



Full wwPDB EM Validation Report ⓘ

Aug 17, 2026 – 11:47 pm BST

PDB ID : 9TLB / pdb_00009tlb
EMDB ID : EMD-56048
Title : Cryo-EM Structure of the oligomeric LPOR:Chlide:NADPH Complexes RF-23
Authors : Gabruk, M.; Desfosses, A.; Estrozi, L.F.; Pintscher, S.; Rawski, M.
Deposited on : 2025-12-10
Resolution : 3.20 Å(reported)
Based on initial model : 7JK9

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

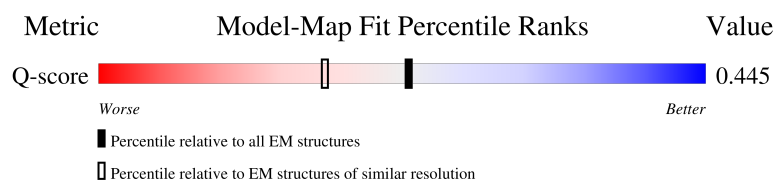
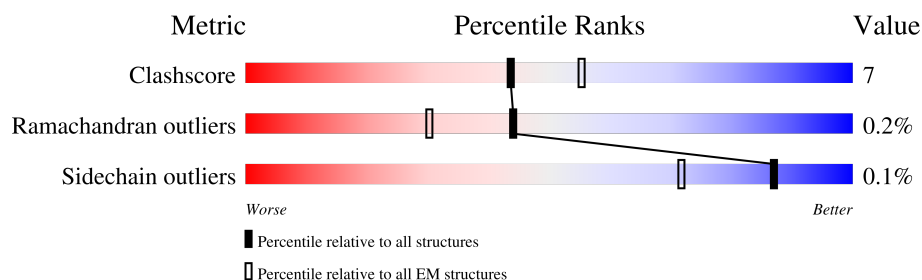
EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15020 (2.70 - 3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	YI	353	
1	YJ	353	
1	YK	353	
1	YL	353	

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Mol	Chain	Length	Quality of chain
1	YM	353	
1	YN	353	
1	YO	353	
1	YP	353	
1	YQ	353	
1	YR	353	
1	YS	353	
1	YT	353	
1	YU	353	
1	YV	353	
1	YW	353	
1	YX	353	
1	YY	353	
1	YZ	353	
1	ZA	353	
1	ZB	353	
1	ZC	353	
1	ZD	353	
1	ZE	353	
1	ZF	353	
1	ZG	353	
1	ZH	353	
1	ZI	353	
1	ZJ	353	
1	ZK	353	

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Mol	Chain	Length	Quality of chain
1	ZL	353	
1	ZM	353	
1	ZN	353	
1	ZO	353	
1	ZP	353	
1	ZQ	353	
1	ZR	353	
1	ZS	353	
1	ZT	353	
1	ZU	353	
1	ZV	353	
1	ZW	353	
1	ZX	353	
1	ZY	353	
1	ZZ	353	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	A1JWG	YI	503	X	-	-	-
4	A1JWG	YJ	503	X	-	-	-
4	A1JWG	YK	503	X	-	-	-
4	A1JWG	YL	503	X	-	-	-
4	A1JWG	YM	503	X	-	-	-
4	A1JWG	YN	503	X	-	-	-
4	A1JWG	YO	503	X	-	-	-
4	A1JWG	YP	503	X	-	-	-
4	A1JWG	YQ	503	X	-	-	-
4	A1JWG	YR	503	X	-	-	-
4	A1JWG	YS	503	X	-	-	-
4	A1JWG	YT	503	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	A1JWG	YU	503	X	-	-	-
4	A1JWG	YV	503	X	-	-	-
4	A1JWG	YW	503	X	-	-	-
4	A1JWG	YX	503	X	-	-	-
4	A1JWG	YY	503	X	-	-	-
4	A1JWG	YZ	503	X	-	-	-
4	A1JWG	ZA	503	X	-	-	-
4	A1JWG	ZB	503	X	-	-	-
4	A1JWG	ZC	503	X	-	-	-
4	A1JWG	ZD	503	X	-	-	-
4	A1JWG	ZE	503	X	-	-	-
4	A1JWG	ZF	503	X	-	-	-
4	A1JWG	ZG	503	X	-	-	-
4	A1JWG	ZH	503	X	-	-	-
4	A1JWG	ZI	503	X	-	-	-
4	A1JWG	ZJ	503	X	-	-	-
4	A1JWG	ZK	503	X	-	-	-
4	A1JWG	ZL	503	X	-	-	-
4	A1JWG	ZM	503	X	-	-	-
4	A1JWG	ZN	503	X	-	-	-
4	A1JWG	ZO	503	X	-	-	-
4	A1JWG	ZP	503	X	-	-	-
4	A1JWG	ZQ	503	X	-	-	-
4	A1JWG	ZR	503	X	-	-	-
4	A1JWG	ZS	503	X	-	-	-
4	A1JWG	ZT	503	X	-	-	-
4	A1JWG	ZU	503	X	-	-	-
4	A1JWG	ZV	503	X	-	-	-
4	A1JWG	ZW	503	X	-	-	-
4	A1JWG	ZX	503	X	-	-	-
4	A1JWG	ZY	503	X	-	-	-
4	A1JWG	ZZ	503	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 111980 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Protochlorophyllide reductase B, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	YI	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YJ	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YK	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YL	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YM	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YN	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YO	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YP	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YQ	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YR	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YS	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YT	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YU	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YV	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YW	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	YX	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	YY	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	YZ	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZA	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZB	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZC	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZD	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZE	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZF	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZG	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZH	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZI	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZJ	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZK	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZL	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZM	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZN	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZO	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZP	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZQ	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZR	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZS	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZT	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		

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Mol	Chain	Residues	Atoms					AltConf	Trace
1	ZU	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZV	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZW	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZX	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		
1	ZY	316	Total	C	N	O	S	0	0
			2421	1534	418	459	10		
1	ZZ	318	Total	C	N	O	S	0	0
			2437	1544	421	462	10		

There are 836 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
YI	49	MET	-	initiating methionine	UNP P21218
YI	50	GLY	-	expression tag	UNP P21218
YI	51	SER	-	expression tag	UNP P21218
YI	52	SER	-	expression tag	UNP P21218
YI	53	HIS	-	expression tag	UNP P21218
YI	54	HIS	-	expression tag	UNP P21218
YI	55	HIS	-	expression tag	UNP P21218
YI	56	HIS	-	expression tag	UNP P21218
YI	57	HIS	-	expression tag	UNP P21218
YI	58	HIS	-	expression tag	UNP P21218
YI	59	SER	-	expression tag	UNP P21218
YI	60	SER	-	expression tag	UNP P21218
YI	61	GLY	-	expression tag	UNP P21218
YI	62	LEU	-	expression tag	UNP P21218
YI	63	VAL	-	expression tag	UNP P21218
YI	64	PRO	-	expression tag	UNP P21218
YI	65	ARG	-	expression tag	UNP P21218
YI	66	GLY	-	expression tag	UNP P21218
YI	67	SER	-	expression tag	UNP P21218
YJ	49	MET	-	initiating methionine	UNP P21218
YJ	50	GLY	-	expression tag	UNP P21218
YJ	51	SER	-	expression tag	UNP P21218
YJ	52	SER	-	expression tag	UNP P21218
YJ	53	HIS	-	expression tag	UNP P21218
YJ	54	HIS	-	expression tag	UNP P21218
YJ	55	HIS	-	expression tag	UNP P21218
YJ	56	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YJ	57	HIS	-	expression tag	UNP P21218
YJ	58	HIS	-	expression tag	UNP P21218
YJ	59	SER	-	expression tag	UNP P21218
YJ	60	SER	-	expression tag	UNP P21218
YJ	61	GLY	-	expression tag	UNP P21218
YJ	62	LEU	-	expression tag	UNP P21218
YJ	63	VAL	-	expression tag	UNP P21218
YJ	64	PRO	-	expression tag	UNP P21218
YJ	65	ARG	-	expression tag	UNP P21218
YJ	66	GLY	-	expression tag	UNP P21218
YJ	67	SER	-	expression tag	UNP P21218
YK	49	MET	-	initiating methionine	UNP P21218
YK	50	GLY	-	expression tag	UNP P21218
YK	51	SER	-	expression tag	UNP P21218
YK	52	SER	-	expression tag	UNP P21218
YK	53	HIS	-	expression tag	UNP P21218
YK	54	HIS	-	expression tag	UNP P21218
YK	55	HIS	-	expression tag	UNP P21218
YK	56	HIS	-	expression tag	UNP P21218
YK	57	HIS	-	expression tag	UNP P21218
YK	58	HIS	-	expression tag	UNP P21218
YK	59	SER	-	expression tag	UNP P21218
YK	60	SER	-	expression tag	UNP P21218
YK	61	GLY	-	expression tag	UNP P21218
YK	62	LEU	-	expression tag	UNP P21218
YK	63	VAL	-	expression tag	UNP P21218
YK	64	PRO	-	expression tag	UNP P21218
YK	65	ARG	-	expression tag	UNP P21218
YK	66	GLY	-	expression tag	UNP P21218
YK	67	SER	-	expression tag	UNP P21218
YL	49	MET	-	initiating methionine	UNP P21218
YL	50	GLY	-	expression tag	UNP P21218
YL	51	SER	-	expression tag	UNP P21218
YL	52	SER	-	expression tag	UNP P21218
YL	53	HIS	-	expression tag	UNP P21218
YL	54	HIS	-	expression tag	UNP P21218
YL	55	HIS	-	expression tag	UNP P21218
YL	56	HIS	-	expression tag	UNP P21218
YL	57	HIS	-	expression tag	UNP P21218
YL	58	HIS	-	expression tag	UNP P21218
YL	59	SER	-	expression tag	UNP P21218
YL	60	SER	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YL	61	GLY	-	expression tag	UNP P21218
YL	62	LEU	-	expression tag	UNP P21218
YL	63	VAL	-	expression tag	UNP P21218
YL	64	PRO	-	expression tag	UNP P21218
YL	65	ARG	-	expression tag	UNP P21218
YL	66	GLY	-	expression tag	UNP P21218
YL	67	SER	-	expression tag	UNP P21218
YM	49	MET	-	initiating methionine	UNP P21218
YM	50	GLY	-	expression tag	UNP P21218
YM	51	SER	-	expression tag	UNP P21218
YM	52	SER	-	expression tag	UNP P21218
YM	53	HIS	-	expression tag	UNP P21218
YM	54	HIS	-	expression tag	UNP P21218
YM	55	HIS	-	expression tag	UNP P21218
YM	56	HIS	-	expression tag	UNP P21218
YM	57	HIS	-	expression tag	UNP P21218
YM	58	HIS	-	expression tag	UNP P21218
YM	59	SER	-	expression tag	UNP P21218
YM	60	SER	-	expression tag	UNP P21218
YM	61	GLY	-	expression tag	UNP P21218
YM	62	LEU	-	expression tag	UNP P21218
YM	63	VAL	-	expression tag	UNP P21218
YM	64	PRO	-	expression tag	UNP P21218
YM	65	ARG	-	expression tag	UNP P21218
YM	66	GLY	-	expression tag	UNP P21218
YM	67	SER	-	expression tag	UNP P21218
YN	49	MET	-	initiating methionine	UNP P21218
YN	50	GLY	-	expression tag	UNP P21218
YN	51	SER	-	expression tag	UNP P21218
YN	52	SER	-	expression tag	UNP P21218
YN	53	HIS	-	expression tag	UNP P21218
YN	54	HIS	-	expression tag	UNP P21218
YN	55	HIS	-	expression tag	UNP P21218
YN	56	HIS	-	expression tag	UNP P21218
YN	57	HIS	-	expression tag	UNP P21218
YN	58	HIS	-	expression tag	UNP P21218
YN	59	SER	-	expression tag	UNP P21218
YN	60	SER	-	expression tag	UNP P21218
YN	61	GLY	-	expression tag	UNP P21218
YN	62	LEU	-	expression tag	UNP P21218
YN	63	VAL	-	expression tag	UNP P21218
YN	64	PRO	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YN	65	ARG	-	expression tag	UNP P21218
YN	66	GLY	-	expression tag	UNP P21218
YN	67	SER	-	expression tag	UNP P21218
YO	49	MET	-	initiating methionine	UNP P21218
YO	50	GLY	-	expression tag	UNP P21218
YO	51	SER	-	expression tag	UNP P21218
YO	52	SER	-	expression tag	UNP P21218
YO	53	HIS	-	expression tag	UNP P21218
YO	54	HIS	-	expression tag	UNP P21218
YO	55	HIS	-	expression tag	UNP P21218
YO	56	HIS	-	expression tag	UNP P21218
YO	57	HIS	-	expression tag	UNP P21218
YO	58	HIS	-	expression tag	UNP P21218
YO	59	SER	-	expression tag	UNP P21218
YO	60	SER	-	expression tag	UNP P21218
YO	61	GLY	-	expression tag	UNP P21218
YO	62	LEU	-	expression tag	UNP P21218
YO	63	VAL	-	expression tag	UNP P21218
YO	64	PRO	-	expression tag	UNP P21218
YO	65	ARG	-	expression tag	UNP P21218
YO	66	GLY	-	expression tag	UNP P21218
YO	67	SER	-	expression tag	UNP P21218
YP	49	MET	-	initiating methionine	UNP P21218
YP	50	GLY	-	expression tag	UNP P21218
YP	51	SER	-	expression tag	UNP P21218
YP	52	SER	-	expression tag	UNP P21218
YP	53	HIS	-	expression tag	UNP P21218
YP	54	HIS	-	expression tag	UNP P21218
YP	55	HIS	-	expression tag	UNP P21218
YP	56	HIS	-	expression tag	UNP P21218
YP	57	HIS	-	expression tag	UNP P21218
YP	58	HIS	-	expression tag	UNP P21218
YP	59	SER	-	expression tag	UNP P21218
YP	60	SER	-	expression tag	UNP P21218
YP	61	GLY	-	expression tag	UNP P21218
YP	62	LEU	-	expression tag	UNP P21218
YP	63	VAL	-	expression tag	UNP P21218
YP	64	PRO	-	expression tag	UNP P21218
YP	65	ARG	-	expression tag	UNP P21218
YP	66	GLY	-	expression tag	UNP P21218
YP	67	SER	-	expression tag	UNP P21218
YQ	49	MET	-	initiating methionine	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YQ	50	GLY	-	expression tag	UNP P21218
YQ	51	SER	-	expression tag	UNP P21218
YQ	52	SER	-	expression tag	UNP P21218
YQ	53	HIS	-	expression tag	UNP P21218
YQ	54	HIS	-	expression tag	UNP P21218
YQ	55	HIS	-	expression tag	UNP P21218
YQ	56	HIS	-	expression tag	UNP P21218
YQ	57	HIS	-	expression tag	UNP P21218
YQ	58	HIS	-	expression tag	UNP P21218
YQ	59	SER	-	expression tag	UNP P21218
YQ	60	SER	-	expression tag	UNP P21218
YQ	61	GLY	-	expression tag	UNP P21218
YQ	62	LEU	-	expression tag	UNP P21218
YQ	63	VAL	-	expression tag	UNP P21218
YQ	64	PRO	-	expression tag	UNP P21218
YQ	65	ARG	-	expression tag	UNP P21218
YQ	66	GLY	-	expression tag	UNP P21218
YQ	67	SER	-	expression tag	UNP P21218
YR	49	MET	-	initiating methionine	UNP P21218
YR	50	GLY	-	expression tag	UNP P21218
YR	51	SER	-	expression tag	UNP P21218
YR	52	SER	-	expression tag	UNP P21218
YR	53	HIS	-	expression tag	UNP P21218
YR	54	HIS	-	expression tag	UNP P21218
YR	55	HIS	-	expression tag	UNP P21218
YR	56	HIS	-	expression tag	UNP P21218
YR	57	HIS	-	expression tag	UNP P21218
YR	58	HIS	-	expression tag	UNP P21218
YR	59	SER	-	expression tag	UNP P21218
YR	60	SER	-	expression tag	UNP P21218
YR	61	GLY	-	expression tag	UNP P21218
YR	62	LEU	-	expression tag	UNP P21218
YR	63	VAL	-	expression tag	UNP P21218
YR	64	PRO	-	expression tag	UNP P21218
YR	65	ARG	-	expression tag	UNP P21218
YR	66	GLY	-	expression tag	UNP P21218
YR	67	SER	-	expression tag	UNP P21218
YS	49	MET	-	initiating methionine	UNP P21218
YS	50	GLY	-	expression tag	UNP P21218
YS	51	SER	-	expression tag	UNP P21218
YS	52	SER	-	expression tag	UNP P21218
YS	53	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YS	54	HIS	-	expression tag	UNP P21218
YS	55	HIS	-	expression tag	UNP P21218
YS	56	HIS	-	expression tag	UNP P21218
YS	57	HIS	-	expression tag	UNP P21218
YS	58	HIS	-	expression tag	UNP P21218
YS	59	SER	-	expression tag	UNP P21218
YS	60	SER	-	expression tag	UNP P21218
YS	61	GLY	-	expression tag	UNP P21218
YS	62	LEU	-	expression tag	UNP P21218
YS	63	VAL	-	expression tag	UNP P21218
YS	64	PRO	-	expression tag	UNP P21218
YS	65	ARG	-	expression tag	UNP P21218
YS	66	GLY	-	expression tag	UNP P21218
YS	67	SER	-	expression tag	UNP P21218
YT	49	MET	-	initiating methionine	UNP P21218
YT	50	GLY	-	expression tag	UNP P21218
YT	51	SER	-	expression tag	UNP P21218
YT	52	SER	-	expression tag	UNP P21218
YT	53	HIS	-	expression tag	UNP P21218
YT	54	HIS	-	expression tag	UNP P21218
YT	55	HIS	-	expression tag	UNP P21218
YT	56	HIS	-	expression tag	UNP P21218
YT	57	HIS	-	expression tag	UNP P21218
YT	58	HIS	-	expression tag	UNP P21218
YT	59	SER	-	expression tag	UNP P21218
YT	60	SER	-	expression tag	UNP P21218
YT	61	GLY	-	expression tag	UNP P21218
YT	62	LEU	-	expression tag	UNP P21218
YT	63	VAL	-	expression tag	UNP P21218
YT	64	PRO	-	expression tag	UNP P21218
YT	65	ARG	-	expression tag	UNP P21218
YT	66	GLY	-	expression tag	UNP P21218
YT	67	SER	-	expression tag	UNP P21218
YU	49	MET	-	initiating methionine	UNP P21218
YU	50	GLY	-	expression tag	UNP P21218
YU	51	SER	-	expression tag	UNP P21218
YU	52	SER	-	expression tag	UNP P21218
YU	53	HIS	-	expression tag	UNP P21218
YU	54	HIS	-	expression tag	UNP P21218
YU	55	HIS	-	expression tag	UNP P21218
YU	56	HIS	-	expression tag	UNP P21218
YU	57	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YU	58	HIS	-	expression tag	UNP P21218
YU	59	SER	-	expression tag	UNP P21218
YU	60	SER	-	expression tag	UNP P21218
YU	61	GLY	-	expression tag	UNP P21218
YU	62	LEU	-	expression tag	UNP P21218
YU	63	VAL	-	expression tag	UNP P21218
YU	64	PRO	-	expression tag	UNP P21218
YU	65	ARG	-	expression tag	UNP P21218
YU	66	GLY	-	expression tag	UNP P21218
YU	67	SER	-	expression tag	UNP P21218
YV	49	MET	-	initiating methionine	UNP P21218
YV	50	GLY	-	expression tag	UNP P21218
YV	51	SER	-	expression tag	UNP P21218
YV	52	SER	-	expression tag	UNP P21218
YV	53	HIS	-	expression tag	UNP P21218
YV	54	HIS	-	expression tag	UNP P21218
YV	55	HIS	-	expression tag	UNP P21218
YV	56	HIS	-	expression tag	UNP P21218
YV	57	HIS	-	expression tag	UNP P21218
YV	58	HIS	-	expression tag	UNP P21218
YV	59	SER	-	expression tag	UNP P21218
YV	60	SER	-	expression tag	UNP P21218
YV	61	GLY	-	expression tag	UNP P21218
YV	62	LEU	-	expression tag	UNP P21218
YV	63	VAL	-	expression tag	UNP P21218
YV	64	PRO	-	expression tag	UNP P21218
YV	65	ARG	-	expression tag	UNP P21218
YV	66	GLY	-	expression tag	UNP P21218
YV	67	SER	-	expression tag	UNP P21218
YW	49	MET	-	initiating methionine	UNP P21218
YW	50	GLY	-	expression tag	UNP P21218
YW	51	SER	-	expression tag	UNP P21218
YW	52	SER	-	expression tag	UNP P21218
YW	53	HIS	-	expression tag	UNP P21218
YW	54	HIS	-	expression tag	UNP P21218
YW	55	HIS	-	expression tag	UNP P21218
YW	56	HIS	-	expression tag	UNP P21218
YW	57	HIS	-	expression tag	UNP P21218
YW	58	HIS	-	expression tag	UNP P21218
YW	59	SER	-	expression tag	UNP P21218
YW	60	SER	-	expression tag	UNP P21218
YW	61	GLY	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YW	62	LEU	-	expression tag	UNP P21218
YW	63	VAL	-	expression tag	UNP P21218
YW	64	PRO	-	expression tag	UNP P21218
YW	65	ARG	-	expression tag	UNP P21218
YW	66	GLY	-	expression tag	UNP P21218
YW	67	SER	-	expression tag	UNP P21218
YX	49	MET	-	initiating methionine	UNP P21218
YX	50	GLY	-	expression tag	UNP P21218
YX	51	SER	-	expression tag	UNP P21218
YX	52	SER	-	expression tag	UNP P21218
YX	53	HIS	-	expression tag	UNP P21218
YX	54	HIS	-	expression tag	UNP P21218
YX	55	HIS	-	expression tag	UNP P21218
YX	56	HIS	-	expression tag	UNP P21218
YX	57	HIS	-	expression tag	UNP P21218
YX	58	HIS	-	expression tag	UNP P21218
YX	59	SER	-	expression tag	UNP P21218
YX	60	SER	-	expression tag	UNP P21218
YX	61	GLY	-	expression tag	UNP P21218
YX	62	LEU	-	expression tag	UNP P21218
YX	63	VAL	-	expression tag	UNP P21218
YX	64	PRO	-	expression tag	UNP P21218
YX	65	ARG	-	expression tag	UNP P21218
YX	66	GLY	-	expression tag	UNP P21218
YX	67	SER	-	expression tag	UNP P21218
YY	49	MET	-	initiating methionine	UNP P21218
YY	50	GLY	-	expression tag	UNP P21218
YY	51	SER	-	expression tag	UNP P21218
YY	52	SER	-	expression tag	UNP P21218
YY	53	HIS	-	expression tag	UNP P21218
YY	54	HIS	-	expression tag	UNP P21218
YY	55	HIS	-	expression tag	UNP P21218
YY	56	HIS	-	expression tag	UNP P21218
YY	57	HIS	-	expression tag	UNP P21218
YY	58	HIS	-	expression tag	UNP P21218
YY	59	SER	-	expression tag	UNP P21218
YY	60	SER	-	expression tag	UNP P21218
YY	61	GLY	-	expression tag	UNP P21218
YY	62	LEU	-	expression tag	UNP P21218
YY	63	VAL	-	expression tag	UNP P21218
YY	64	PRO	-	expression tag	UNP P21218
YY	65	ARG	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
YY	66	GLY	-	expression tag	UNP P21218
YY	67	SER	-	expression tag	UNP P21218
YZ	49	MET	-	initiating methionine	UNP P21218
YZ	50	GLY	-	expression tag	UNP P21218
YZ	51	SER	-	expression tag	UNP P21218
YZ	52	SER	-	expression tag	UNP P21218
YZ	53	HIS	-	expression tag	UNP P21218
YZ	54	HIS	-	expression tag	UNP P21218
YZ	55	HIS	-	expression tag	UNP P21218
YZ	56	HIS	-	expression tag	UNP P21218
YZ	57	HIS	-	expression tag	UNP P21218
YZ	58	HIS	-	expression tag	UNP P21218
YZ	59	SER	-	expression tag	UNP P21218
YZ	60	SER	-	expression tag	UNP P21218
YZ	61	GLY	-	expression tag	UNP P21218
YZ	62	LEU	-	expression tag	UNP P21218
YZ	63	VAL	-	expression tag	UNP P21218
YZ	64	PRO	-	expression tag	UNP P21218
YZ	65	ARG	-	expression tag	UNP P21218
YZ	66	GLY	-	expression tag	UNP P21218
YZ	67	SER	-	expression tag	UNP P21218
ZA	49	MET	-	initiating methionine	UNP P21218
ZA	50	GLY	-	expression tag	UNP P21218
ZA	51	SER	-	expression tag	UNP P21218
ZA	52	SER	-	expression tag	UNP P21218
ZA	53	HIS	-	expression tag	UNP P21218
ZA	54	HIS	-	expression tag	UNP P21218
ZA	55	HIS	-	expression tag	UNP P21218
ZA	56	HIS	-	expression tag	UNP P21218
ZA	57	HIS	-	expression tag	UNP P21218
ZA	58	HIS	-	expression tag	UNP P21218
ZA	59	SER	-	expression tag	UNP P21218
ZA	60	SER	-	expression tag	UNP P21218
ZA	61	GLY	-	expression tag	UNP P21218
ZA	62	LEU	-	expression tag	UNP P21218
ZA	63	VAL	-	expression tag	UNP P21218
ZA	64	PRO	-	expression tag	UNP P21218
ZA	65	ARG	-	expression tag	UNP P21218
ZA	66	GLY	-	expression tag	UNP P21218
ZA	67	SER	-	expression tag	UNP P21218
ZB	49	MET	-	initiating methionine	UNP P21218
ZB	50	GLY	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZB	51	SER	-	expression tag	UNP P21218
ZB	52	SER	-	expression tag	UNP P21218
ZB	53	HIS	-	expression tag	UNP P21218
ZB	54	HIS	-	expression tag	UNP P21218
ZB	55	HIS	-	expression tag	UNP P21218
ZB	56	HIS	-	expression tag	UNP P21218
ZB	57	HIS	-	expression tag	UNP P21218
ZB	58	HIS	-	expression tag	UNP P21218
ZB	59	SER	-	expression tag	UNP P21218
ZB	60	SER	-	expression tag	UNP P21218
ZB	61	GLY	-	expression tag	UNP P21218
ZB	62	LEU	-	expression tag	UNP P21218
ZB	63	VAL	-	expression tag	UNP P21218
ZB	64	PRO	-	expression tag	UNP P21218
ZB	65	ARG	-	expression tag	UNP P21218
ZB	66	GLY	-	expression tag	UNP P21218
ZB	67	SER	-	expression tag	UNP P21218
ZC	49	MET	-	initiating methionine	UNP P21218
ZC	50	GLY	-	expression tag	UNP P21218
ZC	51	SER	-	expression tag	UNP P21218
ZC	52	SER	-	expression tag	UNP P21218
ZC	53	HIS	-	expression tag	UNP P21218
ZC	54	HIS	-	expression tag	UNP P21218
ZC	55	HIS	-	expression tag	UNP P21218
ZC	56	HIS	-	expression tag	UNP P21218
ZC	57	HIS	-	expression tag	UNP P21218
ZC	58	HIS	-	expression tag	UNP P21218
ZC	59	SER	-	expression tag	UNP P21218
ZC	60	SER	-	expression tag	UNP P21218
ZC	61	GLY	-	expression tag	UNP P21218
ZC	62	LEU	-	expression tag	UNP P21218
ZC	63	VAL	-	expression tag	UNP P21218
ZC	64	PRO	-	expression tag	UNP P21218
ZC	65	ARG	-	expression tag	UNP P21218
ZC	66	GLY	-	expression tag	UNP P21218
ZC	67	SER	-	expression tag	UNP P21218
ZD	49	MET	-	initiating methionine	UNP P21218
ZD	50	GLY	-	expression tag	UNP P21218
ZD	51	SER	-	expression tag	UNP P21218
ZD	52	SER	-	expression tag	UNP P21218
ZD	53	HIS	-	expression tag	UNP P21218
ZD	54	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZD	55	HIS	-	expression tag	UNP P21218
ZD	56	HIS	-	expression tag	UNP P21218
ZD	57	HIS	-	expression tag	UNP P21218
ZD	58	HIS	-	expression tag	UNP P21218
ZD	59	SER	-	expression tag	UNP P21218
ZD	60	SER	-	expression tag	UNP P21218
ZD	61	GLY	-	expression tag	UNP P21218
ZD	62	LEU	-	expression tag	UNP P21218
ZD	63	VAL	-	expression tag	UNP P21218
ZD	64	PRO	-	expression tag	UNP P21218
ZD	65	ARG	-	expression tag	UNP P21218
ZD	66	GLY	-	expression tag	UNP P21218
ZD	67	SER	-	expression tag	UNP P21218
ZE	49	MET	-	initiating methionine	UNP P21218
ZE	50	GLY	-	expression tag	UNP P21218
ZE	51	SER	-	expression tag	UNP P21218
ZE	52	SER	-	expression tag	UNP P21218
ZE	53	HIS	-	expression tag	UNP P21218
ZE	54	HIS	-	expression tag	UNP P21218
ZE	55	HIS	-	expression tag	UNP P21218
ZE	56	HIS	-	expression tag	UNP P21218
ZE	57	HIS	-	expression tag	UNP P21218
ZE	58	HIS	-	expression tag	UNP P21218
ZE	59	SER	-	expression tag	UNP P21218
ZE	60	SER	-	expression tag	UNP P21218
ZE	61	GLY	-	expression tag	UNP P21218
ZE	62	LEU	-	expression tag	UNP P21218
ZE	63	VAL	-	expression tag	UNP P21218
ZE	64	PRO	-	expression tag	UNP P21218
ZE	65	ARG	-	expression tag	UNP P21218
ZE	66	GLY	-	expression tag	UNP P21218
ZE	67	SER	-	expression tag	UNP P21218
ZF	49	MET	-	initiating methionine	UNP P21218
ZF	50	GLY	-	expression tag	UNP P21218
ZF	51	SER	-	expression tag	UNP P21218
ZF	52	SER	-	expression tag	UNP P21218
ZF	53	HIS	-	expression tag	UNP P21218
ZF	54	HIS	-	expression tag	UNP P21218
ZF	55	HIS	-	expression tag	UNP P21218
ZF	56	HIS	-	expression tag	UNP P21218
ZF	57	HIS	-	expression tag	UNP P21218
ZF	58	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZF	59	SER	-	expression tag	UNP P21218
ZF	60	SER	-	expression tag	UNP P21218
ZF	61	GLY	-	expression tag	UNP P21218
ZF	62	LEU	-	expression tag	UNP P21218
ZF	63	VAL	-	expression tag	UNP P21218
ZF	64	PRO	-	expression tag	UNP P21218
ZF	65	ARG	-	expression tag	UNP P21218
ZF	66	GLY	-	expression tag	UNP P21218
ZF	67	SER	-	expression tag	UNP P21218
ZG	49	MET	-	initiating methionine	UNP P21218
ZG	50	GLY	-	expression tag	UNP P21218
ZG	51	SER	-	expression tag	UNP P21218
ZG	52	SER	-	expression tag	UNP P21218
ZG	53	HIS	-	expression tag	UNP P21218
ZG	54	HIS	-	expression tag	UNP P21218
ZG	55	HIS	-	expression tag	UNP P21218
ZG	56	HIS	-	expression tag	UNP P21218
ZG	57	HIS	-	expression tag	UNP P21218
ZG	58	HIS	-	expression tag	UNP P21218
ZG	59	SER	-	expression tag	UNP P21218
ZG	60	SER	-	expression tag	UNP P21218
ZG	61	GLY	-	expression tag	UNP P21218
ZG	62	LEU	-	expression tag	UNP P21218
ZG	63	VAL	-	expression tag	UNP P21218
ZG	64	PRO	-	expression tag	UNP P21218
ZG	65	ARG	-	expression tag	UNP P21218
ZG	66	GLY	-	expression tag	UNP P21218
ZG	67	SER	-	expression tag	UNP P21218
ZH	49	MET	-	initiating methionine	UNP P21218
ZH	50	GLY	-	expression tag	UNP P21218
ZH	51	SER	-	expression tag	UNP P21218
ZH	52	SER	-	expression tag	UNP P21218
ZH	53	HIS	-	expression tag	UNP P21218
ZH	54	HIS	-	expression tag	UNP P21218
ZH	55	HIS	-	expression tag	UNP P21218
ZH	56	HIS	-	expression tag	UNP P21218
ZH	57	HIS	-	expression tag	UNP P21218
ZH	58	HIS	-	expression tag	UNP P21218
ZH	59	SER	-	expression tag	UNP P21218
ZH	60	SER	-	expression tag	UNP P21218
ZH	61	GLY	-	expression tag	UNP P21218
ZH	62	LEU	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZH	63	VAL	-	expression tag	UNP P21218
ZH	64	PRO	-	expression tag	UNP P21218
ZH	65	ARG	-	expression tag	UNP P21218
ZH	66	GLY	-	expression tag	UNP P21218
ZH	67	SER	-	expression tag	UNP P21218
ZI	49	MET	-	initiating methionine	UNP P21218
ZI	50	GLY	-	expression tag	UNP P21218
ZI	51	SER	-	expression tag	UNP P21218
ZI	52	SER	-	expression tag	UNP P21218
ZI	53	HIS	-	expression tag	UNP P21218
ZI	54	HIS	-	expression tag	UNP P21218
ZI	55	HIS	-	expression tag	UNP P21218
ZI	56	HIS	-	expression tag	UNP P21218
ZI	57	HIS	-	expression tag	UNP P21218
ZI	58	HIS	-	expression tag	UNP P21218
ZI	59	SER	-	expression tag	UNP P21218
ZI	60	SER	-	expression tag	UNP P21218
ZI	61	GLY	-	expression tag	UNP P21218
ZI	62	LEU	-	expression tag	UNP P21218
ZI	63	VAL	-	expression tag	UNP P21218
ZI	64	PRO	-	expression tag	UNP P21218
ZI	65	ARG	-	expression tag	UNP P21218
ZI	66	GLY	-	expression tag	UNP P21218
ZI	67	SER	-	expression tag	UNP P21218
ZJ	49	MET	-	initiating methionine	UNP P21218
ZJ	50	GLY	-	expression tag	UNP P21218
ZJ	51	SER	-	expression tag	UNP P21218
ZJ	52	SER	-	expression tag	UNP P21218
ZJ	53	HIS	-	expression tag	UNP P21218
ZJ	54	HIS	-	expression tag	UNP P21218
ZJ	55	HIS	-	expression tag	UNP P21218
ZJ	56	HIS	-	expression tag	UNP P21218
ZJ	57	HIS	-	expression tag	UNP P21218
ZJ	58	HIS	-	expression tag	UNP P21218
ZJ	59	SER	-	expression tag	UNP P21218
ZJ	60	SER	-	expression tag	UNP P21218
ZJ	61	GLY	-	expression tag	UNP P21218
ZJ	62	LEU	-	expression tag	UNP P21218
ZJ	63	VAL	-	expression tag	UNP P21218
ZJ	64	PRO	-	expression tag	UNP P21218
ZJ	65	ARG	-	expression tag	UNP P21218
ZJ	66	GLY	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZJ	67	SER	-	expression tag	UNP P21218
ZK	49	MET	-	initiating methionine	UNP P21218
ZK	50	GLY	-	expression tag	UNP P21218
ZK	51	SER	-	expression tag	UNP P21218
ZK	52	SER	-	expression tag	UNP P21218
ZK	53	HIS	-	expression tag	UNP P21218
ZK	54	HIS	-	expression tag	UNP P21218
ZK	55	HIS	-	expression tag	UNP P21218
ZK	56	HIS	-	expression tag	UNP P21218
ZK	57	HIS	-	expression tag	UNP P21218
ZK	58	HIS	-	expression tag	UNP P21218
ZK	59	SER	-	expression tag	UNP P21218
ZK	60	SER	-	expression tag	UNP P21218
ZK	61	GLY	-	expression tag	UNP P21218
ZK	62	LEU	-	expression tag	UNP P21218
ZK	63	VAL	-	expression tag	UNP P21218
ZK	64	PRO	-	expression tag	UNP P21218
ZK	65	ARG	-	expression tag	UNP P21218
ZK	66	GLY	-	expression tag	UNP P21218
ZK	67	SER	-	expression tag	UNP P21218
ZL	49	MET	-	initiating methionine	UNP P21218
ZL	50	GLY	-	expression tag	UNP P21218
ZL	51	SER	-	expression tag	UNP P21218
ZL	52	SER	-	expression tag	UNP P21218
ZL	53	HIS	-	expression tag	UNP P21218
ZL	54	HIS	-	expression tag	UNP P21218
ZL	55	HIS	-	expression tag	UNP P21218
ZL	56	HIS	-	expression tag	UNP P21218
ZL	57	HIS	-	expression tag	UNP P21218
ZL	58	HIS	-	expression tag	UNP P21218
ZL	59	SER	-	expression tag	UNP P21218
ZL	60	SER	-	expression tag	UNP P21218
ZL	61	GLY	-	expression tag	UNP P21218
ZL	62	LEU	-	expression tag	UNP P21218
ZL	63	VAL	-	expression tag	UNP P21218
ZL	64	PRO	-	expression tag	UNP P21218
ZL	65	ARG	-	expression tag	UNP P21218
ZL	66	GLY	-	expression tag	UNP P21218
ZL	67	SER	-	expression tag	UNP P21218
ZM	49	MET	-	initiating methionine	UNP P21218
ZM	50	GLY	-	expression tag	UNP P21218
ZM	51	SER	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZM	52	SER	-	expression tag	UNP P21218
ZM	53	HIS	-	expression tag	UNP P21218
ZM	54	HIS	-	expression tag	UNP P21218
ZM	55	HIS	-	expression tag	UNP P21218
ZM	56	HIS	-	expression tag	UNP P21218
ZM	57	HIS	-	expression tag	UNP P21218
ZM	58	HIS	-	expression tag	UNP P21218
ZM	59	SER	-	expression tag	UNP P21218
ZM	60	SER	-	expression tag	UNP P21218
ZM	61	GLY	-	expression tag	UNP P21218
ZM	62	LEU	-	expression tag	UNP P21218
ZM	63	VAL	-	expression tag	UNP P21218
ZM	64	PRO	-	expression tag	UNP P21218
ZM	65	ARG	-	expression tag	UNP P21218
ZM	66	GLY	-	expression tag	UNP P21218
ZM	67	SER	-	expression tag	UNP P21218
ZN	49	MET	-	initiating methionine	UNP P21218
ZN	50	GLY	-	expression tag	UNP P21218
ZN	51	SER	-	expression tag	UNP P21218
ZN	52	SER	-	expression tag	UNP P21218
ZN	53	HIS	-	expression tag	UNP P21218
ZN	54	HIS	-	expression tag	UNP P21218
ZN	55	HIS	-	expression tag	UNP P21218
ZN	56	HIS	-	expression tag	UNP P21218
ZN	57	HIS	-	expression tag	UNP P21218
ZN	58	HIS	-	expression tag	UNP P21218
ZN	59	SER	-	expression tag	UNP P21218
ZN	60	SER	-	expression tag	UNP P21218
ZN	61	GLY	-	expression tag	UNP P21218
ZN	62	LEU	-	expression tag	UNP P21218
ZN	63	VAL	-	expression tag	UNP P21218
ZN	64	PRO	-	expression tag	UNP P21218
ZN	65	ARG	-	expression tag	UNP P21218
ZN	66	GLY	-	expression tag	UNP P21218
ZN	67	SER	-	expression tag	UNP P21218
ZO	49	MET	-	initiating methionine	UNP P21218
ZO	50	GLY	-	expression tag	UNP P21218
ZO	51	SER	-	expression tag	UNP P21218
ZO	52	SER	-	expression tag	UNP P21218
ZO	53	HIS	-	expression tag	UNP P21218
ZO	54	HIS	-	expression tag	UNP P21218
ZO	55	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZO	56	HIS	-	expression tag	UNP P21218
ZO	57	HIS	-	expression tag	UNP P21218
ZO	58	HIS	-	expression tag	UNP P21218
ZO	59	SER	-	expression tag	UNP P21218
ZO	60	SER	-	expression tag	UNP P21218
ZO	61	GLY	-	expression tag	UNP P21218
ZO	62	LEU	-	expression tag	UNP P21218
ZO	63	VAL	-	expression tag	UNP P21218
ZO	64	PRO	-	expression tag	UNP P21218
ZO	65	ARG	-	expression tag	UNP P21218
ZO	66	GLY	-	expression tag	UNP P21218
ZO	67	SER	-	expression tag	UNP P21218
ZP	49	MET	-	initiating methionine	UNP P21218
ZP	50	GLY	-	expression tag	UNP P21218
ZP	51	SER	-	expression tag	UNP P21218
ZP	52	SER	-	expression tag	UNP P21218
ZP	53	HIS	-	expression tag	UNP P21218
ZP	54	HIS	-	expression tag	UNP P21218
ZP	55	HIS	-	expression tag	UNP P21218
ZP	56	HIS	-	expression tag	UNP P21218
ZP	57	HIS	-	expression tag	UNP P21218
ZP	58	HIS	-	expression tag	UNP P21218
ZP	59	SER	-	expression tag	UNP P21218
ZP	60	SER	-	expression tag	UNP P21218
ZP	61	GLY	-	expression tag	UNP P21218
ZP	62	LEU	-	expression tag	UNP P21218
ZP	63	VAL	-	expression tag	UNP P21218
ZP	64	PRO	-	expression tag	UNP P21218
ZP	65	ARG	-	expression tag	UNP P21218
ZP	66	GLY	-	expression tag	UNP P21218
ZP	67	SER	-	expression tag	UNP P21218
ZQ	49	MET	-	initiating methionine	UNP P21218
ZQ	50	GLY	-	expression tag	UNP P21218
ZQ	51	SER	-	expression tag	UNP P21218
ZQ	52	SER	-	expression tag	UNP P21218
ZQ	53	HIS	-	expression tag	UNP P21218
ZQ	54	HIS	-	expression tag	UNP P21218
ZQ	55	HIS	-	expression tag	UNP P21218
ZQ	56	HIS	-	expression tag	UNP P21218
ZQ	57	HIS	-	expression tag	UNP P21218
ZQ	58	HIS	-	expression tag	UNP P21218
ZQ	59	SER	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZQ	60	SER	-	expression tag	UNP P21218
ZQ	61	GLY	-	expression tag	UNP P21218
ZQ	62	LEU	-	expression tag	UNP P21218
ZQ	63	VAL	-	expression tag	UNP P21218
ZQ	64	PRO	-	expression tag	UNP P21218
ZQ	65	ARG	-	expression tag	UNP P21218
ZQ	66	GLY	-	expression tag	UNP P21218
ZQ	67	SER	-	expression tag	UNP P21218
ZR	49	MET	-	initiating methionine	UNP P21218
ZR	50	GLY	-	expression tag	UNP P21218
ZR	51	SER	-	expression tag	UNP P21218
ZR	52	SER	-	expression tag	UNP P21218
ZR	53	HIS	-	expression tag	UNP P21218
ZR	54	HIS	-	expression tag	UNP P21218
ZR	55	HIS	-	expression tag	UNP P21218
ZR	56	HIS	-	expression tag	UNP P21218
ZR	57	HIS	-	expression tag	UNP P21218
ZR	58	HIS	-	expression tag	UNP P21218
ZR	59	SER	-	expression tag	UNP P21218
ZR	60	SER	-	expression tag	UNP P21218
ZR	61	GLY	-	expression tag	UNP P21218
ZR	62	LEU	-	expression tag	UNP P21218
ZR	63	VAL	-	expression tag	UNP P21218
ZR	64	PRO	-	expression tag	UNP P21218
ZR	65	ARG	-	expression tag	UNP P21218
ZR	66	GLY	-	expression tag	UNP P21218
ZR	67	SER	-	expression tag	UNP P21218
ZS	49	MET	-	initiating methionine	UNP P21218
ZS	50	GLY	-	expression tag	UNP P21218
ZS	51	SER	-	expression tag	UNP P21218
ZS	52	SER	-	expression tag	UNP P21218
ZS	53	HIS	-	expression tag	UNP P21218
ZS	54	HIS	-	expression tag	UNP P21218
ZS	55	HIS	-	expression tag	UNP P21218
ZS	56	HIS	-	expression tag	UNP P21218
ZS	57	HIS	-	expression tag	UNP P21218
ZS	58	HIS	-	expression tag	UNP P21218
ZS	59	SER	-	expression tag	UNP P21218
ZS	60	SER	-	expression tag	UNP P21218
ZS	61	GLY	-	expression tag	UNP P21218
ZS	62	LEU	-	expression tag	UNP P21218
ZS	63	VAL	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZS	64	PRO	-	expression tag	UNP P21218
ZS	65	ARG	-	expression tag	UNP P21218
ZS	66	GLY	-	expression tag	UNP P21218
ZS	67	SER	-	expression tag	UNP P21218
ZT	49	MET	-	initiating methionine	UNP P21218
ZT	50	GLY	-	expression tag	UNP P21218
ZT	51	SER	-	expression tag	UNP P21218
ZT	52	SER	-	expression tag	UNP P21218
ZT	53	HIS	-	expression tag	UNP P21218
ZT	54	HIS	-	expression tag	UNP P21218
ZT	55	HIS	-	expression tag	UNP P21218
ZT	56	HIS	-	expression tag	UNP P21218
ZT	57	HIS	-	expression tag	UNP P21218
ZT	58	HIS	-	expression tag	UNP P21218
ZT	59	SER	-	expression tag	UNP P21218
ZT	60	SER	-	expression tag	UNP P21218
ZT	61	GLY	-	expression tag	UNP P21218
ZT	62	LEU	-	expression tag	UNP P21218
ZT	63	VAL	-	expression tag	UNP P21218
ZT	64	PRO	-	expression tag	UNP P21218
ZT	65	ARG	-	expression tag	UNP P21218
ZT	66	GLY	-	expression tag	UNP P21218
ZT	67	SER	-	expression tag	UNP P21218
ZU	49	MET	-	initiating methionine	UNP P21218
ZU	50	GLY	-	expression tag	UNP P21218
ZU	51	SER	-	expression tag	UNP P21218
ZU	52	SER	-	expression tag	UNP P21218
ZU	53	HIS	-	expression tag	UNP P21218
ZU	54	HIS	-	expression tag	UNP P21218
ZU	55	HIS	-	expression tag	UNP P21218
ZU	56	HIS	-	expression tag	UNP P21218
ZU	57	HIS	-	expression tag	UNP P21218
ZU	58	HIS	-	expression tag	UNP P21218
ZU	59	SER	-	expression tag	UNP P21218
ZU	60	SER	-	expression tag	UNP P21218
ZU	61	GLY	-	expression tag	UNP P21218
ZU	62	LEU	-	expression tag	UNP P21218
ZU	63	VAL	-	expression tag	UNP P21218
ZU	64	PRO	-	expression tag	UNP P21218
ZU	65	ARG	-	expression tag	UNP P21218
ZU	66	GLY	-	expression tag	UNP P21218
ZU	67	SER	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZV	49	MET	-	initiating methionine	UNP P21218
ZV	50	GLY	-	expression tag	UNP P21218
ZV	51	SER	-	expression tag	UNP P21218
ZV	52	SER	-	expression tag	UNP P21218
ZV	53	HIS	-	expression tag	UNP P21218
ZV	54	HIS	-	expression tag	UNP P21218
ZV	55	HIS	-	expression tag	UNP P21218
ZV	56	HIS	-	expression tag	UNP P21218
ZV	57	HIS	-	expression tag	UNP P21218
ZV	58	HIS	-	expression tag	UNP P21218
ZV	59	SER	-	expression tag	UNP P21218
ZV	60	SER	-	expression tag	UNP P21218
ZV	61	GLY	-	expression tag	UNP P21218
ZV	62	LEU	-	expression tag	UNP P21218
ZV	63	VAL	-	expression tag	UNP P21218
ZV	64	PRO	-	expression tag	UNP P21218
ZV	65	ARG	-	expression tag	UNP P21218
ZV	66	GLY	-	expression tag	UNP P21218
ZV	67	SER	-	expression tag	UNP P21218
ZW	49	MET	-	initiating methionine	UNP P21218
ZW	50	GLY	-	expression tag	UNP P21218
ZW	51	SER	-	expression tag	UNP P21218
ZW	52	SER	-	expression tag	UNP P21218
ZW	53	HIS	-	expression tag	UNP P21218
ZW	54	HIS	-	expression tag	UNP P21218
ZW	55	HIS	-	expression tag	UNP P21218
ZW	56	HIS	-	expression tag	UNP P21218
ZW	57	HIS	-	expression tag	UNP P21218
ZW	58	HIS	-	expression tag	UNP P21218
ZW	59	SER	-	expression tag	UNP P21218
ZW	60	SER	-	expression tag	UNP P21218
ZW	61	GLY	-	expression tag	UNP P21218
ZW	62	LEU	-	expression tag	UNP P21218
ZW	63	VAL	-	expression tag	UNP P21218
ZW	64	PRO	-	expression tag	UNP P21218
ZW	65	ARG	-	expression tag	UNP P21218
ZW	66	GLY	-	expression tag	UNP P21218
ZW	67	SER	-	expression tag	UNP P21218
ZX	49	MET	-	initiating methionine	UNP P21218
ZX	50	GLY	-	expression tag	UNP P21218
ZX	51	SER	-	expression tag	UNP P21218
ZX	52	SER	-	expression tag	UNP P21218

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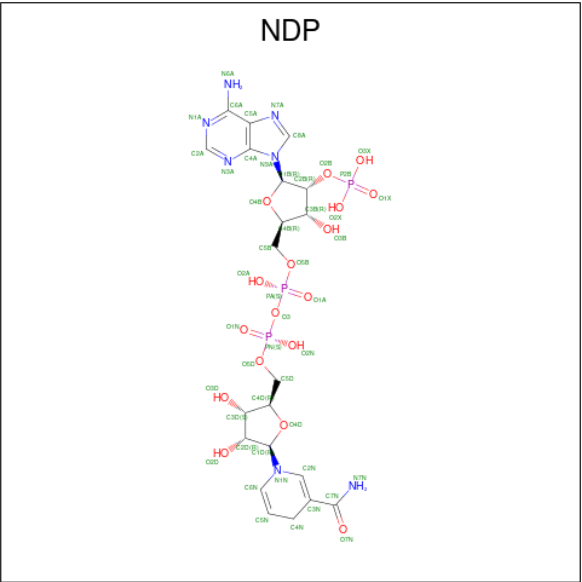
Chain	Residue	Modelled	Actual	Comment	Reference
ZX	53	HIS	-	expression tag	UNP P21218
ZX	54	HIS	-	expression tag	UNP P21218
ZX	55	HIS	-	expression tag	UNP P21218
ZX	56	HIS	-	expression tag	UNP P21218
ZX	57	HIS	-	expression tag	UNP P21218
ZX	58	HIS	-	expression tag	UNP P21218
ZX	59	SER	-	expression tag	UNP P21218
ZX	60	SER	-	expression tag	UNP P21218
ZX	61	GLY	-	expression tag	UNP P21218
ZX	62	LEU	-	expression tag	UNP P21218
ZX	63	VAL	-	expression tag	UNP P21218
ZX	64	PRO	-	expression tag	UNP P21218
ZX	65	ARG	-	expression tag	UNP P21218
ZX	66	GLY	-	expression tag	UNP P21218
ZX	67	SER	-	expression tag	UNP P21218
ZY	49	MET	-	initiating methionine	UNP P21218
ZY	50	GLY	-	expression tag	UNP P21218
ZY	51	SER	-	expression tag	UNP P21218
ZY	52	SER	-	expression tag	UNP P21218
ZY	53	HIS	-	expression tag	UNP P21218
ZY	54	HIS	-	expression tag	UNP P21218
ZY	55	HIS	-	expression tag	UNP P21218
ZY	56	HIS	-	expression tag	UNP P21218
ZY	57	HIS	-	expression tag	UNP P21218
ZY	58	HIS	-	expression tag	UNP P21218
ZY	59	SER	-	expression tag	UNP P21218
ZY	60	SER	-	expression tag	UNP P21218
ZY	61	GLY	-	expression tag	UNP P21218
ZY	62	LEU	-	expression tag	UNP P21218
ZY	63	VAL	-	expression tag	UNP P21218
ZY	64	PRO	-	expression tag	UNP P21218
ZY	65	ARG	-	expression tag	UNP P21218
ZY	66	GLY	-	expression tag	UNP P21218
ZY	67	SER	-	expression tag	UNP P21218
ZZ	49	MET	-	initiating methionine	UNP P21218
ZZ	50	GLY	-	expression tag	UNP P21218
ZZ	51	SER	-	expression tag	UNP P21218
ZZ	52	SER	-	expression tag	UNP P21218
ZZ	53	HIS	-	expression tag	UNP P21218
ZZ	54	HIS	-	expression tag	UNP P21218
ZZ	55	HIS	-	expression tag	UNP P21218
ZZ	56	HIS	-	expression tag	UNP P21218

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Chain	Residue	Modelled	Actual	Comment	Reference
ZZ	57	HIS	-	expression tag	UNP P21218
ZZ	58	HIS	-	expression tag	UNP P21218
ZZ	59	SER	-	expression tag	UNP P21218
ZZ	60	SER	-	expression tag	UNP P21218
ZZ	61	GLY	-	expression tag	UNP P21218
ZZ	62	LEU	-	expression tag	UNP P21218
ZZ	63	VAL	-	expression tag	UNP P21218
ZZ	64	PRO	-	expression tag	UNP P21218
ZZ	65	ARG	-	expression tag	UNP P21218
ZZ	66	GLY	-	expression tag	UNP P21218
ZZ	67	SER	-	expression tag	UNP P21218

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
2	YI	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	YJ	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	YK	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	YL	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	YM	1	Total	C	N	O	P	0
			48	21	7	17	3	

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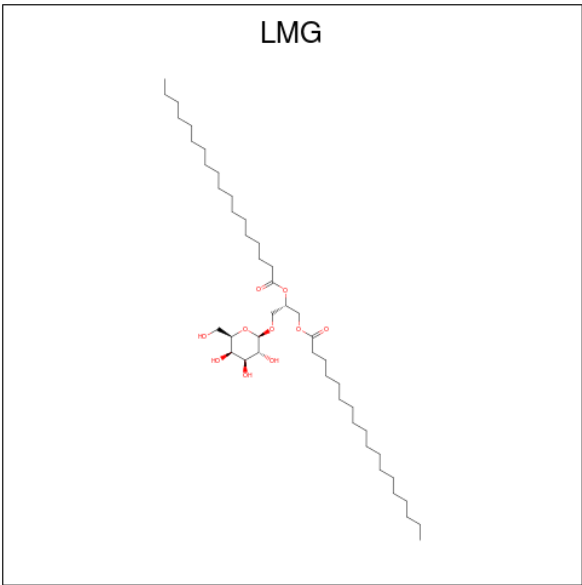
Mol	Chain	Residues	Atoms					AltConf
2	YN	1	Total 48	C 21	N 7	O 17	P 3	0
2	YO	1	Total 48	C 21	N 7	O 17	P 3	0
2	YP	1	Total 48	C 21	N 7	O 17	P 3	0
2	YQ	1	Total 48	C 21	N 7	O 17	P 3	0
2	YR	1	Total 48	C 21	N 7	O 17	P 3	0
2	YS	1	Total 48	C 21	N 7	O 17	P 3	0
2	YT	1	Total 48	C 21	N 7	O 17	P 3	0
2	YU	1	Total 48	C 21	N 7	O 17	P 3	0
2	YV	1	Total 48	C 21	N 7	O 17	P 3	0
2	YW	1	Total 48	C 21	N 7	O 17	P 3	0
2	YX	1	Total 48	C 21	N 7	O 17	P 3	0
2	YY	1	Total 48	C 21	N 7	O 17	P 3	0
2	YZ	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZA	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZB	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZC	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZD	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZE	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZF	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZG	1	Total 48	C 21	N 7	O 17	P 3	0
2	ZH	1	Total 48	C 21	N 7	O 17	P 3	0

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Mol	Chain	Residues	Atoms					AltConf
2	ZI	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZJ	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZK	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZL	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZM	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZN	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZO	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZP	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZQ	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZR	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZS	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZT	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZU	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZV	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZW	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZX	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZY	1	Total	C	N	O	P	0
			48	21	7	17	3	
2	ZZ	1	Total	C	N	O	P	0
			48	21	7	17	3	

- Molecule 3 is 1,2-DISTEAROYL-MONOGALACTOSYL-DIGLYCERIDE (CCD ID: LMG) (formula: C₄₅H₈₆O₁₀) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
3	YI	1	Total	C	O	0
			23	13	10	
3	YJ	1	Total	C	O	0
			23	13	10	
3	YK	1	Total	C	O	0
			23	13	10	
3	YL	1	Total	C	O	0
			23	13	10	
3	YM	1	Total	C	O	0
			23	13	10	
3	YN	1	Total	C	O	0
			23	13	10	
3	YO	1	Total	C	O	0
			23	13	10	
3	YP	1	Total	C	O	0
			23	13	10	
3	YQ	1	Total	C	O	0
			23	13	10	
3	YR	1	Total	C	O	0
			23	13	10	
3	YS	1	Total	C	O	0
			23	13	10	
3	YT	1	Total	C	O	0
			23	13	10	
3	YU	1	Total	C	O	0
			23	13	10	
3	YV	1	Total	C	O	0
			23	13	10	

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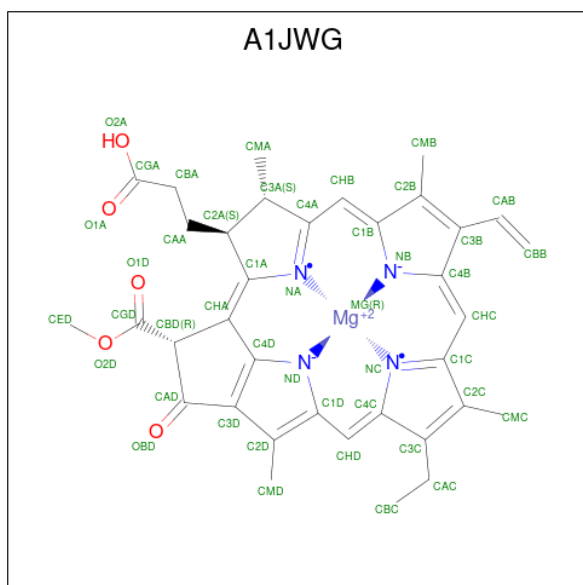
Mol	Chain	Residues	Atoms			AltConf
3	YW	1	Total 23	C 13	O 10	0
3	YX	1	Total 23	C 13	O 10	0
3	YY	1	Total 23	C 13	O 10	0
3	YZ	1	Total 23	C 13	O 10	0
3	ZA	1	Total 23	C 13	O 10	0
3	ZB	1	Total 23	C 13	O 10	0
3	ZC	1	Total 23	C 13	O 10	0
3	ZD	1	Total 23	C 13	O 10	0
3	ZE	1	Total 23	C 13	O 10	0
3	ZF	1	Total 23	C 13	O 10	0
3	ZG	1	Total 23	C 13	O 10	0
3	ZH	1	Total 23	C 13	O 10	0
3	ZI	1	Total 23	C 13	O 10	0
3	ZJ	1	Total 23	C 13	O 10	0
3	ZK	1	Total 23	C 13	O 10	0
3	ZL	1	Total 23	C 13	O 10	0
3	ZM	1	Total 23	C 13	O 10	0
3	ZN	1	Total 23	C 13	O 10	0
3	ZO	1	Total 23	C 13	O 10	0
3	ZP	1	Total 23	C 13	O 10	0
3	ZQ	1	Total 23	C 13	O 10	0

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Mol	Chain	Residues	Atoms			AltConf
3	ZR	1	Total	C	O	0
			23	13	10	
3	ZS	1	Total	C	O	0
			23	13	10	
3	ZT	1	Total	C	O	0
			23	13	10	
3	ZU	1	Total	C	O	0
			23	13	10	
3	ZV	1	Total	C	O	0
			23	13	10	
3	ZW	1	Total	C	O	0
			23	13	10	
3	ZX	1	Total	C	O	0
			23	13	10	
3	ZY	1	Total	C	O	0
			23	13	10	
3	ZZ	1	Total	C	O	0
			23	13	10	

- Molecule 4 is Chlorophyllide a (CCD ID: A1JWG) (formula: $C_{35}H_{34}MgN_4O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
4	YI	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
4	YJ	1	Total	C	Mg	N	O	0
			45	35	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
4	YK	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YL	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YM	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YN	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YO	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YP	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YQ	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YR	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YS	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YT	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YU	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YV	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YW	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YX	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YY	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	YZ	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZA	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZB	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZC	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZD	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZE	1	Total 45	C 35	Mg 1	N 4	O 5	0

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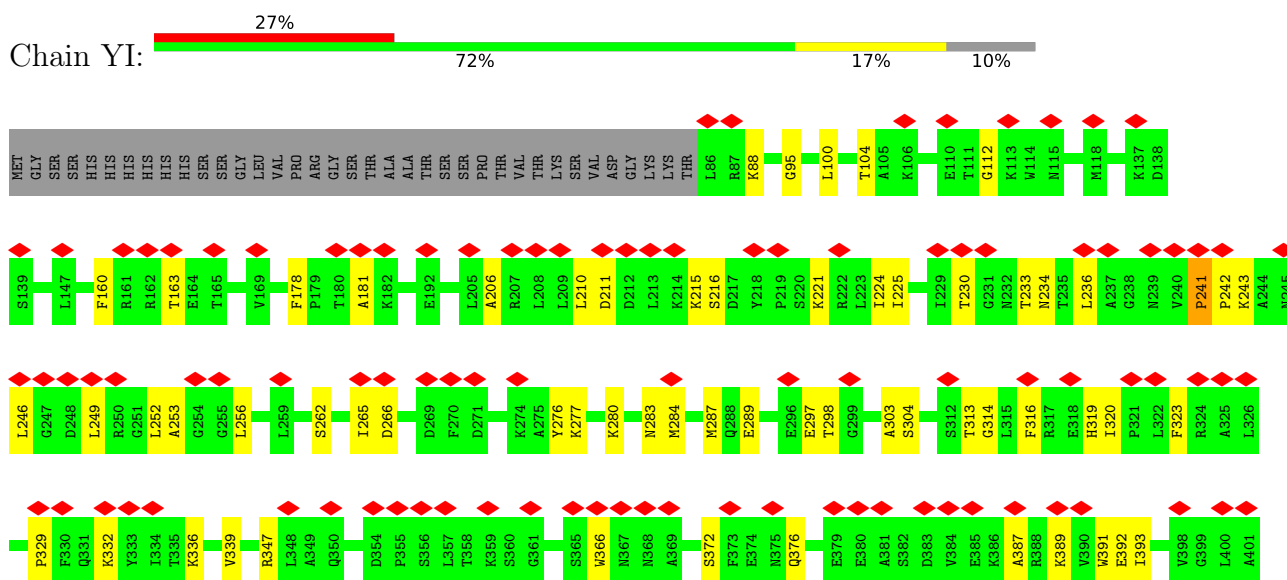
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Mol	Chain	Residues	Atoms					AltConf
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4	ZG	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZH	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZI	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZJ	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZK	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZL	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZM	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZN	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZO	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZP	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZQ	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZR	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZS	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZT	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZU	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZV	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZW	1	Total 45	C 35	Mg 1	N 4	O 5	0
4	ZX	1	Total 45	C 35	Mg 1	N 4	O 5	0
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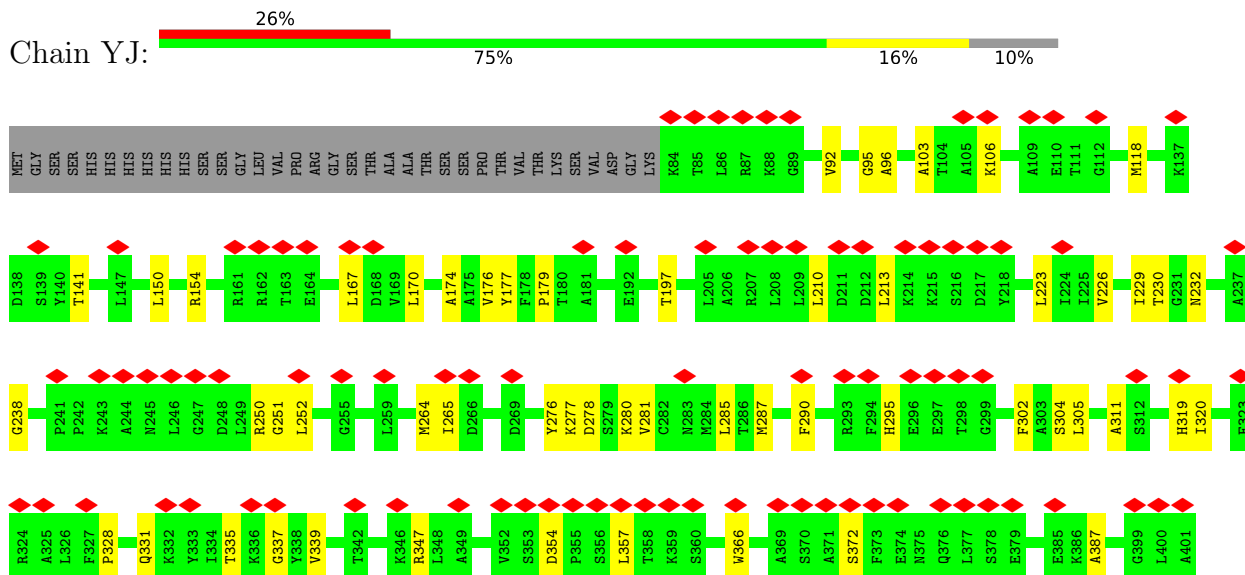
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

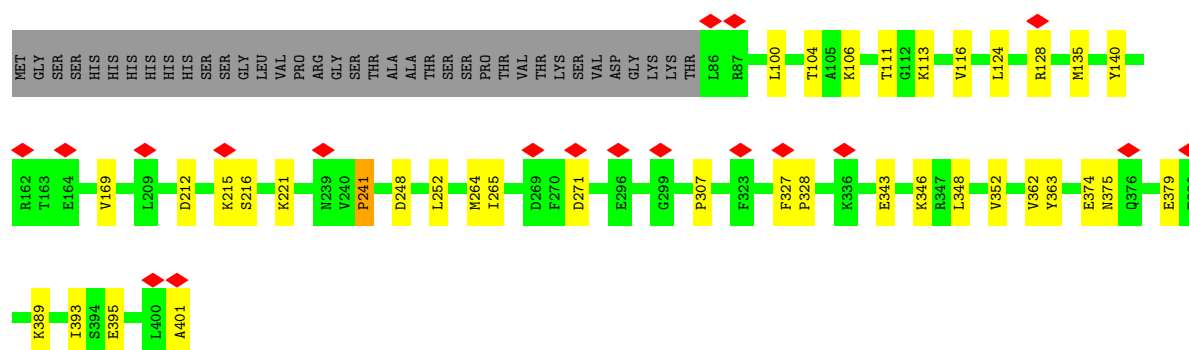
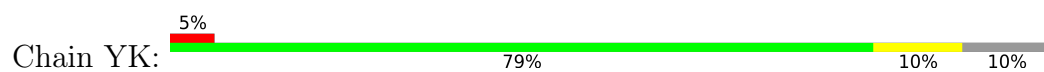
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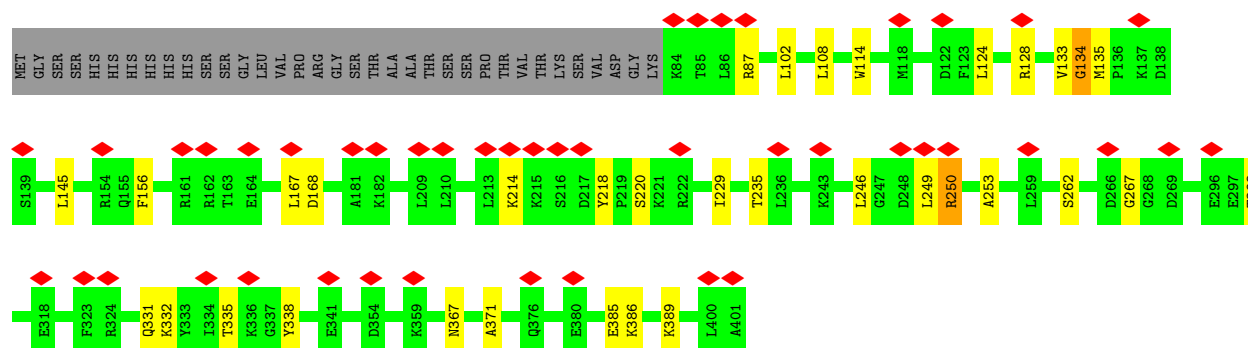
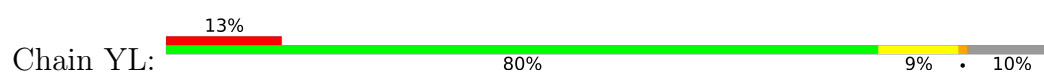
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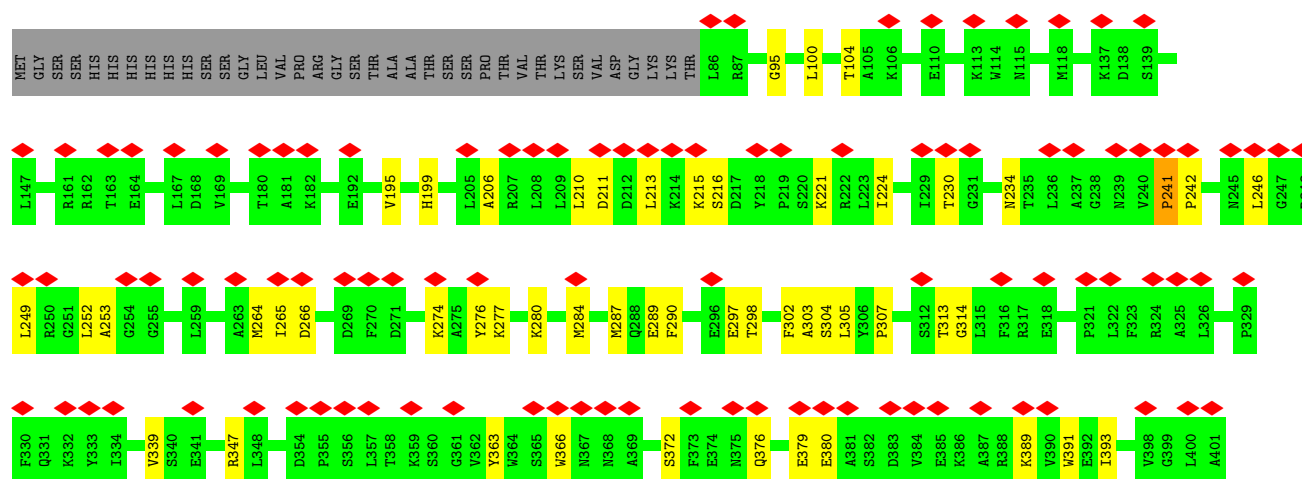
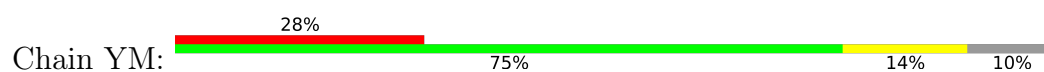
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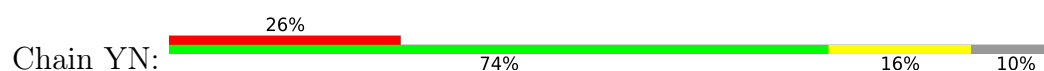
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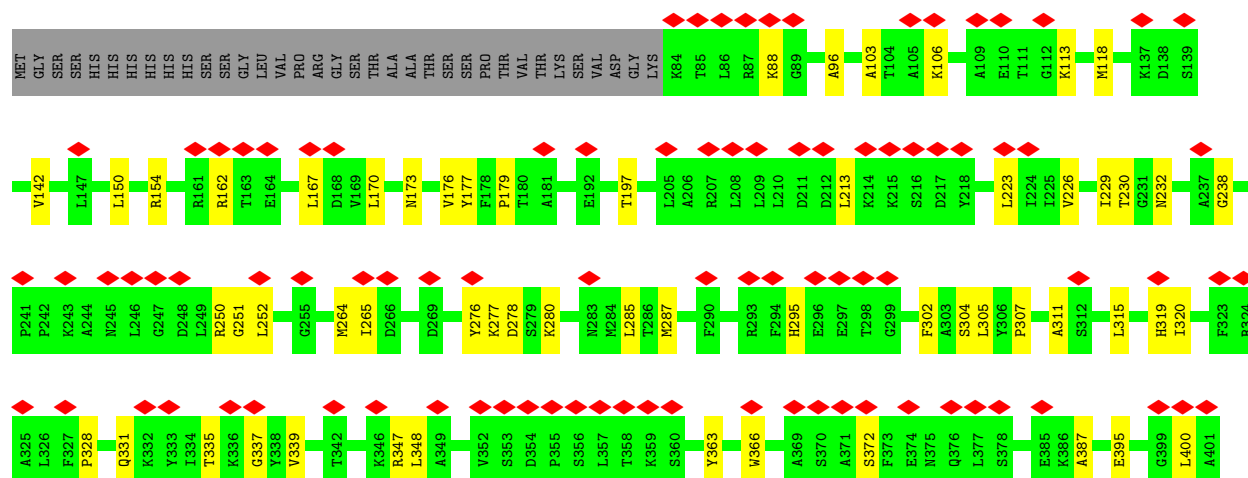


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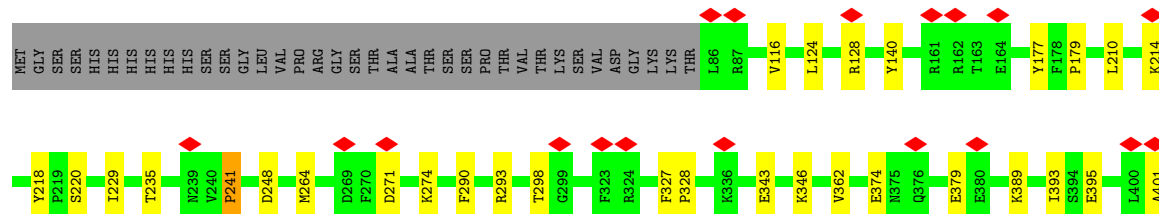
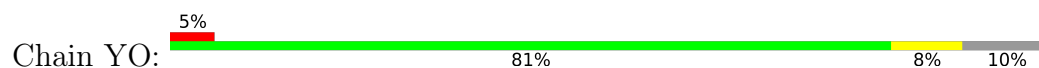


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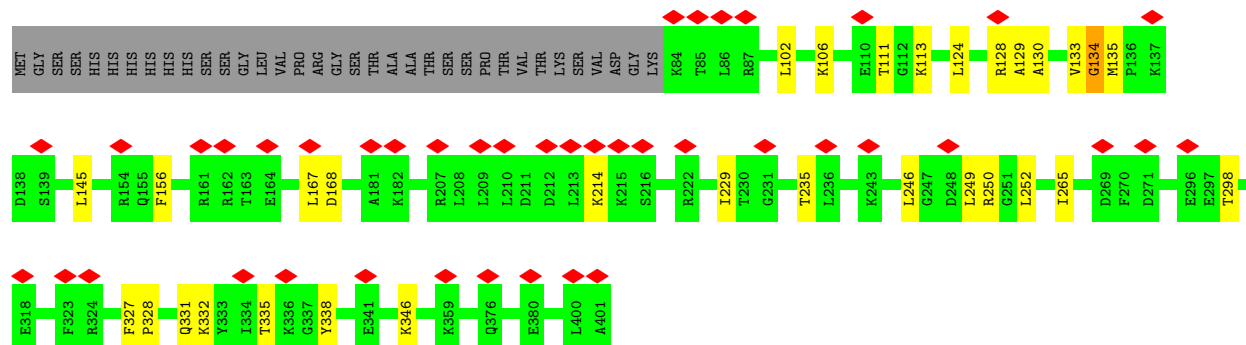
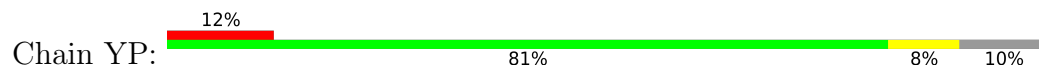




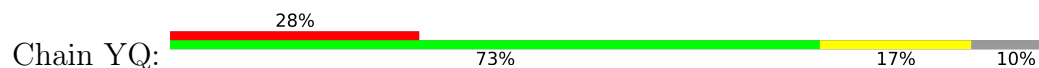
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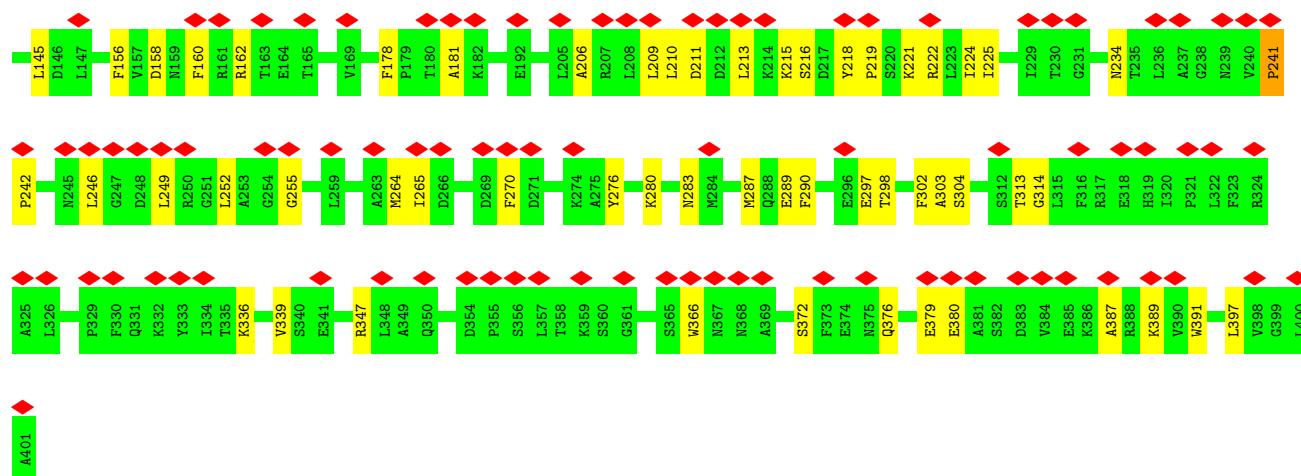


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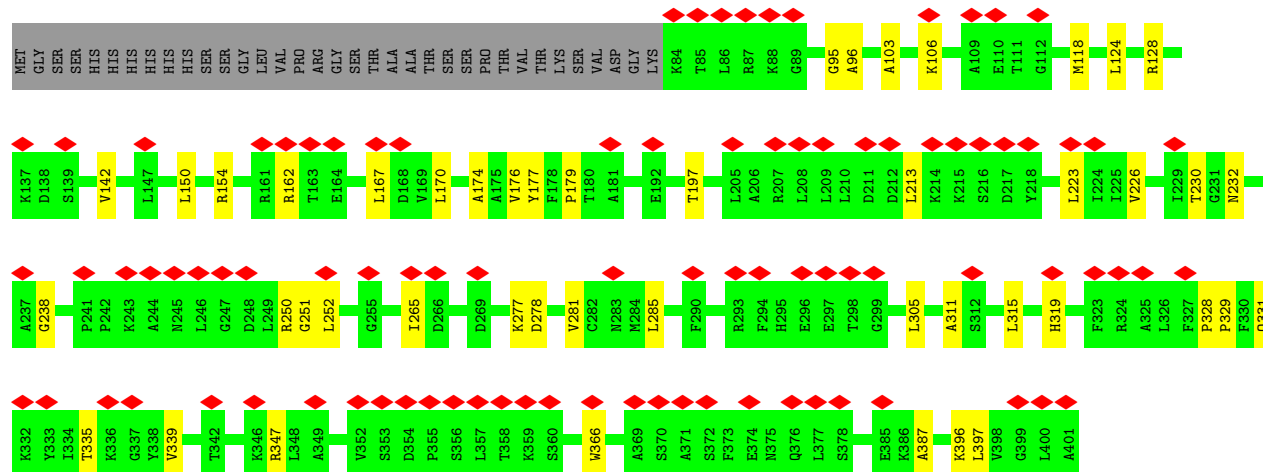
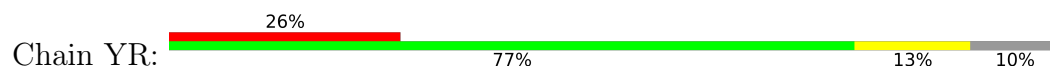


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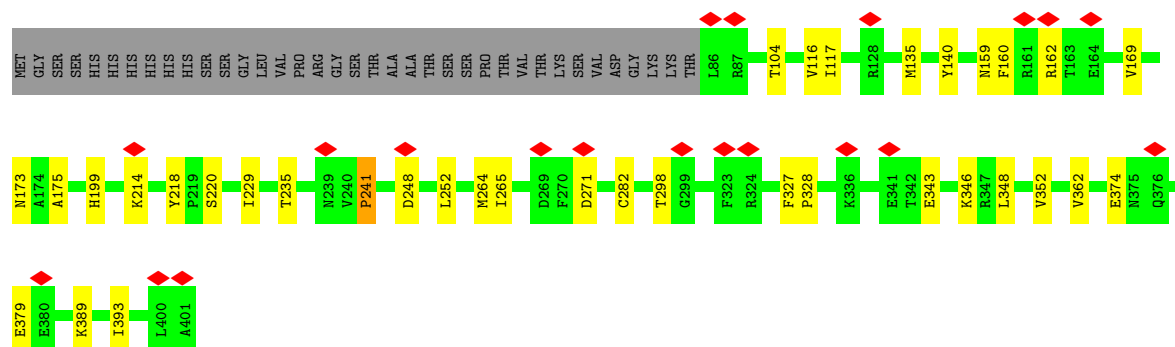
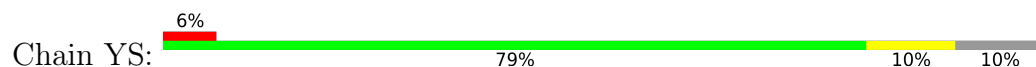




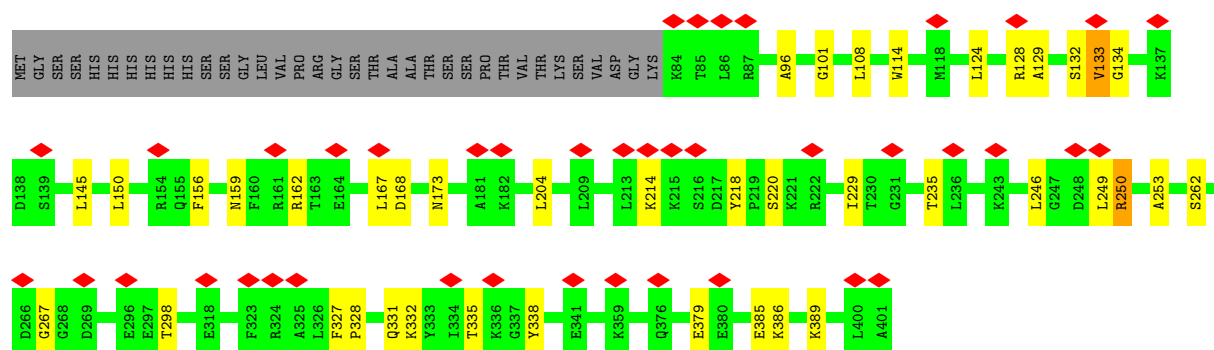
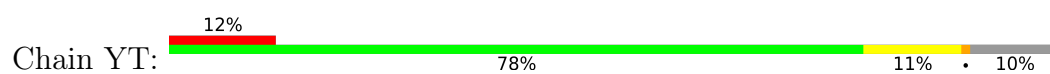
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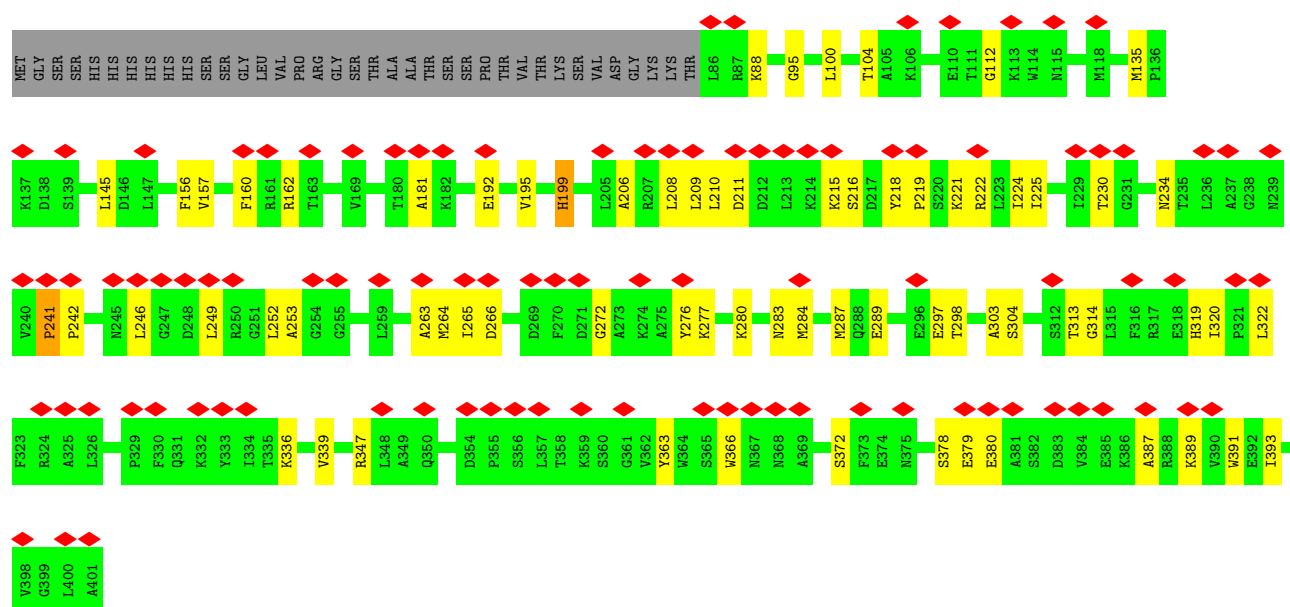
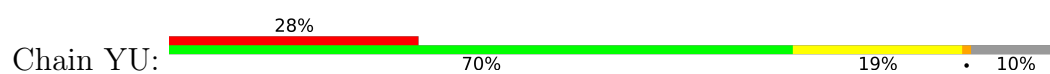
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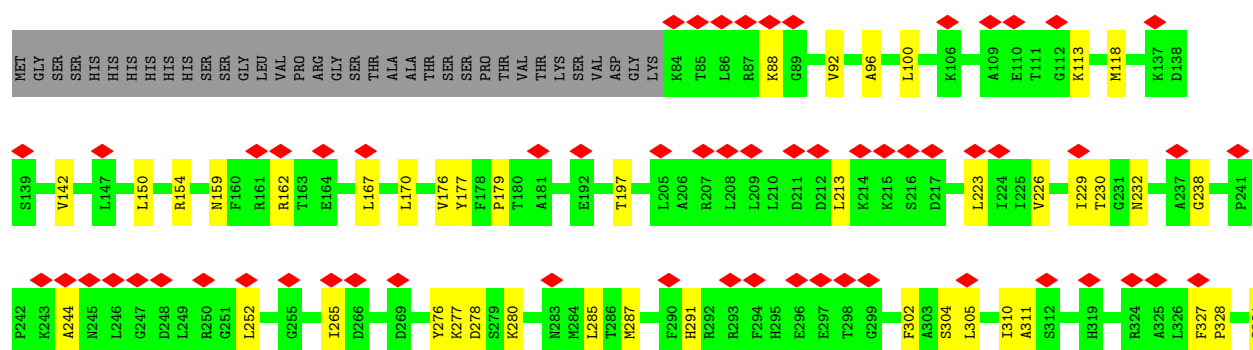
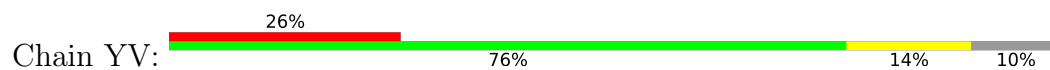
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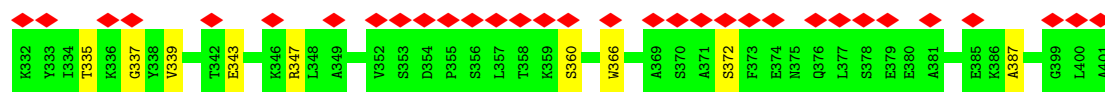


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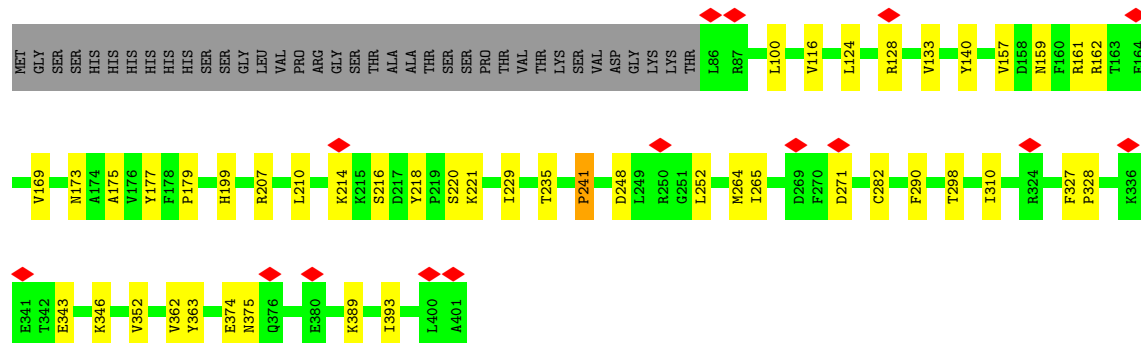
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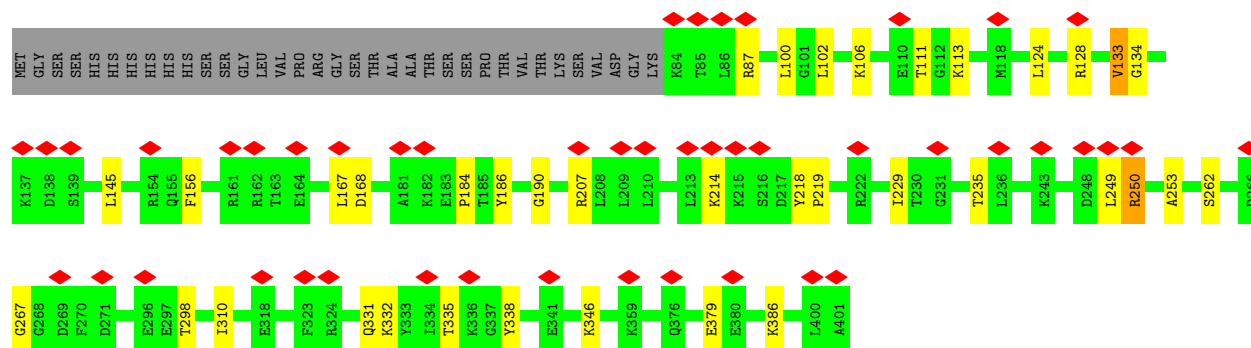
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

Chain YW: 76% 13% 10%



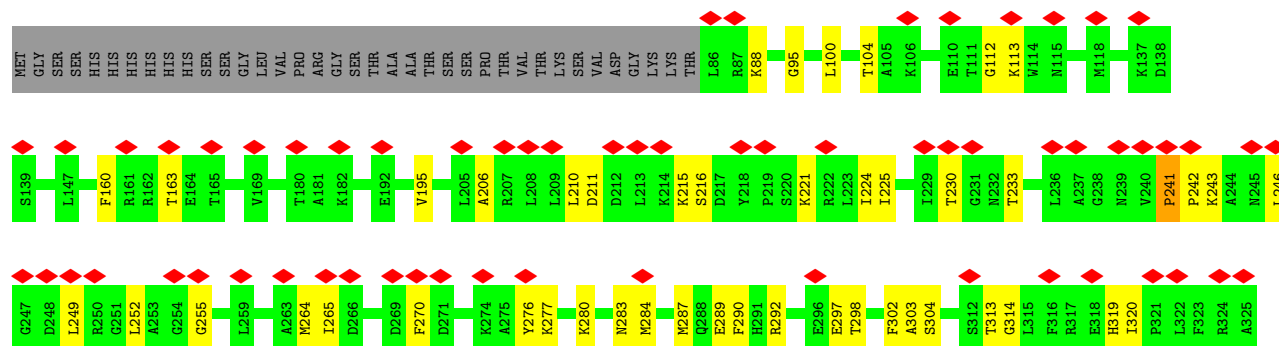
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

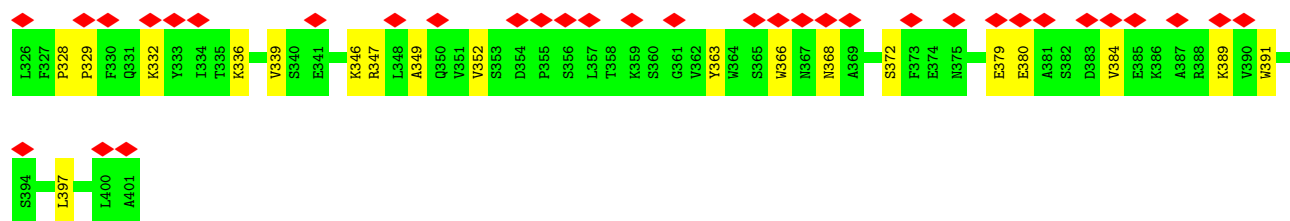
Chain YX: 13% 80% 10% • 10%



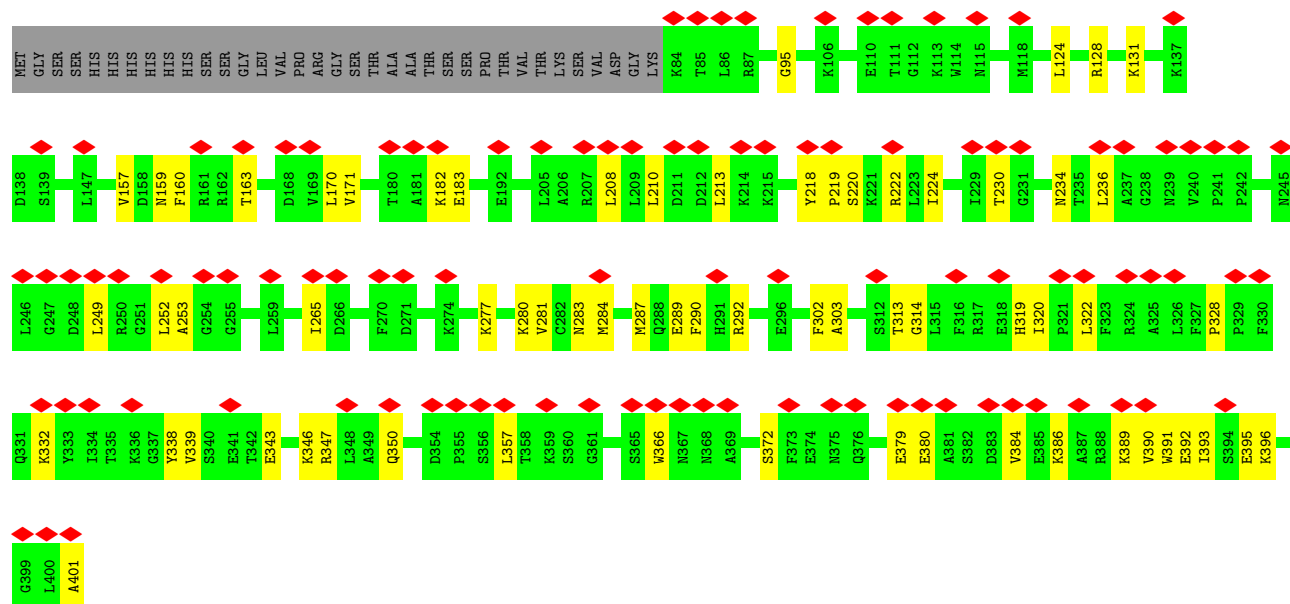
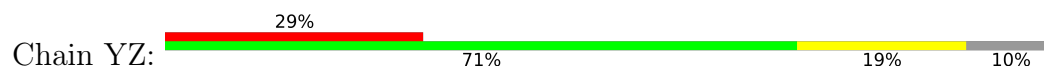
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

Chain YY: 27% 71% 18% 10%

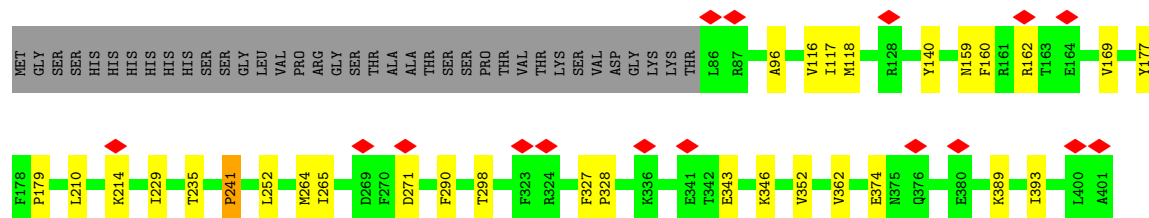
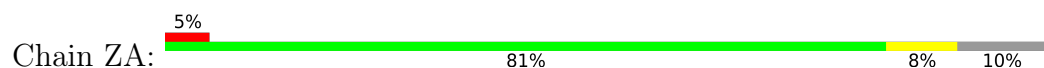




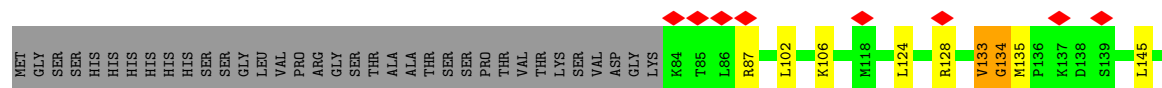
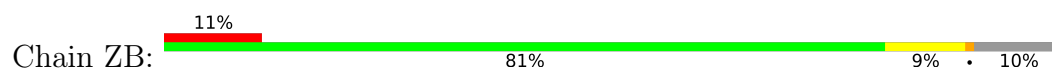
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

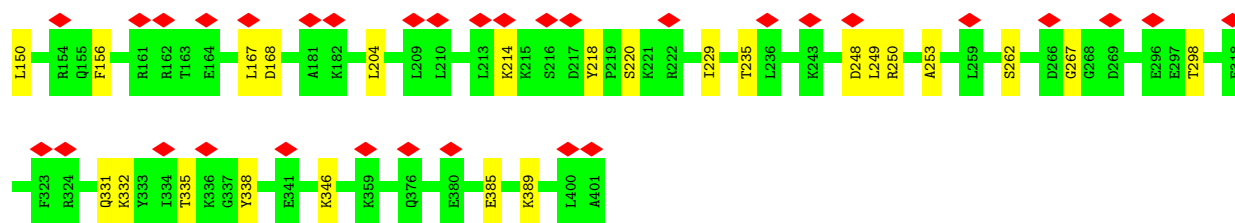


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

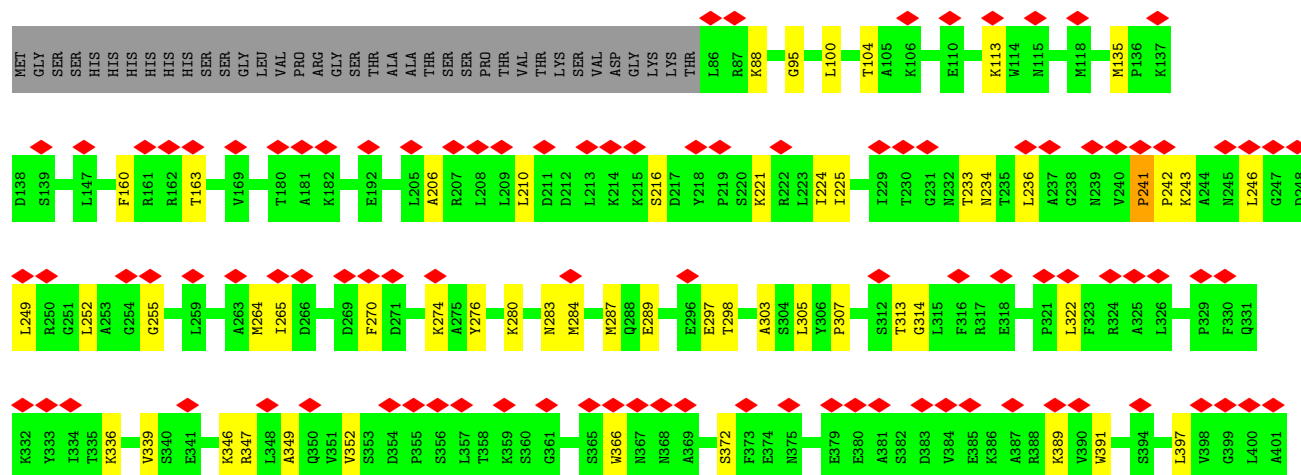
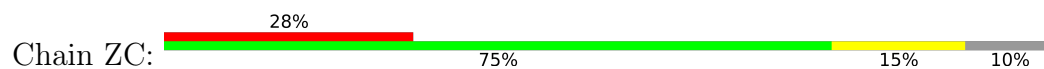


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

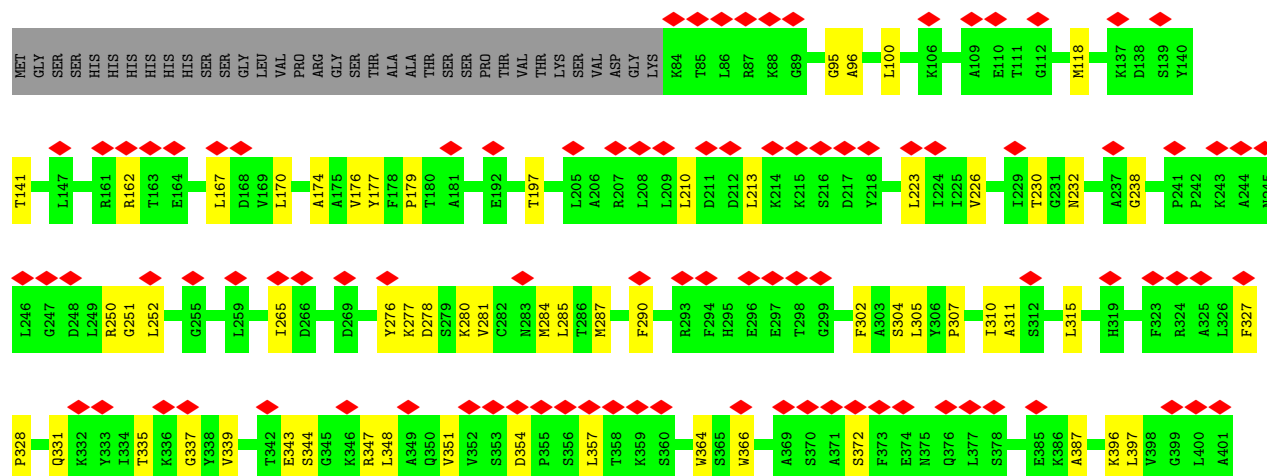
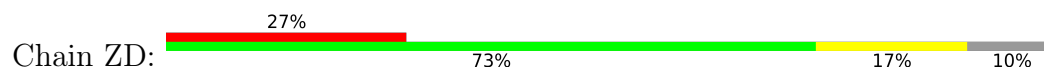




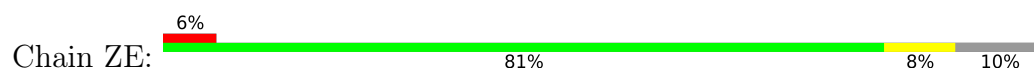
• Molecule 1: Protochlorophyllide reductase B, chloroplastic

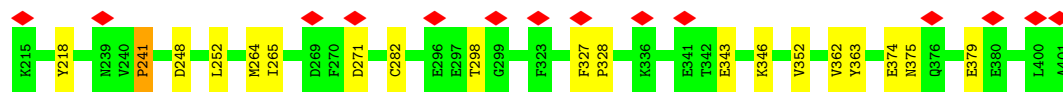


• Molecule 1: Protochlorophyllide reductase B, chloroplastic

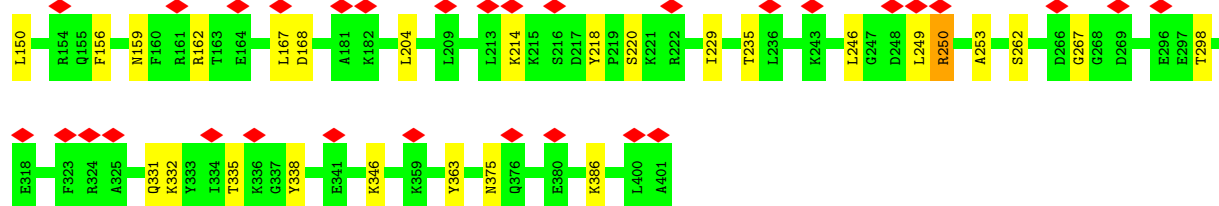


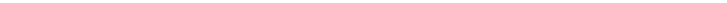
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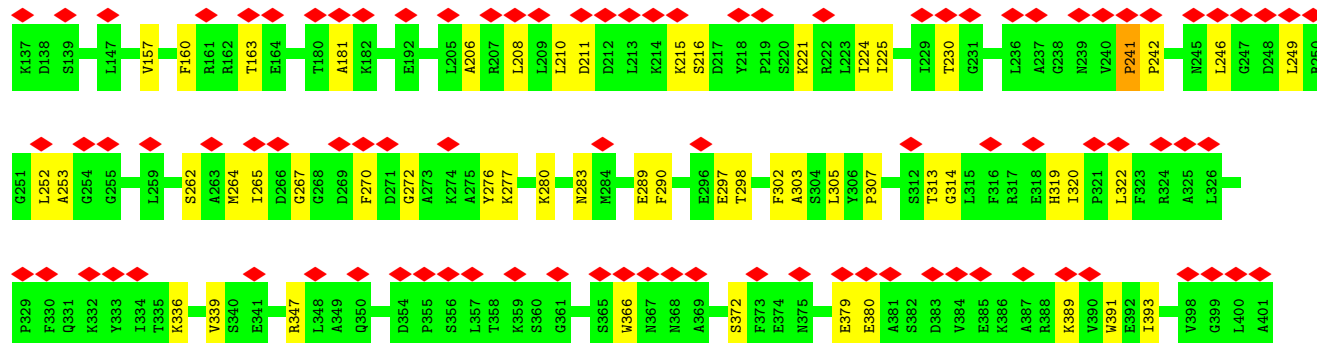


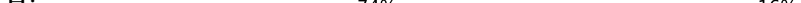


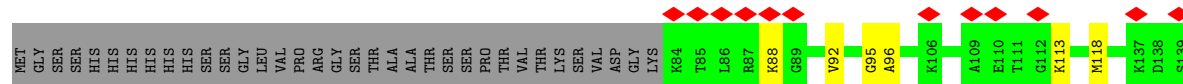
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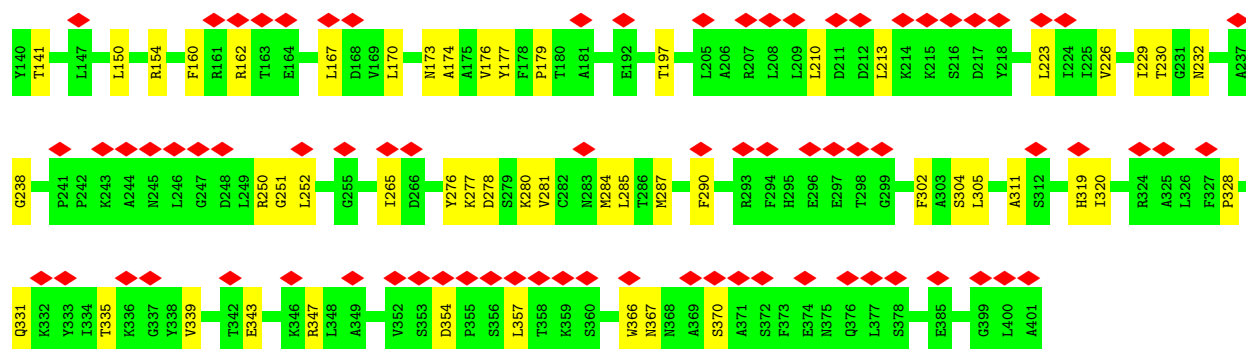


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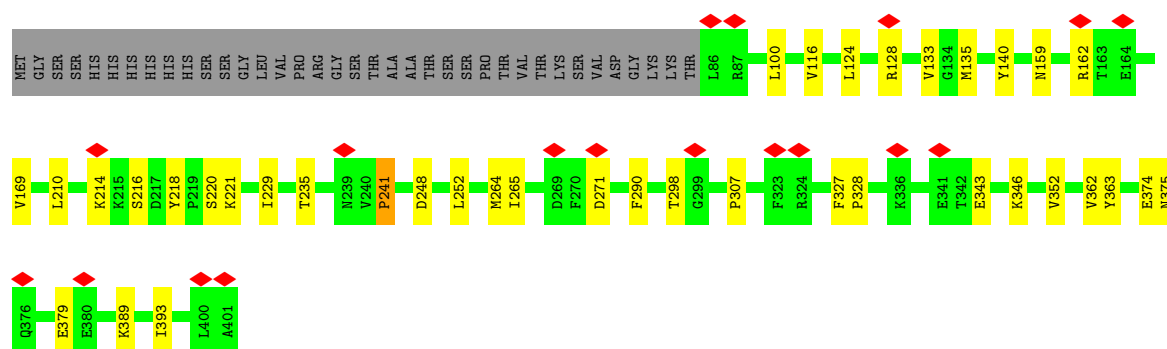
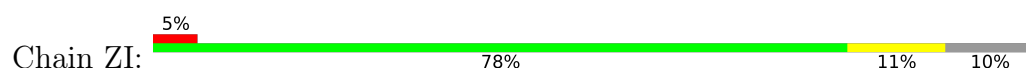


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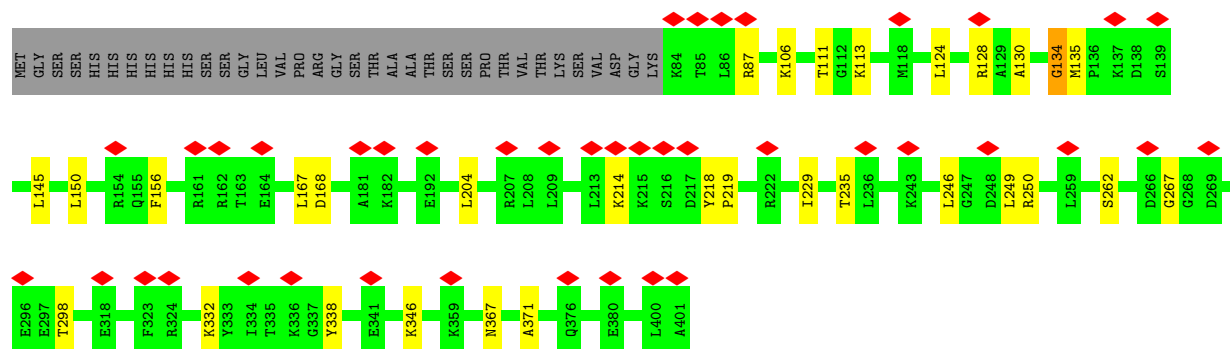
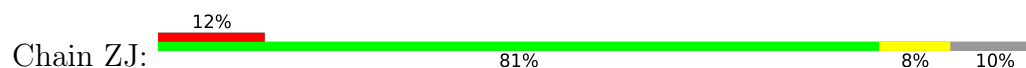




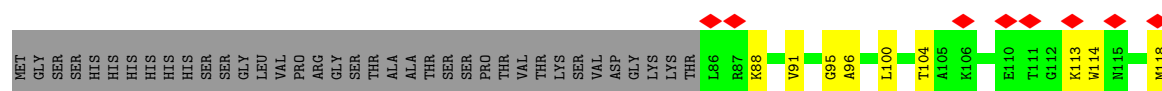
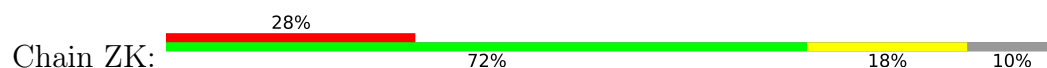
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

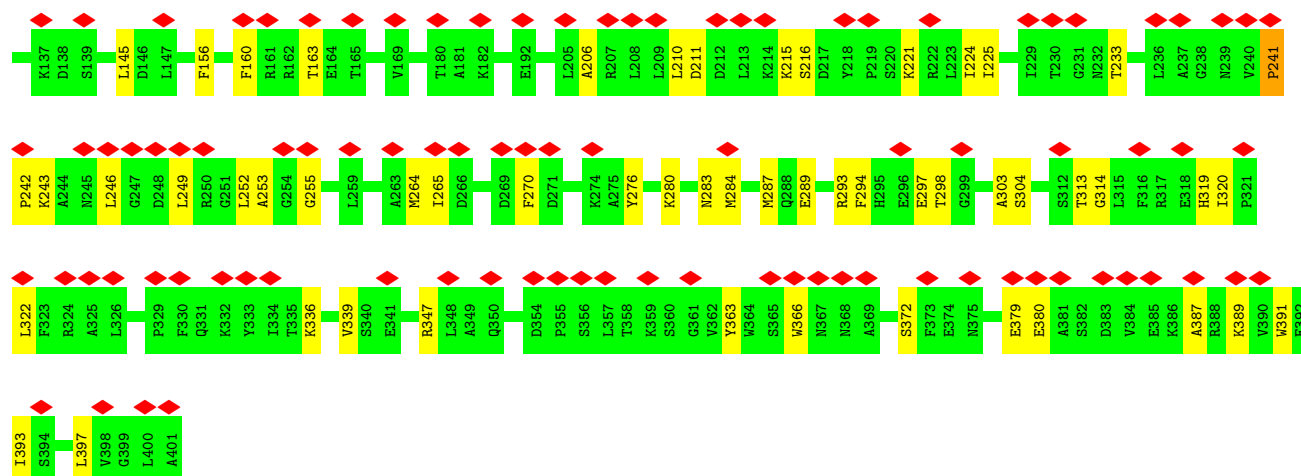


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

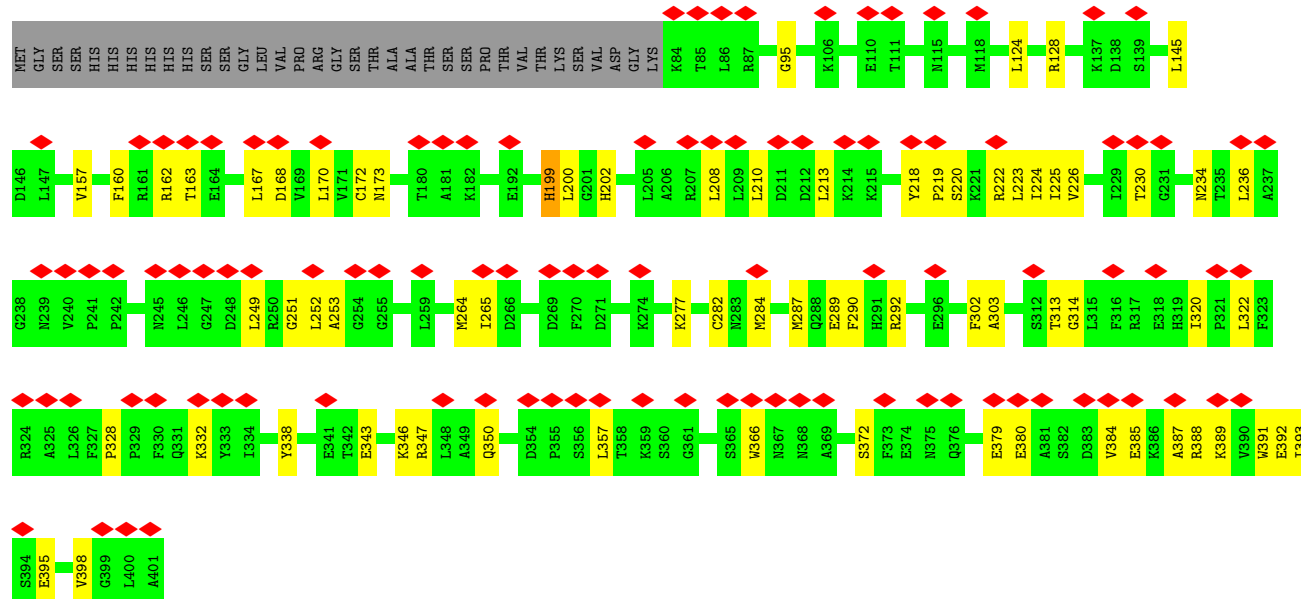


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

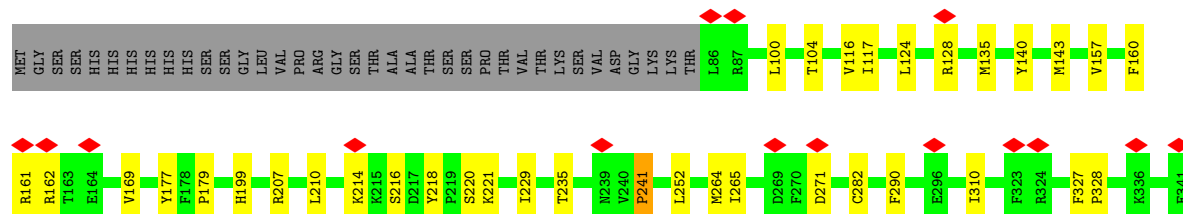
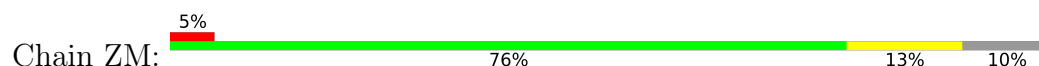


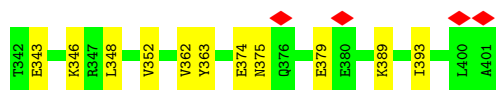


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

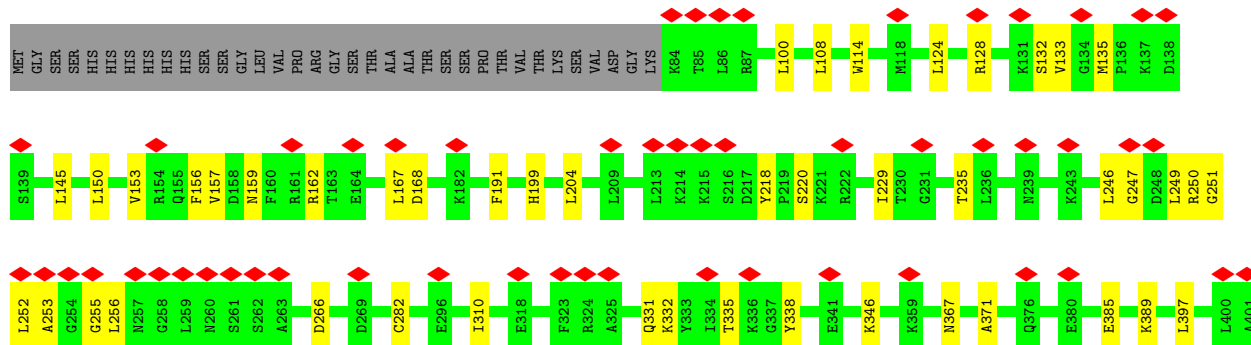
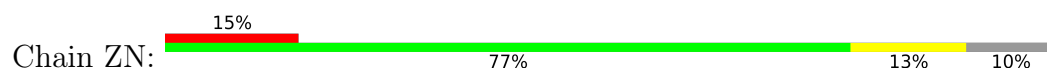


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

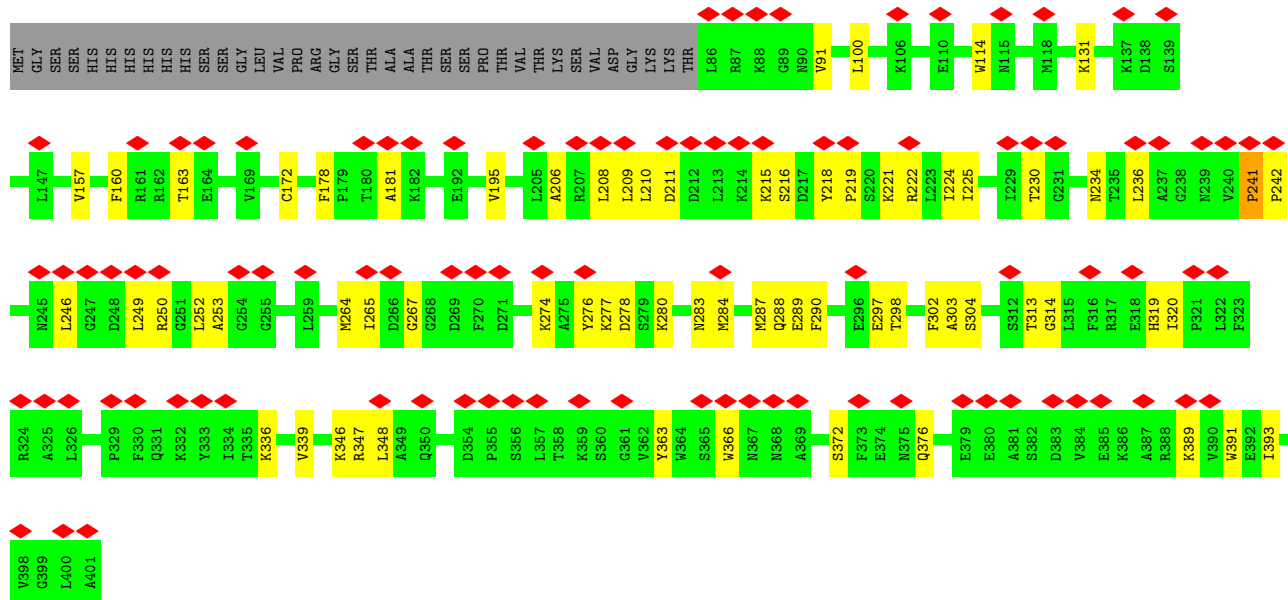
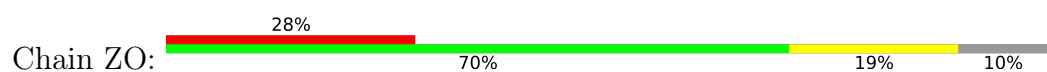




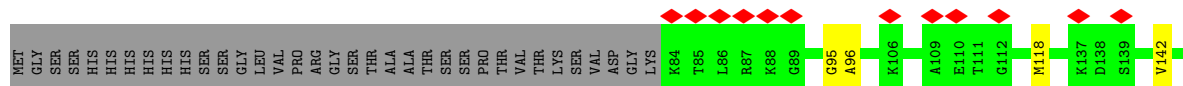
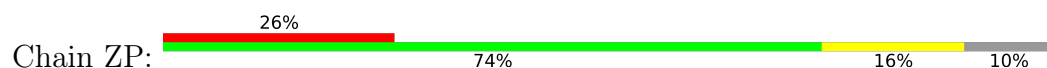
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

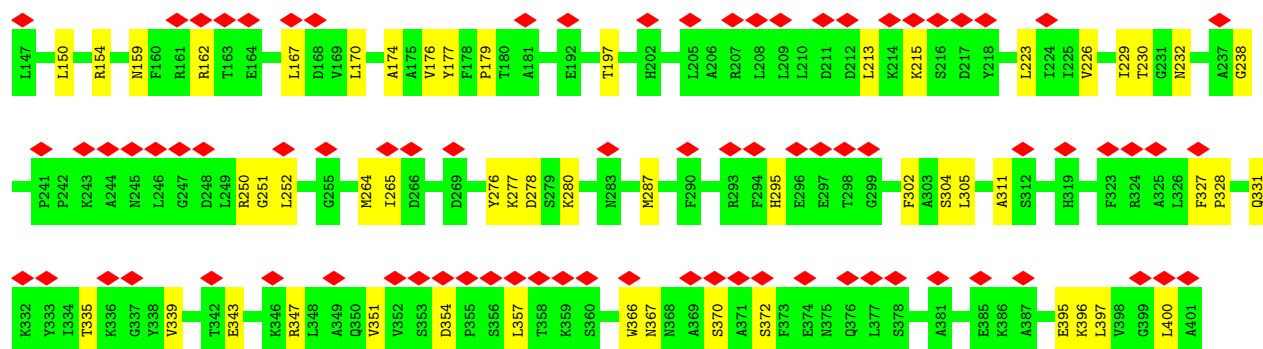


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

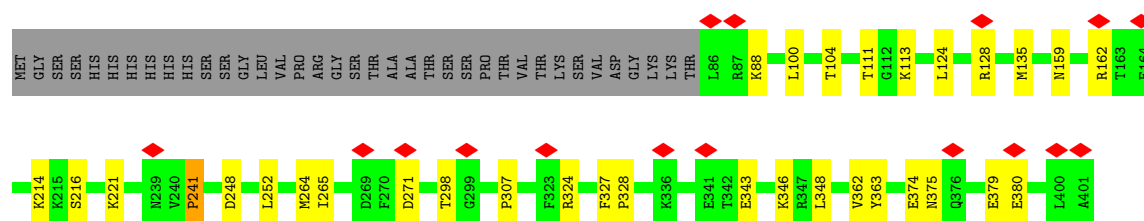
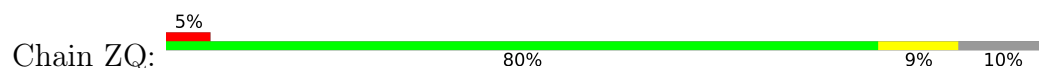


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

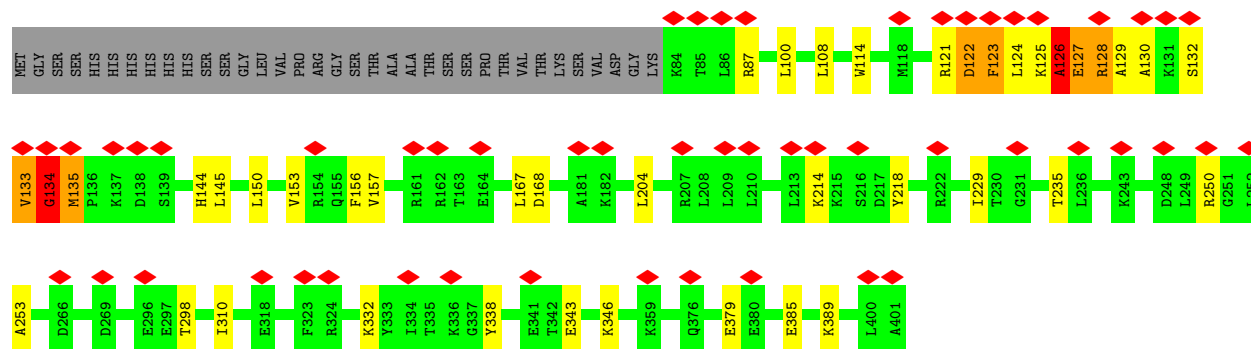
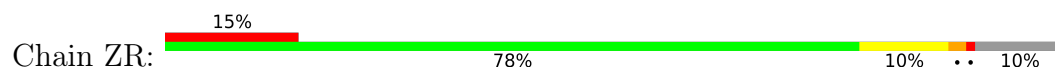




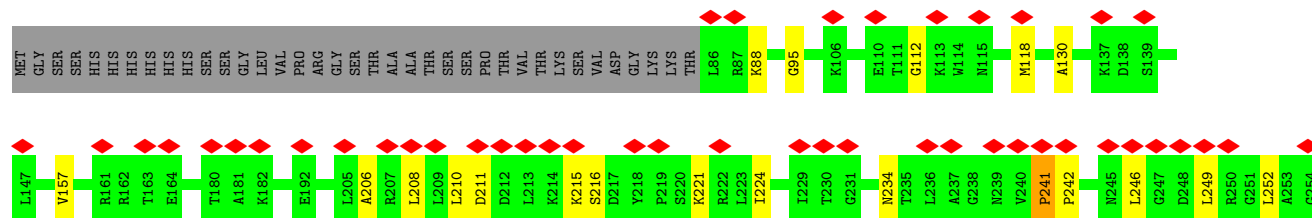
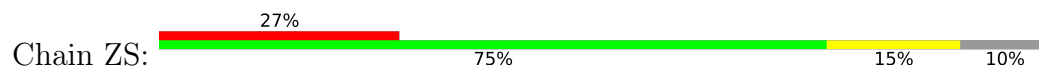
- Molecule 1: Protochlorophyllide reductase B, chloroplastic

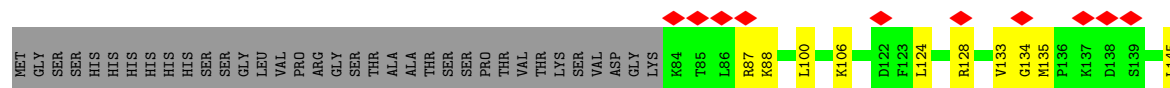


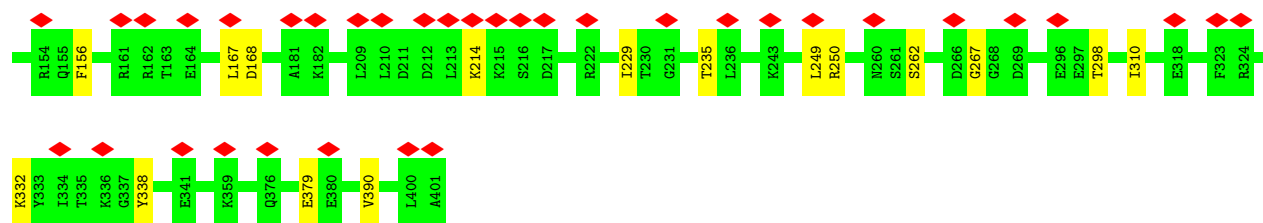
- Molecule 1: Protochlorophyllide reductase B, chloroplastic



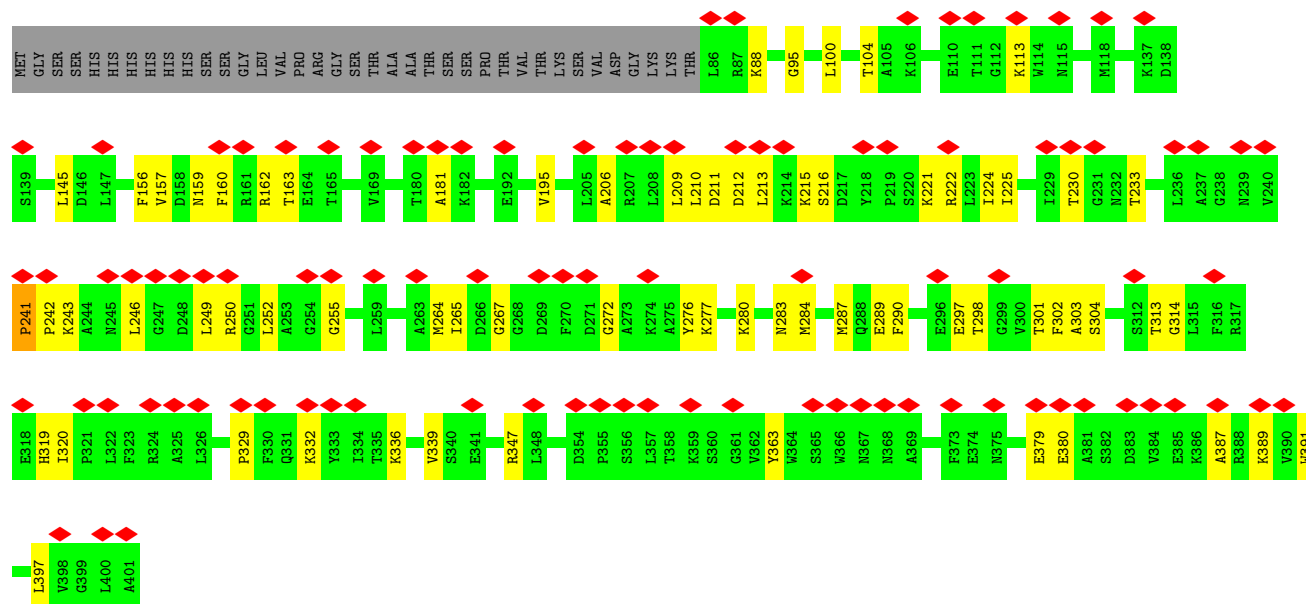
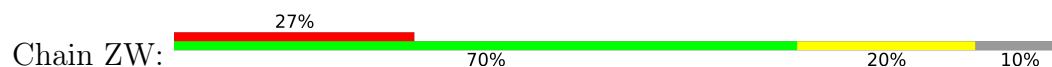
- Molecule 1: Protochlorophyllide reductase B, chloroplastic



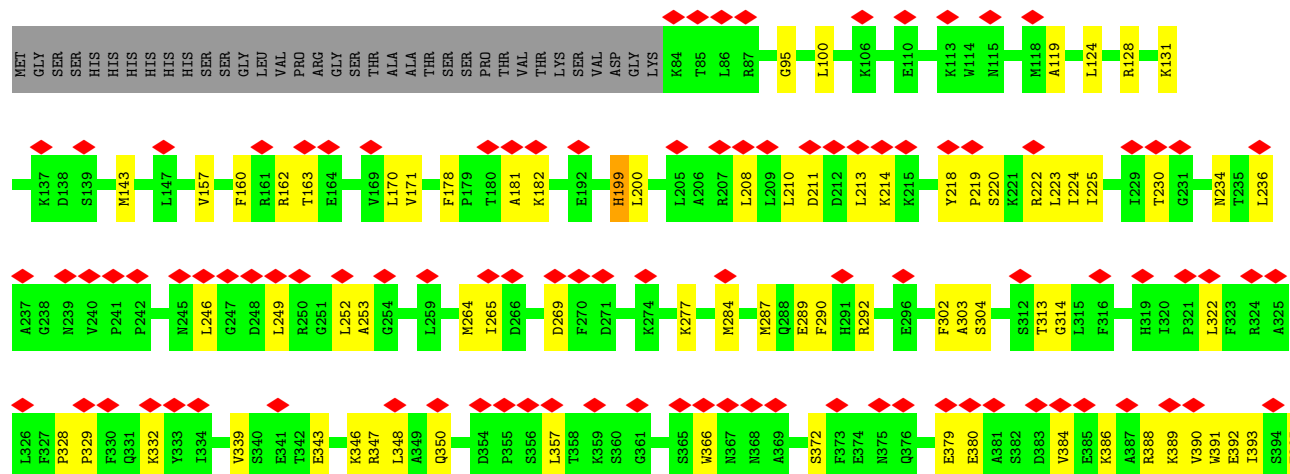


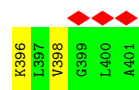


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

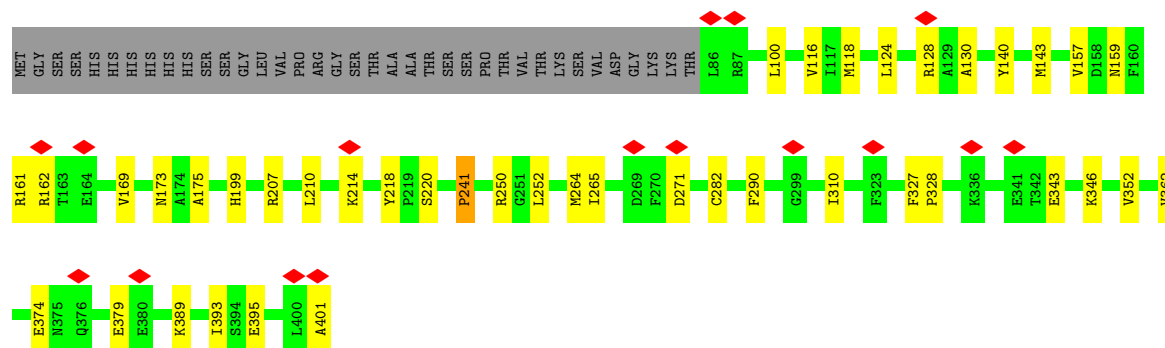
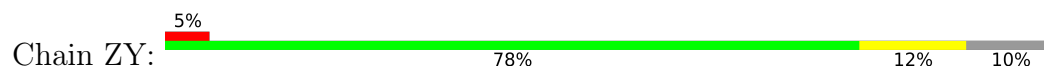


- Molecule 1: Protochlorophyllide reductase B, chloroplastic

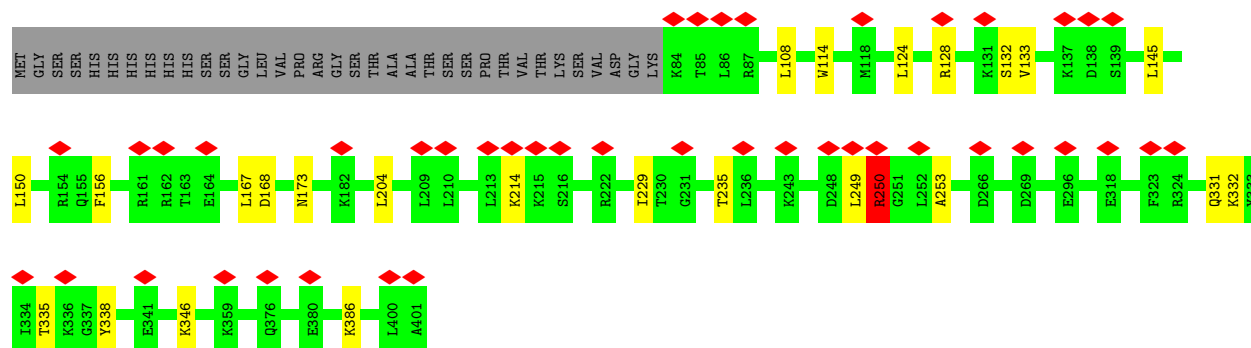
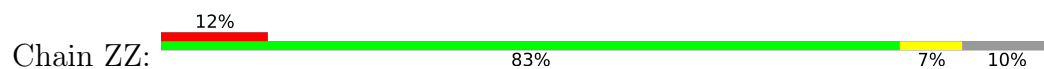




- Molecule 1: Protochlorophyllide reductase B, chloroplastic



- Molecule 1: Protochlorophyllide reductase B, chloroplastic



4 Experimental information

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=-19.328°, rise=153.203 Å, axial sym=D11	Depositor
Number of segments used	25485	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{Å}^2$)	40	Depositor
Minimum defocus (nm)	900	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.334	Depositor
Minimum map value	-0.135	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.017	Depositor
Recommended contour level	0.104	Depositor
Map size (Å)	550.4, 550.4, 550.4	wwPDB
Map dimensions	640, 640, 640	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.86, 0.86, 0.86	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NDP, LMG, A1JWG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	YI	0.20	0/2470	0.44	0/3342
1	YJ	0.19	0/2486	0.39	0/3363
1	YK	0.22	0/2470	0.36	0/3342
1	YL	0.21	0/2486	0.43	2/3363 (0.1%)
1	YM	0.20	0/2470	0.44	0/3342
1	YN	0.19	0/2486	0.39	0/3363
1	YO	0.22	0/2470	0.36	0/3342
1	YP	0.23	1/2486 (0.0%)	0.45	2/3363 (0.1%)
1	YQ	0.20	0/2470	0.44	0/3342
1	YR	0.19	0/2486	0.39	0/3363
1	YS	0.21	0/2470	0.37	0/3342
1	YT	0.21	0/2486	0.40	0/3363
1	YU	0.21	0/2470	0.45	0/3342
1	YV	0.19	0/2486	0.40	0/3363
1	YW	0.21	0/2470	0.36	0/3342
1	YX	0.22	0/2486	0.44	4/3363 (0.1%)
1	YY	0.20	0/2470	0.44	0/3342
1	YZ	0.21	0/2486	0.45	0/3363
1	ZA	0.21	0/2470	0.36	0/3342
1	ZB	0.24	1/2486 (0.0%)	0.48	2/3363 (0.1%)
1	ZC	0.20	0/2470	0.44	0/3342
1	ZD	0.19	0/2486	0.40	0/3363
1	ZE	0.21	0/2470	0.36	0/3342
1	ZF	0.24	1/2486 (0.0%)	0.43	2/3363 (0.1%)
1	ZG	0.20	0/2470	0.44	0/3342
1	ZH	0.19	0/2486	0.40	0/3363
1	ZI	0.21	0/2470	0.37	0/3342
1	ZJ	0.26	1/2486 (0.0%)	0.45	2/3363 (0.1%)
1	ZK	0.20	0/2470	0.44	0/3342
1	ZL	0.21	0/2486	0.45	0/3363
1	ZM	0.22	0/2470	0.37	0/3342
1	ZN	0.21	0/2486	0.44	0/3363

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	ZO	0.21	0/2470	0.44	0/3342
1	ZP	0.19	0/2486	0.40	0/3363
1	ZQ	0.21	0/2470	0.35	0/3342
1	ZR	0.26	1/2486 (0.0%)	0.49	2/3363 (0.1%)
1	ZS	0.20	0/2470	0.44	0/3342
1	ZT	0.19	0/2486	0.41	0/3363
1	ZU	0.21	0/2470	0.37	0/3342
1	ZV	0.23	1/2486 (0.0%)	0.51	2/3363 (0.1%)
1	ZW	0.20	0/2470	0.44	0/3342
1	ZX	0.21	0/2486	0.45	0/3363
1	ZY	0.22	0/2470	0.37	0/3342
1	ZZ	0.20	0/2486	0.41	0/3363
All	All	0.21	6/109032 (0.0%)	0.42	18/147510 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	YI	0	1
1	YK	0	1
1	YL	0	1
1	YM	0	1
1	YO	0	1
1	YQ	0	1
1	YS	0	1
1	YT	0	2
1	YU	0	1
1	YW	0	1
1	YX	0	1
1	YY	0	1
1	ZA	0	1
1	ZC	0	1
1	ZE	0	1
1	ZF	0	1
1	ZG	0	1
1	ZI	0	1
1	ZK	0	1
1	ZM	0	1
1	ZN	0	3
1	ZO	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	ZQ	0	1
1	ZR	0	5
1	ZS	0	1
1	ZU	0	1
1	ZW	0	1
1	ZY	0	1
1	ZZ	0	3
All	All	0	38

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	ZJ	134	GLY	C-N	8.01	1.45	1.33
1	ZB	134	GLY	C-N	6.13	1.42	1.33
1	ZV	134	GLY	C-N	6.13	1.42	1.33
1	ZF	134	GLY	C-N	-5.96	1.25	1.33
1	YP	134	GLY	C-N	5.85	1.42	1.33
1	ZR	134	GLY	C-N	5.05	1.44	1.33

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	ZV	134	GLY	CA-C-N	-13.22	96.22	122.40
1	ZV	134	GLY	C-N-CA	-13.22	96.22	122.40
1	ZB	134	GLY	CA-C-N	-9.80	98.65	122.36
1	ZB	134	GLY	C-N-CA	-9.80	98.65	122.36
1	ZJ	134	GLY	CA-C-N	-9.08	100.40	122.36
1	ZJ	134	GLY	C-N-CA	-9.08	100.40	122.36
1	YP	134	GLY	CA-C-N	-8.88	100.87	122.36
1	YP	134	GLY	C-N-CA	-8.88	100.87	122.36
1	YL	134	GLY	CA-C-N	-6.76	106.00	122.36
1	YL	134	GLY	C-N-CA	-6.76	106.00	122.36
1	ZF	134	GLY	CA-C-N	-6.32	110.18	122.70
1	ZF	134	GLY	C-N-CA	-6.32	110.18	122.70
1	YX	134	GLY	CA-C-N	-6.20	107.80	122.31
1	YX	134	GLY	C-N-CA	-6.20	107.80	122.31
1	ZR	134	GLY	CA-C-N	-6.04	107.05	121.80
1	ZR	134	GLY	C-N-CA	-6.04	107.05	121.80
1	YX	133	VAL	CA-C-N	-5.79	110.07	121.41
1	YX	133	VAL	C-N-CA	-5.79	110.07	121.41

There are no chirality outliers.

All (38) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	YI	241	PRO	Peptide
1	YK	241	PRO	Peptide
1	YL	250	ARG	Peptide
1	YM	241	PRO	Peptide
1	YO	241	PRO	Peptide
1	YQ	241	PRO	Peptide
1	YS	241	PRO	Peptide
1	YT	133	VAL	Peptide
1	YT	250	ARG	Sidechain
1	YU	241	PRO	Peptide
1	YW	241	PRO	Peptide
1	YX	250	ARG	Peptide
1	YY	241	PRO	Peptide
1	ZA	241	PRO	Peptide
1	ZC	241	PRO	Peptide
1	ZE	241	PRO	Peptide
1	ZF	250	ARG	Peptide
1	ZG	241	PRO	Peptide
1	ZI	241	PRO	Peptide
1	ZK	241	PRO	Peptide
1	ZM	241	PRO	Peptide
1	ZN	132	SER	Peptide
1	ZN	251	GLY	Peptide
1	ZN	252	LEU	Peptide
1	ZO	241	PRO	Peptide
1	ZQ	241	PRO	Peptide
1	ZR	122	ASP	Peptide
1	ZR	126	ALA	Peptide
1	ZR	128	ARG	Peptide
1	ZR	133	VAL	Peptide
1	ZR	135	MET	Peptide
1	ZS	241	PRO	Peptide
1	ZU	241	PRO	Peptide
1	ZW	241	PRO	Peptide
1	ZY	241	PRO	Peptide
1	ZZ	132	SER	Peptide
1	ZZ	250	ARG	Peptide,Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	YI	2421	0	2417	37	0
1	YJ	2437	0	2437	42	0
1	YK	2421	0	2417	21	0
1	YL	2437	0	2434	26	0
1	YM	2421	0	2417	32	0
1	YN	2437	0	2437	39	0
1	YO	2421	0	2417	21	0
1	YP	2437	0	2435	23	0
1	YQ	2421	0	2417	36	0
1	YR	2437	0	2437	34	0
1	YS	2421	0	2417	25	0
1	YT	2437	0	2435	28	0
1	YU	2421	0	2417	47	0
1	YV	2437	0	2437	41	0
1	YW	2421	0	2417	33	0
1	YX	2437	0	2436	32	0
1	YY	2421	0	2417	40	0
1	YZ	2437	0	2437	48	0
1	ZA	2421	0	2417	22	0
1	ZB	2437	0	2435	27	0
1	ZC	2421	0	2417	36	0
1	ZD	2437	0	2437	42	0
1	ZE	2421	0	2417	19	0
1	ZF	2437	0	2435	28	0
1	ZG	2421	0	2417	38	0
1	ZH	2437	0	2437	44	0
1	ZI	2421	0	2417	27	0
1	ZJ	2437	0	2435	22	0
1	ZK	2421	0	2417	37	0
1	ZL	2437	0	2435	49	0
1	ZM	2421	0	2417	31	0
1	ZN	2437	0	2436	36	0
1	ZO	2421	0	2417	43	0
1	ZP	2437	0	2437	43	0
1	ZQ	2421	0	2417	21	0
1	ZR	2437	0	2435	41	0
1	ZS	2421	0	2417	34	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	ZT	2437	0	2437	36	0
1	ZU	2421	0	2417	25	0
1	ZV	2437	0	2435	18	0
1	ZW	2421	0	2417	44	0
1	ZX	2437	0	2435	60	0
1	ZY	2421	0	2417	30	0
1	ZZ	2437	0	2435	18	0
2	YI	48	0	26	1	0
2	YJ	48	0	26	0	0
2	YK	48	0	26	1	0
2	YL	48	0	26	1	0
2	YM	48	0	26	2	0
2	YN	48	0	26	2	0
2	YO	48	0	26	0	0
2	YP	48	0	26	1	0
2	YQ	48	0	26	1	0
2	YR	48	0	26	1	0
2	YS	48	0	26	3	0
2	YT	48	0	26	3	0
2	YU	48	0	26	2	0
2	YV	48	0	26	1	0
2	YW	48	0	26	2	0
2	YX	48	0	26	1	0
2	YY	48	0	26	1	0
2	YZ	48	0	26	2	0
2	ZA	48	0	26	0	0
2	ZB	48	0	26	1	0
2	ZC	48	0	26	2	0
2	ZD	48	0	26	1	0
2	ZE	48	0	26	0	0
2	ZF	48	0	26	1	0
2	ZG	48	0	26	1	0
2	ZH	48	0	26	2	0
2	ZI	48	0	26	1	0
2	ZJ	48	0	26	1	0
2	ZK	48	0	26	1	0
2	ZL	48	0	26	4	0
2	ZM	48	0	26	1	0
2	ZN	48	0	26	1	0
2	ZO	48	0	26	0	0
2	ZP	48	0	26	0	0
2	ZQ	48	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	ZR	48	0	26	1	0
2	ZS	48	0	26	1	0
2	ZT	48	0	26	1	0
2	ZU	48	0	26	1	0
2	ZV	48	0	26	0	0
2	ZW	48	0	26	1	0
2	ZX	48	0	26	2	0
2	ZY	48	0	26	2	0
2	ZZ	48	0	26	3	0
3	YI	23	0	16	0	0
3	YJ	23	0	16	0	0
3	YK	23	0	16	1	0
3	YL	23	0	16	0	0
3	YM	23	0	16	0	0
3	YN	23	0	16	0	0
3	YO	23	0	16	1	0
3	YP	23	0	16	0	0
3	YQ	23	0	16	0	0
3	YR	23	0	16	0	0
3	YS	23	0	16	1	0
3	YT	23	0	16	0	0
3	YU	23	0	16	0	0
3	YV	23	0	16	0	0
3	YW	23	0	16	1	0
3	YX	23	0	16	0	0
3	YY	23	0	16	0	0
3	YZ	23	0	16	0	0
3	ZA	23	0	16	1	0
3	ZB	23	0	16	0	0
3	ZC	23	0	16	0	0
3	ZD	23	0	16	0	0
3	ZE	23	0	16	1	0
3	ZF	23	0	16	0	0
3	ZG	23	0	16	0	0
3	ZH	23	0	16	0	0
3	ZI	23	0	16	1	0
3	ZJ	23	0	16	0	0
3	ZK	23	0	16	0	0
3	ZL	23	0	16	0	0
3	ZM	23	0	16	1	0
3	ZN	23	0	16	0	0
3	ZO	23	0	16	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	ZP	23	0	16	0	0
3	ZQ	23	0	16	1	0
3	ZR	23	0	16	0	0
3	ZS	23	0	16	0	0
3	ZT	23	0	16	0	0
3	ZU	23	0	16	0	0
3	ZV	23	0	16	0	0
3	ZW	23	0	16	0	0
3	ZX	23	0	16	1	0
3	ZY	23	0	16	1	0
3	ZZ	23	0	16	0	0
4	YI	45	0	0	0	0
4	YJ	45	0	0	0	0
4	YK	45	0	0	1	0
4	YL	45	0	0	1	0
4	YM	45	0	0	1	0
4	YN	45	0	0	0	0
4	YO	45	0	0	0	0
4	YP	45	0	0	1	0
4	YQ	45	0	0	0	0
4	YR	45	0	0	0	0
4	YS	45	0	0	1	0
4	YT	45	0	0	1	0
4	YU	45	0	0	1	0
4	YV	45	0	0	1	0
4	YW	45	0	0	0	0
4	YX	45	0	0	1	0
4	YY	45	0	0	0	0
4	YZ	45	0	0	1	0
4	ZA	45	0	0	0	0
4	ZB	45	0	0	1	0
4	ZC	45	0	0	1	0
4	ZD	45	0	0	0	0
4	ZE	45	0	0	0	0
4	ZF	45	0	0	1	0
4	ZG	45	0	0	0	0
4	ZH	45	0	0	1	0
4	ZI	45	0	0	1	0
4	ZJ	45	0	0	1	0
4	ZK	45	0	0	0	0
4	ZL	45	0	0	1	0
4	ZM	45	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	ZN	45	0	0	1	0
4	ZO	45	0	0	0	0
4	ZP	45	0	0	0	0
4	ZQ	45	0	0	0	0
4	ZR	45	0	0	1	0
4	ZS	45	0	0	0	0
4	ZT	45	0	0	0	0
4	ZU	45	0	0	1	0
4	ZV	45	0	0	0	0
4	ZW	45	0	0	0	0
4	ZX	45	0	0	1	0
4	ZY	45	0	0	0	0
4	ZZ	45	0	0	1	0
All	All	111980	0	108611	1435	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (1435) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZN:250:ARG:HE	1:ZR:133:VAL:HG13	1.15	1.08
1:ZF:250:ARG:HE	1:ZJ:106:LYS:HZ3	1.14	0.94
1:ZC:206:ALA:O	1:ZC:210:LEU:HB2	1.69	0.92
1:ZI:252:LEU:HD12	1:ZI:265:ILE:HD11	1.52	0.92
1:YU:211:ASP:O	1:YU:215:LYS:HB2	1.70	0.90
1:YK:252:LEU:HD12	1:YK:265:ILE:HD11	1.54	0.89
1:ZU:252:LEU:HD12	1:ZU:265:ILE:HD11	1.53	0.89
1:YM:206:ALA:O	1:YM:210:LEU:HB2	1.74	0.88
1:YS:252:LEU:HD12	1:YS:265:ILE:HD11	1.57	0.87
1:ZR:124:LEU:O	1:ZR:127:GLU:O	1.91	0.87
1:ZG:211:ASP:O	1:ZG:215:LYS:HB2	1.75	0.86
1:ZY:252:LEU:HD12	1:ZY:265:ILE:HD11	1.56	0.86
1:YP:250:ARG:HG3	1:YT:133:VAL:HG22	1.58	0.85
1:ZN:247:GLY:HA3	1:ZN:266:ASP:HA	1.58	0.84
1:ZR:250:ARG:HH22	1:ZV:106:LYS:HD2	1.39	0.84
1:ZU:264:MET:HE2	1:ZU:271:ASP:H	1.43	0.84
1:YK:264:MET:HE2	1:YK:271:ASP:H	1.43	0.83
1:YZ:249:LEU:HD11	1:YZ:386:LYS:HG3	1.57	0.83
1:ZY:264:MET:HE2	1:ZY:271:ASP:H	1.43	0.83
1:ZO:157:VAL:HG11	1:ZO:208:LEU:HD22	1.60	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YS:264:MET:HE2	1:YS:271:ASP:H	1.43	0.82
1:ZO:206:ALA:O	1:ZO:210:LEU:HB2	1.81	0.81
1:ZS:206:ALA:O	1:ZS:210:LEU:HB2	1.80	0.81
1:ZM:264:MET:HE2	1:ZM:271:ASP:H	1.46	0.81
1:ZT:96:ALA:H	1:ZT:118:MET:HE1	1.45	0.81
1:YO:264:MET:HE2	1:YO:271:ASP:H	1.45	0.81
1:ZS:157:VAL:HG11	1:ZS:208:LEU:HD22	1.63	0.81
1:ZI:264:MET:HE2	1:ZI:271:ASP:H	1.44	0.80
1:YY:224:ILE:HG12	1:YY:303:ALA:HB3	1.64	0.80
1:YI:252:LEU:HD12	1:YI:265:ILE:HD12	1.64	0.80
1:YM:211:ASP:O	1:YM:215:LYS:HB2	1.81	0.80
1:ZC:224:ILE:HG12	1:ZC:303:ALA:HB3	1.63	0.80
1:ZG:224:ILE:HG12	1:ZG:303:ALA:HB3	1.64	0.79
1:ZQ:252:LEU:HD12	1:ZQ:265:ILE:HD12	1.64	0.79
1:ZK:224:ILE:HG12	1:ZK:303:ALA:HB3	1.64	0.79
1:ZW:224:ILE:HG12	1:ZW:303:ALA:HB3	1.65	0.79
1:YM:224:ILE:HG12	1:YM:303:ALA:HB3	1.62	0.79
1:ZO:224:ILE:HG12	1:ZO:303:ALA:HB3	1.64	0.79
1:ZD:344:SER:HA	1:ZD:347:ARG:HD2	1.65	0.79
1:ZW:88:LYS:HD3	1:ZW:113:LYS:HZ2	1.47	0.79
1:YW:252:LEU:HD12	1:YW:265:ILE:HD12	1.63	0.79
1:YI:224:ILE:HG12	1:YI:303:ALA:HB3	1.65	0.79
1:YQ:224:ILE:HG12	1:YQ:303:ALA:HB3	1.64	0.79
1:ZM:252:LEU:HD12	1:ZM:265:ILE:HD11	1.64	0.79
1:ZP:150:LEU:HB3	1:ZP:154:ARG:HH12	1.48	0.78
1:YU:224:ILE:HG12	1:YU:303:ALA:HB3	1.63	0.78
1:ZE:252:LEU:HD12	1:ZE:265:ILE:HD12	1.64	0.78
1:ZA:252:LEU:HD12	1:ZA:265:ILE:HD12	1.64	0.78
1:ZZ:250:ARG:HB2	1:ZZ:253:ALA:HB3	1.66	0.78
1:ZH:150:LEU:HB3	1:ZH:154:ARG:HH12	1.49	0.77
1:ZG:252:LEU:HD12	1:ZG:265:ILE:HD12	1.63	0.77
1:ZW:252:LEU:HD12	1:ZW:265:ILE:HD12	1.65	0.77
1:YL:102:LEU:HG	1:YL:133:VAL:HG21	1.65	0.77
1:YQ:252:LEU:HD12	1:YQ:265:ILE:HD12	1.67	0.76
1:YR:150:LEU:HB3	1:YR:154:ARG:HH12	1.51	0.76
1:ZT:150:LEU:HB3	1:ZT:154:ARG:HH12	1.50	0.76
1:ZF:250:ARG:HB2	1:ZF:253:ALA:HB3	1.68	0.76
1:ZA:264:MET:HE2	1:ZA:271:ASP:H	1.51	0.75
1:ZG:206:ALA:O	1:ZG:210:LEU:HB2	1.87	0.75
1:YZ:350:GLN:HE21	1:YZ:357:LEU:HD11	1.52	0.75
1:ZK:252:LEU:HD12	1:ZK:265:ILE:HD12	1.67	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZO:252:LEU:HD12	1:ZO:265:ILE:HD12	1.68	0.75
1:YV:213:LEU:HD13	1:YV:223:LEU:HD11	1.67	0.75
1:ZX:350:GLN:HE21	1:ZX:357:LEU:HD11	1.52	0.75
1:ZR:126:ALA:HB3	1:ZR:130:ALA:HB2	1.68	0.74
1:ZD:213:LEU:HD13	1:ZD:223:LEU:HD11	1.69	0.74
1:ZP:213:LEU:HD13	1:ZP:223:LEU:HD11	1.68	0.74
1:YJ:150:LEU:HB3	1:YJ:154:ARG:HH12	1.52	0.74
1:ZQ:264:MET:HE2	1:ZQ:271:ASP:H	1.53	0.74
1:YL:134:GLY:HA2	1:ZZ:250:ARG:HG3	1.68	0.74
1:YL:214:LYS:HE2	1:YL:298:THR:HG22	1.70	0.74
1:ZL:167:LEU:HD21	1:ZL:170:LEU:HD21	1.70	0.73
1:ZA:210:LEU:HD11	1:ZA:290:PHE:HE1	1.53	0.73
1:YW:210:LEU:HD11	1:YW:290:PHE:HE1	1.54	0.73
1:YM:249:LEU:HD12	1:YM:389:LYS:HE3	1.69	0.73
1:ZR:214:LYS:HE2	1:ZR:298:THR:HG22	1.70	0.73
1:ZS:252:LEU:HB2	1:ZS:265:ILE:HG23	1.71	0.73
1:YY:206:ALA:O	1:YY:210:LEU:HB2	1.88	0.73
1:ZM:210:LEU:HD11	1:ZM:290:PHE:HE1	1.52	0.73
1:YW:264:MET:HE2	1:YW:271:ASP:H	1.53	0.73
1:ZP:96:ALA:H	1:ZP:118:MET:HE1	1.53	0.73
1:ZY:210:LEU:HD11	1:ZY:290:PHE:HE1	1.52	0.73
1:ZD:162:ARG:HE	1:ZI:162:ARG:HH12	1.37	0.73
1:YO:210:LEU:HD11	1:YO:290:PHE:HE1	1.53	0.72
1:YN:213:LEU:HD13	1:YN:223:LEU:HD11	1.70	0.72
1:YY:211:ASP:O	1:YY:215:LYS:HB2	1.89	0.72
1:YI:206:ALA:O	1:YI:210:LEU:HB2	1.88	0.72
1:ZN:250:ARG:HH21	1:ZR:133:VAL:H	1.36	0.72
1:ZJ:214:LYS:HE2	1:ZJ:298:THR:HG22	1.72	0.72
1:ZT:213:LEU:HD13	1:ZT:223:LEU:HD11	1.72	0.72
1:YV:150:LEU:HB3	1:YV:154:ARG:HH12	1.53	0.71
1:YV:162:ARG:HD2	1:ZA:162:ARG:HD2	1.70	0.71
1:ZH:213:LEU:HD13	1:ZH:223:LEU:HD11	1.71	0.71
1:YT:129:ALA:HA	1:YT:132:SER:HB2	1.72	0.71
1:ZL:230:THR:HG22	1:ZL:277:LYS:HE3	1.73	0.71
1:ZW:206:ALA:O	1:ZW:210:LEU:HB2	1.91	0.70
1:YP:214:LYS:HE2	1:YP:298:THR:HG22	1.73	0.70
1:ZS:211:ASP:O	1:ZS:215:LYS:HB2	1.91	0.70
1:YR:213:LEU:HD13	1:YR:223:LEU:HD11	1.72	0.70
1:ZU:210:LEU:HD11	1:ZU:290:PHE:HE1	1.54	0.70
1:ZX:249:LEU:HD11	1:ZX:386:LYS:HG3	1.72	0.69
1:ZE:214:LYS:HE2	1:ZE:298:THR:HG22	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZX:220:SER:HB3	1:ZX:222:ARG:HH12	1.57	0.69
1:ZK:88:LYS:HD2	1:ZK:113:LYS:HA	1.73	0.69
1:YJ:213:LEU:HD13	1:YJ:223:LEU:HD11	1.75	0.69
1:ZS:224:ILE:HG13	1:ZS:303:ALA:HB3	1.75	0.69
1:YV:335:THR:HG23	1:YV:337:GLY:H	1.59	0.68
1:ZV:214:LYS:HE2	1:ZV:298:THR:HG22	1.74	0.68
1:ZX:230:THR:HG22	1:ZX:277:LYS:HE3	1.75	0.68
1:YL:133:VAL:HG13	1:YL:135:MET:HG3	1.75	0.68
1:YY:252:LEU:HD12	1:YY:265:ILE:HD12	1.76	0.68
1:ZL:350:GLN:HE21	1:ZL:357:LEU:HD11	1.58	0.68
1:ZB:218:TYR:HD1	1:ZB:220:SER:H	1.41	0.67
1:YZ:230:THR:HG22	1:YZ:277:LYS:HE3	1.76	0.67
1:YZ:392:GLU:O	1:YZ:395:GLU:HG2	1.94	0.67
1:YX:100:LEU:HD21	1:YX:310:ILE:HD13	1.77	0.67
1:ZC:252:LEU:HD12	1:ZC:265:ILE:HD12	1.75	0.67
1:ZE:241:PRO:HG2	3:ZE:502:LMG:HC8	1.77	0.67
1:YI:336:LYS:O	1:YI:336:LYS:HD3	1.96	0.66
1:YU:252:LEU:HD12	1:YU:265:ILE:HD12	1.76	0.66
1:YU:313:THR:HG22	1:YU:314:GLY:H	1.60	0.66
1:YZ:249:LEU:HD13	1:YZ:390:VAL:HG23	1.77	0.66
1:ZK:206:ALA:O	1:ZK:210:LEU:HB2	1.95	0.66
1:ZP:162:ARG:HG2	1:ZU:162:ARG:HH21	1.60	0.66
1:ZX:292:ARG:HH12	1:ZX:384:VAL:HA	1.60	0.66
1:YL:87:ARG:HH12	1:YL:218:TYR:HD2	1.43	0.66
1:ZF:218:TYR:HD1	1:ZF:220:SER:H	1.41	0.66
1:ZO:313:THR:HG22	1:ZO:314:GLY:H	1.61	0.66
1:ZO:336:LYS:O	1:ZO:336:LYS:HD3	1.96	0.66
1:ZS:336:LYS:O	1:ZS:336:LYS:HD3	1.96	0.66
1:YY:313:THR:HG22	1:YY:314:GLY:H	1.61	0.66
1:ZI:241:PRO:HG2	3:ZI:502:LMG:HC8	1.78	0.66
1:YU:336:LYS:O	1:YU:336:LYS:HD3	1.96	0.66
1:ZO:211:ASP:O	1:ZO:215:LYS:HB2	1.95	0.66
1:YV:252:LEU:HB2	1:YV:265:ILE:HG23	1.77	0.66
1:ZG:336:LYS:O	1:ZG:336:LYS:HD3	1.96	0.66
1:ZW:195:VAL:HG21	1:ZW:265:ILE:HD11	1.78	0.66
1:YM:313:THR:HG22	1:YM:314:GLY:H	1.61	0.65
1:YQ:336:LYS:O	1:YQ:336:LYS:HD3	1.96	0.65
1:ZG:252:LEU:HB2	1:ZG:265:ILE:HG23	1.78	0.65
1:ZK:211:ASP:O	1:ZK:215:LYS:HB2	1.96	0.65
1:ZT:162:ARG:HE	1:ZY:162:ARG:HH12	1.44	0.65
1:ZG:313:THR:HG22	1:ZG:314:GLY:H	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZO:195:VAL:HG21	1:ZO:265:ILE:HD11	1.78	0.65
1:YR:162:ARG:HD2	1:YW:162:ARG:HH21	1.61	0.65
1:ZH:252:LEU:HB2	1:ZH:265:ILE:HG23	1.78	0.65
1:ZL:252:LEU:HB2	1:ZL:265:ILE:HG12	1.78	0.65
1:YN:162:ARG:HE	1:YS:162:ARG:HH12	1.42	0.65
1:ZT:252:LEU:HB2	1:ZT:265:ILE:HG23	1.78	0.65
1:ZT:339:VAL:HG21	1:ZT:347:ARG:HH12	1.61	0.65
1:YQ:297:GLU:HG3	1:YQ:298:THR:HG23	1.79	0.65
1:YZ:220:SER:HB3	1:YZ:222:ARG:HH12	1.60	0.65
1:ZX:210:LEU:HA	1:ZX:213:LEU:HB2	1.79	0.65
1:ZC:313:THR:HG22	1:ZC:314:GLY:H	1.62	0.65
1:ZW:313:THR:HG22	1:ZW:314:GLY:H	1.61	0.65
1:YI:313:THR:HG22	1:YI:314:GLY:H	1.62	0.64
1:YU:206:ALA:O	1:YU:210:LEU:HB2	1.96	0.64
1:ZS:249:LEU:HA	1:ZS:265:ILE:HG22	1.79	0.64
1:YZ:210:LEU:HA	1:YZ:213:LEU:HB2	1.79	0.64
1:ZS:284:MET:HA	1:ZS:287:MET:HE2	1.78	0.64
1:ZX:253:ALA:HA	1:ZX:393:ILE:HG21	1.79	0.64
1:ZC:233:THR:HG23	1:ZC:243:LYS:HD2	1.77	0.64
1:ZS:313:THR:HG22	1:ZS:314:GLY:H	1.62	0.64
1:YN:96:ALA:HB3	1:YN:118:MET:SD	2.38	0.64
1:ZZ:250:ARG:HA	1:ZZ:253:ALA:H	1.62	0.64
1:ZT:335:THR:HG23	1:ZT:337:GLY:H	1.61	0.64
1:YK:241:PRO:HG2	3:YK:502:LMG:HC8	1.78	0.64
1:ZF:250:ARG:HA	1:ZF:253:ALA:H	1.63	0.64
1:YQ:206:ALA:O	1:YQ:210:LEU:HB2	1.98	0.64
1:YV:96:ALA:HB3	1:YV:118:MET:SD	2.38	0.64
1:ZQ:214:LYS:HE2	1:ZQ:298:THR:HA	1.80	0.63
1:YN:339:VAL:HG11	1:YN:347:ARG:HH12	1.63	0.63
1:ZQ:241:PRO:HG2	3:ZQ:502:LMG:HC8	1.79	0.63
1:ZW:95:GLY:HA2	2:ZW:501:NDP:H1B	1.81	0.63
1:ZD:252:LEU:HB2	1:ZD:265:ILE:HG23	1.80	0.63
1:ZR:124:LEU:O	1:ZR:127:GLU:C	2.41	0.63
1:ZC:252:LEU:HB2	1:ZC:265:ILE:HG23	1.79	0.63
1:ZL:289:GLU:HG2	1:ZL:391:TRP:HB2	1.78	0.63
1:YR:96:ALA:HB3	1:YR:118:MET:SD	2.39	0.63
1:YR:252:LEU:HB2	1:YR:265:ILE:HG23	1.80	0.63
1:ZJ:87:ARG:HH12	1:ZJ:218:TYR:HD1	1.44	0.63
1:ZO:297:GLU:HG3	1:ZO:298:THR:HG23	1.80	0.63
1:ZC:297:GLU:HG3	1:ZC:298:THR:HG23	1.81	0.63
1:YJ:252:LEU:HB2	1:YJ:265:ILE:HG23	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YQ:339:VAL:HG21	1:YQ:347:ARG:HH12	1.64	0.62
1:YU:95:GLY:HA2	2:YU:501:NDP:H1B	1.82	0.62
1:ZX:199:HIS:HD1	1:ZX:199:HIS:C	2.07	0.62
1:ZX:252:LEU:HB2	1:ZX:265:ILE:HG23	1.81	0.62
1:ZC:95:GLY:HA2	2:ZC:501:NDP:H1B	1.82	0.62
1:YJ:230:THR:HG22	1:YJ:277:LYS:HD3	1.82	0.62
1:YY:297:GLU:HG3	1:YY:298:THR:HG23	1.82	0.62
1:ZG:297:GLU:HG3	1:ZG:298:THR:HG23	1.82	0.62
1:ZR:124:LEU:HB3	1:ZR:127:GLU:HA	1.81	0.62
1:ZL:249:LEU:HA	1:ZL:265:ILE:HG22	1.81	0.62
1:ZK:313:THR:HG22	1:ZK:314:GLY:H	1.63	0.61
2:ZU:501:NDP:H41N	4:ZU:503:A1JWG:O1A	2.00	0.61
1:YT:250:ARG:HG3	1:YT:253:ALA:HB3	1.81	0.61
1:ZL:220:SER:HB3	1:ZL:222:ARG:HH12	1.65	0.61
1:ZX:289:GLU:HG2	1:ZX:391:TRP:HB2	1.83	0.61
1:YQ:313:THR:HG22	1:YQ:314:GLY:H	1.64	0.61
1:ZN:100:LEU:HD21	1:ZN:310:ILE:HD13	1.82	0.61
1:ZX:200:LEU:HD23	1:ZX:398:VAL:HG12	1.82	0.61
1:YX:111:THR:HG22	1:YX:113:LYS:HG3	1.81	0.61
1:ZG:95:GLY:HA2	2:ZG:501:NDP:H1B	1.81	0.61
1:ZN:250:ARG:NE	1:ZR:133:VAL:HG13	2.01	0.61
1:ZL:210:LEU:HA	1:ZL:213:LEU:HB2	1.82	0.61
1:YX:250:ARG:HB2	1:YX:253:ALA:HB3	1.83	0.61
1:ZA:241:PRO:HG2	3:ZA:502:LMG:HC8	1.83	0.61
1:YL:250:ARG:HG3	1:YP:134:GLY:HA2	1.81	0.61
1:ZL:385:GLU:O	1:ZL:389:LYS:HG2	2.01	0.61
1:ZD:339:VAL:HG11	1:ZD:347:ARG:NH1	2.15	0.61
1:ZR:250:ARG:HH12	1:ZV:106:LYS:NZ	1.99	0.61
1:YP:111:THR:HG22	1:YP:113:LYS:HG3	1.83	0.60
1:YX:250:ARG:HG3	1:ZB:134:GLY:HA2	1.83	0.60
1:YN:252:LEU:HB2	1:YN:265:ILE:HG23	1.82	0.60
1:YR:339:VAL:HG11	1:YR:347:ARG:HH22	1.66	0.60
1:YX:250:ARG:HH12	1:ZB:133:VAL:HA	1.66	0.60
1:ZX:329:PRO:HA	1:ZX:332:LYS:HG2	1.83	0.60
1:YJ:285:LEU:HD12	1:YJ:387:ALA:HB2	1.84	0.60
1:ZJ:150:LEU:HD12	1:ZJ:204:LEU:HD22	1.83	0.60
1:YK:395:GLU:HG2	1:YK:401:ALA:HB2	1.83	0.60
1:YV:285:LEU:HD12	1:YV:387:ALA:HB2	1.84	0.60
2:YT:501:NDP:H41N	4:YT:503:A1JWG:O1A	2.01	0.60
1:ZD:96:ALA:HB3	1:ZD:118:MET:SD	2.42	0.60
2:ZF:501:NDP:H41N	4:ZF:503:A1JWG:O1A	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZS:249:LEU:HD12	1:ZS:389:LYS:HE3	1.84	0.60
1:YV:230:THR:HG22	1:YV:277:LYS:HD3	1.84	0.60
1:YI:297:GLU:HG3	1:YI:298:THR:HG23	1.82	0.59
1:YP:252:LEU:HB2	1:YP:265:ILE:HG23	1.83	0.59
1:ZZ:150:LEU:HD12	1:ZZ:204:LEU:HD22	1.85	0.59
1:YI:339:VAL:HG21	1:YI:347:ARG:HH12	1.67	0.59
1:YM:252:LEU:HD12	1:YM:265:ILE:HD12	1.83	0.59
1:YX:229:ILE:HD11	1:YX:235:THR:HG21	1.84	0.59
1:ZF:150:LEU:HD12	1:ZF:204:LEU:HD13	1.84	0.59
1:ZP:230:THR:HG22	1:ZP:277:LYS:HD3	1.85	0.59
1:YN:335:THR:HG23	1:YN:337:GLY:H	1.66	0.59
1:YI:211:ASP:O	1:YI:215:LYS:HB2	2.02	0.59
1:ZN:229:ILE:HD11	1:ZN:235:THR:HG21	1.84	0.59
2:ZZ:501:NDP:H51N	2:ZZ:501:NDP:H2N	1.84	0.59
1:YU:199:HIS:HD1	1:YU:199:HIS:C	2.11	0.59
1:ZK:339:VAL:HG21	1:ZK:347:ARG:HH12	1.68	0.59
1:YR:285:LEU:HD12	1:YR:387:ALA:HB2	1.85	0.59
1:ZY:241:PRO:HG2	3:ZY:502:LMG:HC8	1.84	0.59
1:YI:249:LEU:HD12	1:YI:389:LYS:HE3	1.83	0.59
2:YT:501:NDP:H51N	2:YT:501:NDP:H2N	1.83	0.59
1:YZ:289:GLU:HG2	1:YZ:391:TRP:HB2	1.84	0.59
1:YI:233:THR:HG23	1:YI:243:LYS:HD2	1.83	0.58
1:YU:249:LEU:HD12	1:YU:389:LYS:HE3	1.84	0.58
1:ZK:297:GLU:HG3	1:ZK:298:THR:HG23	1.83	0.58
1:ZL:253:ALA:HA	1:ZL:393:ILE:HG21	1.84	0.58
1:ZB:250:ARG:HB2	1:ZB:253:ALA:HB3	1.85	0.58
1:ZI:159:ASN:O	1:ZI:162:ARG:HG2	2.02	0.58
1:ZL:313:THR:HG22	1:ZL:314:GLY:H	1.68	0.58
1:YY:95:GLY:HA2	2:YY:501:NDP:H1B	1.86	0.58
1:ZP:252:LEU:HB2	1:ZP:265:ILE:HG23	1.84	0.58
1:ZT:230:THR:HG22	1:ZT:277:LYS:HD3	1.85	0.58
2:YX:501:NDP:H41N	4:YX:503:A1JWG:O1A	2.03	0.58
1:ZO:264:MET:HE1	1:ZO:278:ASP:HB3	1.84	0.58
1:ZV:229:ILE:HD11	1:ZV:235:THR:HG21	1.85	0.58
1:ZZ:331:GLN:HA	1:ZZ:335:THR:HG22	1.85	0.58
1:YT:229:ILE:HD11	1:YT:235:THR:HG21	1.85	0.58
2:ZB:501:NDP:H41N	4:ZB:503:A1JWG:O1A	2.04	0.58
1:ZG:339:VAL:HG21	1:ZG:347:ARG:HH12	1.67	0.58
1:ZX:313:THR:HG22	1:ZX:314:GLY:H	1.69	0.58
1:YQ:290:PHE:HB3	1:YQ:302:PHE:CE2	2.39	0.58
1:ZJ:111:THR:HG22	1:ZJ:113:LYS:HG3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZR:229:ILE:HD11	1:ZR:235:THR:HG21	1.86	0.58
1:ZZ:249:LEU:HD11	1:ZZ:386:LYS:HG3	1.86	0.58
2:ZR:501:NDP:H41N	4:ZR:503:A1JWG:O1A	2.04	0.58
1:YU:339:VAL:HG21	1:YU:347:ARG:HH12	1.68	0.58
1:YM:339:VAL:HG21	1:YM:347:ARG:HH12	1.69	0.57
1:YN:285:LEU:HD12	1:YN:387:ALA:HB2	1.86	0.57
1:YV:226:VAL:HG22	1:YV:305:LEU:HD21	1.86	0.57
1:ZW:297:GLU:HG3	1:ZW:298:THR:HG23	1.85	0.57
1:YL:246:LEU:HD23	1:YL:249:LEU:HD21	1.85	0.57
1:ZC:246:LEU:HD23	1:ZC:249:LEU:HD21	1.84	0.57
2:ZJ:501:NDP:H41N	4:ZJ:503:A1JWG:O1A	2.04	0.57
1:ZW:290:PHE:HB3	1:ZW:302:PHE:CE2	2.39	0.57
1:YO:343:GLU:O	1:YO:346:LYS:HG2	2.03	0.57
1:ZD:162:ARG:HG2	1:ZI:162:ARG:HH22	1.69	0.57
1:ZF:331:GLN:HA	1:ZF:335:THR:HG22	1.85	0.57
1:ZL:95:GLY:HA2	2:ZL:501:NDP:H1B	1.86	0.57
1:ZO:290:PHE:HB3	1:ZO:302:PHE:CE2	2.40	0.57
1:ZW:339:VAL:HG21	1:ZW:347:ARG:HH12	1.69	0.57
1:ZN:250:ARG:O	1:ZN:253:ALA:O	2.22	0.57
1:ZN:331:GLN:HA	1:ZN:335:THR:HG22	1.86	0.57
1:ZU:343:GLU:O	1:ZU:346:LYS:HG2	2.05	0.57
1:ZW:246:LEU:HD23	1:ZW:249:LEU:HD21	1.87	0.57
2:YL:501:NDP:H41N	4:YL:503:A1JWG:O1A	2.04	0.57
1:YY:246:LEU:HD23	1:YY:249:LEU:HD21	1.86	0.57
1:ZO:157:VAL:HG13	1:ZO:209:LEU:HD13	1.87	0.57
1:YJ:287:MET:HE2	1:YJ:304:SER:HB3	1.85	0.57
1:YM:95:GLY:HA2	2:YM:501:NDP:H1B	1.86	0.57
1:YN:230:THR:HG22	1:YN:277:LYS:HD3	1.84	0.57
1:ZL:384:VAL:O	1:ZL:388:ARG:HG2	2.05	0.57
2:ZZ:501:NDP:H41N	4:ZZ:503:A1JWG:O1A	2.05	0.57
1:YR:331:GLN:HA	1:YR:335:THR:HG22	1.87	0.57
1:YQ:252:LEU:HB2	1:YQ:265:ILE:HG23	1.87	0.57
1:ZO:339:VAL:HG21	1:ZO:347:ARG:HH12	1.70	0.57
1:YJ:331:GLN:HA	1:YJ:335:THR:HG22	1.87	0.57
1:YQ:249:LEU:HD12	1:YQ:389:LYS:HE3	1.85	0.57
1:YU:389:LYS:O	1:YU:393:ILE:HG13	2.05	0.57
1:YZ:313:THR:HG22	1:YZ:314:GLY:H	1.70	0.57
1:ZS:95:GLY:HA2	2:ZS:501:NDP:H1B	1.86	0.57
1:YO:241:PRO:HG2	3:YO:502:LMG:HC8	1.86	0.56
1:ZH:230:THR:HG22	1:ZH:277:LYS:HD3	1.86	0.56
1:YZ:236:LEU:HD13	1:ZC:236:LEU:HD13	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZW:249:LEU:HD12	1:ZW:389:LYS:HE3	1.87	0.56
1:YR:95:GLY:HA3	1:YR:174:ALA:HB2	1.87	0.56
1:YZ:95:GLY:HA2	2:YZ:501:NDP:H1B	1.87	0.56
1:ZP:287:MET:HE2	1:ZP:304:SER:HB3	1.86	0.56
1:YV:159:ASN:HD21	1:ZA:159:ASN:HD21	1.53	0.56
1:ZK:252:LEU:HB2	1:ZK:265:ILE:HG23	1.87	0.56
1:ZN:250:ARG:NH2	1:ZR:133:VAL:H	2.03	0.56
1:ZT:226:VAL:HG22	1:ZT:305:LEU:HD21	1.88	0.56
2:YP:501:NDP:H41N	4:YP:503:A1JWG:O1A	2.04	0.56
1:ZT:162:ARG:HG2	1:ZY:162:ARG:HH22	1.69	0.56
1:ZV:100:LEU:HD21	1:ZV:310:ILE:HD13	1.87	0.56
1:ZW:250:ARG:HB3	1:ZW:267:GLY:HA3	1.88	0.56
1:ZY:343:GLU:O	1:ZY:346:LYS:HG2	2.06	0.56
1:YT:331:GLN:HA	1:YT:335:THR:HG22	1.88	0.56
1:YU:297:GLU:HG3	1:YU:298:THR:HG23	1.86	0.56
1:YY:290:PHE:HB3	1:YY:302:PHE:CE2	2.41	0.56
1:ZL:392:GLU:O	1:ZL:395:GLU:HG2	2.05	0.56
1:ZG:249:LEU:HB2	1:ZG:389:LYS:HE3	1.88	0.56
1:YW:343:GLU:O	1:YW:346:LYS:HG2	2.05	0.56
1:ZB:331:GLN:HA	1:ZB:335:THR:HG22	1.87	0.56
1:ZR:125:LYS:C	1:ZR:127:GLU:O	2.49	0.56
1:ZB:150:LEU:HD12	1:ZB:204:LEU:HD22	1.86	0.56
1:ZD:331:GLN:HA	1:ZD:335:THR:HG22	1.86	0.56
1:YR:162:ARG:HD2	1:YW:162:ARG:NH2	2.21	0.56
1:ZG:289:GLU:HG2	1:ZG:391:TRP:HB2	1.88	0.56
1:YQ:95:GLY:HA2	2:YQ:501:NDP:H1B	1.88	0.55
1:YQ:246:LEU:HD23	1:YQ:249:LEU:HD21	1.88	0.55
1:ZX:246:LEU:HD13	1:ZX:386:LYS:HD3	1.87	0.55
1:YZ:234:ASN:HB2	1:ZC:234:ASN:HB2	1.88	0.55
1:YJ:339:VAL:HG21	1:YJ:347:ARG:HH12	1.71	0.55
1:YU:216:SER:OG	1:YU:221:LYS:HD3	2.06	0.55
1:YY:336:LYS:HD2	1:YY:336:LYS:O	2.06	0.55
1:ZX:95:GLY:HA2	2:ZX:501:NDP:H1B	1.87	0.55
1:ZX:211:ASP:HA	1:ZX:214:LYS:HE3	1.88	0.55
1:YL:229:ILE:HD11	1:YL:235:THR:HG21	1.87	0.55
1:ZK:246:LEU:HD23	1:ZK:249:LEU:HD21	1.87	0.55
1:ZS:118:MET:HE1	1:ZS:130:ALA:HB2	1.88	0.55
1:YI:246:LEU:HD23	1:YI:249:LEU:HD21	1.88	0.55
1:YM:246:LEU:HD23	1:YM:249:LEU:HD21	1.88	0.55
1:YR:230:THR:HG22	1:YR:277:LYS:HD3	1.87	0.55
1:YX:250:ARG:HA	1:YX:253:ALA:H	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YX:331:GLN:HA	1:YX:335:THR:HG22	1.87	0.55
1:ZQ:343:GLU:O	1:ZQ:346:LYS:HG2	2.07	0.55
1:ZX:287:MET:HE3	1:ZX:304:SER:HB2	1.87	0.55
1:YP:229:ILE:HD11	1:YP:235:THR:HG21	1.88	0.55
1:ZB:229:ILE:HD11	1:ZB:235:THR:HG21	1.89	0.55
1:ZJ:229:ILE:HD11	1:ZJ:235:THR:HG21	1.89	0.55
1:ZK:336:LYS:HD2	1:ZK:336:LYS:O	2.07	0.55
1:ZN:150:LEU:HD12	1:ZN:204:LEU:HD22	1.89	0.55
1:ZN:218:TYR:HD1	1:ZN:220:SER:H	1.53	0.55
1:YM:289:GLU:HG2	1:YM:391:TRP:HB2	1.89	0.55
1:ZF:229:ILE:HD11	1:ZF:235:THR:HG21	1.88	0.55
1:YJ:96:ALA:HB3	1:YJ:118:MET:SD	2.46	0.55
1:YL:331:GLN:HA	1:YL:335:THR:HG22	1.87	0.55
1:YM:290:PHE:HB3	1:YM:302:PHE:CE1	2.42	0.55
1:YN:162:ARG:HE	1:YS:162:ARG:NH1	2.05	0.55
1:ZT:118:MET:HE3	1:ZT:120:CYS:HB2	1.89	0.55
1:ZD:230:THR:HG22	1:ZD:277:LYS:HD3	1.88	0.55
1:ZH:331:GLN:HA	1:ZH:335:THR:HG22	1.88	0.55
1:YY:249:LEU:HD12	1:YY:389:LYS:HE3	1.90	0.54
1:ZC:339:VAL:HG21	1:ZC:347:ARG:HH12	1.71	0.54
1:ZJ:130:ALA:O	1:ZJ:135:MET:HB2	2.07	0.54
1:ZS:252:LEU:HD12	1:ZS:265:ILE:HD12	1.88	0.54
1:YZ:182:LYS:HG3	1:YZ:183:GLU:CD	2.32	0.54
1:ZH:319:HIS:ND1	1:ZH:320:ILE:HG12	2.22	0.54
1:ZT:331:GLN:HA	1:ZT:335:THR:HG22	1.89	0.54
1:YK:343:GLU:O	1:YK:346:LYS:HG2	2.08	0.54
1:ZK:95:GLY:HA2	2:ZK:501:NDP:H1B	1.88	0.54
2:ZN:501:NDP:H41N	4:ZN:503:A1JWG:O1A	2.07	0.54
1:ZS:297:GLU:HG3	1:ZS:298:THR:HG23	1.88	0.54
1:YZ:249:LEU:HB2	1:YZ:389:LYS:HE2	1.89	0.54
1:ZO:246:LEU:HD23	1:ZO:249:LEU:HD21	1.89	0.54
1:ZP:331:GLN:HA	1:ZP:335:THR:HG22	1.88	0.54
1:ZC:216:SER:OG	1:ZC:221:LYS:HD3	2.07	0.54
1:ZX:292:ARG:HH11	1:ZX:388:ARG:NH1	2.06	0.54
1:YM:297:GLU:HG3	1:YM:298:THR:HG23	1.90	0.54
1:YT:385:GLU:O	1:YT:389:LYS:HG2	2.08	0.54
1:ZL:234:ASN:HB2	1:ZO:234:ASN:HB2	1.89	0.54
1:YQ:160:PHE:CD2	1:YQ:209:LEU:HD21	2.42	0.54
1:YR:347:ARG:HD2	1:YR:366:TRP:CD2	2.43	0.54
1:YT:150:LEU:HD12	1:YT:204:LEU:HD22	1.90	0.54
1:YZ:218:TYR:CD1	1:YZ:219:PRO:HD2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YO:214:LYS:NZ	1:YO:298:THR:HA	2.23	0.54
1:ZT:339:VAL:HG11	1:ZT:347:ARG:NH1	2.23	0.54
1:ZX:366:TRP:NE1	1:ZX:372:SER:HB3	2.23	0.54
1:YJ:339:VAL:HG11	1:YJ:347:ARG:NH1	2.22	0.54
1:YZ:284:MET:HA	1:YZ:287:MET:HE2	1.89	0.54
1:ZH:347:ARG:HB3	1:ZH:366:TRP:CZ2	2.42	0.54
1:ZO:249:LEU:HD12	1:ZO:389:LYS:HE3	1.90	0.53
1:ZQ:124:LEU:HD21	1:ZQ:128:ARG:HH21	1.72	0.53
1:ZX:249:LEU:HD13	1:ZX:390:VAL:HG23	1.91	0.53
1:ZG:290:PHE:HB3	1:ZG:302:PHE:CE1	2.43	0.53
1:ZT:162:ARG:HE	1:ZY:162:ARG:NH1	2.06	0.53
1:YM:284:MET:O	1:YM:287:MET:HG2	2.08	0.53
1:YR:339:VAL:HG11	1:YR:347:ARG:NH2	2.23	0.53
1:ZN:250:ARG:O	1:ZN:253:ALA:C	2.50	0.53
1:ZP:226:VAL:HG22	1:ZP:305:LEU:HD21	1.91	0.53
1:YP:331:GLN:HA	1:YP:335:THR:HG22	1.91	0.53
1:YU:246:LEU:HD23	1:YU:249:LEU:HD21	1.90	0.53
1:ZF:250:ARG:HG3	1:ZJ:134:GLY:HA2	1.91	0.53
1:ZZ:229:ILE:HD11	1:ZZ:235:THR:HG21	1.89	0.53
1:YU:264:MET:HE3	1:YU:265:ILE:H	1.74	0.53
1:ZL:290:PHE:HB3	1:ZL:302:PHE:CZ	2.43	0.53
1:ZL:366:TRP:NE1	1:ZL:372:SER:HB3	2.23	0.53
1:ZW:336:LYS:HD2	1:ZW:336:LYS:O	2.09	0.53
1:ZX:100:LEU:HD21	1:ZX:348:LEU:HD13	1.90	0.53
1:YM:216:SER:OG	1:YM:221:LYS:HD3	2.09	0.53
1:YP:130:ALA:O	1:YP:135:MET:HB2	2.08	0.53
1:YS:229:ILE:HD11	1:YS:235:THR:HG21	1.91	0.53
1:YX:250:ARG:HH22	1:ZB:133:VAL:CG2	2.22	0.53
1:ZS:339:VAL:HG21	1:ZS:347:ARG:HH12	1.73	0.53
1:ZT:319:HIS:ND1	1:ZT:320:ILE:HG12	2.24	0.53
1:YV:331:GLN:HA	1:YV:335:THR:HG22	1.89	0.53
1:ZC:336:LYS:HD2	1:ZC:336:LYS:O	2.09	0.53
1:ZW:225:ILE:HG21	1:ZW:283:ASN:HB3	1.91	0.53
1:YJ:226:VAL:HG22	1:YJ:305:LEU:HD21	1.91	0.53
1:YL:250:ARG:CG	1:YP:134:GLY:HA2	2.39	0.53
1:YY:216:SER:OG	1:YY:221:LYS:HD3	2.09	0.53
1:YY:255:GLY:HA2	1:YY:397:LEU:HD21	1.91	0.53
1:ZE:124:LEU:HD21	1:ZE:128:ARG:HH21	1.73	0.53
1:ZL:199:HIS:ND1	1:ZL:199:HIS:C	2.66	0.53
1:ZQ:362:VAL:HG13	1:ZQ:374:GLU:HG2	1.91	0.53
1:ZS:290:PHE:HB3	1:ZS:302:PHE:CE1	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YM:284:MET:HA	1:YM:287:MET:HE2	1.89	0.53
1:ZW:216:SER:OG	1:ZW:221:LYS:HD3	2.09	0.53
1:ZX:218:TYR:CD1	1:ZX:219:PRO:HD2	2.44	0.53
1:YS:214:LYS:NZ	1:YS:298:THR:HA	2.24	0.52
1:YV:339:VAL:HG11	1:YV:347:ARG:NH1	2.24	0.52
1:YV:347:ARG:HB3	1:YV:366:TRP:CZ2	2.44	0.52
1:ZH:226:VAL:HG22	1:ZH:305:LEU:HD21	1.91	0.52
1:ZO:100:LEU:HD23	1:ZO:348:LEU:HD12	1.91	0.52
1:ZF:87:ARG:HH12	1:ZF:218:TYR:HD2	1.57	0.52
1:ZH:96:ALA:HB3	1:ZH:118:MET:SD	2.50	0.52
1:ZH:162:ARG:HE	1:ZM:162:ARG:HD2	1.74	0.52
1:ZW:255:GLY:HA2	1:ZW:397:LEU:HD21	1.91	0.52
2:ZH:501:NDP:H41N	4:ZH:503:A1JWG:O1A	2.09	0.52
1:ZO:289:GLU:HG2	1:ZO:391:TRP:HB2	1.90	0.52
1:YK:362:VAL:HG13	1:YK:374:GLU:HG2	1.92	0.52
1:YN:331:GLN:HA	1:YN:335:THR:HG22	1.91	0.52
1:YP:250:ARG:NE	1:YT:134:GLY:C	2.65	0.52
1:YU:289:GLU:HG2	1:YU:391:TRP:HB2	1.91	0.52
1:ZH:339:VAL:HG11	1:ZH:347:ARG:NH2	2.25	0.52
1:ZL:226:VAL:HG11	2:ZL:501:NDP:H1D	1.92	0.52
1:YW:159:ASN:HA	1:YW:162:ARG:HH11	1.74	0.52
1:ZC:289:GLU:HG2	1:ZC:391:TRP:HB2	1.92	0.52
1:ZL:172:CYS:HB3	1:ZL:202:HIS:CE1	2.45	0.52
1:ZM:362:VAL:HG13	1:ZM:374:GLU:HG2	1.92	0.52
1:ZX:199:HIS:C	1:ZX:199:HIS:ND1	2.66	0.52
1:YI:236:LEU:HD13	1:ZX:236:LEU:HD13	1.91	0.52
1:ZG:216:SER:OG	1:ZG:221:LYS:HD3	2.10	0.52
1:ZK:255:GLY:HA2	1:ZK:397:LEU:HD21	1.90	0.52
1:YY:252:LEU:HB2	1:YY:265:ILE:HG23	1.92	0.52
1:ZC:249:LEU:HB2	1:ZC:389:LYS:HE3	1.92	0.52
1:ZK:264:MET:SD	1:ZK:270:PHE:HA	2.50	0.52
1:YJ:223:LEU:HB2	1:YJ:302:PHE:HD1	1.75	0.52
1:YQ:255:GLY:HA2	1:YQ:397:LEU:HD21	1.92	0.52
1:YU:195:VAL:HG21	1:YU:265:ILE:HD11	1.91	0.52
1:ZA:343:GLU:HA	1:ZA:346:LYS:HG2	1.91	0.52
1:ZI:389:LYS:O	1:ZI:393:ILE:HG13	2.09	0.52
1:ZL:236:LEU:HD13	1:ZO:236:LEU:HD13	1.92	0.52
1:ZX:252:LEU:HD12	1:ZX:265:ILE:HG12	1.92	0.52
1:YS:241:PRO:HG2	3:YS:502:LMG:HC8	1.92	0.52
1:ZB:102:LEU:HD21	1:ZB:133:VAL:HG11	1.92	0.52
1:ZE:362:VAL:HG13	1:ZE:374:GLU:HG2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZH:281:VAL:HA	1:ZH:284:MET:HE3	1.92	0.52
1:ZK:284:MET:O	1:ZK:287:MET:HG2	2.10	0.52
1:ZW:211:ASP:O	1:ZW:215:LYS:HB2	2.09	0.52
1:YI:234:ASN:HB2	1:ZX:234:ASN:HB2	1.92	0.51
1:YN:232:ASN:O	1:YN:238:GLY:HA3	2.10	0.51
1:YQ:289:GLU:HG2	1:YQ:391:TRP:HB2	1.91	0.51
1:ZL:284:MET:O	1:ZL:287:MET:HG2	2.09	0.51
1:ZX:392:GLU:O	1:ZX:395:GLU:HG2	2.09	0.51
1:YI:216:SER:OG	1:YI:221:LYS:HD3	2.10	0.51
1:YI:319:HIS:CG	1:YI:320:ILE:H	2.28	0.51
1:YN:276:TYR:HE1	1:YN:280:LYS:HE2	1.75	0.51
1:YU:199:HIS:C	1:YU:199:HIS:ND1	2.69	0.51
1:ZB:248:ASP:C	1:ZB:249:LEU:HD23	2.35	0.51
1:ZR:125:LYS:O	1:ZR:127:GLU:O	2.28	0.51
1:ZV:124:LEU:HD21	1:ZV:128:ARG:HH21	1.75	0.51
1:ZD:100:LEU:HD12	1:ZD:310:ILE:HD12	1.92	0.51
1:ZF:124:LEU:HD21	1:ZF:128:ARG:NH2	2.26	0.51
1:YZ:366:TRP:NE1	1:YZ:372:SER:HB3	2.25	0.51
1:ZK:320:ILE:HG22	1:ZK:322:LEU:H	1.76	0.51
1:ZQ:88:LYS:HG2	1:ZQ:113:LYS:HA	1.92	0.51
1:ZS:284:MET:O	1:ZS:287:MET:HG2	2.10	0.51
1:YM:195:VAL:HG21	1:YM:265:ILE:HD11	1.93	0.51
1:YR:226:VAL:HG22	1:YR:305:LEU:HD21	1.90	0.51
1:ZL:230:THR:CG2	1:ZL:277:LYS:HE3	2.41	0.51
1:ZS:255:GLY:HA2	1:ZS:397:LEU:HD21	1.92	0.51
1:YI:289:GLU:HG2	1:YI:391:TRP:HB2	1.92	0.51
1:ZD:281:VAL:HA	1:ZD:284:MET:HE3	1.92	0.51
1:ZD:343:GLU:O	1:ZD:347:ARG:HG3	2.11	0.51
1:ZO:160:PHE:O	1:ZO:163:THR:HG22	2.11	0.51
1:YU:246:LEU:O	1:YU:266:ASP:HB3	2.11	0.51
1:ZC:284:MET:O	1:ZC:287:MET:HG2	2.10	0.51
1:ZH:177:TYR:CE2	1:ZH:179:PRO:HG3	2.45	0.51
1:ZR:332:LYS:HA	1:ZR:338:TYR:HB2	1.92	0.51
1:ZP:347:ARG:HB3	1:ZP:366:TRP:CZ2	2.46	0.51
1:ZV:332:LYS:HA	1:ZV:338:TYR:HB2	1.93	0.51
1:ZW:264:MET:HE3	1:ZW:265:ILE:H	1.75	0.51
1:YJ:232:ASN:O	1:YJ:238:GLY:HA3	2.11	0.51
1:YN:319:HIS:ND1	1:YN:320:ILE:HG12	2.26	0.51
1:YR:176:VAL:HG12	1:YR:197:THR:HG21	1.93	0.51
1:YY:225:ILE:HG21	1:YY:283:ASN:HB3	1.92	0.51
1:ZK:249:LEU:HD12	1:ZK:389:LYS:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZR:121:ARG:HG3	1:ZR:122:ASP:N	2.25	0.51
1:ZR:150:LEU:HD12	1:ZR:204:LEU:HD22	1.93	0.51
1:YI:316:PHE:HD2	1:YI:323:PHE:HE2	1.57	0.50
1:YX:214:LYS:HE2	1:YX:298:THR:HG22	1.94	0.50
1:ZC:160:PHE:O	1:ZC:163:THR:HG22	2.12	0.50
1:ZM:241:PRO:HG2	3:ZM:502:LMG:HC8	1.93	0.50
1:ZW:284:MET:O	1:ZW:287:MET:HG2	2.11	0.50
1:YJ:176:VAL:HG12	1:YJ:197:THR:HG21	1.93	0.50
1:YJ:319:HIS:ND1	1:YJ:320:ILE:HG12	2.26	0.50
1:YR:265:ILE:HD12	1:YR:278:ASP:HB3	1.93	0.50
1:YU:192:GLU:HG3	1:YU:263:ALA:O	2.12	0.50
1:ZB:87:ARG:HH12	1:ZB:218:TYR:HD2	1.59	0.50
1:ZN:332:LYS:HA	1:ZN:338:TYR:HB2	1.93	0.50
1:ZO:216:SER:OG	1:ZO:221:LYS:HD3	2.10	0.50
1:ZP:223:LEU:HB2	1:ZP:302:PHE:HD1	1.76	0.50
1:ZY:210:LEU:HD11	1:ZY:290:PHE:CE1	2.40	0.50
1:ZD:354:ASP:HB3	1:ZD:357:LEU:HG	1.93	0.50
1:ZQ:111:THR:HG22	1:ZQ:113:LYS:HG3	1.93	0.50
1:ZR:250:ARG:HA	1:ZR:253:ALA:H	1.76	0.50
1:ZV:145:LEU:HD12	1:ZV:156:PHE:CG	2.46	0.50
2:YV:501:NDP:H41N	4:YV:503:A1JWG:CBA	2.42	0.50
1:YZ:157:VAL:HG11	1:YZ:208:LEU:HB3	1.92	0.50
1:ZI:362:VAL:HG13	1:ZI:374:GLU:HG2	1.94	0.50
1:ZN:108:LEU:HD12	1:ZN:114:TRP:CD1	2.47	0.50
1:ZZ:108:LEU:HD12	1:ZZ:114:TRP:CD1	2.45	0.50
1:YO:214:LYS:HZ3	1:YO:298:THR:HA	1.77	0.50
1:YT:145:LEU:HD12	1:YT:156:PHE:CG	2.46	0.50
1:YW:159:ASN:HA	1:YW:162:ARG:NH1	2.27	0.50
1:ZD:167:LEU:HD21	1:ZD:170:LEU:HD21	1.93	0.50
1:ZH:167:LEU:HD21	1:ZH:170:LEU:HD21	1.93	0.50
1:ZI:214:LYS:NZ	1:ZI:298:THR:HA	2.26	0.50
1:ZW:289:GLU:HG2	1:ZW:391:TRP:HB2	1.94	0.50
1:YQ:216:SER:OG	1:YQ:221:LYS:HD3	2.12	0.50
1:ZP:395:GLU:HG2	1:ZP:400:LEU:HD23	1.94	0.50
1:ZZ:332:LYS:HA	1:ZZ:338:TYR:HB2	1.94	0.50
1:YV:118:MET:HB3	1:YV:142:VAL:HA	1.93	0.50
1:YU:284:MET:HA	1:YU:287:MET:HE2	1.94	0.50
1:YY:319:HIS:CG	1:YY:320:ILE:H	2.29	0.50
1:ZL:200:LEU:HD23	1:ZL:398:VAL:HG12	1.94	0.50
1:YN:173:ASN:HD21	2:YN:501:NDP:H4D	1.76	0.50
1:YQ:211:ASP:O	1:YQ:215:LYS:HB2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YQ:264:MET:SD	1:YQ:270:PHE:HA	2.52	0.50
1:YR:167:LEU:HD21	1:YR:170:LEU:HD21	1.93	0.50
1:ZD:232:ASN:O	1:ZD:238:GLY:HA3	2.12	0.50
1:ZK:249:LEU:HB2	1:ZK:389:LYS:HE3	1.94	0.50
1:YM:304:SER:HB3	1:YM:363:TYR:HA	1.94	0.49
1:YN:223:LEU:HB2	1:YN:302:PHE:HD1	1.77	0.49
1:YY:160:PHE:O	1:YY:163:THR:HG22	2.11	0.49
1:YY:284:MET:O	1:YY:287:MET:HG2	2.11	0.49
1:YY:339:VAL:HG21	1:YY:347:ARG:HH12	1.77	0.49
1:ZH:276:TYR:HE1	1:ZH:280:LYS:HE2	1.75	0.49
1:ZM:229:ILE:HD11	1:ZM:235:THR:HG21	1.93	0.49
1:YI:88:LYS:HE2	1:YI:112:GLY:O	2.12	0.49
1:YV:177:TYR:CE2	1:YV:179:PRO:HG3	2.47	0.49
1:YZ:230:THR:CG2	1:YZ:277:LYS:HE3	2.43	0.49
1:ZH:281:VAL:O	1:ZH:285:LEU:HD23	2.13	0.49
1:ZL:284:MET:HA	1:ZL:287:MET:HE2	1.93	0.49
1:YP:129:ALA:O	1:YP:133:VAL:HG12	2.12	0.49
1:YR:250:ARG:HG3	1:YR:251:GLY:N	2.27	0.49
1:YT:124:LEU:HD21	1:YT:128:ARG:HH21	1.78	0.49
1:YW:218:TYR:HD2	1:YW:220:SER:H	1.59	0.49
1:YY:264:MET:SD	1:YY:270:PHE:HA	2.52	0.49
1:ZI:343:GLU:HA	1:ZI:346:LYS:HG2	1.94	0.49
1:ZN:385:GLU:O	1:ZN:389:LYS:HG2	2.12	0.49
1:ZO:304:SER:HB3	1:ZO:363:TYR:HA	1.94	0.49
1:ZP:232:ASN:O	1:ZP:238:GLY:HA3	2.12	0.49
1:ZP:339:VAL:HG11	1:ZP:347:ARG:NH2	2.28	0.49
1:ZP:396:LYS:HD2	1:ZP:397:LEU:HD22	1.95	0.49
1:YI:160:PHE:O	1:YI:163:THR:HG22	2.13	0.49
1:YQ:225:ILE:HG21	1:YQ:283:ASN:HB3	1.94	0.49
1:YV:232:ASN:O	1:YV:238:GLY:HA3	2.13	0.49
1:YX:145:LEU:HD12	1:YX:156:PHE:CG	2.47	0.49
1:ZD:162:ARG:HE	1:ZI:162:ARG:NH1	2.05	0.49
1:ZD:226:VAL:HG22	1:ZD:305:LEU:HD21	1.94	0.49
1:ZJ:145:LEU:HD12	1:ZJ:156:PHE:CG	2.47	0.49
1:ZK:233:THR:HG23	1:ZK:243:LYS:HD2	1.94	0.49
1:ZT:167:LEU:HD21	1:ZT:170:LEU:HD21	1.94	0.49
1:ZW:319:HIS:CG	1:ZW:320:ILE:H	2.30	0.49
1:YT:246:LEU:HD23	1:YT:249:LEU:HD21	1.94	0.49
1:ZL:218:TYR:CD1	1:ZL:219:PRO:HD2	2.48	0.49
1:ZO:250:ARG:HB3	1:ZO:267:GLY:HA3	1.95	0.49
1:ZO:284:MET:O	1:ZO:287:MET:HG2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZY:362:VAL:HG13	1:ZY:374:GLU:HG2	1.95	0.49
1:YN:177:TYR:CE2	1:YN:179:PRO:HG3	2.48	0.49
1:YQ:276:TYR:O	1:YQ:280:LYS:HG2	2.12	0.49
1:YS:159:ASN:O	1:YS:162:ARG:HG2	2.11	0.49
1:YU:221:LYS:O	1:YU:222:ARG:HD3	2.12	0.49
1:YU:230:THR:HG22	1:YU:277:LYS:HD3	1.95	0.49
1:YV:88:LYS:HA	1:YV:113:LYS:HZ3	1.78	0.49
1:ZM:389:LYS:O	1:ZM:393:ILE:HG13	2.13	0.49
1:YJ:177:TYR:CE2	1:YJ:179:PRO:HG3	2.48	0.49
1:YO:229:ILE:HD11	1:YO:235:THR:HG21	1.95	0.49
1:ZA:362:VAL:HG13	1:ZA:374:GLU:HG2	1.94	0.49
1:ZB:385:GLU:O	1:ZB:389:LYS:HG2	2.13	0.49
1:ZK:319:HIS:CG	1:ZK:320:ILE:H	2.31	0.49
1:ZL:162:ARG:HD2	1:ZQ:162:ARG:HH12	1.77	0.49
1:ZP:276:TYR:CE1	1:ZP:280:LYS:HE2	2.47	0.49
1:YJ:95:GLY:HA3	1:YJ:174:ALA:HB2	1.94	0.49
1:YO:389:LYS:O	1:YO:393:ILE:HG13	2.13	0.49
1:YX:250:ARG:HE	1:ZB:106:LYS:HZ3	1.61	0.49
1:YY:289:GLU:HG2	1:YY:391:TRP:HB2	1.94	0.49
1:ZC:284:MET:HA	1:ZC:287:MET:HE2	1.94	0.49
1:ZR:123:PHE:O	1:ZR:126:ALA:N	2.46	0.49
1:ZS:304:SER:HB3	1:ZS:363:TYR:HA	1.94	0.49
1:ZU:389:LYS:O	1:ZU:393:ILE:HG13	2.12	0.49
1:ZW:181:ALA:HB3	1:ZW:272:GLY:HA3	1.94	0.49
1:ZD:265:ILE:HD12	1:ZD:278:ASP:HB3	1.95	0.49
1:ZG:246:LEU:HD23	1:ZG:249:LEU:HD21	1.95	0.49
1:ZK:216:SER:OG	1:ZK:221:LYS:HD3	2.13	0.49
1:ZP:176:VAL:HG12	1:ZP:197:THR:HG21	1.95	0.49
1:ZP:339:VAL:HG11	1:ZP:347:ARG:NH1	2.28	0.49
1:ZW:160:PHE:CD2	1:ZW:209:LEU:HD21	2.48	0.49
1:ZT:265:ILE:HD12	1:ZT:278:ASP:HB3	1.95	0.49
1:ZX:249:LEU:HD12	1:ZX:389:LYS:HB2	1.95	0.49
1:YK:106:LYS:HE2	1:ZY:250:ARG:HE	1.77	0.48
1:YO:362:VAL:HG13	1:YO:374:GLU:HG2	1.95	0.48
1:ZU:210:LEU:HD11	1:ZU:290:PHE:CE1	2.43	0.48
1:YI:95:GLY:HA2	2:YI:501:NDP:H1B	1.95	0.48
1:YN:226:VAL:HG22	1:YN:305:LEU:HD21	1.95	0.48
1:YN:287:MET:HE2	1:YN:304:SER:N	2.28	0.48
1:YU:225:ILE:HG21	1:YU:283:ASN:HB3	1.94	0.48
1:YW:214:LYS:HE2	1:YW:298:THR:HG22	1.95	0.48
1:ZH:343:GLU:O	1:ZH:347:ARG:HG3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZP:276:TYR:HE1	1:ZP:280:LYS:HE2	1.78	0.48
1:ZT:177:TYR:CE2	1:ZT:179:PRO:HG3	2.47	0.48
1:ZT:287:MET:HE2	1:ZT:304:SER:N	2.28	0.48
1:ZX:160:PHE:O	1:ZX:163:THR:HG22	2.13	0.48
1:YR:311:ALA:O	1:YR:328:PRO:HG3	2.13	0.48
1:YU:319:HIS:CG	1:YU:320:ILE:H	2.31	0.48
1:YV:276:TYR:HE1	1:YV:280:LYS:HE2	1.77	0.48
1:YX:184:PRO:HB2	1:YX:186:TYR:CE1	2.49	0.48
1:ZH:176:VAL:HG12	1:ZH:197:THR:HG21	1.95	0.48
1:ZN:250:ARG:HH12	1:ZR:134:GLY:HA2	1.43	0.48
1:ZP:250:ARG:HG3	1:ZP:251:GLY:N	2.27	0.48
1:ZT:176:VAL:HG12	1:ZT:197:THR:HG21	1.95	0.48
1:YN:177:TYR:CD2	1:YN:179:PRO:HG3	2.49	0.48
1:YR:232:ASN:O	1:YR:238:GLY:HA3	2.12	0.48
1:YS:362:VAL:HG13	1:YS:374:GLU:HG2	1.95	0.48
1:YZ:220:SER:HB3	1:YZ:222:ARG:NH1	2.27	0.48
1:ZA:214:LYS:HD2	1:ZA:214:LYS:HA	1.63	0.48
1:ZD:276:TYR:HE1	1:ZD:280:LYS:HE2	1.78	0.48
1:ZH:223:LEU:HB2	1:ZH:302:PHE:HD1	1.77	0.48
1:ZH:265:ILE:HD12	1:ZH:278:ASP:HB3	1.95	0.48
1:YN:150:LEU:HB3	1:YN:154:ARG:HH12	1.78	0.48
1:YN:162:ARG:HG2	1:YS:162:ARG:HH22	1.79	0.48
1:YV:291:HIS:CD2	1:YV:360:SER:HB3	2.49	0.48
1:ZH:160:PHE:CZ	1:ZH:167:LEU:HB2	2.49	0.48
1:ZH:250:ARG:HG3	1:ZH:251:GLY:N	2.27	0.48
1:ZJ:332:LYS:HA	1:ZJ:338:TYR:HB2	1.95	0.48
1:ZO:225:ILE:HG21	1:ZO:283:ASN:HB3	1.96	0.48
1:YJ:177:TYR:CD2	1:YJ:179:PRO:HG3	2.48	0.48
1:YN:176:VAL:HG12	1:YN:197:THR:HG21	1.96	0.48
1:ZD:287:MET:HE2	1:ZD:304:SER:N	2.29	0.48
1:ZE:214:LYS:NZ	1:ZE:298:THR:HA	2.27	0.48
1:ZE:343:GLU:HA	1:ZE:346:LYS:HG2	1.96	0.48
1:ZJ:246:LEU:HB3	1:ZJ:249:LEU:HD11	1.96	0.48
1:YO:210:LEU:HD11	1:YO:290:PHE:CE1	2.42	0.48
1:YV:167:LEU:HD21	1:YV:170:LEU:HD21	1.95	0.48
1:ZD:177:TYR:CE2	1:ZD:179:PRO:HG3	2.49	0.48
1:ZT:177:TYR:CD2	1:ZT:179:PRO:HG3	2.49	0.48
1:ZW:233:THR:HG23	1:ZW:243:LYS:HD2	1.95	0.48
1:ZX:220:SER:HB3	1:ZX:222:ARG:NH1	2.25	0.48
1:ZY:389:LYS:O	1:ZY:393:ILE:HG13	2.14	0.48
1:YI:225:ILE:HG21	1:YI:283:ASN:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YU:88:LYS:HE3	1:YU:112:GLY:C	2.39	0.48
1:YX:250:ARG:CG	1:ZB:134:GLY:HA2	2.43	0.48
1:YZ:160:PHE:O	1:YZ:163:THR:HG22	2.14	0.48
1:ZD:281:VAL:O	1:ZD:285:LEU:HD23	2.14	0.48
1:YN:305:LEU:HD11	1:YN:348:LEU:HG	1.94	0.48
1:YR:162:ARG:HD2	1:YW:162:ARG:HE	1.78	0.48
1:YW:362:VAL:HG13	1:YW:374:GLU:HG2	1.95	0.48
1:ZH:311:ALA:O	1:ZH:328:PRO:HG3	2.14	0.48
1:YO:264:MET:HE2	1:YO:271:ASP:N	2.21	0.48
1:YW:241:PRO:HG2	3:YW:502:LMG:HC8	1.96	0.48
1:ZD:276:TYR:CE1	1:ZD:280:LYS:HE2	2.49	0.48
1:ZF:332:LYS:HA	1:ZF:338:TYR:HB2	1.95	0.48
1:ZN:145:LEU:HD12	1:ZN:156:PHE:CG	2.49	0.48
1:ZS:88:LYS:HE3	1:ZS:112:GLY:C	2.39	0.48
1:YJ:250:ARG:HG3	1:YJ:251:GLY:N	2.29	0.47
1:ZD:250:ARG:HG3	1:ZD:251:GLY:N	2.28	0.47
1:ZM:135:MET:HE3	1:ZM:135:MET:HB3	1.76	0.47
1:YL:250:ARG:HE	1:YP:106:LYS:HZ3	1.61	0.47
1:YN:311:ALA:O	1:YN:328:PRO:HG3	2.14	0.47
1:YU:157:VAL:HG11	1:YU:208:LEU:HG	1.94	0.47
1:YU:276:TYR:O	1:YU:280:LYS:HG2	2.13	0.47
1:ZB:124:LEU:HD21	1:ZB:128:ARG:HH21	1.79	0.47
1:ZH:210:LEU:HD13	1:ZH:290:PHE:HZ	1.79	0.47
1:ZH:347:ARG:HD2	1:ZH:366:TRP:CD2	2.49	0.47
1:ZL:220:SER:HB3	1:ZL:222:ARG:NH1	2.28	0.47
1:ZN:124:LEU:HD21	1:ZN:128:ARG:HH21	1.79	0.47
1:ZS:379:GLU:HG2	1:ZS:380:GLU:N	2.29	0.47
1:ZU:214:LYS:HD2	1:ZU:214:LYS:HA	1.63	0.47
1:ZV:249:LEU:HD21	1:ZV:390:VAL:HG23	1.97	0.47
1:ZW:284:MET:HA	1:ZW:287:MET:HE2	1.96	0.47
1:ZW:329:PRO:HA	1:ZW:332:LYS:HB2	1.96	0.47
1:YM:199:HIS:CD2	1:YM:199:HIS:C	2.92	0.47
1:YT:96:ALA:HA	1:YT:101:GLY:HA3	1.97	0.47
1:YV:347:ARG:HD2	1:YV:366:TRP:CD2	2.49	0.47
1:YX:250:ARG:HH22	1:ZB:133:VAL:HG23	1.79	0.47
1:ZF:145:LEU:HD12	1:ZF:156:PHE:CG	2.48	0.47
1:ZH:95:GLY:HA3	1:ZH:174:ALA:HB2	1.96	0.47
1:ZN:250:ARG:HD3	1:ZR:135:MET:HG2	1.96	0.47
1:ZT:250:ARG:HG3	1:ZT:251:GLY:N	2.29	0.47
1:ZY:199:HIS:CE1	1:ZY:282:CYS:HB3	2.49	0.47
1:YI:287:MET:HE3	1:YI:304:SER:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YK:212:ASP:HA	1:YK:215:LYS:HE2	1.95	0.47
1:YY:284:MET:HA	1:YY:287:MET:HE2	1.96	0.47
1:ZB:145:LEU:HD12	1:ZB:156:PHE:CG	2.49	0.47
1:ZD:311:ALA:O	1:ZD:328:PRO:HG3	2.15	0.47
1:ZM:214:LYS:HD2	1:ZM:214:LYS:HA	1.58	0.47
1:ZS:246:LEU:HD23	1:ZS:249:LEU:HD21	1.96	0.47
1:ZZ:173:ASN:OD1	2:ZZ:501:NDP:H4D	2.14	0.47
1:YJ:319:HIS:HD1	1:YJ:320:ILE:HG12	1.80	0.47
1:YO:395:GLU:HG2	1:YO:401:ALA:HB2	1.96	0.47
1:YP:124:LEU:HD21	1:YP:128:ARG:HH21	1.77	0.47
1:YS:389:LYS:O	1:YS:393:ILE:HG13	2.14	0.47
1:YX:250:ARG:HH12	1:ZB:133:VAL:CA	2.27	0.47
1:ZC:225:ILE:HG21	1:ZC:283:ASN:HB3	1.96	0.47
1:ZO:276:TYR:O	1:ZO:280:LYS:HG2	2.14	0.47
1:ZP:265:ILE:HD12	1:ZP:278:ASP:HB3	1.96	0.47
1:ZP:311:ALA:O	1:ZP:328:PRO:HG3	2.14	0.47
1:YM:246:LEU:O	1:YM:266:ASP:HB3	2.15	0.47
1:ZH:339:VAL:HG11	1:ZH:347:ARG:NH1	2.30	0.47
1:ZJ:250:ARG:NH1	1:ZN:133:VAL:N	2.63	0.47
1:ZP:167:LEU:HD21	1:ZP:170:LEU:HD21	1.97	0.47
1:YK:124:LEU:HD21	1:YK:128:ARG:HH21	1.80	0.47
1:YL:124:LEU:HD21	1:YL:128:ARG:HH21	1.79	0.47
1:YL:134:GLY:HA2	1:ZZ:250:ARG:CG	2.41	0.47
1:YR:118:MET:HB3	1:YR:142:VAL:HA	1.97	0.47
1:YY:379:GLU:HG2	1:YY:380:GLU:N	2.29	0.47
1:YZ:159:ASN:HA	1:ZE:162:ARG:HH12	1.80	0.47
1:YZ:252:LEU:HB2	1:YZ:265:ILE:HG23	1.97	0.47
1:ZA:214:LYS:NZ	1:ZA:298:THR:HA	2.29	0.47
1:ZC:255:GLY:HA2	1:ZC:397:LEU:HD21	1.96	0.47
1:ZC:349:ALA:HA	1:ZC:352:VAL:HG22	1.96	0.47
1:ZF:124:LEU:HD21	1:ZF:128:ARG:HH21	1.80	0.47
1:ZH:160:PHE:HZ	1:ZH:167:LEU:HB2	1.79	0.47
1:ZH:232:ASN:O	1:ZH:238:GLY:HA3	2.14	0.47
1:ZH:287:MET:HE2	1:ZH:304:SER:N	2.29	0.47
1:ZK:289:GLU:HG2	1:ZK:391:TRP:HB2	1.96	0.47
1:ZK:379:GLU:HG2	1:ZK:380:GLU:N	2.29	0.47
1:ZL:251:GLY:HA3	1:ZL:264:MET:O	2.15	0.47
1:ZL:343:GLU:O	1:ZL:346:LYS:HG2	2.14	0.47
1:ZR:100:LEU:HD11	1:ZR:310:ILE:HD13	1.97	0.47
1:ZS:320:ILE:HG22	1:ZS:322:LEU:H	1.80	0.47
1:ZT:232:ASN:O	1:ZT:238:GLY:HA3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZX:230:THR:CG2	1:ZX:277:LYS:HE3	2.42	0.47
1:YM:276:TYR:O	1:YM:280:LYS:HG2	2.14	0.47
1:YN:276:TYR:CE1	1:YN:280:LYS:HE2	2.50	0.47
1:YQ:158:ASP:O	1:YQ:162:ARG:HG3	2.15	0.47
1:YQ:379:GLU:HG2	1:YQ:380:GLU:N	2.30	0.47
1:YT:173:ASN:OD1	2:YT:501:NDP:H4D	2.14	0.47
1:YU:320:ILE:HG22	1:YU:322:LEU:H	1.79	0.47
1:YZ:280:LYS:HD3	1:YZ:283:ASN:HD22	1.79	0.47
1:YZ:280:LYS:HD3	1:YZ:283:ASN:ND2	2.30	0.47
1:ZM:264:MET:HE2	1:ZM:271:ASP:N	2.23	0.47
1:ZP:96:ALA:HB3	1:ZP:118:MET:SD	2.55	0.47
1:YL:145:LEU:HD12	1:YL:156:PHE:CG	2.50	0.47
1:YN:395:GLU:HG2	1:YN:400:LEU:HD23	1.97	0.47
1:YR:177:TYR:CE2	1:YR:179:PRO:HG3	2.50	0.47
1:YV:265:ILE:HD12	1:YV:278:ASP:HB3	1.95	0.47
1:YW:229:ILE:HD11	1:YW:235:THR:HG21	1.97	0.47
1:YX:106:LYS:HD2	1:YX:133:VAL:HG22	1.97	0.47
1:ZH:173:ASN:OD1	2:ZH:501:NDP:H4D	2.15	0.47
1:ZN:250:ARG:CD	1:ZR:135:MET:HG2	2.44	0.47
1:ZO:252:LEU:HB2	1:ZO:265:ILE:HG23	1.97	0.47
1:ZX:292:ARG:HD2	1:ZX:388:ARG:HH12	1.79	0.47
1:ZZ:145:LEU:HD12	1:ZZ:156:PHE:CG	2.50	0.47
1:YJ:311:ALA:O	1:YJ:328:PRO:HG3	2.14	0.47
1:YP:332:LYS:HA	1:YP:338:TYR:HB2	1.96	0.47
1:YT:246:LEU:HB3	1:YT:249:LEU:HD21	1.97	0.47
1:YW:199:HIS:CE1	1:YW:282:CYS:HB3	2.50	0.47
1:YY:304:SER:HB3	1:YY:363:TYR:HA	1.96	0.47
1:ZE:88:LYS:HG2	1:ZE:113:LYS:HA	1.95	0.47
1:ZN:246:LEU:HD21	1:ZN:249:LEU:HD11	1.96	0.47
1:ZX:290:PHE:HB3	1:ZX:302:PHE:CE2	2.50	0.47
1:YL:250:ARG:HA	1:YL:253:ALA:H	1.80	0.46
1:YY:233:THR:HG23	1:YY:243:LYS:HD2	1.97	0.46
1:ZD:339:VAL:HG21	1:ZD:347:ARG:HH12	1.80	0.46
1:ZE:264:MET:HE2	1:ZE:271:ASP:H	1.79	0.46
1:ZF:135:MET:HE3	1:ZF:135:MET:HB3	1.77	0.46
1:ZP:354:ASP:HB3	1:ZP:357:LEU:HG	1.97	0.46
1:ZS:289:GLU:HG2	1:ZS:391:TRP:HB2	1.96	0.46
1:ZW:212:ASP:HA	1:ZW:215:LYS:HB3	1.97	0.46
1:YL:332:LYS:HA	1:YL:338:TYR:HB2	1.96	0.46
1:ZN:124:LEU:HD21	1:ZN:128:ARG:NH2	2.30	0.46
1:ZS:276:TYR:O	1:ZS:280:LYS:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZU:214:LYS:NZ	1:ZU:298:THR:HA	2.30	0.46
1:YJ:210:LEU:HD13	1:YJ:290:PHE:HZ	1.79	0.46
1:YM:230:THR:HG22	1:YM:277:LYS:HD3	1.97	0.46
1:YN:250:ARG:HG3	1:YN:251:GLY:N	2.29	0.46
1:YN:307:PRO:HG3	1:YN:348:LEU:HD11	1.98	0.46
1:YY:349:ALA:HA	1:YY:352:VAL:HG22	1.96	0.46
1:YZ:290:PHE:HB3	1:YZ:302:PHE:CZ	2.50	0.46
2:YZ:501:NDP:H41N	4:YZ:503:A1JWG:O1A	2.15	0.46
1:ZD:347:ARG:HB3	1:ZD:366:TRP:CZ2	2.49	0.46
1:ZH:177:TYR:CD2	1:ZH:179:PRO:HG3	2.50	0.46
1:ZK:276:TYR:O	1:ZK:280:LYS:HG2	2.15	0.46
1:ZP:177:TYR:CE2	1:ZP:179:PRO:HG3	2.51	0.46
1:ZQ:159:ASN:O	1:ZQ:162:ARG:HG2	2.15	0.46
1:ZR:385:GLU:O	1:ZR:389:LYS:HG2	2.16	0.46
1:YN:118:MET:HB3	1:YN:142:VAL:HA	1.97	0.46
1:YS:199:HIS:CE1	1:YS:282:CYS:HB3	2.50	0.46
1:YS:343:GLU:HA	1:YS:346:LYS:HG2	1.97	0.46
1:YX:249:LEU:HD11	1:YX:386:LYS:HG3	1.95	0.46
1:ZG:276:TYR:O	1:ZG:280:LYS:HG2	2.14	0.46
1:ZM:199:HIS:CE1	1:ZM:282:CYS:HB3	2.50	0.46
1:YI:276:TYR:O	1:YI:280:LYS:HG2	2.15	0.46
1:YJ:276:TYR:HE1	1:YJ:280:LYS:HE2	1.81	0.46
1:YR:396:LYS:HD2	1:YR:397:LEU:HD22	1.96	0.46
1:YT:214:LYS:HZ1	1:YT:298:THR:HA	1.79	0.46
1:YU:289:GLU:HB2	1:YU:387:ALA:HB1	1.97	0.46
1:ZA:210:LEU:HD11	1:ZA:290:PHE:CE1	2.41	0.46
1:ZM:210:LEU:HD11	1:ZM:290:PHE:CE1	2.41	0.46
1:ZR:250:ARG:NH2	1:ZV:133:VAL:CG1	2.79	0.46
1:ZW:276:TYR:O	1:ZW:280:LYS:HG2	2.15	0.46
1:ZZ:124:LEU:HD21	1:ZZ:128:ARG:HH21	1.81	0.46
1:YQ:145:LEU:HD22	1:YQ:156:PHE:CD2	2.51	0.46
1:YU:379:GLU:HG2	1:YU:380:GLU:N	2.30	0.46
1:YW:210:LEU:HD11	1:YW:290:PHE:CE1	2.42	0.46
1:YX:124:LEU:HD21	1:YX:128:ARG:NH2	2.31	0.46
1:ZF:150:LEU:HD12	1:ZF:204:LEU:CD1	2.46	0.46
1:ZL:292:ARG:HH12	1:ZL:384:VAL:HA	1.80	0.46
1:ZZ:346:LYS:HE3	1:ZZ:346:LYS:HB2	1.82	0.46
1:YQ:366:TRP:NE1	1:YQ:372:SER:HB3	2.31	0.46
1:YU:160:PHE:CD2	1:YU:209:LEU:HD21	2.51	0.46
1:YY:368:ASN:HD22	1:YY:368:ASN:H	1.63	0.46
1:ZB:214:LYS:NZ	1:ZB:298:THR:HA	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZG:181:ALA:HB3	1:ZG:272:GLY:HA3	1.97	0.46
1:ZT:96:ALA:HB3	1:ZT:118:MET:SD	2.55	0.46
1:ZT:311:ALA:O	1:ZT:328:PRO:HG3	2.15	0.46
1:ZZ:214:LYS:HD2	1:ZZ:214:LYS:HA	1.73	0.46
1:YM:274:LYS:HE3	1:YM:274:LYS:HB3	1.78	0.46
1:YN:265:ILE:HD12	1:YN:278:ASP:HB3	1.97	0.46
1:YO:116:VAL:O	1:YO:140:TYR:HA	2.15	0.46
1:ZA:96:ALA:O	1:ZA:118:MET:HE3	2.16	0.46
1:ZB:133:VAL:O	1:ZB:135:MET:HG2	2.15	0.46
1:ZN:191:PHE:HA	1:ZN:256:LEU:HD21	1.97	0.46
1:ZR:87:ARG:HH21	1:ZR:218:TYR:HD2	1.64	0.46
1:ZS:249:LEU:HB2	1:ZS:389:LYS:HE3	1.96	0.46
1:ZW:160:PHE:O	1:ZW:163:THR:HG22	2.16	0.46
1:ZA:169:VAL:HG11	1:ZA:352:VAL:HG13	1.98	0.46
1:ZO:91:VAL:HG21	1:ZO:114:TRP:CE3	2.51	0.46
1:ZO:253:ALA:HA	1:ZO:393:ILE:HG21	1.98	0.46
1:ZS:319:HIS:CG	1:ZS:320:ILE:H	2.34	0.46
1:ZU:362:VAL:HG13	1:ZU:374:GLU:HG2	1.98	0.46
1:ZX:343:GLU:HA	1:ZX:346:LYS:HG3	1.98	0.46
1:YM:264:MET:HE3	1:YM:265:ILE:H	1.81	0.46
1:YN:167:LEU:HD21	1:YN:170:LEU:HD21	1.98	0.46
1:YY:328:PRO:O	1:YY:332:LYS:HG3	2.16	0.46
1:YZ:339:VAL:HG13	1:YZ:343:GLU:OE2	2.16	0.46
1:ZM:157:VAL:O	1:ZM:161:ARG:HG3	2.16	0.46
1:ZR:343:GLU:HA	1:ZR:346:LYS:HG2	1.98	0.46
1:YU:234:ASN:ND2	1:YU:378:SER:HA	2.31	0.45
1:YY:112:GLY:O	1:YY:113:LYS:HD2	2.16	0.45
1:YY:276:TYR:O	1:YY:280:LYS:HG2	2.15	0.45
1:ZX:343:GLU:O	1:ZX:346:LYS:HG3	2.15	0.45
1:YJ:265:ILE:HD12	1:YJ:278:ASP:HB3	1.98	0.45
1:YQ:218:TYR:CD1	1:YQ:219:PRO:HD2	2.51	0.45
1:YU:145:LEU:HD22	1:YU:156:PHE:CD2	2.52	0.45
1:YY:195:VAL:HG21	1:YY:265:ILE:HD11	1.98	0.45
1:ZC:276:TYR:O	1:ZC:280:LYS:HG2	2.15	0.45
1:ZF:249:LEU:HD11	1:ZF:386:LYS:HG3	1.99	0.45
1:ZK:284:MET:HA	1:ZK:287:MET:HE2	1.96	0.45
1:YM:379:GLU:HG2	1:YM:380:GLU:N	2.31	0.45
1:YV:343:GLU:O	1:YV:347:ARG:HG3	2.17	0.45
1:YX:218:TYR:CD1	1:YX:219:PRO:HD2	2.51	0.45
1:YZ:213:LEU:HD23	1:YZ:213:LEU:HA	1.81	0.45
1:ZK:289:GLU:HB2	1:ZK:387:ALA:HB1	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZS:216:SER:OG	1:ZS:221:LYS:HD3	2.17	0.45
2:ZX:501:NDP:H41N	4:ZX:503:A1JWG:O1A	2.15	0.45
1:ZY:207:ARG:HE	1:ZY:207:ARG:HB3	1.64	0.45
1:YM:100:LEU:O	1:YM:104:THR:HG22	2.16	0.45
1:YX:262:SER:HB3	1:YX:267:GLY:HA2	1.99	0.45
1:YZ:322:LEU:HD23	1:YZ:322:LEU:HA	1.82	0.45
1:ZA:229:ILE:HD11	1:ZA:235:THR:HG21	1.97	0.45
1:ZJ:346:LYS:HE3	1:ZJ:346:LYS:HB2	1.85	0.45
1:ZN:191:PHE:CE2	1:ZN:397:LEU:HD21	2.50	0.45
1:YP:145:LEU:HD12	1:YP:156:PHE:CG	2.51	0.45
1:YY:329:PRO:HA	1:YY:332:LYS:HD3	1.99	0.45
1:ZC:264:MET:SD	1:ZC:270:PHE:HA	2.56	0.45
1:ZM:218:TYR:HD1	1:ZM:220:SER:H	1.64	0.45
1:ZO:274:LYS:HE3	1:ZO:274:LYS:HB3	1.78	0.45
1:ZR:145:LEU:HD12	1:ZR:156:PHE:CG	2.51	0.45
1:ZS:274:LYS:HE3	1:ZS:274:LYS:HB3	1.79	0.45
1:ZX:264:MET:HG3	1:ZX:269:ASP:O	2.17	0.45
1:YL:218:TYR:HD1	1:YL:220:SER:H	1.62	0.45
1:YL:385:GLU:O	1:YL:389:LYS:HG2	2.16	0.45
1:YV:100:LEU:HD12	1:YV:310:ILE:HD12	1.98	0.45
1:YX:332:LYS:HA	1:YX:338:TYR:HB2	1.97	0.45
1:ZB:214:LYS:HE2	1:ZB:298:THR:HG22	1.97	0.45
1:ZF:250:ARG:NE	1:ZJ:106:LYS:HZ3	1.97	0.45
1:ZL:160:PHE:O	1:ZL:163:THR:HG22	2.16	0.45
1:ZU:104:THR:HG21	1:ZU:348:LEU:HD22	1.99	0.45
1:YP:102:LEU:HD21	1:YP:133:VAL:HG11	1.97	0.45
1:YQ:88:LYS:HG3	1:YQ:113:LYS:C	2.41	0.45
1:ZA:214:LYS:HE2	1:ZA:298:THR:HG22	1.99	0.45
1:ZP:215:LYS:HB2	1:ZP:215:LYS:HE2	1.86	0.45
1:ZP:343:GLU:O	1:ZP:347:ARG:HG3	2.17	0.45
1:YL:133:VAL:HG13	1:YL:133:VAL:O	2.17	0.45
1:YL:367:ASN:OD1	1:YL:371:ALA:HB3	2.17	0.45
1:YW:116:VAL:O	1:YW:140:TYR:HA	2.17	0.45
1:ZL:320:ILE:HG22	1:ZL:322:LEU:H	1.82	0.45
1:ZN:346:LYS:HE3	1:ZN:346:LYS:HB2	1.76	0.45
1:YJ:354:ASP:HB3	1:YJ:357:LEU:HG	1.99	0.45
1:YX:214:LYS:NZ	1:YX:298:THR:HA	2.31	0.45
1:ZK:366:TRP:NE1	1:ZK:372:SER:HB3	2.32	0.45
1:ZO:221:LYS:C	1:ZO:222:ARG:HG3	2.42	0.45
1:ZR:127:GLU:HG3	1:ZR:128:ARG:N	2.32	0.45
1:YP:167:LEU:HD12	1:YP:168:ASP:H	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YP:246:LEU:HB3	1:YP:249:LEU:HD21	1.97	0.45
1:YZ:319:HIS:CG	1:YZ:320:ILE:H	2.35	0.45
1:ZG:366:TRP:NE1	1:ZG:372:SER:HB3	2.32	0.45
1:ZJ:167:LEU:HD12	1:ZJ:168:ASP:H	1.81	0.45
1:ZU:169:VAL:HG11	1:ZU:352:VAL:HG13	1.99	0.45
1:ZX:284:MET:O	1:ZX:287:MET:HG3	2.16	0.45
1:YU:253:ALA:HA	1:YU:393:ILE:HG21	1.99	0.44
1:YV:176:VAL:HG12	1:YV:197:THR:HG21	1.98	0.44
1:ZD:176:VAL:HG12	1:ZD:197:THR:HG21	1.98	0.44
1:ZT:354:ASP:HB3	1:ZT:357:LEU:HG	1.99	0.44
1:ZT:366:TRP:HA	1:ZT:372:SER:HA	1.99	0.44
1:ZW:304:SER:HB3	1:ZW:363:TYR:HA	1.99	0.44
1:YJ:167:LEU:HD21	1:YJ:170:LEU:HD21	2.00	0.44
1:YT:108:LEU:HD12	1:YT:114:TRP:CD1	2.51	0.44
1:YV:276:TYR:CE1	1:YV:280:LYS:HE2	2.51	0.44
1:ZC:100:LEU:O	1:ZC:104:THR:HG22	2.18	0.44
1:ZY:214:LYS:HD2	1:ZY:214:LYS:HA	1.58	0.44
1:YJ:276:TYR:CE1	1:YJ:280:LYS:HE2	2.52	0.44
1:YY:366:TRP:NE1	1:YY:372:SER:HB3	2.33	0.44
1:YZ:292:ARG:HH12	1:YZ:384:VAL:HA	1.81	0.44
1:ZD:396:LYS:HD2	1:ZD:397:LEU:HD22	1.98	0.44
1:ZG:88:LYS:HD2	1:ZG:113:LYS:HA	1.98	0.44
1:ZJ:124:LEU:HD21	1:ZJ:128:ARG:HH21	1.83	0.44
1:ZL:167:LEU:HD12	1:ZL:168:ASP:H	1.83	0.44
1:ZO:178:PHE:HB3	1:ZO:181:ALA:HB2	1.99	0.44
1:ZY:327:PHE:HB3	1:ZY:328:PRO:HD3	2.00	0.44
1:YL:250:ARG:HB2	1:YL:253:ALA:HB3	2.00	0.44
1:YY:346:LYS:HA	1:YY:346:LYS:HD3	1.83	0.44
1:YZ:395:GLU:HB2	1:YZ:401:ALA:HB3	1.99	0.44
1:ZA:116:VAL:O	1:ZA:140:TYR:HA	2.18	0.44
1:ZC:249:LEU:HD12	1:ZC:389:LYS:HE3	2.00	0.44
1:ZH:210:LEU:HD13	1:ZH:290:PHE:CZ	2.53	0.44
1:ZO:319:HIS:CG	1:ZO:320:ILE:H	2.36	0.44
1:ZY:218:TYR:HD2	1:ZY:220:SER:H	1.65	0.44
1:YI:249:LEU:HB2	1:YI:389:LYS:HE3	2.00	0.44
1:YL:135:MET:HB3	1:YL:135:MET:HE2	1.55	0.44
1:ZC:366:TRP:NE1	1:ZC:372:SER:HB3	2.33	0.44
1:ZE:87:ARG:HH22	1:ZE:218:TYR:HB2	1.82	0.44
1:ZK:96:ALA:O	1:ZK:118:MET:HE3	2.17	0.44
1:ZV:250:ARG:NH1	1:ZZ:133:VAL:N	2.65	0.44
1:ZY:175:ALA:HB3	2:ZY:501:NDP:H3D	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YO:379:GLU:H	1:YO:379:GLU:CD	2.25	0.44
1:YU:181:ALA:HB3	1:YU:272:GLY:HA3	2.00	0.44
1:YV:366:TRP:HA	1:YV:372:SER:HA	2.00	0.44
1:YW:157:VAL:O	1:YW:161:ARG:HG3	2.18	0.44
1:YX:167:LEU:HD12	1:YX:168:ASP:H	1.82	0.44
1:YX:346:LYS:HE3	1:YX:346:LYS:HB2	1.85	0.44
1:YZ:253:ALA:HB2	1:YZ:393:ILE:HD13	2.00	0.44
1:ZE:199:HIS:CD2	1:ZE:282:CYS:HB3	2.53	0.44
1:ZH:276:TYR:CE1	1:ZH:280:LYS:HE2	2.52	0.44
1:ZO:131:LYS:HD3	1:ZO:131:LYS:HA	1.86	0.44
1:ZP:347:ARG:HD2	1:ZP:366:TRP:CD2	2.52	0.44
1:ZP:367:ASN:HB2	1:ZP:370:SER:O	2.17	0.44
1:ZX:157:VAL:HG11	1:ZX:208:LEU:HB3	1.98	0.44
1:ZB:332:LYS:HA	1:ZB:338:TYR:HB2	2.00	0.44
1:ZI:379:GLU:H	1:ZI:379:GLU:CD	2.25	0.44
1:ZO:230:THR:HG22	1:ZO:277:LYS:HD3	1.98	0.44
1:ZQ:135:MET:HE2	1:ZQ:135:MET:HB3	1.79	0.44
1:YK:389:LYS:O	1:YK:393:ILE:HG13	2.18	0.44
1:YU:366:TRP:NE1	1:YU:372:SER:HB3	2.32	0.44
1:YV:339:VAL:HG11	1:YV:347:ARG:CZ	2.48	0.44
1:YX:207:ARG:HE	1:YX:207:ARG:HB3	1.59	0.44
1:YY:100:LEU:O	1:YY:104:THR:HG22	2.17	0.44
1:ZA:117:ILE:HD12	1:ZA:160:PHE:HE2	1.83	0.44
1:ZI:327:PHE:HB3	1:ZI:328:PRO:HD3	2.00	0.44
1:ZK:160:PHE:O	1:ZK:163:THR:HG22	2.17	0.44
1:ZO:346:LYS:HA	1:ZO:346:LYS:HD3	1.85	0.44
1:ZV:167:LEU:HD12	1:ZV:168:ASP:H	1.83	0.44
1:ZX:211:ASP:O	1:ZX:214:LYS:HG2	2.18	0.44
1:ZY:159:ASN:O	1:ZY:162:ARG:HG2	2.17	0.44
1:YJ:229:ILE:HD12	1:YJ:229:ILE:HA	1.88	0.44
1:YJ:319:HIS:CE1	1:YJ:320:ILE:HG12	2.53	0.44
1:YK:100:LEU:HD21	1:YK:307:PRO:HG2	2.00	0.44
1:YL:167:LEU:HD12	1:YL:168:ASP:H	1.82	0.44
1:YV:311:ALA:O	1:YV:328:PRO:HG3	2.18	0.44
1:YV:339:VAL:HG11	1:YV:347:ARG:NH2	2.33	0.44
1:YX:124:LEU:HD21	1:YX:128:ARG:HH21	1.83	0.44
1:YY:88:LYS:HG3	1:YY:113:LYS:C	2.42	0.44
1:ZC:135:MET:HE3	1:ZC:135:MET:HB3	1.91	0.44
1:ZF:214:LYS:HZ1	1:ZF:298:THR:HA	1.83	0.44
1:ZN:167:LEU:HD12	1:ZN:168:ASP:H	1.82	0.44
1:ZU:229:ILE:HD11	1:ZU:235:THR:HG21	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZU:327:PHE:HB3	1:ZU:328:PRO:HD3	2.00	0.44
1:ZW:222:ARG:HG2	1:ZW:301:THR:HB	1.99	0.44
1:YR:347:ARG:HB3	1:YR:366:TRP:CZ2	2.53	0.43
2:YS:501:NDP:H41N	4:YS:503:A1JWG:O1A	2.18	0.43
1:YT:167:LEU:HD12	1:YT:168:ASP:H	1.82	0.43
1:YV:223:LEU:HB2	1:YV:302:PHE:HD1	1.83	0.43
1:ZF:250:ARG:HH21	1:ZJ:106:LYS:HD2	1.83	0.43
1:ZG:211:ASP:O	1:ZG:215:LYS:CB	2.58	0.43
2:ZI:501:NDP:H41N	4:ZI:503:A1JWG:O1A	2.16	0.43
1:ZT:223:LEU:HB2	1:ZT:302:PHE:HD1	1.83	0.43
1:YI:100:LEU:O	1:YI:104:THR:HG22	2.18	0.43
1:YN:315:LEU:HB2	2:YN:501:NDP:O1N	2.18	0.43
1:YS:214:LYS:HZ2	1:YS:298:THR:HA	1.83	0.43
1:ZJ:262:SER:HB3	1:ZJ:267:GLY:HA2	2.00	0.43
2:ZL:501:NDP:H41N	4:ZL:503:A1JWG:O1A	2.17	0.43
1:ZP:339:VAL:HG11	1:ZP:347:ARG:CZ	2.48	0.43
1:ZR:108:LEU:HD12	1:ZR:114:TRP:CD1	2.52	0.43
1:YP:346:LYS:HE3	1:YP:346:LYS:HB2	1.89	0.43
1:YT:159:ASN:O	1:YT:162:ARG:HG2	2.18	0.43
1:YU:249:LEU:HB2	1:YU:389:LYS:HE3	2.01	0.43
1:YU:253:ALA:HA	1:YU:393:ILE:HD13	2.00	0.43
1:ZE:363:TYR:CD2	1:ZE:375:ASN:HB3	2.53	0.43
1:ZG:230:THR:HG22	1:ZG:277:LYS:HD3	2.00	0.43
1:ZK:100:LEU:O	1:ZK:104:THR:HG22	2.18	0.43
1:ZN:367:ASN:OD1	1:ZN:371:ALA:HB3	2.19	0.43
1:ZR:124:LEU:O	1:ZR:124:LEU:HD23	2.19	0.43
1:ZW:289:GLU:HB2	1:ZW:387:ALA:HB1	1.99	0.43
1:ZW:379:GLU:HG2	1:ZW:380:GLU:N	2.32	0.43
1:ZX:182:LYS:H	1:ZX:182:LYS:HD2	1.82	0.43
1:ZY:379:GLU:H	1:ZY:379:GLU:CD	2.26	0.43
1:YI:246:LEU:O	1:YI:266:ASP:HB3	2.19	0.43
1:YI:366:TRP:NE1	1:YI:372:SER:HB3	2.33	0.43
1:YJ:366:TRP:HA	1:YJ:372:SER:HA	1.99	0.43
1:YM:213:LEU:HD12	1:YM:221:LYS:HD2	1.99	0.43
1:YR:177:TYR:CD2	1:YR:179:PRO:HG3	2.53	0.43
1:YS:327:PHE:HB3	1:YS:328:PRO:HD3	2.00	0.43
1:YW:264:MET:CE	1:YW:271:ASP:H	2.28	0.43
1:YZ:343:GLU:O	1:YZ:346:LYS:HG2	2.18	0.43
1:ZB:124:LEU:HD21	1:ZB:128:ARG:NH2	2.34	0.43
1:ZB:346:LYS:HE3	1:ZB:346:LYS:HB2	1.90	0.43
1:ZC:280:LYS:HA	1:ZC:280:LYS:HD2	1.87	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZD:285:LEU:HD12	1:ZD:387:ALA:HB2	2.00	0.43
1:ZH:339:VAL:HG11	1:ZH:347:ARG:CZ	2.48	0.43
1:ZI:218:TYR:HD2	1:ZI:220:SER:H	1.66	0.43
1:ZM:327:PHE:HB3	1:ZM:328:PRO:HD3	2.00	0.43
1:ZU:199:HIS:CE1	1:ZU:282:CYS:HB3	2.53	0.43
1:ZX:224:ILE:HD12	1:ZX:303:ALA:HB3	2.01	0.43
1:YO:248:ASP:CG	1:YO:248:ASP:O	2.61	0.43
1:YQ:249:LEU:HB2	1:YQ:389:LYS:HE3	2.01	0.43
1:YS:169:VAL:HG11	1:YS:352:VAL:HG13	2.01	0.43
1:YS:175:ALA:HB3	2:YS:501:NDP:H3D	1.99	0.43
1:YS:379:GLU:H	1:YS:379:GLU:CD	2.25	0.43
1:YX:186:TYR:HD2	1:YX:190:GLY:O	2.02	0.43
1:ZB:167:LEU:HD12	1:ZB:168:ASP:H	1.83	0.43
1:ZM:117:ILE:HD12	1:ZM:160:PHE:HE2	1.83	0.43
1:ZM:265:ILE:HD12	1:ZM:265:ILE:HA	1.85	0.43
1:ZT:92:VAL:HB	1:ZT:170:LEU:HD13	2.00	0.43
1:ZV:87:ARG:HG2	1:ZV:88:LYS:N	2.33	0.43
1:ZV:100:LEU:HD21	1:ZV:310:ILE:CD1	2.49	0.43
1:ZX:178:PHE:HB3	1:ZX:181:ALA:HB2	2.00	0.43
1:ZX:289:GLU:HG2	1:ZX:391:TRP:CB	2.48	0.43
1:ZY:124:LEU:HD21	1:ZY:128:ARG:HH21	1.83	0.43
1:YK:248:ASP:O	1:YK:248:ASP:CG	2.62	0.43
1:YO:327:PHE:HB3	1:YO:328:PRO:HD3	2.00	0.43
1:YS:116:VAL:O	1:YS:140:TYR:HA	2.19	0.43
1:YS:218:TYR:HD2	1:YS:220:SER:H	1.66	0.43
1:YT:332:LYS:HA	1:YT:338:TYR:HB2	1.99	0.43
1:YU:100:LEU:O	1:YU:104:THR:HG22	2.19	0.43
1:YZ:253:ALA:HA	1:YZ:393:ILE:HG21	2.00	0.43
1:ZG:290:PHE:HB3	1:ZG:302:PHE:CZ	2.53	0.43
1:ZI:116:VAL:O	1:ZI:140:TYR:HA	2.19	0.43
1:ZI:214:LYS:HZ3	1:ZI:298:THR:HA	1.83	0.43
1:ZM:116:VAL:O	1:ZM:140:TYR:HA	2.19	0.43
1:ZM:379:GLU:H	1:ZM:379:GLU:CD	2.26	0.43
1:ZO:249:LEU:HB2	1:ZO:389:LYS:HE3	2.01	0.43
1:ZW:252:LEU:HB2	1:ZW:265:ILE:HG23	2.01	0.43
1:ZX:322:LEU:HD23	1:ZX:322:LEU:HA	1.86	0.43
1:ZY:116:VAL:O	1:ZY:140:TYR:HA	2.18	0.43
1:YJ:92:VAL:HB	1:YJ:170:LEU:HD13	1.99	0.43
1:YJ:335:THR:HG23	1:YJ:337:GLY:H	1.82	0.43
1:YQ:100:LEU:O	1:YQ:104:THR:HG22	2.18	0.43
1:YS:104:THR:HG21	1:YS:348:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YW:169:VAL:HG11	1:YW:352:VAL:HG13	2.01	0.43
1:ZD:366:TRP:HA	1:ZD:372:SER:HA	2.00	0.43
1:ZG:264:MET:SD	1:ZG:270:PHE:HA	2.58	0.43
1:ZI:210:LEU:HD22	1:ZI:290:PHE:HE1	1.83	0.43
1:ZI:229:ILE:HD11	1:ZI:235:THR:HG21	2.00	0.43
1:YW:175:ALA:HB3	2:YW:501:NDP:H3D	1.99	0.43
1:ZF:246:LEU:HB3	1:ZF:249:LEU:HD21	2.00	0.43
1:ZL:332:LYS:HG2	1:ZL:338:TYR:CG	2.54	0.43
1:ZN:199:HIS:CE1	1:ZN:282:CYS:HB3	2.54	0.43
3:ZX:502:LMG:HC1	3:ZX:502:LMG:HC8	1.65	0.43
1:YT:218:TYR:HD1	1:YT:220:SER:H	1.65	0.43
1:ZL:347:ARG:O	1:ZL:350:GLN:HB3	2.19	0.43
1:ZR:379:GLU:CD	1:ZR:379:GLU:H	2.27	0.43
1:ZT:315:LEU:HD13	2:ZT:501:NDP:H71N	1.84	0.43
1:ZW:230:THR:HG22	1:ZW:277:LYS:HD3	2.00	0.43
1:YM:366:TRP:NE1	1:YM:372:SER:HB3	2.34	0.43
1:YO:218:TYR:HD2	1:YO:220:SER:H	1.67	0.43
1:YQ:289:GLU:HB2	1:YQ:387:ALA:HB1	1.99	0.43
1:YY:280:LYS:HA	1:YY:280:LYS:HD2	1.88	0.43
1:ZE:169:VAL:HG11	1:ZE:352:VAL:HG13	2.01	0.43
1:ZE:327:PHE:HB3	1:ZE:328:PRO:HD3	2.00	0.43
1:ZF:167:LEU:HD12	1:ZF:168:ASP:H	1.83	0.43
1:ZJ:367:ASN:OD1	1:ZJ:371:ALA:HB3	2.19	0.43
1:ZL:157:VAL:HG11	1:ZL:208:LEU:HB3	2.00	0.43
1:ZL:289:GLU:HB2	1:ZL:387:ALA:HB1	2.00	0.43
1:ZQ:216:SER:OG	1:ZQ:221:LYS:HD3	2.19	0.43
1:ZV:135:MET:HE2	1:ZV:135:MET:HB2	1.86	0.43
1:ZW:100:LEU:O	1:ZW:104:THR:HG22	2.18	0.43
1:ZW:145:LEU:HD22	1:ZW:156:PHE:CD2	2.54	0.43
1:ZW:221:LYS:C	1:ZW:222:ARG:HG3	2.43	0.43
1:ZX:100:LEU:CD2	1:ZX:348:LEU:HD13	2.49	0.43
1:YL:249:LEU:HD11	1:YL:386:LYS:HG3	2.01	0.42
1:YM:253:ALA:HA	1:YM:393:ILE:HG21	2.01	0.42
1:YZ:277:LYS:O	1:YZ:281:VAL:HG23	2.19	0.42
1:ZD:315:LEU:HD13	2:ZD:501:NDP:H71N	1.84	0.42
1:ZK:304:SER:HB3	1:ZK:363:TYR:HA	2.01	0.42
1:ZL:264:MET:HE2	1:ZL:264:MET:HB2	1.86	0.42
1:ZM:104:THR:HG21	1:ZM:348:LEU:HD22	2.01	0.42
1:ZR:250:ARG:HH12	1:ZV:106:LYS:HZ2	1.67	0.42
1:YJ:210:LEU:HD13	1:YJ:290:PHE:CZ	2.54	0.42
1:YJ:347:ARG:HB3	1:YJ:366:TRP:CZ2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YW:327:PHE:HB3	1:YW:328:PRO:HD3	2.01	0.42
1:ZI:248:ASP:CG	1:ZI:248:ASP:O	2.61	0.42
1:ZJ:246:LEU:HB3	1:ZJ:249:LEU:CD1	2.49	0.42
1:ZM:124:LEU:HD21	1:ZM:128:ARG:HH21	1.84	0.42
1:ZP:229:ILE:HD12	1:ZP:229:ILE:HA	1.89	0.42
1:ZU:116:VAL:O	1:ZU:140:TYR:HA	2.19	0.42
1:ZY:157:VAL:O	1:ZY:161:ARG:HG3	2.19	0.42
1:YO:124:LEU:HD21	1:YO:128:ARG:HH21	1.84	0.42
1:YQ:280:LYS:HA	1:YQ:280:LYS:HD2	1.85	0.42
1:YT:250:ARG:NH2	1:YX:133:VAL:N	2.66	0.42
2:YU:501:NDP:H41N	4:YU:503:A1JWG:O1A	2.19	0.42
1:YV:88:LYS:HA	1:YV:113:LYS:NZ	2.34	0.42
1:ZC:274:LYS:HE3	1:ZC:274:LYS:HB3	1.79	0.42
1:ZE:248:ASP:CG	1:ZE:248:ASP:O	2.62	0.42
1:ZF:159:ASN:O	1:ZF:162:ARG:HG2	2.19	0.42
1:ZH:88:LYS:HA	1:ZH:113:LYS:NZ	2.34	0.42
1:ZK:91:VAL:HG21	1:ZK:114:TRP:HE3	1.83	0.42
1:ZQ:327:PHE:HB3	1:ZQ:328:PRO:HD3	2.01	0.42
1:ZR:167:LEU:HD12	1:ZR:168:ASP:H	1.83	0.42
1:ZZ:167:LEU:HD12	1:ZZ:168:ASP:H	1.83	0.42
1:YI:289:GLU:HB2	1:YI:387:ALA:HB1	2.00	0.42
1:YK:363:TYR:CD2	1:YK:375:ASN:HB3	2.54	0.42
1:YM:280:LYS:HA	1:YM:280:LYS:HD2	1.94	0.42
1:YW:248:ASP:CG	1:YW:248:ASP:O	2.62	0.42
1:YZ:347:ARG:O	1:YZ:350:GLN:HB3	2.20	0.42
1:ZR:127:GLU:C	1:ZR:129:ALA:H	2.27	0.42
1:ZU:248:ASP:CG	1:ZU:248:ASP:O	2.62	0.42
1:ZY:118:MET:HE1	1:ZY:130:ALA:HB2	2.02	0.42
1:YK:111:THR:HG22	1:YK:113:LYS:HG3	2.02	0.42
1:YK:379:GLU:H	1:YK:379:GLU:CD	2.28	0.42
1:YQ:221:LYS:C	1:YQ:222:ARG:HG3	2.45	0.42
1:ZA:327:PHE:HB3	1:ZA:328:PRO:HD3	2.01	0.42
1:ZB:250:ARG:NH2	1:ZF:133:VAL:N	2.67	0.42
1:ZD:357:LEU:HD22	1:ZD:364:TRP:CH2	2.54	0.42
1:ZG:100:LEU:O	1:ZG:104:THR:HG22	2.20	0.42
1:ZI:169:VAL:HG11	1:ZI:352:VAL:HG13	2.02	0.42
1:ZK:225:ILE:HG21	1:ZK:283:ASN:HB3	2.02	0.42
1:ZL:213:LEU:HD23	1:ZL:213:LEU:HA	1.84	0.42
1:ZS:366:TRP:NE1	1:ZS:372:SER:HB3	2.34	0.42
1:YK:327:PHE:HB3	1:YK:328:PRO:HD3	2.01	0.42
2:YK:501:NDP:H41N	4:YK:503:A1JWG:O1A	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YP:250:ARG:CD	1:YT:133:VAL:HG13	2.50	0.42
1:YQ:234:ASN:HD21	1:YQ:376:GLN:HE22	1.68	0.42
1:YS:248:ASP:O	1:YS:248:ASP:CG	2.63	0.42
1:YW:214:LYS:HD2	1:YW:214:LYS:HA	1.70	0.42
1:YZ:332:LYS:HG2	1:YZ:338:TYR:CG	2.54	0.42
1:ZG:249:LEU:HD12	1:ZG:389:LYS:HE3	2.01	0.42
1:ZH:367:ASN:HB2	1:ZH:370:SER:O	2.19	0.42
1:ZM:169:VAL:HG11	1:ZM:352:VAL:HG13	2.02	0.42
2:ZM:501:NDP:H41N	4:ZM:503:A1JWG:O1A	2.20	0.42
1:ZR:214:LYS:NZ	1:ZR:298:THR:HA	2.34	0.42
1:ZX:379:GLU:HG3	1:ZX:380:GLU:N	2.34	0.42
1:YI:234:ASN:HD21	1:YI:376:GLN:HE22	1.67	0.42
1:YU:210:LEU:HD23	1:YU:210:LEU:O	2.20	0.42
1:YU:304:SER:HB3	1:YU:363:TYR:HA	2.00	0.42
1:ZD:95:GLY:HA3	1:ZD:174:ALA:HB2	2.01	0.42
1:ZG:210:LEU:HD23	1:ZG:210:LEU:O	2.19	0.42
1:ZG:225:ILE:HG21	1:ZG:283:ASN:HB3	2.02	0.42
1:ZJ:87:ARG:NH1	1:ZJ:218:TYR:HD1	2.15	0.42
1:ZM:343:GLU:O	1:ZM:346:LYS:HG2	2.20	0.42
1:ZP:118:MET:HB3	1:ZP:142:VAL:HA	2.00	0.42
1:YL:262:SER:HB3	1:YL:267:GLY:HA2	2.01	0.42
1:YM:305:LEU:HD12	1:YM:307:PRO:HD3	2.02	0.42
1:YR:315:LEU:HB2	2:YR:501:NDP:O1N	2.20	0.42
1:YT:249:LEU:HD11	1:YT:386:LYS:HG3	2.01	0.42
1:YU:135:MET:HE3	1:YU:135:MET:HB3	1.94	0.42
1:YZ:379:GLU:HG3	1:YZ:380:GLU:N	2.34	0.42
1:ZD:210:LEU:HD13	1:ZD:290:PHE:HZ	1.85	0.42
1:ZG:131:LYS:HA	1:ZG:131:LYS:HD3	1.85	0.42
1:ZG:157:VAL:HG11	1:ZG:208:LEU:HG	2.02	0.42
1:ZK:293:ARG:HB2	1:ZK:294:PHE:CD1	2.54	0.42
1:ZM:216:SER:OG	1:ZM:221:LYS:HD3	2.19	0.42
1:ZO:172:CYS:HB2	1:ZO:225:ILE:HD13	2.02	0.42
1:ZQ:248:ASP:O	1:ZQ:248:ASP:CG	2.62	0.42
1:ZQ:363:TYR:CD2	1:ZQ:375:ASN:HB3	2.55	0.42
1:ZU:100:LEU:HD21	1:ZU:307:PRO:HG2	2.02	0.42
1:YR:281:VAL:O	1:YR:285:LEU:HD23	2.20	0.42
1:YV:291:HIS:HD2	1:YV:360:SER:HB3	1.85	0.42
1:ZF:346:LYS:HB2	1:ZF:346:LYS:HE3	1.78	0.42
1:ZH:92:VAL:HB	1:ZH:170:LEU:HD13	2.02	0.42
1:ZS:280:LYS:HA	1:ZS:280:LYS:HD2	1.84	0.42
1:YJ:118:MET:HB2	1:YJ:141:THR:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YM:234:ASN:HD21	1:YM:376:GLN:HE22	1.68	0.42
1:YQ:213:LEU:HD12	1:YQ:221:LYS:HD2	2.02	0.42
1:YV:159:ASN:HD21	1:ZA:159:ASN:ND2	2.17	0.42
1:ZF:363:TYR:CD2	1:ZF:375:ASN:HB3	2.54	0.42
1:ZG:280:LYS:HA	1:ZG:280:LYS:HD2	1.88	0.42
1:ZJ:218:TYR:CD1	1:ZJ:219:PRO:HD2	2.55	0.42
1:ZK:145:LEU:HD22	1:ZK:156:PHE:CD2	2.54	0.42
1:ZL:223:LEU:HD22	1:ZL:225:ILE:HD11	2.01	0.42
1:ZW:157:VAL:HG13	1:ZW:209:LEU:HD13	2.02	0.42
1:YI:178:PHE:HB3	1:YI:181:ALA:HB2	2.02	0.41
1:YJ:287:MET:HE2	1:YJ:304:SER:CB	2.50	0.41
1:YK:169:VAL:HG11	1:YK:352:VAL:HG13	2.02	0.41
1:YP:246:LEU:HD23	1:YP:249:LEU:HD21	2.02	0.41
1:YS:117:ILE:HD12	1:YS:160:PHE:HE2	1.85	0.41
1:YT:262:SER:HB3	1:YT:267:GLY:HA2	2.00	0.41
1:YV:287:MET:HE2	1:YV:304:SER:N	2.34	0.41
1:YX:379:GLU:H	1:YX:379:GLU:CD	2.28	0.41
1:YZ:131:LYS:HA	1:YZ:131:LYS:HD3	1.80	0.41
1:YZ:182:LYS:HE2	1:YZ:183:GLU:OE1	2.20	0.41
1:ZA:389:LYS:O	1:ZA:393:ILE:HG13	2.20	0.41
1:ZD:210:LEU:HD13	1:ZD:290:PHE:CZ	2.56	0.41
1:ZE:135:MET:HE2	1:ZE:135:MET:HB3	1.76	0.41
1:ZG:264:MET:HE2	1:ZG:264:MET:HB2	1.84	0.41
1:ZL:265:ILE:HD12	1:ZL:282:CYS:SG	2.60	0.41
1:ZQ:264:MET:CE	1:ZQ:271:ASP:H	2.29	0.41
1:YU:252:LEU:HD12	1:YU:265:ILE:HG23	2.02	0.41
1:YZ:124:LEU:HD21	1:YZ:128:ARG:NH2	2.35	0.41
1:ZH:88:LYS:HA	1:ZH:113:LYS:HZ2	1.85	0.41
1:ZI:100:LEU:HD21	1:ZI:307:PRO:HG2	2.02	0.41
1:ZT:347:ARG:HB3	1:ZT:366:TRP:CZ2	2.55	0.41
1:ZU:143:MET:HE2	1:ZU:143:MET:HB3	1.85	0.41
1:YI:253:ALA:HA	1:YI:393:ILE:HG21	2.02	0.41
1:YO:177:TYR:CE2	1:YO:179:PRO:HG3	2.55	0.41
1:YO:293:ARG:HE	1:YO:293:ARG:HB2	1.75	0.41
1:YR:162:ARG:HD2	1:YW:162:ARG:NE	2.35	0.41
1:ZF:214:LYS:HA	1:ZF:214:LYS:HD2	1.93	0.41
1:ZH:319:HIS:CE1	1:ZH:320:ILE:HG12	2.54	0.41
1:ZL:379:GLU:HG3	1:ZL:380:GLU:N	2.35	0.41
1:ZN:133:VAL:C	1:ZN:135:MET:N	2.78	0.41
1:ZQ:100:LEU:HD21	1:ZQ:307:PRO:HG2	2.01	0.41
1:ZR:153:VAL:O	1:ZR:157:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZT:295:HIS:HB2	1:ZT:302:PHE:CE2	2.55	0.41
1:ZY:143:MET:HE2	1:ZY:143:MET:HB3	1.99	0.41
1:ZC:88:LYS:HD2	1:ZC:113:LYS:HA	2.03	0.41
1:ZD:223:LEU:HB2	1:ZD:302:PHE:HD1	1.85	0.41
1:ZN:256:LEU:HD12	1:ZN:256:LEU:HA	1.77	0.41
1:ZQ:379:GLU:HG2	1:ZQ:380:GLU:N	2.35	0.41
1:ZT:244:ALA:HB2	1:ZT:277:LYS:HB3	2.02	0.41
1:ZV:379:GLU:CD	1:ZV:379:GLU:H	2.28	0.41
1:YL:389:LYS:O	1:YL:392:GLU:HG2	2.21	0.41
1:YJ:103:ALA:O	1:YJ:106:LYS:HG2	2.21	0.41
1:YJ:281:VAL:O	1:YJ:285:LEU:HD23	2.20	0.41
1:YN:103:ALA:O	1:YN:106:LYS:HG2	2.20	0.41
1:YN:319:HIS:CE1	1:YN:320:ILE:HG12	2.55	0.41
1:YR:103:ALA:O	1:YR:106:LYS:HG2	2.21	0.41
1:YR:179:PRO:HG2	1:YR:319:HIS:HB2	2.02	0.41
1:YR:339:VAL:HG11	1:YR:347:ARG:HH12	1.85	0.41
1:YT:250:ARG:CG	1:YT:253:ALA:HB3	2.49	0.41
1:YV:92:VAL:HB	1:YV:170:LEU:HD13	2.03	0.41
1:YZ:393:ILE:O	1:YZ:396:LYS:HG2	2.20	0.41
1:ZC:346:LYS:HD3	1:ZC:346:LYS:HA	1.81	0.41
1:ZK:280:LYS:HA	1:ZK:280:LYS:HD2	1.89	0.41
1:ZM:363:TYR:CD2	1:ZM:375:ASN:HB3	2.56	0.41
1:ZN:153:VAL:O	1:ZN:157:VAL:HG23	2.20	0.41
1:ZO:366:TRP:NE1	1:ZO:372:SER:HB3	2.35	0.41
1:ZP:287:MET:HE2	1:ZP:304:SER:CB	2.50	0.41
1:ZT:319:HIS:CE1	1:ZT:320:ILE:HG12	2.55	0.41
1:ZU:216:SER:OG	1:ZU:221:LYS:HD3	2.21	0.41
1:ZV:262:SER:HB3	1:ZV:267:GLY:HA2	2.02	0.41
1:ZX:253:ALA:HA	1:ZX:393:ILE:CG2	2.46	0.41
1:YL:108:LEU:HD12	1:YL:114:TRP:CD1	2.56	0.41
1:YS:173:ASN:OD1	2:YS:501:NDP:H4D	2.20	0.41
1:YZ:170:LEU:HD23	1:YZ:171:VAL:N	2.35	0.41
1:YZ:224:ILE:CD1	1:YZ:303:ALA:HB3	2.51	0.41
1:ZE:214:LYS:HZ3	1:ZE:298:THR:HA	1.85	0.41
1:ZH:118:MET:HB2	1:ZH:141:THR:O	2.21	0.41
1:ZI:216:SER:OG	1:ZI:221:LYS:HD3	2.21	0.41
1:ZI:363:TYR:CD2	1:ZI:375:ASN:HB3	2.55	0.41
1:ZL:124:LEU:HD21	1:ZL:128:ARG:NH2	2.36	0.41
1:ZL:249:LEU:HD12	1:ZL:389:LYS:HB2	2.03	0.41
1:ZM:177:TYR:CE2	1:ZM:179:PRO:HG3	2.56	0.41
1:ZM:343:GLU:HA	1:ZM:346:LYS:HG2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZN:255:GLY:CA	1:ZR:133:VAL:HG11	2.51	0.41
1:ZO:284:MET:O	1:ZO:288:GLN:HG3	2.21	0.41
1:ZP:162:ARG:HG2	1:ZU:162:ARG:NH2	2.33	0.41
1:ZU:218:TYR:HD1	1:ZU:220:SER:H	1.69	0.41
1:ZW:313:THR:HG22	1:ZW:314:GLY:N	2.33	0.41
1:ZX:131:LYS:HA	1:ZX:131:LYS:HD3	1.80	0.41
1:ZX:292:ARG:HH11	1:ZX:388:ARG:HH11	1.67	0.41
1:YJ:229:ILE:HD12	1:YJ:232:ASN:HB2	2.02	0.41
2:YM:501:NDP:H41N	4:YM:503:A1JWG:O1A	2.19	0.41
1:YY:230:THR:HG22	1:YY:277:LYS:HD3	2.03	0.41
1:ZO:218:TYR:CD1	1:ZO:219:PRO:HD2	2.56	0.41
1:ZO:234:ASN:HD21	1:ZO:376:GLN:HE22	1.69	0.41
1:ZQ:104:THR:HG21	1:ZQ:348:LEU:HD22	2.03	0.41
1:ZR:121:ARG:CB	1:ZR:144:HIS:CE1	3.04	0.41
1:ZT:118:MET:CE	1:ZT:120:CYS:HB2	2.49	0.41
1:YJ:179:PRO:HB2	1:YJ:319:HIS:CD2	2.56	0.41
1:YJ:251:GLY:HA3	1:YJ:264:MET:O	2.20	0.41
1:YK:116:VAL:O	1:YK:140:TYR:HA	2.21	0.41
1:YN:295:HIS:HB2	1:YN:302:PHE:CE2	2.56	0.41
1:YR:124:LEU:HD21	1:YR:128:ARG:NH2	2.36	0.41
1:YW:177:TYR:CE2	1:YW:179:PRO:HG3	2.56	0.41
1:ZD:327:PHE:HB3	1:ZD:328:PRO:HD3	2.03	0.41
1:ZL:328:PRO:O	1:ZL:332:LYS:HG3	2.20	0.41
1:ZN:255:GLY:HA2	1:ZR:133:VAL:HG21	2.03	0.41
1:ZP:162:ARG:HE	1:ZU:162:ARG:HE	1.67	0.41
1:ZR:250:ARG:HB2	1:ZR:253:ALA:HB3	2.01	0.41
1:ZU:124:LEU:HD21	1:ZU:128:ARG:HH21	1.84	0.41
1:ZW:159:ASN:HA	1:ZW:162:ARG:CZ	2.51	0.41
1:ZX:223:LEU:HD22	1:ZX:225:ILE:HD11	2.03	0.41
1:ZY:100:LEU:HD11	1:ZY:310:ILE:CD1	2.51	0.41
1:YI:230:THR:HG22	1:YI:277:LYS:HD3	2.03	0.41
1:YI:256:LEU:HD12	1:YI:262:SER:O	2.21	0.41
1:YJ:295:HIS:HB2	1:YJ:302:PHE:CE2	2.55	0.41
1:YN:366:TRP:HA	1:YN:372:SER:HA	2.03	0.41
1:YP:250:ARG:NH2	1:YT:133:VAL:H	2.18	0.41
1:YQ:313:THR:HG22	1:YQ:314:GLY:N	2.34	0.41
1:YV:339:VAL:HG11	1:YV:347:ARG:HH12	1.85	0.41
1:YY:292:ARG:HH12	1:YY:384:VAL:HA	1.86	0.41
1:YZ:328:PRO:O	1:YZ:332:LYS:HG3	2.21	0.41
1:ZA:252:LEU:HB2	1:ZA:265:ILE:HG23	2.02	0.41
2:ZC:501:NDP:H41N	4:ZC:503:A1JWG:O1A	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZD:307:PRO:HG3	1:ZD:348:LEU:HD11	2.03	0.41
1:ZD:335:THR:HG23	1:ZD:337:GLY:H	1.85	0.41
1:ZE:379:GLU:H	1:ZE:379:GLU:CD	2.28	0.41
1:ZG:319:HIS:CG	1:ZG:320:ILE:H	2.38	0.41
1:ZG:379:GLU:HG2	1:ZG:380:GLU:N	2.35	0.41
1:ZI:124:LEU:HD21	1:ZI:128:ARG:HH21	1.85	0.41
1:ZI:135:MET:HE3	1:ZI:135:MET:HB3	1.84	0.41
1:ZM:143:MET:HE2	1:ZM:143:MET:HB3	1.91	0.41
1:ZP:366:TRP:HA	1:ZP:372:SER:HA	2.02	0.41
1:ZS:305:LEU:HD12	1:ZS:307:PRO:HD3	2.03	0.41
1:ZX:162:ARG:HD2	1:ZX:162:ARG:O	2.21	0.41
1:ZZ:249:LEU:HD11	1:ZZ:386:LYS:CG	2.50	0.41
1:YI:284:MET:O	1:YI:287:MET:HG3	2.21	0.41
1:YK:135:MET:HE3	1:YK:135:MET:HB3	1.99	0.41
1:YK:216:SER:OG	1:YK:221:LYS:HD3	2.20	0.41
1:YT:327:PHE:HB3	1:YT:328:PRO:HD3	2.03	0.41
1:ZA:177:TYR:CE2	1:ZA:179:PRO:HG3	2.56	0.41
1:ZC:322:LEU:HD23	1:ZC:322:LEU:HA	1.86	0.41
1:ZG:305:LEU:HD12	1:ZG:307:PRO:HD3	2.03	0.41
1:ZH:354:ASP:HB3	1:ZH:357:LEU:HG	2.03	0.41
1:ZM:207:ARG:HE	1:ZM:207:ARG:HB3	1.63	0.41
1:ZN:133:VAL:C	1:ZN:135:MET:H	2.29	0.41
1:ZS:290:PHE:HB3	1:ZS:302:PHE:CZ	2.56	0.41
1:ZW:213:LEU:HD12	1:ZW:221:LYS:HD2	2.02	0.41
1:ZY:169:VAL:HG11	1:ZY:352:VAL:HG13	2.03	0.41
1:YM:389:LYS:O	1:YM:393:ILE:HG13	2.22	0.40
1:YT:379:GLU:H	1:YT:379:GLU:CD	2.29	0.40
1:YW:216:SER:OG	1:YW:221:LYS:HD3	2.21	0.40
1:YX:102:LEU:HD12	1:YX:102:LEU:HA	1.92	0.40
1:ZB:262:SER:HB3	1:ZB:267:GLY:HA2	2.03	0.40
1:ZD:118:MET:HB2	1:ZD:141:THR:O	2.20	0.40
1:ZG:160:PHE:O	1:ZG:163:THR:HG22	2.21	0.40
1:ZH:339:VAL:HG11	1:ZH:347:ARG:HH22	1.85	0.40
1:ZI:346:LYS:HB2	1:ZI:346:LYS:HE2	1.89	0.40
1:ZI:362:VAL:CG1	1:ZI:374:GLU:HG2	2.51	0.40
1:ZL:224:ILE:HD12	1:ZL:303:ALA:HB3	2.03	0.40
1:ZN:159:ASN:O	1:ZN:162:ARG:HG2	2.20	0.40
1:ZP:95:GLY:HA3	1:ZP:174:ALA:HB2	2.03	0.40
1:ZY:173:ASN:OD1	2:ZY:501:NDP:H4D	2.20	0.40
1:ZY:264:MET:HE2	1:ZY:271:ASP:N	2.24	0.40
1:YK:104:THR:HG21	1:YK:348:LEU:HD22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:YN:88:LYS:HA	1:YN:113:LYS:NZ	2.36	0.40
1:YP:327:PHE:HB3	1:YP:328:PRO:HD3	2.03	0.40
1:YV:229:ILE:HD12	1:YV:229:ILE:HA	1.90	0.40
1:YV:327:PHE:HB3	1:YV:328:PRO:HD3	2.03	0.40
1:YW:173:ASN:OD1	2:YW:501:NDP:H4D	2.20	0.40
1:YX:87:ARG:NH2	1:YX:218:TYR:HB2	2.36	0.40
1:YY:224:ILE:HD12	1:YY:352:VAL:HG12	2.03	0.40
1:ZF:262:SER:HB3	1:ZF:267:GLY:HA2	2.01	0.40
1:ZG:320:ILE:HG22	1:ZG:322:LEU:H	1.86	0.40
1:ZH:229:ILE:HD12	1:ZH:232:ASN:HD22	1.86	0.40
1:ZP:251:GLY:HA3	1:ZP:264:MET:O	2.21	0.40
1:ZP:295:HIS:HB2	1:ZP:302:PHE:CE2	2.56	0.40
1:ZS:234:ASN:ND2	1:ZS:378:SER:HA	2.36	0.40
1:ZX:170:LEU:HD23	1:ZX:171:VAL:N	2.36	0.40
1:YI:225:ILE:CG2	1:YI:283:ASN:HB3	2.51	0.40
1:YU:218:TYR:CD1	1:YU:219:PRO:HD2	2.56	0.40
1:YW:100:LEU:HD11	1:YW:310:ILE:CD1	2.51	0.40
1:ZC:305:LEU:HD12	1:ZC:307:PRO:HD3	2.03	0.40
1:ZC:339:VAL:HG11	1:ZC:347:ARG:NH1	2.37	0.40
1:ZG:253:ALA:HA	1:ZG:393:ILE:HG21	2.03	0.40
1:ZP:327:PHE:HB3	1:ZP:328:PRO:HD3	2.04	0.40
1:ZQ:324:ARG:HE	1:ZQ:324:ARG:HB2	1.73	0.40
1:ZU:214:LYS:HE2	1:ZU:298:THR:HG22	2.03	0.40
1:ZW:280:LYS:HA	1:ZW:280:LYS:HD2	1.86	0.40
1:ZX:124:LEU:HD21	1:ZX:128:ARG:NH2	2.36	0.40
1:ZX:339:VAL:HG13	1:ZX:343:GLU:OE2	2.21	0.40
1:ZX:347:ARG:O	1:ZX:350:GLN:HB3	2.21	0.40
1:ZX:380:GLU:H	1:ZX:380:GLU:HG2	1.73	0.40
1:ZY:395:GLU:HG2	1:ZY:401:ALA:HB2	2.03	0.40
1:YI:329:PRO:HA	1:YI:332:LYS:HD3	2.03	0.40
1:YN:229:ILE:HD13	1:YN:363:TYR:OH	2.22	0.40
1:YO:264:MET:HE1	1:YO:274:LYS:HB3	2.02	0.40
1:YQ:287:MET:HE3	1:YQ:304:SER:HB2	2.03	0.40
1:YU:162:ARG:HD2	1:YU:162:ARG:O	2.22	0.40
1:YV:244:ALA:HB2	1:YV:277:LYS:HB3	2.04	0.40
1:YW:124:LEU:HD21	1:YW:128:ARG:HH21	1.87	0.40
1:YW:363:TYR:CD2	1:YW:375:ASN:HB3	2.57	0.40
1:YW:389:LYS:O	1:YW:393:ILE:HG13	2.22	0.40
1:YZ:234:ASN:CA	1:ZC:234:ASN:HB2	2.51	0.40
1:ZF:108:LEU:HD12	1:ZF:114:TRP:CD1	2.55	0.40
1:ZK:253:ALA:HA	1:ZK:393:ILE:HG21	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:ZL:145:LEU:HD12	1:ZL:145:LEU:HA	1.84	0.40
1:ZL:173:ASN:OD1	2:ZL:501:NDP:H4D	2.22	0.40
1:ZM:100:LEU:HD11	1:ZM:310:ILE:CD1	2.51	0.40
1:ZO:264:MET:HE1	1:ZO:278:ASP:CB	2.50	0.40
1:ZP:96:ALA:H	1:ZP:118:MET:CE	2.28	0.40
1:ZP:351:VAL:HA	1:ZP:357:LEU:HD12	2.02	0.40
1:ZT:103:ALA:O	1:ZT:106:LYS:HG2	2.21	0.40
1:ZX:119:ALA:HA	1:ZX:143:MET:O	2.22	0.40
1:ZX:328:PRO:N	1:ZX:329:PRO:HD2	2.37	0.40
1:ZX:393:ILE:O	1:ZX:396:LYS:HG2	2.21	0.40
1:YN:251:GLY:HA3	1:YN:264:MET:O	2.21	0.40
1:YQ:178:PHE:HB3	1:YQ:181:ALA:HB2	2.02	0.40
1:YR:328:PRO:N	1:YR:329:PRO:HD2	2.37	0.40
1:YS:135:MET:HE2	1:YS:135:MET:HB3	1.77	0.40
1:YU:313:THR:HG22	1:YU:314:GLY:N	2.33	0.40
1:YV:177:TYR:CD2	1:YV:179:PRO:HG3	2.56	0.40
1:YW:207:ARG:HE	1:YW:207:ARG:HB3	1.63	0.40
1:YZ:249:LEU:HA	1:YZ:265:ILE:HG22	2.04	0.40
1:ZD:347:ARG:O	1:ZD:351:VAL:HG23	2.22	0.40
1:ZG:262:SER:CB	1:ZG:267:GLY:HA2	2.52	0.40
1:ZX:343:GLU:HA	1:ZX:346:LYS:CG	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	YI	314/353 (89%)	299 (95%)	13 (4%)	2 (1%)	21	56
1	YJ	316/353 (90%)	302 (96%)	14 (4%)	0	100	100
1	YK	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	YL	316/353 (90%)	302 (96%)	14 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	YM	314/353 (89%)	298 (95%)	14 (4%)	2 (1%)	21	56
1	YN	316/353 (90%)	302 (96%)	14 (4%)	0	100	100
1	YO	314/353 (89%)	301 (96%)	13 (4%)	0	100	100
1	YP	316/353 (90%)	297 (94%)	19 (6%)	0	100	100
1	YQ	314/353 (89%)	297 (95%)	15 (5%)	2 (1%)	21	56
1	YR	316/353 (90%)	304 (96%)	12 (4%)	0	100	100
1	YS	314/353 (89%)	301 (96%)	13 (4%)	0	100	100
1	YT	316/353 (90%)	295 (93%)	21 (7%)	0	100	100
1	YU	314/353 (89%)	300 (96%)	12 (4%)	2 (1%)	21	56
1	YV	316/353 (90%)	304 (96%)	12 (4%)	0	100	100
1	YW	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	YX	316/353 (90%)	301 (95%)	15 (5%)	0	100	100
1	YY	314/353 (89%)	300 (96%)	12 (4%)	2 (1%)	21	56
1	YZ	316/353 (90%)	298 (94%)	18 (6%)	0	100	100
1	ZA	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	ZB	316/353 (90%)	298 (94%)	18 (6%)	0	100	100
1	ZC	314/353 (89%)	300 (96%)	12 (4%)	2 (1%)	21	56
1	ZD	316/353 (90%)	304 (96%)	12 (4%)	0	100	100
1	ZE	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	ZF	316/353 (90%)	296 (94%)	20 (6%)	0	100	100
1	ZG	314/353 (89%)	298 (95%)	14 (4%)	2 (1%)	21	56
1	ZH	316/353 (90%)	304 (96%)	12 (4%)	0	100	100
1	ZI	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	ZJ	316/353 (90%)	301 (95%)	15 (5%)	0	100	100
1	ZK	314/353 (89%)	301 (96%)	11 (4%)	2 (1%)	21	56
1	ZL	316/353 (90%)	297 (94%)	19 (6%)	0	100	100
1	ZM	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	ZN	316/353 (90%)	288 (91%)	28 (9%)	0	100	100
1	ZO	314/353 (89%)	295 (94%)	17 (5%)	2 (1%)	21	56
1	ZP	316/353 (90%)	303 (96%)	13 (4%)	0	100	100
1	ZQ	314/353 (89%)	297 (95%)	17 (5%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	ZR	316/353 (90%)	292 (92%)	19 (6%)	5 (2%)	7	36
1	ZS	314/353 (89%)	297 (95%)	15 (5%)	2 (1%)	21	56
1	ZT	316/353 (90%)	303 (96%)	13 (4%)	0	100	100
1	ZU	314/353 (89%)	301 (96%)	13 (4%)	0	100	100
1	ZV	316/353 (90%)	295 (93%)	21 (7%)	0	100	100
1	ZW	314/353 (89%)	299 (95%)	13 (4%)	2 (1%)	21	56
1	ZX	316/353 (90%)	298 (94%)	18 (6%)	0	100	100
1	ZY	314/353 (89%)	300 (96%)	14 (4%)	0	100	100
1	ZZ	316/353 (90%)	296 (94%)	20 (6%)	0	100	100
All	All	13860/15532 (89%)	13164 (95%)	669 (5%)	27 (0%)	44	73

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	ZR	127	GLU
1	ZR	123	PHE
1	ZR	126	ALA
1	ZR	132	SER
1	ZR	134	GLY
1	YI	241	PRO
1	YI	242	PRO
1	YY	242	PRO
1	ZC	242	PRO
1	ZK	242	PRO
1	ZW	242	PRO
1	YY	241	PRO
1	ZC	241	PRO
1	ZK	241	PRO
1	ZW	241	PRO
1	YM	241	PRO
1	YQ	241	PRO
1	YU	241	PRO
1	ZG	241	PRO
1	ZO	241	PRO
1	ZS	241	PRO
1	YM	242	PRO
1	YQ	242	PRO
1	YU	242	PRO
1	ZG	242	PRO

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Mol	Chain	Res	Type
1	ZO	242	PRO
1	ZS	242	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	YI	259/290 (89%)	259 (100%)	0	100	100
1	YJ	261/290 (90%)	261 (100%)	0	100	100
1	YK	259/290 (89%)	259 (100%)	0	100	100
1	YL	261/290 (90%)	261 (100%)	0	100	100
1	YM	259/290 (89%)	259 (100%)	0	100	100
1	YN	261/290 (90%)	261 (100%)	0	100	100
1	YO	259/290 (89%)	259 (100%)	0	100	100
1	YP	261/290 (90%)	261 (100%)	0	100	100
1	YQ	259/290 (89%)	259 (100%)	0	100	100
1	YR	261/290 (90%)	261 (100%)	0	100	100
1	YS	259/290 (89%)	259 (100%)	0	100	100
1	YT	261/290 (90%)	261 (100%)	0	100	100
1	YU	259/290 (89%)	258 (100%)	1 (0%)	84	86
1	YV	261/290 (90%)	261 (100%)	0	100	100
1	YW	259/290 (89%)	258 (100%)	1 (0%)	84	86
1	YX	261/290 (90%)	261 (100%)	0	100	100
1	YY	259/290 (89%)	259 (100%)	0	100	100
1	YZ	261/290 (90%)	261 (100%)	0	100	100
1	ZA	259/290 (89%)	259 (100%)	0	100	100
1	ZB	261/290 (90%)	260 (100%)	1 (0%)	84	86
1	ZC	259/290 (89%)	259 (100%)	0	100	100
1	ZD	261/290 (90%)	261 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	ZE	259/290 (89%)	259 (100%)	0	100	100
1	ZF	261/290 (90%)	261 (100%)	0	100	100
1	ZG	259/290 (89%)	259 (100%)	0	100	100
1	ZH	261/290 (90%)	261 (100%)	0	100	100
1	ZI	259/290 (89%)	258 (100%)	1 (0%)	84	86
1	ZJ	261/290 (90%)	261 (100%)	0	100	100
1	ZK	259/290 (89%)	259 (100%)	0	100	100
1	ZL	261/290 (90%)	260 (100%)	1 (0%)	84	86
1	ZM	259/290 (89%)	259 (100%)	0	100	100
1	ZN	261/290 (90%)	261 (100%)	0	100	100
1	ZO	259/290 (89%)	259 (100%)	0	100	100
1	ZP	261/290 (90%)	260 (100%)	1 (0%)	84	86
1	ZQ	259/290 (89%)	259 (100%)	0	100	100
1	ZR	261/290 (90%)	261 (100%)	0	100	100
1	ZS	259/290 (89%)	259 (100%)	0	100	100
1	ZT	261/290 (90%)	261 (100%)	0	100	100
1	ZU	259/290 (89%)	259 (100%)	0	100	100
1	ZV	261/290 (90%)	261 (100%)	0	100	100
1	ZW	259/290 (89%)	259 (100%)	0	100	100
1	ZX	261/290 (90%)	260 (100%)	1 (0%)	84	86
1	ZY	259/290 (89%)	259 (100%)	0	100	100
1	ZZ	261/290 (90%)	260 (100%)	1 (0%)	84	86
All	All	11440/12760 (90%)	11432 (100%)	8 (0%)	87	91

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	YU	199	HIS
1	YW	133	VAL
1	ZB	133	VAL
1	ZI	133	VAL
1	ZL	199	HIS
1	ZP	159	ASN
1	ZX	199	HIS

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Mol	Chain	Res	Type
1	ZZ	250	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (92) such sidechains are listed below:

Mol	Chain	Res	Type
1	YI	234	ASN
1	YI	288	GLN
1	YI	331	GLN
1	YJ	331	GLN
1	YK	115	ASN
1	YK	199	HIS
1	YL	90	ASN
1	YL	115	ASN
1	YL	239	ASN
1	YL	283	ASN
1	YM	234	ASN
1	YM	283	ASN
1	YM	331	GLN
1	YN	155	GLN
1	YO	115	ASN
1	YO	159	ASN
1	YP	90	ASN
1	YQ	234	ASN
1	YS	115	ASN
1	YS	199	HIS
1	YT	90	ASN
1	YV	159	ASN
1	YV	291	HIS
1	YV	331	GLN
1	YW	115	ASN
1	YW	199	HIS
1	YW	288	GLN
1	YW	331	GLN
1	YX	90	ASN
1	YY	199	HIS
1	YY	234	ASN
1	YZ	198	ASN
1	YZ	283	ASN
1	YZ	350	GLN
1	ZA	115	ASN
1	ZB	90	ASN
1	ZC	199	HIS

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Mol	Chain	Res	Type
1	ZC	234	ASN
1	ZD	202	HIS
1	ZD	331	GLN
1	ZF	90	ASN
1	ZF	115	ASN
1	ZF	295	HIS
1	ZF	331	GLN
1	ZG	234	ASN
1	ZI	115	ASN
1	ZI	155	GLN
1	ZJ	90	ASN
1	ZJ	331	GLN
1	ZK	234	ASN
1	ZK	288	GLN
1	ZK	295	HIS
1	ZK	331	GLN
1	ZL	159	ASN
1	ZL	198	ASN
1	ZL	202	HIS
1	ZL	291	HIS
1	ZL	295	HIS
1	ZM	115	ASN
1	ZM	155	GLN
1	ZM	199	HIS
1	ZN	90	ASN
1	ZN	199	HIS
1	ZO	234	ASN
1	ZO	295	HIS
1	ZP	331	GLN
1	ZQ	159	ASN
1	ZQ	199	HIS
1	ZQ	331	GLN
1	ZR	90	ASN
1	ZR	239	ASN
1	ZR	331	GLN
1	ZS	234	ASN
1	ZT	331	GLN
1	ZU	115	ASN
1	ZU	199	HIS
1	ZU	239	ASN
1	ZV	90	ASN
1	ZV	115	ASN

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Mol	Chain	Res	Type
1	ZV	239	ASN
1	ZW	199	HIS
1	ZW	283	ASN
1	ZX	198	ASN
1	ZX	202	HIS
1	ZX	331	GLN
1	ZX	350	GLN
1	ZY	115	ASN
1	ZY	159	ASN
1	ZY	199	HIS
1	ZY	331	GLN
1	ZZ	90	ASN
1	ZZ	239	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

132 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	A1JWG	ZT	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.33	3 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NDP	YT	501	-	49,52,52	2.18	6 (12%)	66,80,80	1.71	14 (21%)
3	LMG	ZS	502	-	23,23,55	0.69	0	31,31,63	0.83	0
4	A1JWG	YK	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.21	4 (6%)
2	NDP	YY	501	-	49,52,52	2.13	4 (8%)	66,80,80	1.72	14 (21%)
4	A1JWG	ZQ	503	-	48,53,53	1.39	5 (10%)	62,89,89	1.09	2 (3%)
4	A1JWG	YO	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.18	2 (3%)
4	A1JWG	ZZ	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.09	3 (4%)
4	A1JWG	ZY	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.17	3 (4%)
3	LMG	ZB	502	-	23,23,55	0.72	0	31,31,63	0.83	1 (3%)
3	LMG	ZG	502	-	23,23,55	0.69	0	31,31,63	0.86	0
3	LMG	ZJ	502	-	23,23,55	0.71	0	31,31,63	0.81	0
3	LMG	ZL	502	-	23,23,55	0.69	0	31,31,63	0.83	0
4	A1JWG	YZ	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.18	3 (4%)
2	NDP	ZU	501	-	49,52,52	2.14	8 (16%)	66,80,80	1.68	15 (22%)
3	LMG	YX	502	-	23,23,55	0.71	0	31,31,63	0.83	0
4	A1JWG	ZR	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.10	3 (4%)
3	LMG	ZV	502	-	23,23,55	0.71	0	31,31,63	0.79	0
2	NDP	YS	501	-	49,52,52	2.14	8 (16%)	66,80,80	1.68	15 (22%)
2	NDP	ZP	501	-	49,52,52	2.20	6 (12%)	66,80,80	1.70	17 (25%)
3	LMG	ZQ	502	-	23,23,55	0.71	0	31,31,63	0.83	0
3	LMG	YY	502	-	23,23,55	0.69	0	31,31,63	0.85	1 (3%)
3	LMG	YI	502	-	23,23,55	0.68	0	31,31,63	0.83	0
4	A1JWG	ZN	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.09	3 (4%)
3	LMG	ZA	502	-	23,23,55	0.73	0	31,31,63	0.80	0
3	LMG	YS	502	-	23,23,55	0.73	0	31,31,63	0.78	0
4	A1JWG	YT	503	-	48,53,53	1.40	4 (8%)	62,89,89	1.14	3 (4%)
4	A1JWG	ZL	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.16	3 (4%)
2	NDP	YL	501	-	49,52,52	2.19	4 (8%)	66,80,80	1.73	15 (22%)
2	NDP	ZX	501	-	49,52,52	2.17	4 (8%)	66,80,80	1.71	15 (22%)
3	LMG	ZK	502	-	23,23,55	0.69	0	31,31,63	0.84	0
3	LMG	ZY	502	-	23,23,55	0.71	0	31,31,63	0.79	0
4	A1JWG	YV	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.23	5 (8%)
2	NDP	ZA	501	-	49,52,52	2.11	8 (16%)	66,80,80	1.70	15 (22%)
3	LMG	ZF	502	-	23,23,55	0.71	0	31,31,63	0.77	0
2	NDP	ZZ	501	-	49,52,52	2.17	6 (12%)	66,80,80	1.70	14 (21%)
3	LMG	ZW	502	-	23,23,55	0.68	0	31,31,63	0.87	0
2	NDP	ZK	501	-	49,52,52	2.15	5 (10%)	66,80,80	1.72	15 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	A1JWG	YI	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.15	2 (3%)
4	A1JWG	YJ	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.31	3 (4%)
3	LMG	ZX	502	-	23,23,55	0.66	0	31,31,63	1.03	2 (6%)
4	A1JWG	ZF	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.09	3 (4%)
4	A1JWG	ZB	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.10	3 (4%)
2	NDP	ZO	501	-	49,52,52	2.18	5 (10%)	66,80,80	1.71	14 (21%)
4	A1JWG	ZX	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.18	3 (4%)
4	A1JWG	ZJ	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.09	3 (4%)
4	A1JWG	ZW	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.18	3 (4%)
2	NDP	YM	501	-	49,52,52	2.14	4 (8%)	66,80,80	1.71	14 (21%)
2	NDP	YQ	501	-	49,52,52	2.14	4 (8%)	66,80,80	1.71	15 (22%)
3	LMG	YK	502	-	23,23,55	0.72	0	31,31,63	0.80	0
3	LMG	YN	502	-	23,23,55	0.70	0	31,31,63	0.81	0
3	LMG	ZD	502	-	23,23,55	0.69	0	31,31,63	0.87	1 (3%)
4	A1JWG	ZG	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.16	2 (3%)
2	NDP	YK	501	-	49,52,52	2.13	8 (16%)	66,80,80	1.66	13 (19%)
2	NDP	ZH	501	-	49,52,52	2.20	6 (12%)	66,80,80	1.71	15 (22%)
3	LMG	ZM	502	-	23,23,55	0.71	0	31,31,63	0.81	0
3	LMG	YP	502	-	23,23,55	0.71	0	31,31,63	0.81	0
3	LMG	YM	502	-	23,23,55	0.68	0	31,31,63	0.82	0
2	NDP	ZY	501	-	49,52,52	2.13	7 (14%)	66,80,80	1.68	14 (21%)
3	LMG	YR	502	-	23,23,55	0.67	0	31,31,63	0.92	0
4	A1JWG	ZA	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.18	2 (3%)
4	A1JWG	YS	503	-	48,53,53	1.39	5 (10%)	62,89,89	1.08	2 (3%)
3	LMG	ZH	502	-	23,23,55	0.71	0	31,31,63	0.80	0
4	A1JWG	ZU	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.08	2 (3%)
2	NDP	ZB	501	-	49,52,52	2.18	5 (10%)	66,80,80	1.72	14 (21%)
4	A1JWG	ZI	503	-	48,53,53	1.39	5 (10%)	62,89,89	1.08	4 (6%)
3	LMG	YJ	502	-	23,23,55	0.70	0	31,31,63	0.85	0
3	LMG	ZE	502	-	23,23,55	0.71	0	31,31,63	0.83	0
3	LMG	ZN	502	-	23,23,55	0.72	0	31,31,63	0.85	0
2	NDP	YR	501	-	49,52,52	2.19	6 (12%)	66,80,80	1.73	17 (25%)
4	A1JWG	ZD	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.31	4 (6%)
4	A1JWG	ZC	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.16	3 (4%)
3	LMG	ZO	502	-	23,23,55	0.69	0	31,31,63	0.76	0
2	NDP	YN	501	-	49,52,52	2.20	6 (12%)	66,80,80	1.71	15 (22%)
4	A1JWG	YQ	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.17	2 (3%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NDP	ZI	501	-	49,52,52	2.15	8 (16%)	66,80,80	1.67	13 (19%)
2	NDP	ZL	501	-	49,52,52	2.14	4 (8%)	66,80,80	1.71	15 (22%)
4	A1JWG	ZE	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.19	3 (4%)
3	LMG	YU	502	-	23,23,55	0.68	0	31,31,63	0.80	0
4	A1JWG	ZP	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.29	4 (6%)
2	NDP	YP	501	-	49,52,52	2.19	6 (12%)	66,80,80	1.70	14 (21%)
4	A1JWG	YN	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.28	4 (6%)
4	A1JWG	YX	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.12	3 (4%)
4	A1JWG	ZH	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.31	5 (8%)
2	NDP	ZS	501	-	49,52,52	2.15	4 (8%)	66,80,80	1.70	15 (22%)
2	NDP	ZW	501	-	49,52,52	2.13	4 (8%)	66,80,80	1.70	13 (19%)
4	A1JWG	YU	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.14	2 (3%)
2	NDP	ZR	501	-	49,52,52	2.21	6 (12%)	66,80,80	1.73	14 (21%)
3	LMG	YQ	502	-	23,23,55	0.71	0	31,31,63	0.85	0
3	LMG	ZU	502	-	23,23,55	0.71	0	31,31,63	0.83	0
4	A1JWG	YP	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.11	3 (4%)
2	NDP	ZG	501	-	49,52,52	2.15	4 (8%)	66,80,80	1.71	15 (22%)
2	NDP	ZM	501	-	49,52,52	2.14	7 (14%)	66,80,80	1.70	15 (22%)
4	A1JWG	YL	503	-	48,53,53	1.40	5 (10%)	62,89,89	1.11	3 (4%)
4	A1JWG	YY	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.17	2 (3%)
3	LMG	YV	502	-	23,23,55	0.71	0	31,31,63	0.82	0
2	NDP	YW	501	-	49,52,52	2.15	8 (16%)	66,80,80	1.68	14 (21%)
4	A1JWG	ZV	503	-	48,53,53	1.39	5 (10%)	62,89,89	1.12	3 (4%)
4	A1JWG	YR	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.32	5 (8%)
3	LMG	ZZ	502	-	23,23,55	0.71	0	31,31,63	0.83	0
3	LMG	ZP	502	-	23,23,55	0.72	0	31,31,63	0.85	0
4	A1JWG	ZM	503	-	48,53,53	1.39	5 (10%)	62,89,89	1.09	4 (6%)
4	A1JWG	YM	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.13	3 (4%)
3	LMG	ZT	502	-	23,23,55	0.68	0	31,31,63	0.88	1 (3%)
2	NDP	YZ	501	-	49,52,52	2.16	5 (10%)	66,80,80	1.72	16 (24%)
4	A1JWG	YW	503	-	48,53,53	1.38	5 (10%)	62,89,89	1.09	3 (4%)
2	NDP	ZE	501	-	49,52,52	2.11	7 (14%)	66,80,80	1.71	15 (22%)
4	A1JWG	ZO	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.18	2 (3%)
3	LMG	YO	502	-	23,23,55	0.71	0	31,31,63	0.80	0
3	LMG	YL	502	-	23,23,55	0.71	0	31,31,63	0.76	0
3	LMG	YW	502	-	23,23,55	0.72	0	31,31,63	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	A1JWG	ZS	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.14	2 (3%)
3	LMG	YZ	502	-	23,23,55	0.69	0	31,31,63	0.84	0
2	NDP	ZQ	501	-	49,52,52	2.14	8 (16%)	66,80,80	1.68	15 (22%)
3	LMG	ZR	502	-	23,23,55	0.70	0	31,31,63	0.78	0
2	NDP	ZF	501	-	49,52,52	2.19	4 (8%)	66,80,80	1.72	14 (21%)
3	LMG	ZC	502	-	23,23,55	0.69	0	31,31,63	0.84	0
4	A1JWG	ZK	503	-	48,53,53	1.37	5 (10%)	62,89,89	1.14	2 (3%)
2	NDP	ZJ	501	-	49,52,52	2.18	6 (12%)	66,80,80	1.71	13 (19%)
2	NDP	YV	501	-	49,52,52	2.21	5 (10%)	66,80,80	1.71	15 (22%)
2	NDP	YX	501	-	49,52,52	2.19	5 (10%)	66,80,80	1.72	14 (21%)
2	NDP	YU	501	-	49,52,52	2.16	4 (8%)	66,80,80	1.71	15 (22%)
2	NDP	ZC	501	-	49,52,52	2.14	4 (8%)	66,80,80	1.70	14 (21%)
2	NDP	ZN	501	-	49,52,52	2.19	5 (10%)	66,80,80	1.72	13 (19%)
2	NDP	ZV	501	-	49,52,52	2.18	6 (12%)	66,80,80	1.71	14 (21%)
3	LMG	ZI	502	-	23,23,55	0.72	0	31,31,63	0.80	0
3	LMG	YT	502	-	23,23,55	0.74	0	31,31,63	0.83	0
2	NDP	ZT	501	-	49,52,52	2.20	5 (10%)	66,80,80	1.71	17 (25%)
2	NDP	ZD	501	-	49,52,52	2.20	5 (10%)	66,80,80	1.72	17 (25%)
2	NDP	YO	501	-	49,52,52	2.16	8 (16%)	66,80,80	1.68	13 (19%)
2	NDP	YI	501	-	49,52,52	2.15	4 (8%)	66,80,80	1.72	15 (22%)
2	NDP	YJ	501	-	49,52,52	2.20	5 (10%)	66,80,80	1.71	17 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	A1JWG	ZT	503	-	3/3/11/15	8/15/91/91	-
2	NDP	YT	501	-	-	13/34/77/77	0/5/5/5
4	A1JWG	YK	503	-	3/3/11/15	7/15/91/91	-
3	LMG	ZS	502	-	-	2/16/36/70	0/1/1/1
2	NDP	YY	501	-	-	9/34/77/77	0/5/5/5
4	A1JWG	ZQ	503	-	3/3/11/15	10/15/91/91	-
4	A1JWG	YO	503	-	3/3/11/15	8/15/91/91	-
4	A1JWG	ZZ	503	-	3/3/11/15	9/15/91/91	-
4	A1JWG	ZY	503	-	3/3/11/15	6/15/91/91	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	LMG	ZB	502	-	-	2/16/36/70	0/1/1/1
3	LMG	ZG	502	-	-	5/16/36/70	0/1/1/1
3	LMG	ZJ	502	-	-	4/16/36/70	0/1/1/1
3	LMG	ZL	502	-	-	2/16/36/70	0/1/1/1
4	A1JWG	YZ	503	-	3/3/11/15	5/15/91/91	-
2	NDP	ZU	501	-	-	11/34/77/77	0/5/5/5
3	LMG	YX	502	-	-	6/16/36/70	0/1/1/1
4	A1JWG	ZR	503	-	3/3/11/15	9/15/91/91	-
3	LMG	ZV	502	-	-	8/16/36/70	0/1/1/1
2	NDP	YS	501	-	-	12/34/77/77	0/5/5/5
2	NDP	ZP	501	-	-	11/34/77/77	0/5/5/5
3	LMG	ZQ	502	-	-	2/16/36/70	0/1/1/1
3	LMG	YY	502	-	-	2/16/36/70	0/1/1/1
3	LMG	YI	502	-	-	4/16/36/70	0/1/1/1
4	A1JWG	ZN	503	-	3/3/11/15	9/15/91/91	-
3	LMG	ZA	502	-	-	3/16/36/70	0/1/1/1
3	LMG	YS	502	-	-	4/16/36/70	0/1/1/1
4	A1JWG	YT	503	-	3/3/11/15	7/15/91/91	-
4	A1JWG	ZL	503	-	3/3/11/15	10/15/91/91	-
2	NDP	YL	501	-	-	5/34/77/77	0/5/5/5
2	NDP	ZX	501	-	-	10/34/77/77	0/5/5/5
3	LMG	ZK	502	-	-	2/16/36/70	0/1/1/1
3	LMG	ZY	502	-	-	7/16/36/70	0/1/1/1
4	A1JWG	YV	503	-	3/3/11/15	7/15/91/91	-
2	NDP	ZA	501	-	-	10/34/77/77	0/5/5/5
3	LMG	ZF	502	-	-	3/16/36/70	0/1/1/1
2	NDP	ZZ	501	-	-	13/34/77/77	0/5/5/5
3	LMG	ZW	502	-	-	6/16/36/70	0/1/1/1
2	NDP	ZK	501	-	-	10/34/77/77	0/5/5/5
4	A1JWG	YI	503	-	3/3/11/15	9/15/91/91	-
4	A1JWG	YJ	503	-	3/3/11/15	8/15/91/91	-
4	A1JWG	ZF	503	-	3/3/11/15	8/15/91/91	-
3	LMG	ZX	502	-	-	6/16/36/70	0/1/1/1
4	A1JWG	ZB	503	-	2/2/11/15	8/15/91/91	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NDP	ZO	501	-	-	9/34/77/77	0/5/5/5
4	A1JWG	ZX	503	-	3/3/11/15	7/15/91/91	-
4	A1JWG	ZJ	503	-	3/3/11/15	8/15/91/91	-
4	A1JWG	ZW	503	-	3/3/11/15	9/15/91/91	-
2	NDP	YM	501	-	-	7/34/77/77	0/5/5/5
2	NDP	YQ	501	-	-	12/34/77/77	0/5/5/5
3	LMG	YK	502	-	-	2/16/36/70	0/1/1/1
3	LMG	YN	502	-	-	6/16/36/70	0/1/1/1
4	A1JWG	ZG	503	-	3/3/11/15	8/15/91/91	-
3	LMG	ZD	502	-	-	3/16/36/70	0/1/1/1
2	NDP	YK	501	-	-	12/34/77/77	0/5/5/5
2	NDP	ZH	501	-	-	15/34/77/77	0/5/5/5
3	LMG	ZM	502	-	-	7/16/36/70	0/1/1/1
3	LMG	YP	502	-	-	4/16/36/70	0/1/1/1
3	LMG	YM	502	-	-	4/16/36/70	0/1/1/1
2	NDP	ZY	501	-	-	13/34/77/77	0/5/5/5
3	LMG	YR	502	-	-	7/16/36/70	0/1/1/1
4	A1JWG	ZA	503	-	3/3/11/15	8/15/91/91	-
4	A1JWG	YS	503	-	3/3/11/15	10/15/91/91	-
3	LMG	ZH	502	-	-	4/16/36/70	0/1/1/1
4	A1JWG	ZU	503	-	3/3/11/15	10/15/91/91	-
2	NDP	ZB	501	-	-	5/34/77/77	0/5/5/5
4	A1JWG	ZI	503	-	3/3/11/15	7/15/91/91	-
3	LMG	YJ	502	-	-	7/16/36/70	0/1/1/1
3	LMG	ZE	502	-	-	3/16/36/70	0/1/1/1
4	A1JWG	ZD	503	-	3/3/11/15	9/15/91/91	-
2	NDP	YR	501	-	-	6/34/77/77	0/5/5/5
3	LMG	ZN	502	-	-	4/16/36/70	0/1/1/1
4	A1JWG	ZC	503	-	3/3/11/15	10/15/91/91	-
3	LMG	ZO	502	-	-	3/16/36/70	0/1/1/1
4	A1JWG	YQ	503	-	3/3/11/15	9/15/91/91	-
2	NDP	YN	501	-	-	10/34/77/77	0/5/5/5
2	NDP	ZI	501	-	-	10/34/77/77	0/5/5/5
4	A1JWG	ZE	503	-	3/3/11/15	7/15/91/91	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NDP	ZL	501	-	-	11/34/77/77	0/5/5/5
3	LMG	YU	502	-	-	2/16/36/70	0/1/1/1
4	A1JWG	ZP	503	-	3/3/11/15	9/15/91/91	-
2	NDP	YP	501	-	-	11/34/77/77	0/5/5/5
4	A1JWG	YN	503	-	3/3/11/15	9/15/91/91	-
4	A1JWG	YX	503	-	3/3/11/15	7/15/91/91	-
4	A1JWG	ZH	503	-	3/3/11/15	5/15/91/91	-
2	NDP	ZS	501	-	-	12/34/77/77	0/5/5/5
4	A1JWG	YU	503	-	3/3/11/15	9/15/91/91	-
2	NDP	ZW	501	-	-	11/34/77/77	0/5/5/5
2	NDP	ZR	501	-	-	7/34/77/77	0/5/5/5
3	LMG	YQ	502	-	-	2/16/36/70	0/1/1/1
3	LMG	ZU	502	-	-	6/16/36/70	0/1/1/1
4	A1JWG	YP	503	-	3/3/11/15	7/15/91/91	-
2	NDP	ZG	501	-	-	9/34/77/77	0/5/5/5
4	A1JWG	YY	503	-	3/3/11/15	9/15/91/91	-
4	A1JWG	YL	503	-	3/3/11/15	7/15/91/91	-
2	NDP	ZM	501	-	-	10/34/77/77	0/5/5/5
3	LMG	YV	502	-	-	8/16/36/70	0/1/1/1
4	A1JWG	ZV	503	-	3/3/11/15	6/15/91/91	-
2	NDP	YW	501	-	-	15/34/77/77	0/5/5/5
4	A1JWG	YR	503	-	3/3/11/15	7/15/91/91	-
3	LMG	ZZ	502	-	-	9/16/36/70	0/1/1/1
3	LMG	ZP	502	-	-	7/16/36/70	0/1/1/1
4	A1JWG	ZM	503	-	3/3/11/15	8/15/91/91	-
4	A1JWG	YM	503	-	3/3/11/15	8/15/91/91	-
3	LMG	ZT	502	-	-	12/16/36/70	0/1/1/1
4	A1JWG	YW	503	-	3/3/11/15	8/15/91/91	-
2	NDP	YZ	501	-	-	10/34/77/77	0/5/5/5
2	NDP	ZE	501	-	-	8/34/77/77	0/5/5/5
4	A1JWG	ZO	503	-	3/3/11/15	9/15/91/91	-
3	LMG	YO	502	-	-	6/16/36/70	0/1/1/1
3	LMG	YL	502	-	-	0/16/36/70	0/1/1/1
3	LMG	YW	502	-	-	10/16/36/70	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	A1JWG	ZS	503	-	3/3/11/15	9/15/91/91	-
3	LMG	YZ	502	-	-	4/16/36/70	0/1/1/1
2	NDP	ZQ	501	-	-	10/34/77/77	0/5/5/5
3	LMG	ZR	502	-	-	3/16/36/70	0/1/1/1
2	NDP	ZF	501	-	-	6/34/77/77	0/5/5/5
3	LMG	ZC	502	-	-	3/16/36/70	0/1/1/1
4	A1JWG	ZK	503	-	3/3/11/15	9/15/91/91	-
2	NDP	ZJ	501	-	-	11/34/77/77	0/5/5/5
2	NDP	YV	501	-	-	13/34/77/77	0/5/5/5
2	NDP	YX	501	-	-	5/34/77/77	0/5/5/5
2	NDP	YU	501	-	-	12/34/77/77	0/5/5/5
2	NDP	ZC	501	-	-	9/34/77/77	0/5/5/5
2	NDP	ZN	501	-	-	5/34/77/77	0/5/5/5
2	NDP	ZV	501	-	-	11/34/77/77	0/5/5/5
3	LMG	ZI	502	-	-	0/16/36/70	0/1/1/1
3	LMG	YT	502	-	-	8/16/36/70	0/1/1/1
2	NDP	ZT	501	-	-	11/34/77/77	0/5/5/5
2	NDP	ZD	501	-	-	11/34/77/77	0/5/5/5
2	NDP	YO	501	-	-	11/34/77/77	0/5/5/5
2	NDP	YI	501	-	-	10/34/77/77	0/5/5/5
2	NDP	YJ	501	-	-	12/34/77/77	0/5/5/5

All (466) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	ZR	501	NDP	P2B-O2B	12.63	1.83	1.59
2	ZH	501	NDP	P2B-O2B	12.63	1.83	1.59
2	ZO	501	NDP	P2B-O2B	12.61	1.83	1.59
2	YJ	501	NDP	P2B-O2B	12.60	1.83	1.59
2	ZP	501	NDP	P2B-O2B	12.57	1.83	1.59
2	YN	501	NDP	P2B-O2B	12.57	1.83	1.59
2	YR	501	NDP	P2B-O2B	12.56	1.83	1.59
2	YV	501	NDP	P2B-O2B	12.55	1.83	1.59
2	ZD	501	NDP	P2B-O2B	12.54	1.83	1.59
2	ZT	501	NDP	P2B-O2B	12.54	1.83	1.59
2	ZX	501	NDP	P2B-O2B	12.49	1.82	1.59
2	YL	501	NDP	P2B-O2B	12.48	1.82	1.59
2	ZN	501	NDP	P2B-O2B	12.46	1.82	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	ZF	501	NDP	P2B-O2B	12.46	1.82	1.59
2	YX	501	NDP	P2B-O2B	12.42	1.82	1.59
2	YP	501	NDP	P2B-O2B	12.41	1.82	1.59
2	YZ	501	NDP	P2B-O2B	12.40	1.82	1.59
2	ZB	501	NDP	P2B-O2B	12.40	1.82	1.59
2	YU	501	NDP	P2B-O2B	12.37	1.82	1.59
2	YT	501	NDP	P2B-O2B	12.35	1.82	1.59
2	ZK	501	NDP	P2B-O2B	12.35	1.82	1.59
2	ZJ	501	NDP	P2B-O2B	12.34	1.82	1.59
2	ZS	501	NDP	P2B-O2B	12.34	1.82	1.59
2	ZV	501	NDP	P2B-O2B	12.33	1.82	1.59
2	ZG	501	NDP	P2B-O2B	12.32	1.82	1.59
2	ZC	501	NDP	P2B-O2B	12.30	1.82	1.59
2	YO	501	NDP	P2B-O2B	12.30	1.82	1.59
2	YQ	501	NDP	P2B-O2B	12.29	1.82	1.59
2	ZL	501	NDP	P2B-O2B	12.27	1.82	1.59
2	YI	501	NDP	P2B-O2B	12.27	1.82	1.59
2	ZZ	501	NDP	P2B-O2B	12.27	1.82	1.59
2	YM	501	NDP	P2B-O2B	12.24	1.82	1.59
2	YW	501	NDP	P2B-O2B	12.20	1.82	1.59
2	ZW	501	NDP	P2B-O2B	12.19	1.82	1.59
2	ZQ	501	NDP	P2B-O2B	12.19	1.82	1.59
2	ZI	501	NDP	P2B-O2B	12.18	1.82	1.59
2	YY	501	NDP	P2B-O2B	12.14	1.82	1.59
2	ZM	501	NDP	P2B-O2B	12.14	1.82	1.59
2	YS	501	NDP	P2B-O2B	12.13	1.82	1.59
2	ZU	501	NDP	P2B-O2B	12.11	1.82	1.59
2	YK	501	NDP	P2B-O2B	12.06	1.82	1.59
2	ZY	501	NDP	P2B-O2B	12.01	1.82	1.59
2	ZE	501	NDP	P2B-O2B	11.91	1.81	1.59
2	ZA	501	NDP	P2B-O2B	11.87	1.81	1.59
4	ZF	503	A1JWG	MG-NA	4.84	2.17	2.06
4	YP	503	A1JWG	MG-NA	4.84	2.17	2.06
4	ZJ	503	A1JWG	MG-NA	4.84	2.17	2.06
4	YX	503	A1JWG	MG-NA	4.84	2.17	2.06
4	ZB	503	A1JWG	MG-NA	4.83	2.17	2.06
4	ZR	503	A1JWG	MG-NA	4.83	2.17	2.06
4	ZN	503	A1JWG	MG-NA	4.83	2.17	2.06
4	YL	503	A1JWG	MG-NA	4.82	2.17	2.06
4	YS	503	A1JWG	MG-NA	4.82	2.17	2.06
4	ZU	503	A1JWG	MG-NA	4.81	2.17	2.06
4	ZQ	503	A1JWG	MG-NA	4.80	2.17	2.06

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	ZC	503	A1JWG	MG-NA	4.80	2.17	2.06
4	YT	503	A1JWG	MG-NA	4.79	2.17	2.06
4	YQ	503	A1JWG	MG-NA	4.79	2.17	2.06
4	YI	503	A1JWG	MG-NA	4.79	2.17	2.06
4	YW	503	A1JWG	MG-NA	4.79	2.17	2.06
4	ZV	503	A1JWG	MG-NA	4.78	2.17	2.06
4	ZZ	503	A1JWG	MG-NA	4.77	2.17	2.06
4	ZW	503	A1JWG	MG-NA	4.77	2.17	2.06
4	ZM	503	A1JWG	MG-NA	4.77	2.17	2.06
4	YY	503	A1JWG	MG-NA	4.77	2.17	2.06
4	ZI	503	A1JWG	MG-NA	4.77	2.17	2.06
4	YU	503	A1JWG	MG-NA	4.77	2.17	2.06
4	ZL	503	A1JWG	MG-NA	4.76	2.17	2.06
4	ZG	503	A1JWG	MG-NA	4.76	2.17	2.06
4	ZS	503	A1JWG	MG-NA	4.76	2.17	2.06
4	ZK	503	A1JWG	MG-NA	4.76	2.17	2.06
4	YM	503	A1JWG	MG-NA	4.76	2.17	2.06
4	ZX	503	A1JWG	MG-NA	4.74	2.17	2.06
4	ZP	503	A1JWG	MG-NA	4.74	2.17	2.06
4	ZD	503	A1JWG	MG-NA	4.72	2.17	2.06
4	ZY	503	A1JWG	MG-NA	4.72	2.17	2.06
4	ZO	503	A1JWG	MG-NA	4.72	2.17	2.06
4	YN	503	A1JWG	MG-NA	4.72	2.17	2.06
4	YO	503	A1JWG	MG-NA	4.72	2.17	2.06
4	ZA	503	A1JWG	MG-NA	4.72	2.17	2.06
4	YJ	503	A1JWG	MG-NA	4.71	2.17	2.06
4	YZ	503	A1JWG	MG-NA	4.70	2.17	2.06
4	YV	503	A1JWG	MG-NA	4.70	2.17	2.06
4	ZT	503	A1JWG	MG-NA	4.70	2.17	2.06
4	ZE	503	A1JWG	MG-NA	4.70	2.17	2.06
4	YK	503	A1JWG	MG-NA	4.69	2.17	2.06
4	YR	503	A1JWG	MG-NA	4.66	2.17	2.06
4	ZH	503	A1JWG	MG-NA	4.65	2.17	2.06
4	YT	503	A1JWG	C4D-ND	-4.31	1.31	1.37
4	ZB	503	A1JWG	C4D-ND	-4.28	1.31	1.37
4	YL	503	A1JWG	C4D-ND	-4.28	1.31	1.37
4	YX	503	A1JWG	C4D-ND	-4.25	1.31	1.37
4	YP	503	A1JWG	C4D-ND	-4.25	1.31	1.37
4	ZR	503	A1JWG	C4D-ND	-4.24	1.31	1.37
4	ZZ	503	A1JWG	C4D-ND	-4.24	1.31	1.37
4	ZV	503	A1JWG	C4D-ND	-4.23	1.31	1.37
4	ZJ	503	A1JWG	C4D-ND	-4.23	1.31	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	ZN	503	A1JWG	C4D-ND	-4.21	1.31	1.37
4	ZF	503	A1JWG	C4D-ND	-4.20	1.31	1.37
4	ZU	503	A1JWG	C4D-ND	-4.19	1.31	1.37
4	YV	503	A1JWG	C4D-ND	-4.19	1.31	1.37
4	ZH	503	A1JWG	C4D-ND	-4.19	1.31	1.37
4	ZQ	503	A1JWG	C4D-ND	-4.18	1.32	1.37
4	ZI	503	A1JWG	C4D-ND	-4.16	1.32	1.37
4	ZM	503	A1JWG	C4D-ND	-4.15	1.32	1.37
4	YR	503	A1JWG	C4D-ND	-4.14	1.32	1.37
4	YW	503	A1JWG	C4D-ND	-4.12	1.32	1.37
4	YS	503	A1JWG	C4D-ND	-4.12	1.32	1.37
4	ZD	503	A1JWG	C4D-ND	-4.08	1.32	1.37
4	ZP	503	A1JWG	C4D-ND	-4.08	1.32	1.37
4	YJ	503	A1JWG	C4D-ND	-4.06	1.32	1.37
4	YN	503	A1JWG	C4D-ND	-4.05	1.32	1.37
4	ZT	503	A1JWG	C4D-ND	-4.02	1.32	1.37
4	YO	503	A1JWG	C4D-ND	-3.97	1.32	1.37
2	ZD	501	NDP	PN-O5D	3.95	1.75	1.59
4	ZY	503	A1JWG	C4D-ND	-3.95	1.32	1.37
2	ZZ	501	NDP	PN-O5D	3.94	1.75	1.59
2	YT	501	NDP	PN-O5D	3.94	1.75	1.59
4	ZE	503	A1JWG	C4D-ND	-3.94	1.32	1.37
4	YK	503	A1JWG	C4D-ND	-3.94	1.32	1.37
4	ZA	503	A1JWG	C4D-ND	-3.93	1.32	1.37
2	ZT	501	NDP	PN-O5D	3.93	1.75	1.59
2	YP	501	NDP	PN-O5D	3.92	1.75	1.59
2	ZP	501	NDP	PN-O5D	3.91	1.75	1.59
4	ZL	503	A1JWG	C4D-ND	-3.88	1.32	1.37
4	YZ	503	A1JWG	C4D-ND	-3.88	1.32	1.37
2	YL	501	NDP	PN-O5D	3.87	1.75	1.59
2	ZB	501	NDP	PN-O5D	3.87	1.75	1.59
2	YJ	501	NDP	PN-O5D	3.87	1.75	1.59
2	ZJ	501	NDP	PN-O5D	3.86	1.74	1.59
4	YQ	503	A1JWG	C4D-ND	-3.86	1.32	1.37
2	ZY	501	NDP	PN-O5D	3.85	1.74	1.59
4	ZO	503	A1JWG	C4D-ND	-3.85	1.32	1.37
4	YU	503	A1JWG	C4D-ND	-3.85	1.32	1.37
2	ZA	501	NDP	PN-O5D	3.85	1.74	1.59
4	ZX	503	A1JWG	C4D-ND	-3.85	1.32	1.37
2	ZR	501	NDP	PN-O5D	3.85	1.74	1.59
2	ZV	501	NDP	PN-O5D	3.84	1.74	1.59
2	ZF	501	NDP	PN-O5D	3.84	1.74	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	ZC	503	A1JWG	C4D-ND	-3.84	1.32	1.37
2	ZU	501	NDP	PN-O5D	3.84	1.74	1.59
2	YV	501	NDP	PN-O5D	3.84	1.74	1.59
4	ZK	503	A1JWG	C4D-ND	-3.83	1.32	1.37
2	ZI	501	NDP	PN-O5D	3.83	1.74	1.59
2	YK	501	NDP	PN-O5D	3.83	1.74	1.59
4	YM	503	A1JWG	C4D-ND	-3.83	1.32	1.37
4	YY	503	A1JWG	C4D-ND	-3.83	1.32	1.37
4	ZW	503	A1JWG	C4D-ND	-3.83	1.32	1.37
2	YX	501	NDP	PN-O5D	3.83	1.74	1.59
4	ZG	503	A1JWG	C4D-ND	-3.82	1.32	1.37
4	ZS	503	A1JWG	C4D-ND	-3.82	1.32	1.37
2	ZN	501	NDP	PN-O5D	3.81	1.74	1.59
2	YS	501	NDP	PN-O5D	3.81	1.74	1.59
2	YR	501	NDP	PN-O5D	3.81	1.74	1.59
4	YI	503	A1JWG	C4D-ND	-3.80	1.32	1.37
2	YO	501	NDP	PN-O5D	3.79	1.74	1.59
2	YW	501	NDP	PN-O5D	3.78	1.74	1.59
2	ZH	501	NDP	PN-O5D	3.78	1.74	1.59
2	YM	501	NDP	PN-O5D	3.76	1.74	1.59
2	YY	501	NDP	PN-O5D	3.76	1.74	1.59
2	ZM	501	NDP	PN-O5D	3.75	1.74	1.59
2	YZ	501	NDP	PN-O5D	3.74	1.74	1.59
2	ZE	501	NDP	PN-O5D	3.74	1.74	1.59
2	YU	501	NDP	PN-O5D	3.73	1.74	1.59
2	YI	501	NDP	PN-O5D	3.72	1.74	1.59
2	YQ	501	NDP	PN-O5D	3.71	1.74	1.59
2	ZO	501	NDP	PN-O5D	3.70	1.74	1.59
2	ZG	501	NDP	PN-O5D	3.70	1.74	1.59
2	ZC	501	NDP	PN-O5D	3.70	1.74	1.59
2	ZQ	501	NDP	PN-O5D	3.69	1.74	1.59
2	ZW	501	NDP	PN-O5D	3.69	1.74	1.59
2	ZS	501	NDP	PN-O5D	3.66	1.74	1.59
2	ZX	501	NDP	PN-O5D	3.66	1.74	1.59
2	ZK	501	NDP	PN-O5D	3.64	1.74	1.59
2	YN	501	NDP	PN-O5D	3.63	1.74	1.59
2	ZL	501	NDP	PN-O5D	3.56	1.73	1.59
2	ZA	501	NDP	O2B-C2B	-3.21	1.32	1.44
2	ZM	501	NDP	O2B-C2B	-3.20	1.32	1.44
2	ZE	501	NDP	O2B-C2B	-3.20	1.32	1.44
2	YS	501	NDP	O2B-C2B	-3.17	1.32	1.44
2	YW	501	NDP	O2B-C2B	-3.16	1.32	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	ZU	501	NDP	O2B-C2B	-3.14	1.32	1.44
2	YK	501	NDP	O2B-C2B	-3.14	1.32	1.44
2	ZY	501	NDP	O2B-C2B	-3.14	1.32	1.44
2	ZI	501	NDP	O2B-C2B	-3.14	1.32	1.44
2	ZV	501	NDP	O2B-C2B	-3.13	1.32	1.44
2	ZJ	501	NDP	O2B-C2B	-3.12	1.32	1.44
2	YO	501	NDP	O2B-C2B	-3.11	1.32	1.44
2	YP	501	NDP	O2B-C2B	-3.11	1.32	1.44
2	YX	501	NDP	O2B-C2B	-3.10	1.32	1.44
2	ZZ	501	NDP	O2B-C2B	-3.10	1.32	1.44
2	ZB	501	NDP	O2B-C2B	-3.09	1.32	1.44
2	YY	501	NDP	O2B-C2B	-3.09	1.32	1.44
2	ZQ	501	NDP	O2B-C2B	-3.09	1.32	1.44
2	YI	501	NDP	O2B-C2B	-3.08	1.32	1.44
2	YM	501	NDP	O2B-C2B	-3.08	1.32	1.44
2	ZF	501	NDP	O2B-C2B	-3.08	1.32	1.44
2	YL	501	NDP	O2B-C2B	-3.08	1.32	1.44
2	YT	501	NDP	O2B-C2B	-3.07	1.32	1.44
2	ZN	501	NDP	O2B-C2B	-3.06	1.32	1.44
2	ZR	501	NDP	O2B-C2B	-3.04	1.33	1.44
2	ZW	501	NDP	O2B-C2B	-3.03	1.33	1.44
2	ZG	501	NDP	O2B-C2B	-3.02	1.33	1.44
2	ZL	501	NDP	O2B-C2B	-3.02	1.33	1.44
2	ZS	501	NDP	O2B-C2B	-3.02	1.33	1.44
2	ZC	501	NDP	O2B-C2B	-3.02	1.33	1.44
2	YQ	501	NDP	O2B-C2B	-3.01	1.33	1.44
2	ZK	501	NDP	O2B-C2B	-3.01	1.33	1.44
2	YU	501	NDP	O2B-C2B	-3.01	1.33	1.44
2	YZ	501	NDP	O2B-C2B	-3.00	1.33	1.44
2	ZX	501	NDP	O2B-C2B	-2.99	1.33	1.44
2	YR	501	NDP	O2B-C2B	-2.99	1.33	1.44
2	ZO	501	NDP	O2B-C2B	-2.99	1.33	1.44
2	ZH	501	NDP	O2B-C2B	-2.98	1.33	1.44
2	ZT	501	NDP	O2B-C2B	-2.97	1.33	1.44
2	YV	501	NDP	O2B-C2B	-2.97	1.33	1.44
2	ZD	501	NDP	O2B-C2B	-2.97	1.33	1.44
2	YJ	501	NDP	O2B-C2B	-2.96	1.33	1.44
2	ZP	501	NDP	O2B-C2B	-2.96	1.33	1.44
2	YN	501	NDP	O2B-C2B	-2.96	1.33	1.44
2	YN	501	NDP	O5D-C5D	-2.43	1.35	1.44
4	ZC	503	A1JWG	C1D-ND	2.41	1.40	1.37
4	ZX	503	A1JWG	C1D-ND	2.39	1.40	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	YQ	503	A1JWG	C1D-ND	2.38	1.40	1.37
4	YZ	503	A1JWG	C1D-ND	2.38	1.40	1.37
4	YI	503	A1JWG	C1D-ND	2.38	1.40	1.37
4	ZL	503	A1JWG	C1D-ND	2.38	1.40	1.37
2	ZH	501	NDP	O3D-C3D	-2.37	1.37	1.43
4	YM	503	A1JWG	C1D-ND	2.37	1.40	1.37
4	ZW	503	A1JWG	C1D-ND	2.36	1.40	1.37
4	YY	503	A1JWG	C1D-ND	2.36	1.40	1.37
4	ZG	503	A1JWG	C1D-ND	2.36	1.40	1.37
4	ZS	503	A1JWG	C1D-ND	2.35	1.40	1.37
4	ZK	503	A1JWG	C1D-ND	2.35	1.40	1.37
4	YU	503	A1JWG	C1D-ND	2.34	1.40	1.37
4	ZO	503	A1JWG	C1D-ND	2.32	1.40	1.37
2	YW	501	NDP	O5D-C5D	-2.32	1.35	1.44
2	ZP	501	NDP	O5D-C5D	-2.30	1.35	1.44
2	ZB	501	NDP	C7N-C3N	-2.29	1.43	1.48
2	YL	501	NDP	C7N-C3N	-2.29	1.43	1.48
2	ZY	501	NDP	O5D-C5D	-2.28	1.36	1.44
2	ZJ	501	NDP	C7N-C3N	-2.28	1.43	1.48
2	YJ	501	NDP	O5D-C5D	-2.28	1.36	1.44
2	ZD	501	NDP	O5D-C5D	-2.27	1.36	1.44
2	YS	501	NDP	O5D-C5D	-2.27	1.36	1.44
2	ZE	501	NDP	O3B-C3B	-2.27	1.37	1.43
4	YM	503	A1JWG	C3D-C4D	2.27	1.45	1.39
2	YN	501	NDP	O3D-C3D	-2.27	1.37	1.43
2	ZR	501	NDP	C7N-C3N	-2.27	1.43	1.48
2	ZT	501	NDP	O5D-C5D	-2.27	1.36	1.44
2	ZN	501	NDP	C7N-C3N	-2.26	1.43	1.48
2	YX	501	NDP	C7N-C3N	-2.26	1.43	1.48
2	ZF	501	NDP	C7N-C3N	-2.26	1.43	1.48
4	YI	503	A1JWG	C3D-C4D	2.25	1.45	1.39
2	YS	501	NDP	O3B-C3B	-2.25	1.37	1.43
2	ZY	501	NDP	O3B-C3B	-2.25	1.37	1.43
2	ZD	501	NDP	O3D-C3D	-2.24	1.37	1.43
4	ZG	503	A1JWG	C3D-C4D	2.24	1.45	1.39
2	YP	501	NDP	C7N-C3N	-2.24	1.43	1.48
4	YQ	503	A1JWG	C3D-C4D	2.24	1.45	1.39
4	ZC	503	A1JWG	C3D-C4D	2.23	1.45	1.39
4	ZC	503	A1JWG	C4B-NB	2.23	1.40	1.37
4	YY	503	A1JWG	C3D-C4D	2.23	1.45	1.39
2	YT	501	NDP	C7N-C3N	-2.23	1.43	1.48
4	ZL	503	A1JWG	C3D-C4D	2.23	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	YZ	503	A1JWG	C3D-C4D	2.23	1.45	1.39
4	ZX	503	A1JWG	C3D-C4D	2.23	1.45	1.39
2	YW	501	NDP	O3B-C3B	-2.22	1.37	1.43
4	YS	503	A1JWG	C4B-NB	2.22	1.40	1.37
2	YK	501	NDP	O3B-C3B	-2.22	1.37	1.43
4	ZS	503	A1JWG	C3D-C4D	2.22	1.45	1.39
2	ZU	501	NDP	O3B-C3B	-2.22	1.37	1.43
4	YU	503	A1JWG	C3D-C4D	2.22	1.45	1.39
4	ZA	503	A1JWG	C1D-ND	2.22	1.40	1.37
4	ZK	503	A1JWG	C3D-C4D	2.22	1.45	1.39
4	ZY	503	A1JWG	C1D-ND	2.22	1.40	1.37
2	ZV	501	NDP	C7N-C3N	-2.22	1.43	1.48
4	ZW	503	A1JWG	C3D-C4D	2.22	1.45	1.39
2	ZZ	501	NDP	C7N-C3N	-2.22	1.43	1.48
4	ZO	503	A1JWG	C3D-C4D	2.21	1.45	1.39
4	YM	503	A1JWG	C4B-NB	2.21	1.40	1.37
4	ZU	503	A1JWG	C4B-NB	2.21	1.40	1.37
4	YW	503	A1JWG	C4B-NB	2.21	1.40	1.37
2	ZA	501	NDP	O3B-C3B	-2.21	1.37	1.43
2	ZP	501	NDP	O3D-C3D	-2.21	1.37	1.43
4	ZN	503	A1JWG	C1D-ND	2.21	1.40	1.37
4	ZX	503	A1JWG	C4B-NB	2.21	1.40	1.37
4	YO	503	A1JWG	C3D-C4D	2.21	1.45	1.39
4	ZA	503	A1JWG	C3D-C4D	2.21	1.45	1.39
4	ZF	503	A1JWG	C1D-ND	2.21	1.40	1.37
4	ZY	503	A1JWG	C4B-NB	2.21	1.40	1.37
2	ZI	501	NDP	O3B-C3B	-2.20	1.37	1.43
4	ZE	503	A1JWG	C3D-C4D	2.20	1.45	1.39
4	ZY	503	A1JWG	C3D-C4D	2.20	1.45	1.39
4	YK	503	A1JWG	C3D-C4D	2.20	1.45	1.39
2	ZQ	501	NDP	O3B-C3B	-2.20	1.37	1.43
2	ZT	501	NDP	O3D-C3D	-2.20	1.37	1.43
4	YZ	503	A1JWG	C4B-NB	2.20	1.40	1.37
4	ZA	503	A1JWG	C4B-NB	2.20	1.40	1.37
4	YO	503	A1JWG	C4B-NB	2.20	1.40	1.37
4	YP	503	A1JWG	C1D-ND	2.20	1.40	1.37
4	ZL	503	A1JWG	C4B-NB	2.20	1.40	1.37
4	ZJ	503	A1JWG	C1D-ND	2.20	1.40	1.37
4	ZS	503	A1JWG	C4B-NB	2.20	1.40	1.37
4	YJ	503	A1JWG	C1D-ND	2.20	1.40	1.37
4	YO	503	A1JWG	C1D-ND	2.20	1.40	1.37
4	YY	503	A1JWG	C4B-NB	2.20	1.40	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	YU	503	A1JWG	C4B-NB	2.19	1.40	1.37
4	ZB	503	A1JWG	C1D-ND	2.19	1.40	1.37
4	ZK	503	A1JWG	C4B-NB	2.19	1.40	1.37
4	ZQ	503	A1JWG	C4B-NB	2.19	1.40	1.37
2	YR	501	NDP	O3D-C3D	-2.18	1.37	1.43
4	YX	503	A1JWG	C1D-ND	2.18	1.40	1.37
4	ZM	503	A1JWG	C4B-NB	2.18	1.40	1.37
2	ZM	501	NDP	O3B-C3B	-2.18	1.37	1.43
4	YK	503	A1JWG	C4B-NB	2.18	1.40	1.37
4	YK	503	A1JWG	C1D-ND	2.18	1.40	1.37
4	ZE	503	A1JWG	C1D-ND	2.18	1.40	1.37
4	ZI	503	A1JWG	C4B-NB	2.18	1.40	1.37
2	YJ	501	NDP	O3D-C3D	-2.18	1.37	1.43
4	YN	503	A1JWG	C1D-ND	2.18	1.40	1.37
4	ZG	503	A1JWG	C4B-NB	2.18	1.40	1.37
4	ZE	503	A1JWG	C4B-NB	2.17	1.40	1.37
2	YR	501	NDP	O5D-C5D	-2.17	1.36	1.44
4	ZR	503	A1JWG	C1D-ND	2.17	1.40	1.37
4	ZU	503	A1JWG	C1D-ND	2.17	1.40	1.37
4	ZT	503	A1JWG	C1D-ND	2.17	1.40	1.37
4	ZD	503	A1JWG	C1D-ND	2.17	1.40	1.37
4	YI	503	A1JWG	C4B-NB	2.17	1.40	1.37
4	ZO	503	A1JWG	C4B-NB	2.17	1.40	1.37
4	YS	503	A1JWG	C1D-ND	2.17	1.40	1.37
2	YO	501	NDP	O3B-C3B	-2.16	1.37	1.43
4	ZP	503	A1JWG	C1D-ND	2.16	1.40	1.37
4	ZZ	503	A1JWG	C1D-ND	2.16	1.40	1.37
4	YQ	503	A1JWG	C4B-NB	2.16	1.40	1.37
4	YT	503	A1JWG	C1D-ND	2.16	1.40	1.37
4	ZN	503	A1JWG	C4B-NB	2.16	1.40	1.37
4	ZQ	503	A1JWG	C1D-ND	2.15	1.40	1.37
4	YJ	503	A1JWG	C3D-C4D	2.15	1.45	1.39
4	ZI	503	A1JWG	C1D-ND	2.15	1.40	1.37
4	ZT	503	A1JWG	C3D-C4D	2.15	1.45	1.39
4	ZF	503	A1JWG	C4B-NB	2.15	1.40	1.37
4	YL	503	A1JWG	C1D-ND	2.15	1.40	1.37
2	YQ	501	NDP	O5D-C5D	-2.15	1.36	1.44
4	ZV	503	A1JWG	C1D-ND	2.14	1.40	1.37
2	ZK	501	NDP	O5D-C5D	-2.14	1.36	1.44
4	YN	503	A1JWG	C3D-C4D	2.14	1.45	1.39
4	ZD	503	A1JWG	C3D-C4D	2.14	1.45	1.39
4	ZM	503	A1JWG	C1D-ND	2.14	1.40	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	ZG	501	NDP	O5D-C5D	-2.14	1.36	1.44
4	YX	503	A1JWG	C4B-NB	2.14	1.40	1.37
4	ZJ	503	A1JWG	C4B-NB	2.14	1.40	1.37
4	ZP	503	A1JWG	C3D-C4D	2.14	1.45	1.39
4	ZR	503	A1JWG	C4B-NB	2.14	1.40	1.37
2	ZO	501	NDP	O5D-C5D	-2.14	1.36	1.44
4	ZW	503	A1JWG	C4B-NB	2.14	1.40	1.37
2	ZZ	501	NDP	O5D-C5D	-2.13	1.36	1.44
4	YV	503	A1JWG	C1D-ND	2.13	1.40	1.37
4	ZP	503	A1JWG	C4B-NB	2.13	1.40	1.37
4	YP	503	A1JWG	C4B-NB	2.13	1.40	1.37
4	YV	503	A1JWG	C3D-C4D	2.13	1.45	1.39
2	ZC	501	NDP	O5D-C5D	-2.13	1.36	1.44
2	YS	501	NDP	O5B-C5B	-2.13	1.36	1.44
4	ZV	503	A1JWG	C4B-NB	2.12	1.40	1.37
4	ZD	503	A1JWG	C4B-NB	2.12	1.40	1.37
4	ZZ	503	A1JWG	C3D-C4D	2.12	1.45	1.39
4	YN	503	A1JWG	C4B-NB	2.12	1.40	1.37
4	YR	503	A1JWG	C1D-ND	2.12	1.40	1.37
4	YR	503	A1JWG	C3D-C4D	2.12	1.45	1.39
4	ZB	503	A1JWG	C4B-NB	2.12	1.40	1.37
4	ZR	503	A1JWG	C3D-C4D	2.12	1.45	1.39
2	YO	501	NDP	O5D-C5D	-2.12	1.36	1.44
2	YU	501	NDP	O5D-C5D	-2.12	1.36	1.44
2	YY	501	NDP	O5D-C5D	-2.12	1.36	1.44
4	YV	503	A1JWG	C4B-NB	2.12	1.40	1.37
4	ZH	503	A1JWG	C3D-C4D	2.11	1.45	1.39
4	ZZ	503	A1JWG	C4B-NB	2.11	1.40	1.37
2	YW	501	NDP	O5B-C5B	-2.11	1.36	1.44
2	ZQ	501	NDP	O5D-C5D	-2.11	1.36	1.44
4	YJ	503	A1JWG	C4B-NB	2.11	1.40	1.37
4	YL	503	A1JWG	C4B-NB	2.11	1.40	1.37
4	YW	503	A1JWG	C1D-ND	2.11	1.40	1.37
2	YK	501	NDP	O5D-C5D	-2.11	1.36	1.44
2	ZS	501	NDP	O5D-C5D	-2.11	1.36	1.44
2	ZL	501	NDP	O5D-C5D	-2.11	1.36	1.44
2	ZU	501	NDP	O5B-C5B	-2.11	1.36	1.44
2	YV	501	NDP	O5D-C5D	-2.11	1.36	1.44
4	ZH	503	A1JWG	C1D-ND	2.10	1.40	1.37
4	YL	503	A1JWG	C3D-C4D	2.10	1.45	1.39
2	YT	501	NDP	O5D-C5D	-2.10	1.36	1.44
4	ZN	503	A1JWG	C3D-C4D	2.10	1.45	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	YI	501	NDP	O5D-C5D	-2.10	1.36	1.44
4	ZF	503	A1JWG	C3D-C4D	2.10	1.45	1.39
2	ZU	501	NDP	O5D-C5D	-2.10	1.36	1.44
2	YM	501	NDP	O5D-C5D	-2.10	1.36	1.44
2	ZE	501	NDP	O5D-C5D	-2.10	1.36	1.44
4	ZT	503	A1JWG	C4B-NB	2.10	1.40	1.37
2	ZM	501	NDP	O5D-C5D	-2.10	1.36	1.44
4	YP	503	A1JWG	C3D-C4D	2.10	1.45	1.39
4	YR	503	A1JWG	C4B-NB	2.10	1.40	1.37
4	ZM	503	A1JWG	C3D-C4D	2.10	1.45	1.39
2	ZW	501	NDP	O5D-C5D	-2.10	1.36	1.44
4	YS	503	A1JWG	C3D-C4D	2.10	1.45	1.39
4	ZV	503	A1JWG	C3D-C4D	2.10	1.45	1.39
2	ZI	501	NDP	O5D-C5D	-2.09	1.36	1.44
2	YP	501	NDP	O3D-C3D	-2.09	1.38	1.43
2	ZX	501	NDP	O5D-C5D	-2.09	1.36	1.44
4	YX	503	A1JWG	C3D-C4D	2.09	1.45	1.39
4	ZU	503	A1JWG	C3D-C4D	2.09	1.45	1.39
2	ZQ	501	NDP	C7N-C3N	-2.09	1.44	1.48
2	ZA	501	NDP	O5B-C5B	-2.09	1.36	1.44
4	ZB	503	A1JWG	C3D-C4D	2.09	1.45	1.39
4	ZI	503	A1JWG	C3D-C4D	2.08	1.45	1.39
4	ZJ	503	A1JWG	C3D-C4D	2.08	1.45	1.39
2	ZY	501	NDP	O5B-C5B	-2.08	1.36	1.44
4	YW	503	A1JWG	C3D-C4D	2.08	1.45	1.39
4	YT	503	A1JWG	C3D-C4D	2.08	1.45	1.39
2	ZV	501	NDP	O3D-C3D	-2.08	1.38	1.43
2	YO	501	NDP	O5B-C5B	-2.07	1.36	1.44
2	ZA	501	NDP	O5D-C5D	-2.07	1.36	1.44
2	YW	501	NDP	O3D-C3D	-2.07	1.38	1.43
4	ZH	503	A1JWG	C4B-NB	2.07	1.40	1.37
4	ZQ	503	A1JWG	C3D-C4D	2.07	1.45	1.39
2	ZJ	501	NDP	O3D-C3D	-2.07	1.38	1.43
2	ZE	501	NDP	C7N-C3N	-2.07	1.44	1.48
2	YS	501	NDP	O3D-C3D	-2.07	1.38	1.43
2	ZI	501	NDP	O5B-C5B	-2.06	1.36	1.44
2	YZ	501	NDP	O5D-C5D	-2.06	1.36	1.44
2	ZA	501	NDP	C7N-C3N	-2.06	1.44	1.48
2	ZY	501	NDP	O3D-C3D	-2.05	1.38	1.43
2	YK	501	NDP	C7N-C3N	-2.05	1.44	1.48
2	ZU	501	NDP	O3D-C3D	-2.05	1.38	1.43
2	YS	501	NDP	C7N-C3N	-2.05	1.44	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	ZH	501	NDP	O3B-C3B	-2.04	1.38	1.43
2	ZO	501	NDP	O3B-C3B	-2.04	1.38	1.43
2	ZJ	501	NDP	O5D-C5D	-2.04	1.36	1.44
2	YR	501	NDP	O3B-C3B	-2.03	1.38	1.43
2	YO	501	NDP	O3D-C3D	-2.03	1.38	1.43
2	ZI	501	NDP	O3D-C3D	-2.03	1.38	1.43
2	ZV	501	NDP	O5D-C5D	-2.03	1.37	1.44
2	YN	501	NDP	O3B-C3B	-2.03	1.38	1.43
2	YK	501	NDP	O3D-C3D	-2.03	1.38	1.43
2	ZQ	501	NDP	O5B-C5B	-2.03	1.37	1.44
2	YV	501	NDP	O3B-C3B	-2.03	1.38	1.43
2	YK	501	NDP	O5B-C5B	-2.02	1.37	1.44
2	ZZ	501	NDP	O3D-C3D	-2.02	1.38	1.43
2	YT	501	NDP	O3D-C3D	-2.02	1.38	1.43
2	ZH	501	NDP	O5B-C5B	-2.02	1.37	1.44
2	YZ	501	NDP	C7N-C3N	-2.02	1.44	1.48
2	YP	501	NDP	O5D-C5D	-2.02	1.37	1.44
2	ZA	501	NDP	O3D-C3D	-2.01	1.38	1.43
2	ZR	501	NDP	O3D-C3D	-2.01	1.38	1.43
2	ZP	501	NDP	O3B-C3B	-2.01	1.38	1.43
2	YO	501	NDP	C7N-C3N	-2.01	1.44	1.48
2	ZB	501	NDP	O5D-C5D	-2.01	1.37	1.44
2	ZK	501	NDP	O3B-C3B	-2.01	1.38	1.43
2	ZU	501	NDP	C7N-C3N	-2.01	1.44	1.48
2	ZE	501	NDP	O3D-C3D	-2.01	1.38	1.43
2	YW	501	NDP	C7N-C3N	-2.01	1.44	1.48
2	ZR	501	NDP	O5D-C5D	-2.01	1.37	1.44
2	ZQ	501	NDP	O3D-C3D	-2.00	1.38	1.43
2	ZN	501	NDP	O5D-C5D	-2.00	1.37	1.44
2	ZI	501	NDP	C7N-C3N	-2.00	1.44	1.48
2	ZM	501	NDP	C7N-C3N	-2.00	1.44	1.48
2	ZM	501	NDP	O3D-C3D	-2.00	1.38	1.43
2	YX	501	NDP	O5D-C5D	-2.00	1.37	1.44

All (782) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZK	501	NDP	PN-O3-PA	-7.23	108.02	132.83
2	YM	501	NDP	PN-O3-PA	-7.22	108.05	132.83
2	YQ	501	NDP	PN-O3-PA	-7.20	108.13	132.83
2	YI	501	NDP	PN-O3-PA	-7.19	108.14	132.83
2	YR	501	NDP	PN-O3-PA	-7.17	108.22	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	YU	501	NDP	PN-O3-PA	-7.14	108.31	132.83
2	ZW	501	NDP	PN-O3-PA	-7.14	108.33	132.83
2	YY	501	NDP	PN-O3-PA	-7.12	108.38	132.83
2	ZO	501	NDP	PN-O3-PA	-7.11	108.44	132.83
2	YJ	501	NDP	PN-O3-PA	-7.10	108.46	132.83
2	ZX	501	NDP	PN-O3-PA	-7.09	108.48	132.83
2	ZS	501	NDP	PN-O3-PA	-7.09	108.49	132.83
2	ZB	501	NDP	PN-O3-PA	-7.09	108.51	132.83
2	ZP	501	NDP	PN-O3-PA	-7.09	108.51	132.83
2	YX	501	NDP	PN-O3-PA	-7.09	108.51	132.83
2	ZF	501	NDP	PN-O3-PA	-7.09	108.51	132.83
2	YL	501	NDP	PN-O3-PA	-7.08	108.52	132.83
2	ZC	501	NDP	PN-O3-PA	-7.08	108.54	132.83
2	ZR	501	NDP	PN-O3-PA	-7.07	108.55	132.83
2	ZT	501	NDP	PN-O3-PA	-7.07	108.56	132.83
2	ZN	501	NDP	PN-O3-PA	-7.07	108.56	132.83
2	ZG	501	NDP	PN-O3-PA	-7.07	108.57	132.83
2	YN	501	NDP	PN-O3-PA	-7.06	108.59	132.83
2	ZD	501	NDP	PN-O3-PA	-7.04	108.66	132.83
2	YV	501	NDP	PN-O3-PA	-7.04	108.67	132.83
2	ZL	501	NDP	PN-O3-PA	-7.03	108.69	132.83
2	YZ	501	NDP	PN-O3-PA	-7.03	108.70	132.83
2	ZZ	501	NDP	PN-O3-PA	-7.01	108.78	132.83
2	ZH	501	NDP	PN-O3-PA	-7.01	108.78	132.83
2	YT	501	NDP	PN-O3-PA	-6.98	108.86	132.83
4	ZT	503	A1JWG	C1A-NA-C4A	6.89	109.80	106.71
2	ZV	501	NDP	PN-O3-PA	-6.89	109.19	132.83
2	YW	501	NDP	PN-O3-PA	-6.84	109.36	132.83
2	ZJ	501	NDP	PN-O3-PA	-6.83	109.38	132.83
2	YS	501	NDP	PN-O3-PA	-6.82	109.42	132.83
2	ZY	501	NDP	PN-O3-PA	-6.82	109.42	132.83
4	YJ	503	A1JWG	C1A-NA-C4A	6.82	109.77	106.71
2	YP	501	NDP	PN-O3-PA	-6.79	109.53	132.83
4	YR	503	A1JWG	C1A-NA-C4A	6.74	109.74	106.71
2	ZQ	501	NDP	PN-O3-PA	-6.65	110.02	132.83
2	ZE	501	NDP	PN-O3-PA	-6.63	110.06	132.83
2	ZM	501	NDP	PN-O3-PA	-6.63	110.06	132.83
4	ZH	503	A1JWG	C1A-NA-C4A	6.63	109.69	106.71
4	ZD	503	A1JWG	C1A-NA-C4A	6.58	109.66	106.71
4	ZP	503	A1JWG	C1A-NA-C4A	6.55	109.65	106.71
2	YO	501	NDP	PN-O3-PA	-6.53	110.42	132.83
2	ZI	501	NDP	PN-O3-PA	-6.49	110.55	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZU	501	NDP	PN-O3-PA	-6.49	110.55	132.83
2	ZA	501	NDP	PN-O3-PA	-6.46	110.65	132.83
2	YK	501	NDP	PN-O3-PA	-6.42	110.80	132.83
4	YN	503	A1JWG	C1A-NA-C4A	6.42	109.59	106.71
4	ZO	503	A1JWG	C1A-NA-C4A	6.07	109.44	106.71
4	YY	503	A1JWG	C1A-NA-C4A	6.01	109.41	106.71
4	YK	503	A1JWG	C1A-NA-C4A	5.99	109.40	106.71
4	YQ	503	A1JWG	C1A-NA-C4A	5.99	109.40	106.71
4	ZE	503	A1JWG	C1A-NA-C4A	5.95	109.38	106.71
4	YV	503	A1JWG	C1A-NA-C4A	5.93	109.37	106.71
4	YO	503	A1JWG	C1A-NA-C4A	5.91	109.36	106.71
4	ZW	503	A1JWG	C1A-NA-C4A	5.88	109.35	106.71
4	ZG	503	A1JWG	C1A-NA-C4A	5.88	109.35	106.71
4	ZA	503	A1JWG	C1A-NA-C4A	5.83	109.33	106.71
4	ZS	503	A1JWG	C1A-NA-C4A	5.81	109.32	106.71
4	ZX	503	A1JWG	C1A-NA-C4A	5.79	109.31	106.71
4	YU	503	A1JWG	C1A-NA-C4A	5.77	109.30	106.71
4	YZ	503	A1JWG	C1A-NA-C4A	5.77	109.30	106.71
4	ZL	503	A1JWG	C1A-NA-C4A	5.74	109.29	106.71
4	YI	503	A1JWG	C1A-NA-C4A	5.73	109.28	106.71
4	ZK	503	A1JWG	C1A-NA-C4A	5.68	109.26	106.71
4	ZY	503	A1JWG	C1A-NA-C4A	5.66	109.25	106.71
4	ZC	503	A1JWG	C1A-NA-C4A	5.64	109.24	106.71
4	YM	503	A1JWG	C1A-NA-C4A	5.62	109.23	106.71
4	YT	503	A1JWG	C1A-NA-C4A	5.40	109.13	106.71
4	ZV	503	A1JWG	C1A-NA-C4A	5.32	109.10	106.71
4	YW	503	A1JWG	C1A-NA-C4A	5.20	109.05	106.71
4	ZQ	503	A1JWG	C1A-NA-C4A	5.12	109.01	106.71
4	YL	503	A1JWG	C1A-NA-C4A	5.08	108.99	106.71
4	ZR	503	A1JWG	C1A-NA-C4A	5.07	108.99	106.71
4	ZZ	503	A1JWG	C1A-NA-C4A	5.04	108.97	106.71
4	YS	503	A1JWG	C1A-NA-C4A	5.02	108.97	106.71
4	YX	503	A1JWG	C1A-NA-C4A	5.01	108.96	106.71
4	YP	503	A1JWG	C1A-NA-C4A	4.93	108.92	106.71
4	ZN	503	A1JWG	C1A-NA-C4A	4.93	108.92	106.71
4	ZU	503	A1JWG	C1A-NA-C4A	4.92	108.92	106.71
4	ZF	503	A1JWG	C1A-NA-C4A	4.92	108.92	106.71
4	ZM	503	A1JWG	C1A-NA-C4A	4.89	108.90	106.71
4	ZJ	503	A1JWG	C1A-NA-C4A	4.82	108.87	106.71
4	ZI	503	A1JWG	C1A-NA-C4A	4.78	108.85	106.71
4	ZB	503	A1JWG	C1A-NA-C4A	4.76	108.85	106.71
2	ZE	501	NDP	O2B-P2B-O1X	-3.71	95.06	109.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZA	501	NDP	O2B-P2B-O1X	-3.69	95.13	109.39
2	ZR	501	NDP	O2B-P2B-O1X	-3.56	95.66	109.39
2	YK	501	NDP	O2B-P2B-O1X	-3.51	95.83	109.39
2	YI	501	NDP	O2B-P2B-O1X	-3.51	95.83	109.39
2	YO	501	NDP	O2B-P2B-O1X	-3.51	95.86	109.39
2	ZQ	501	NDP	O2B-P2B-O1X	-3.50	95.89	109.39
2	ZM	501	NDP	C3B-C2B-C1B	-3.49	96.32	102.89
2	YY	501	NDP	O2B-P2B-O1X	-3.48	95.94	109.39
2	YQ	501	NDP	O2B-P2B-O1X	-3.47	95.98	109.39
2	ZK	501	NDP	O2B-P2B-O1X	-3.47	96.01	109.39
2	ZC	501	NDP	O2B-P2B-O1X	-3.47	96.02	109.39
2	ZG	501	NDP	O2B-P2B-O1X	-3.45	96.08	109.39
2	YZ	501	NDP	O2B-P2B-O1X	-3.44	96.10	109.39
2	ZX	501	NDP	O2B-P2B-O1X	-3.44	96.10	109.39
2	ZO	501	NDP	O2B-P2B-O1X	-3.44	96.12	109.39
2	ZA	501	NDP	C3B-C2B-C1B	-3.44	96.43	102.89
2	YO	501	NDP	C3B-C2B-C1B	-3.38	96.54	102.89
4	ZD	503	A1JWG	C3A-C4A-CHB	-3.36	118.54	122.46
2	ZU	501	NDP	C3B-C2B-C1B	-3.35	96.59	102.89
2	YT	501	NDP	O2B-P2B-O1X	-3.34	96.48	109.39
2	ZM	501	NDP	O2B-P2B-O1X	-3.33	96.52	109.39
2	ZJ	501	NDP	C3B-C2B-C1B	-3.33	96.63	102.89
2	ZW	501	NDP	O2B-P2B-O1X	-3.31	96.60	109.39
2	YP	501	NDP	C3B-C2B-C1B	-3.31	96.67	102.89
2	ZH	501	NDP	O2B-P2B-O1X	-3.31	96.62	109.39
2	YV	501	NDP	O2B-P2B-O1X	-3.30	96.65	109.39
4	ZT	503	A1JWG	C3A-C4A-CHB	-3.30	118.62	122.46
2	ZT	501	NDP	O2B-P2B-O1X	-3.30	96.67	109.39
2	YW	501	NDP	C3B-C2B-C1B	-3.28	96.72	102.89
2	ZD	501	NDP	O2B-P2B-O1X	-3.28	96.73	109.39
2	ZB	501	NDP	O2B-P2B-O1X	-3.27	96.75	109.39
2	ZN	501	NDP	C3B-C2B-C1B	-3.27	96.73	102.89
2	ZV	501	NDP	O2B-P2B-O1X	-3.27	96.77	109.39
2	YL	501	NDP	C3B-C2B-C1B	-3.27	96.74	102.89
2	YR	501	NDP	O2B-P2B-O1X	-3.26	96.80	109.39
2	YN	501	NDP	O2B-P2B-O1X	-3.26	96.80	109.39
2	ZL	501	NDP	O2B-P2B-O1X	-3.26	96.80	109.39
2	ZZ	501	NDP	O2B-P2B-O1X	-3.26	96.81	109.39
2	YX	501	NDP	O2B-P2B-O1X	-3.25	96.84	109.39
2	ZP	501	NDP	O2B-P2B-O1X	-3.25	96.84	109.39
2	YJ	501	NDP	O2B-P2B-O1X	-3.25	96.86	109.39
2	ZJ	501	NDP	O2B-P2B-O1X	-3.24	96.87	109.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZR	501	NDP	C3B-C2B-C1B	-3.24	96.79	102.89
2	ZF	501	NDP	O2B-P2B-O1X	-3.24	96.89	109.39
2	ZN	501	NDP	O2B-P2B-O1X	-3.24	96.89	109.39
2	YT	501	NDP	C3B-C2B-C1B	-3.23	96.81	102.89
2	ZS	501	NDP	O2B-P2B-O1X	-3.23	96.92	109.39
2	YU	501	NDP	O2B-P2B-O1X	-3.22	96.95	109.39
2	ZF	501	NDP	C3B-C2B-C1B	-3.22	96.84	102.89
2	YS	501	NDP	O2B-P2B-O1X	-3.22	96.97	109.39
4	ZP	503	A1JWG	C3A-C4A-CHB	-3.22	118.71	122.46
2	YS	501	NDP	C3B-C2B-C1B	-3.20	96.87	102.89
2	ZB	501	NDP	C3B-C2B-C1B	-3.20	96.87	102.89
2	ZI	501	NDP	O2B-P2B-O1X	-3.20	97.04	109.39
2	YX	501	NDP	C3B-C2B-C1B	-3.20	96.87	102.89
2	ZV	501	NDP	C3B-C2B-C1B	-3.20	96.88	102.89
2	YW	501	NDP	O2B-P2B-O1X	-3.19	97.07	109.39
2	YP	501	NDP	O2B-P2B-O1X	-3.19	97.07	109.39
4	YJ	503	A1JWG	C3A-C4A-CHB	-3.16	118.78	122.46
2	ZI	501	NDP	C3B-C2B-C1B	-3.15	96.96	102.89
2	YM	501	NDP	O2B-P2B-O1X	-3.14	97.25	109.39
4	YN	503	A1JWG	C3A-C4A-CHB	-3.14	118.80	122.46
2	ZD	501	NDP	PA-O5B-C5B	-3.14	103.26	121.68
2	YL	501	NDP	O2B-P2B-O1X	-3.13	97.31	109.39
2	ZP	501	NDP	PA-O5B-C5B	-3.12	103.36	121.68
2	ZT	501	NDP	PA-O5B-C5B	-3.12	103.37	121.68
2	ZY	501	NDP	O2B-P2B-O1X	-3.12	97.34	109.39
2	ZY	501	NDP	PA-O5B-C5B	-3.11	103.45	121.68
2	YS	501	NDP	PA-O5B-C5B	-3.11	103.47	121.68
2	YJ	501	NDP	PA-O5B-C5B	-3.10	103.50	121.68
2	ZI	501	NDP	PA-O5B-C5B	-3.10	103.50	121.68
2	ZU	501	NDP	PA-O5B-C5B	-3.10	103.51	121.68
2	YO	501	NDP	PA-O5B-C5B	-3.10	103.53	121.68
2	ZE	501	NDP	C3B-C2B-C1B	-3.09	97.07	102.89
2	ZZ	501	NDP	C3B-C2B-C1B	-3.09	97.07	102.89
2	YM	501	NDP	PA-O5B-C5B	-3.09	103.56	121.68
2	YW	501	NDP	PA-O5B-C5B	-3.08	103.61	121.68
2	ZU	501	NDP	O2B-P2B-O1X	-3.08	97.51	109.39
2	ZZ	501	NDP	PA-O5B-C5B	-3.08	103.64	121.68
2	YN	501	NDP	PA-O5B-C5B	-3.07	103.67	121.68
2	YT	501	NDP	PA-O5B-C5B	-3.07	103.67	121.68
2	YQ	501	NDP	PA-O5B-C5B	-3.07	103.68	121.68
2	YI	501	NDP	PA-O5B-C5B	-3.07	103.70	121.68
2	ZA	501	NDP	PA-O5B-C5B	-3.07	103.70	121.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	YY	501	NDP	PA-O5B-C5B	-3.06	103.71	121.68
4	ZH	503	A1JWG	C3A-C4A-CHB	-3.06	118.89	122.46
2	ZK	501	NDP	PA-O5B-C5B	-3.06	103.72	121.68
2	YV	501	NDP	PA-O5B-C5B	-3.06	103.72	121.68
2	YR	501	NDP	PA-O5B-C5B	-3.06	103.73	121.68
2	YK	501	NDP	PA-O5B-C5B	-3.06	103.74	121.68
2	ZV	501	NDP	PA-O5B-C5B	-3.06	103.75	121.68
2	ZW	501	NDP	PA-O5B-C5B	-3.05	103.78	121.68
2	ZY	501	NDP	C3B-C2B-C1B	-3.05	97.16	102.89
2	YP	501	NDP	PA-O5B-C5B	-3.04	103.83	121.68
2	ZJ	501	NDP	PA-O5B-C5B	-3.04	103.83	121.68
2	YK	501	NDP	C3B-C2B-C1B	-3.04	97.17	102.89
2	ZL	501	NDP	PA-O5B-C5B	-3.04	103.85	121.68
2	ZX	501	NDP	PA-O5B-C5B	-3.04	103.85	121.68
2	ZE	501	NDP	PA-O5B-C5B	-3.04	103.85	121.68
2	ZG	501	NDP	PA-O5B-C5B	-3.04	103.86	121.68
2	ZQ	501	NDP	PA-O5B-C5B	-3.03	103.90	121.68
2	ZS	501	NDP	PN-O5D-C5D	-3.03	103.92	121.68
2	ZM	501	NDP	PA-O5B-C5B	-3.03	103.92	121.68
2	ZO	501	NDP	PA-O5B-C5B	-3.03	103.92	121.68
2	ZC	501	NDP	PA-O5B-C5B	-3.03	103.94	121.68
2	ZK	501	NDP	PN-O5D-C5D	-3.02	103.96	121.68
2	YU	501	NDP	PA-O5B-C5B	-3.02	103.96	121.68
2	ZS	501	NDP	PA-O5B-C5B	-3.02	104.00	121.68
2	YZ	501	NDP	PA-O5B-C5B	-3.01	104.01	121.68
2	ZH	501	NDP	PA-O5B-C5B	-3.01	104.04	121.68
2	ZR	501	NDP	PA-O5B-C5B	-3.00	104.06	121.68
2	ZZ	501	NDP	PN-O5D-C5D	-3.00	104.07	121.68
2	YN	501	NDP	PN-O5D-C5D	-3.00	104.07	121.68
2	ZF	501	NDP	PA-O5B-C5B	-3.00	104.09	121.68
2	ZN	501	NDP	PA-O5B-C5B	-3.00	104.10	121.68
2	YQ	501	NDP	PN-O5D-C5D	-2.99	104.14	121.68
4	ZO	503	A1JWG	O2D-CGD-O1D	-2.99	117.99	123.84
2	YT	501	NDP	PN-O5D-C5D	-2.99	104.16	121.68
2	ZB	501	NDP	PA-O5B-C5B	-2.99	104.17	121.68
2	YL	501	NDP	PA-O5B-C5B	-2.98	104.18	121.68
2	YX	501	NDP	PA-O5B-C5B	-2.98	104.19	121.68
2	YU	501	NDP	PN-O5D-C5D	-2.98	104.20	121.68
4	ZK	503	A1JWG	O2D-CGD-O1D	-2.98	118.02	123.84
2	ZH	501	NDP	C3B-C2B-C1B	-2.97	97.30	102.89
2	ZO	501	NDP	C3B-C2B-C1B	-2.97	97.31	102.89
2	YW	501	NDP	PN-O5D-C5D	-2.95	104.37	121.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZO	501	NDP	PN-O5D-C5D	-2.95	104.40	121.68
2	ZW	501	NDP	PN-O5D-C5D	-2.94	104.46	121.68
4	YY	503	A1JWG	O2D-CGD-O1D	-2.93	118.10	123.84
2	ZG	501	NDP	PN-O5D-C5D	-2.92	104.56	121.68
2	ZC	501	NDP	PN-O5D-C5D	-2.92	104.57	121.68
2	YS	501	NDP	PN-O5D-C5D	-2.92	104.57	121.68
2	ZY	501	NDP	PN-O5D-C5D	-2.92	104.57	121.68
4	ZG	503	A1JWG	O2D-CGD-O1D	-2.92	118.13	123.84
2	YI	501	NDP	PN-O5D-C5D	-2.92	104.58	121.68
2	YY	501	NDP	PN-O5D-C5D	-2.91	104.59	121.68
2	ZH	501	NDP	PN-O5D-C5D	-2.91	104.62	121.68
2	YM	501	NDP	PN-O5D-C5D	-2.91	104.63	121.68
2	YR	501	NDP	C3B-C2B-C1B	-2.90	97.43	102.89
2	YZ	501	NDP	PN-O5D-C5D	-2.90	104.68	121.68
2	ZL	501	NDP	PN-O5D-C5D	-2.89	104.71	121.68
2	ZX	501	NDP	PN-O5D-C5D	-2.89	104.74	121.68
2	YV	501	NDP	C3B-C2B-C1B	-2.88	97.47	102.89
4	YI	503	A1JWG	O2D-CGD-O1D	-2.88	118.21	123.84
2	ZQ	501	NDP	C3B-C2B-C1B	-2.88	97.48	102.89
4	ZX	503	A1JWG	O2D-CGD-O1D	-2.85	118.26	123.84
2	YM	501	NDP	C3B-C2B-C1B	-2.85	97.53	102.89
2	YJ	501	NDP	C3B-C2B-C1B	-2.83	97.57	102.89
2	ZL	501	NDP	C3B-C2B-C1B	-2.82	97.59	102.89
2	ZV	501	NDP	PN-O5D-C5D	-2.80	105.28	121.68
4	ZS	503	A1JWG	O2D-CGD-O1D	-2.79	118.39	123.84
2	YY	501	NDP	C3B-C2B-C1B	-2.79	97.65	102.89
2	YZ	501	NDP	C3B-C2B-C1B	-2.78	97.66	102.89
4	YR	503	A1JWG	C3A-C4A-CHB	-2.78	119.22	122.46
2	ZP	501	NDP	PN-O5D-C5D	-2.78	105.40	121.68
2	ZN	501	NDP	PN-O5D-C5D	-2.77	105.44	121.68
4	ZW	503	A1JWG	O2D-CGD-O1D	-2.77	118.42	123.84
2	ZF	501	NDP	PN-O5D-C5D	-2.77	105.46	121.68
2	ZI	501	NDP	PN-O5D-C5D	-2.76	105.48	121.68
2	YJ	501	NDP	PN-O5D-C5D	-2.76	105.49	121.68
2	YO	501	NDP	PN-O5D-C5D	-2.76	105.49	121.68
2	YR	501	NDP	PN-O5D-C5D	-2.76	105.49	121.68
2	ZJ	501	NDP	PN-O5D-C5D	-2.76	105.51	121.68
2	ZB	501	NDP	PN-O5D-C5D	-2.76	105.52	121.68
2	ZT	501	NDP	PN-O5D-C5D	-2.75	105.54	121.68
2	YL	501	NDP	PN-O5D-C5D	-2.75	105.54	121.68
2	YP	501	NDP	PN-O5D-C5D	-2.74	105.59	121.68
2	YX	501	NDP	PN-O5D-C5D	-2.74	105.61	121.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	YU	501	NDP	C3B-C2B-C1B	-2.74	97.74	102.89
2	YV	501	NDP	PN-O5D-C5D	-2.74	105.61	121.68
2	YV	501	NDP	C4D-O4D-C1D	-2.74	103.43	109.47
2	ZD	501	NDP	PN-O5D-C5D	-2.73	105.68	121.68
2	ZG	501	NDP	C3B-C2B-C1B	-2.72	97.77	102.89
2	ZR	501	NDP	PN-O5D-C5D	-2.72	105.72	121.68
2	ZM	501	NDP	PN-O5D-C5D	-2.71	105.79	121.68
2	ZQ	501	NDP	PN-O5D-C5D	-2.71	105.80	121.68
2	ZT	501	NDP	C3B-C2B-C1B	-2.71	97.80	102.89
2	YK	501	NDP	PN-O5D-C5D	-2.71	105.81	121.68
2	ZU	501	NDP	PN-O5D-C5D	-2.70	105.84	121.68
2	ZA	501	NDP	PN-O5D-C5D	-2.69	105.89	121.68
2	ZD	501	NDP	C3B-C2B-C1B	-2.69	97.83	102.89
2	ZE	501	NDP	PN-O5D-C5D	-2.68	105.96	121.68
2	YI	501	NDP	C3B-C2B-C1B	-2.67	97.86	102.89
2	ZP	501	NDP	C3B-C2B-C1B	-2.66	97.89	102.89
4	ZC	503	A1JWG	O2D-CGD-O1D	-2.65	118.65	123.84
2	ZY	501	NDP	O3X-P2B-O2X	2.65	117.78	107.64
2	YN	501	NDP	C3B-C2B-C1B	-2.65	97.90	102.89
2	ZX	501	NDP	C3B-C2B-C1B	-2.65	97.90	102.89
2	ZU	501	NDP	O3X-P2B-O2X	2.65	117.77	107.64
2	ZE	501	NDP	O5D-PN-O1N	-2.64	98.74	109.07
2	ZQ	501	NDP	O5D-PN-O1N	-2.64	98.75	109.07
2	ZM	501	NDP	O5D-PN-O1N	-2.63	98.78	109.07
2	ZW	501	NDP	C3B-C2B-C1B	-2.63	97.94	102.89
2	ZC	501	NDP	C3B-C2B-C1B	-2.63	97.95	102.89
2	YN	501	NDP	C1D-N1N-C2N	-2.63	116.74	121.11
4	YR	503	A1JWG	O2D-CGD-O1D	-2.62	118.72	123.84
2	ZK	501	NDP	C3B-C2B-C1B	-2.62	97.97	102.89
2	YL	501	NDP	O3X-P2B-O2X	2.62	117.64	107.64
4	YP	503	A1JWG	O2D-CGD-O1D	-2.62	118.72	123.84
2	YP	501	NDP	O3X-P2B-O2X	2.61	117.60	107.64
4	YX	503	A1JWG	O2D-CGD-O1D	-2.60	118.75	123.84
2	YT	501	NDP	C5B-C4B-C3B	-2.60	105.45	115.18
4	ZJ	503	A1JWG	O2D-CGD-O1D	-2.60	118.76	123.84
2	ZS	501	NDP	C3B-C2B-C1B	-2.59	98.03	102.89
2	YQ	501	NDP	C3B-C2B-C1B	-2.58	98.03	102.89
2	YK	501	NDP	O3X-P2B-O2X	2.58	117.50	107.64
2	YR	501	NDP	C2D-C1D-N1N	-2.58	106.84	113.30
2	ZM	501	NDP	O3X-P2B-O2X	2.58	117.50	107.64
2	ZL	501	NDP	O3X-P2B-O2X	2.58	117.48	107.64
4	ZN	503	A1JWG	O2D-CGD-O1D	-2.57	118.81	123.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZP	501	NDP	O3X-P2B-O2X	2.57	117.45	107.64
4	YU	503	A1JWG	O2D-CGD-O1D	-2.57	118.82	123.84
2	YJ	501	NDP	O3X-P2B-O2X	2.57	117.44	107.64
2	YO	501	NDP	O3X-P2B-O2X	2.57	117.44	107.64
2	YN	501	NDP	O3X-P2B-O2X	2.56	117.44	107.64
2	ZD	501	NDP	O3X-P2B-O2X	2.56	117.44	107.64
2	ZW	501	NDP	O3X-P2B-O2X	2.56	117.43	107.64
2	YR	501	NDP	O3X-P2B-O2X	2.56	117.43	107.64
4	YZ	503	A1JWG	O2D-CGD-O1D	-2.56	118.83	123.84
2	YQ	501	NDP	O3X-P2B-O2X	2.56	117.42	107.64
2	ZZ	501	NDP	C5B-C4B-C3B	-2.56	105.60	115.18
2	YT	501	NDP	O3X-P2B-O2X	2.56	117.41	107.64
3	ZX	502	LMG	C7-O1-C1	2.56	118.73	113.74
2	ZG	501	NDP	O3X-P2B-O2X	2.56	117.41	107.64
2	ZQ	501	NDP	O3X-P2B-O2X	2.56	117.41	107.64
2	ZC	501	NDP	O3X-P2B-O2X	2.56	117.41	107.64
2	YZ	501	NDP	O3X-P2B-O2X	2.55	117.39	107.64
2	ZK	501	NDP	O3X-P2B-O2X	2.55	117.39	107.64
2	ZX	501	NDP	O3X-P2B-O2X	2.55	117.37	107.64
2	ZZ	501	NDP	O3X-P2B-O2X	2.54	117.36	107.64
2	YV	501	NDP	O3X-P2B-O2X	2.54	117.35	107.64
2	ZV	501	NDP	O3X-P2B-O2X	2.54	117.34	107.64
2	YM	501	NDP	O3X-P2B-O2X	2.54	117.33	107.64
2	ZN	501	NDP	O3X-P2B-O2X	2.54	117.33	107.64
2	ZJ	501	NDP	O3X-P2B-O2X	2.54	117.33	107.64
2	YS	501	NDP	O3X-P2B-O2X	2.54	117.33	107.64
2	YJ	501	NDP	C1D-N1N-C2N	-2.54	116.89	121.11
4	ZL	503	A1JWG	O2D-CGD-O1D	-2.54	118.88	123.84
2	ZH	501	NDP	O3X-P2B-O2X	2.54	117.33	107.64
2	ZO	501	NDP	O3X-P2B-O2X	2.53	117.33	107.64
2	ZB	501	NDP	O3X-P2B-O2X	2.53	117.32	107.64
2	ZT	501	NDP	O3X-P2B-O2X	2.53	117.32	107.64
2	ZP	501	NDP	C1D-N1N-C2N	-2.53	116.89	121.11
2	YX	501	NDP	O3X-P2B-O2X	2.53	117.31	107.64
2	ZP	501	NDP	C5B-C4B-C3B	-2.53	105.70	115.18
2	YJ	501	NDP	C5B-C4B-C3B	-2.53	105.71	115.18
2	ZF	501	NDP	O3X-P2B-O2X	2.52	117.28	107.64
2	YN	501	NDP	C5B-C4B-C3B	-2.52	105.72	115.18
2	YR	501	NDP	C5B-C4B-C3B	-2.52	105.73	115.18
2	YW	501	NDP	O3X-P2B-O2X	2.52	117.27	107.64
2	ZI	501	NDP	O5D-PN-O1N	-2.52	99.23	109.07
2	YV	501	NDP	C5B-C4B-C3B	-2.52	105.75	115.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	ZB	503	A1JWG	O2D-CGD-O1D	-2.52	118.92	123.84
2	ZV	501	NDP	C5B-C4B-C3B	-2.51	105.76	115.18
2	ZI	501	NDP	O3X-P2B-O2X	2.51	117.23	107.64
2	ZA	501	NDP	O5D-PN-O1N	-2.51	99.26	109.07
2	YZ	501	NDP	C3N-C2N-N1N	-2.51	119.51	123.10
2	YK	501	NDP	O5D-PN-O1N	-2.51	99.27	109.07
2	ZJ	501	NDP	C5B-C4B-C3B	-2.51	105.79	115.18
2	YO	501	NDP	O5D-PN-O1N	-2.50	99.29	109.07
2	ZT	501	NDP	C5B-C4B-C3B	-2.50	105.82	115.18
2	ZA	501	NDP	C5B-C4B-C3B	-2.50	105.82	115.18
2	ZR	501	NDP	O3X-P2B-O2X	2.50	117.18	107.64
2	ZS	501	NDP	O3X-P2B-O2X	2.50	117.18	107.64
2	YP	501	NDP	C5B-C4B-C3B	-2.49	105.83	115.18
2	YU	501	NDP	O3X-P2B-O2X	2.49	117.16	107.64
2	YI	501	NDP	O3X-P2B-O2X	2.49	117.15	107.64
2	ZN	501	NDP	C5B-C4B-C3B	-2.49	105.86	115.18
2	ZD	501	NDP	C5B-C4B-C3B	-2.49	105.86	115.18
2	YS	501	NDP	C5B-C4B-C3B	-2.49	105.87	115.18
2	YI	501	NDP	O2N-PN-O1N	2.48	124.53	112.24
2	ZR	501	NDP	O5D-PN-O1N	-2.48	99.36	109.07
2	ZU	501	NDP	O5D-PN-O1N	-2.48	99.36	109.07
2	YV	501	NDP	O5D-PN-O1N	-2.48	99.36	109.07
2	ZY	501	NDP	C5B-C4B-C3B	-2.48	105.87	115.18
2	YY	501	NDP	O3X-P2B-O2X	2.48	117.12	107.64
2	ZE	501	NDP	C5B-C4B-C3B	-2.48	105.88	115.18
4	ZF	503	A1JWG	O2D-CGD-O1D	-2.48	118.99	123.84
2	YW	501	NDP	C5B-C4B-C3B	-2.48	105.89	115.18
2	YI	501	NDP	C5B-C4B-C3B	-2.48	105.90	115.18
2	ZN	501	NDP	O5D-PN-O1N	-2.47	99.43	109.07
2	YY	501	NDP	C5B-C4B-C3B	-2.46	105.95	115.18
2	YV	501	NDP	O7N-C7N-C3N	2.46	125.54	120.90
4	YV	503	A1JWG	C3A-C4A-CHB	-2.46	119.59	122.46
2	YR	501	NDP	O2N-PN-O1N	2.46	124.42	112.24
2	YX	501	NDP	O5D-PN-O1N	-2.46	99.45	109.07
4	ZH	503	A1JWG	O2D-CGD-O1D	-2.46	119.03	123.84
2	ZB	501	NDP	O5D-PN-O1N	-2.45	99.48	109.07
2	ZH	501	NDP	C5B-C4B-C3B	-2.45	106.00	115.18
2	YM	501	NDP	O2N-PN-O1N	2.45	124.35	112.24
2	YM	501	NDP	C5B-C4B-C3B	-2.45	106.00	115.18
2	ZE	501	NDP	O3X-P2B-O2X	2.45	116.99	107.64
2	ZL	501	NDP	C5B-C4B-C3B	-2.45	106.01	115.18
2	ZF	501	NDP	O5D-PN-O1N	-2.45	99.51	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	YK	503	A1JWG	O2D-CGD-O1D	-2.44	119.06	123.84
2	ZA	501	NDP	O3X-P2B-O2X	2.44	116.98	107.64
4	ZA	503	A1JWG	O2D-CGD-O1D	-2.44	119.07	123.84
2	ZS	501	NDP	O5D-PN-O1N	-2.44	99.54	109.07
2	YY	501	NDP	O2N-PN-O1N	2.44	124.29	112.24
2	ZW	501	NDP	O5D-PN-O1N	-2.43	99.56	109.07
2	ZQ	501	NDP	C5B-C4B-C3B	-2.43	106.07	115.18
2	ZK	501	NDP	O2N-PN-O1N	2.43	124.26	112.24
2	YK	501	NDP	C2A-N1A-C6A	-2.43	114.59	118.77
2	ZX	501	NDP	O2N-PN-O1N	2.43	124.26	112.24
2	YL	501	NDP	C5B-C4B-C3B	-2.43	106.08	115.18
2	YL	501	NDP	O5D-PN-O1N	-2.43	99.58	109.07
2	ZF	501	NDP	C5B-C4B-C3B	-2.43	106.09	115.18
2	YQ	501	NDP	O2N-PN-O1N	2.42	124.22	112.24
2	ZE	501	NDP	O2N-PN-O1N	2.42	124.19	112.24
2	ZL	501	NDP	O5D-PN-O1N	-2.42	99.62	109.07
2	ZX	501	NDP	C3N-C2N-N1N	-2.42	119.65	123.10
4	YT	503	A1JWG	O2D-CGD-O1D	-2.42	119.11	123.84
2	ZO	501	NDP	O2N-PN-O1N	2.42	124.18	112.24
2	ZQ	501	NDP	O2N-PN-O1N	2.41	124.18	112.24
2	ZN	501	NDP	O2N-PN-O1N	2.41	124.17	112.24
2	ZI	501	NDP	C2A-N1A-C6A	-2.41	114.63	118.77
2	ZP	501	NDP	C2D-C1D-N1N	-2.41	107.26	113.30
2	ZX	501	NDP	O5D-PN-O1N	-2.41	99.65	109.07
4	YO	503	A1JWG	O2D-CGD-O1D	-2.41	119.13	123.84
2	YP	501	NDP	C2A-N1A-C6A	-2.41	114.63	118.77
2	ZG	501	NDP	C2A-N1A-C6A	-2.41	114.63	118.77
2	ZM	501	NDP	C5B-C4B-C3B	-2.41	106.15	115.18
2	YU	501	NDP	C2A-N1A-C6A	-2.41	114.64	118.77
2	ZZ	501	NDP	C2A-N1A-C6A	-2.41	114.64	118.77
2	ZI	501	NDP	C5B-C4B-C3B	-2.41	106.16	115.18
2	ZF	501	NDP	O2N-PN-O1N	2.41	124.14	112.24
2	ZV	501	NDP	C2A-N1A-C6A	-2.41	114.64	118.77
2	ZG	501	NDP	O2N-PN-O1N	2.41	124.14	112.24
2	YI	501	NDP	O5D-PN-O1N	-2.41	99.67	109.07
2	YK	501	NDP	C5B-C4B-C3B	-2.41	106.17	115.18
2	ZW	501	NDP	O2N-PN-O1N	2.41	124.13	112.24
2	ZK	501	NDP	C2A-N1A-C6A	-2.40	114.64	118.77
2	ZC	501	NDP	O5D-PN-O1N	-2.40	99.68	109.07
2	ZR	501	NDP	O2N-PN-O1N	2.40	124.12	112.24
2	ZC	501	NDP	C2A-N1A-C6A	-2.40	114.64	118.77
2	ZO	501	NDP	C2A-N1A-C6A	-2.40	114.64	118.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZS	501	NDP	C2A-N1A-C6A	-2.40	114.64	118.77
2	ZH	501	NDP	O5D-PN-O1N	-2.40	99.69	109.07
2	YT	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	ZR	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	ZJ	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	YX	501	NDP	C5B-C4B-C3B	-2.40	106.19	115.18
2	ZW	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	YO	501	NDP	C5B-C4B-C3B	-2.40	106.20	115.18
2	YI	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	YO	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	ZE	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	ZM	501	NDP	O2N-PN-O1N	2.40	124.09	112.24
2	ZN	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	ZG	501	NDP	O5D-PN-O1N	-2.40	99.70	109.07
2	YX	501	NDP	C2A-N1A-C6A	-2.40	114.65	118.77
2	ZC	501	NDP	O2N-PN-O1N	2.40	124.09	112.24
2	YQ	501	NDP	O5D-PN-O1N	-2.40	99.70	109.07
2	YL	501	NDP	O2N-PN-O1N	2.40	124.08	112.24
2	YL	501	NDP	C2A-N1A-C6A	-2.39	114.66	118.77
2	YQ	501	NDP	C2A-N1A-C6A	-2.39	114.66	118.77
2	YY	501	NDP	C2A-N1A-C6A	-2.39	114.66	118.77
2	ZK	501	NDP	O5D-PN-O1N	-2.39	99.72	109.07
4	YQ	503	A1JWG	O2D-CGD-O1D	-2.39	119.16	123.84
2	ZB	501	NDP	C5B-C4B-C3B	-2.39	106.22	115.18
2	ZY	501	NDP	C2A-N1A-C6A	-2.39	114.66	118.77
2	YY	501	NDP	N3A-C4A-N9A	2.39	131.02	127.08
2	YX	501	NDP	O2N-PN-O1N	2.39	124.06	112.24
2	ZR	501	NDP	C5B-C4B-C3B	-2.39	106.23	115.18
2	ZA	501	NDP	C2A-N1A-C6A	-2.39	114.67	118.77
2	YZ	501	NDP	C2A-N1A-C6A	-2.39	114.67	118.77
2	ZB	501	NDP	O2N-PN-O1N	2.39	124.04	112.24
2	ZF	501	NDP	C2A-N1A-C6A	-2.39	114.67	118.77
2	ZL	501	NDP	C2A-N1A-C6A	-2.39	114.67	118.77
2	ZM	501	NDP	C2A-N1A-C6A	-2.39	114.67	118.77
2	YZ	501	NDP	N3A-C4A-N9A	2.39	131.01	127.08
2	ZB	501	NDP	C2A-N1A-C6A	-2.39	114.67	118.77
2	YR	501	NDP	O5D-PN-O1N	-2.38	99.75	109.07
2	ZC	501	NDP	N3A-C4A-N9A	2.38	131.01	127.08
2	ZS	501	NDP	C5B-C4B-C3B	-2.38	106.25	115.18
2	YZ	501	NDP	O5D-PN-O1N	-2.38	99.76	109.07
2	ZU	501	NDP	C2A-N1A-C6A	-2.38	114.68	118.77
2	ZO	501	NDP	O5D-PN-O1N	-2.38	99.76	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZV	501	NDP	O5D-PN-O1N	-2.38	99.77	109.07
2	ZG	501	NDP	N3A-C4A-N9A	2.38	131.00	127.08
2	ZQ	501	NDP	C2A-N1A-C6A	-2.38	114.68	118.77
2	YM	501	NDP	C2A-N1A-C6A	-2.38	114.68	118.77
2	ZX	501	NDP	C2A-N1A-C6A	-2.38	114.69	118.77
2	YZ	501	NDP	O2N-PN-O1N	2.38	123.99	112.24
2	YS	501	NDP	C2A-N1A-C6A	-2.38	114.69	118.77
2	ZX	501	NDP	N3A-C4A-N9A	2.38	131.00	127.08
2	ZZ	501	NDP	C3N-C2N-N1N	-2.37	119.71	123.10
2	ZU	501	NDP	C5B-C4B-C3B	-2.37	106.30	115.18
2	ZW	501	NDP	C5B-C4B-C3B	-2.37	106.30	115.18
4	YV	503	A1JWG	O2D-CGD-O1D	-2.37	119.20	123.84
2	YY	501	NDP	O5D-PN-O1N	-2.37	99.81	109.07
2	ZD	501	NDP	C2A-N1A-C6A	-2.37	114.70	118.77
2	YU	501	NDP	O2N-PN-O1N	2.37	123.94	112.24
2	YW	501	NDP	C2A-N1A-C6A	-2.36	114.71	118.77
2	YJ	501	NDP	C2D-C1D-N1N	-2.36	107.39	113.30
4	ZE	503	A1JWG	O2D-CGD-O1D	-2.36	119.23	123.84
2	YU	501	NDP	C5B-C4B-C3B	-2.35	106.36	115.18
2	YU	501	NDP	O5D-PN-O1N	-2.35	99.87	109.07
2	YN	501	NDP	C2A-N1A-C6A	-2.35	114.73	118.77
2	YP	501	NDP	O5D-PN-O1N	-2.35	99.89	109.07
2	YQ	501	NDP	C5B-C4B-C3B	-2.35	106.38	115.18
2	ZJ	501	NDP	O5D-PN-O1N	-2.35	99.89	109.07
2	YJ	501	NDP	C2A-N1A-C6A	-2.35	114.74	118.77
2	ZL	501	NDP	N3A-C4A-N9A	2.35	130.95	127.08
2	ZP	501	NDP	C2A-N1A-C6A	-2.35	114.74	118.77
2	ZS	501	NDP	O2N-PN-O1N	2.35	123.84	112.24
2	ZK	501	NDP	N3A-C4A-N9A	2.34	130.94	127.08
2	ZO	501	NDP	N3A-C4A-N9A	2.34	130.94	127.08
2	YI	501	NDP	N3A-C4A-N9A	2.34	130.94	127.08
2	ZT	501	NDP	C2A-N1A-C6A	-2.34	114.75	118.77
2	ZH	501	NDP	C2A-N1A-C6A	-2.34	114.75	118.77
2	YR	501	NDP	C2A-N1A-C6A	-2.34	114.76	118.77
2	YJ	501	NDP	O2N-PN-O1N	2.34	123.79	112.24
2	ZL	501	NDP	O2N-PN-O1N	2.34	123.78	112.24
2	YM	501	NDP	O5D-PN-O1N	-2.33	99.95	109.07
2	ZS	501	NDP	N3A-C4A-N9A	2.33	130.93	127.08
2	ZK	501	NDP	C5B-C4B-C3B	-2.33	106.44	115.18
2	ZP	501	NDP	O2N-PN-O1N	2.33	123.77	112.24
2	YV	501	NDP	C2A-N1A-C6A	-2.33	114.77	118.77
2	ZD	501	NDP	C1D-N1N-C2N	-2.33	117.24	121.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	YN	501	NDP	O2N-PN-O1N	2.33	123.74	112.24
3	ZB	502	LMG	O1-C7-C8	-2.32	105.29	110.90
4	ZI	503	A1JWG	C3A-C4A-CHB	-2.32	119.76	122.46
2	YZ	501	NDP	C5B-C4B-C3B	-2.32	106.50	115.18
2	ZD	501	NDP	O2N-PN-O1N	2.31	123.67	112.24
2	ZT	501	NDP	O2N-PN-O1N	2.31	123.65	112.24
2	ZG	501	NDP	C5B-C4B-C3B	-2.30	106.54	115.18
4	ZM	503	A1JWG	C3A-C4A-CHB	-2.30	119.78	122.46
2	YK	501	NDP	N3A-C4A-N9A	2.30	130.88	127.08
2	YU	501	NDP	N3A-C4A-N9A	2.30	130.88	127.08
2	ZZ	501	NDP	N3A-C4A-N9A	2.30	130.88	127.08
2	YY	501	NDP	C5A-C4A-N3A	-2.30	123.75	126.75
2	YM	501	NDP	N3A-C4A-N9A	2.30	130.87	127.08
2	ZH	501	NDP	O2N-PN-O1N	2.30	123.61	112.24
2	ZO	501	NDP	C5B-C4B-C3B	-2.30	106.57	115.18
2	YN	501	NDP	O5D-PN-O1N	-2.30	100.09	109.07
2	ZI	501	NDP	N3A-C4A-N9A	2.29	130.86	127.08
2	ZC	501	NDP	C5B-C4B-C3B	-2.29	106.58	115.18
2	ZT	501	NDP	C1D-N1N-C2N	-2.29	117.30	121.11
2	ZV	501	NDP	N3A-C4A-N9A	2.29	130.86	127.08
2	ZJ	501	NDP	N3A-C4A-N9A	2.29	130.85	127.08
2	ZN	501	NDP	N3A-C4A-N9A	2.29	130.85	127.08
2	YT	501	NDP	N3A-C4A-N9A	2.28	130.85	127.08
2	ZE	501	NDP	N3A-C4A-N9A	2.28	130.85	127.08
2	YP	501	NDP	N3A-C4A-N9A	2.28	130.84	127.08
2	ZW	501	NDP	N3A-C4A-N9A	2.28	130.84	127.08
2	YL	501	NDP	N3A-C4A-N9A	2.28	130.84	127.08
2	ZQ	501	NDP	N3A-C4A-N9A	2.28	130.84	127.08
2	ZX	501	NDP	C5A-C4A-N3A	-2.27	123.79	126.75
4	ZZ	503	A1JWG	O2D-CGD-O1D	-2.27	119.40	123.84
2	YQ	501	NDP	N3A-C4A-N9A	2.27	130.82	127.08
2	ZU	501	NDP	O2N-PN-O1N	2.27	123.46	112.24
2	ZM	501	NDP	N3A-C4A-N9A	2.27	130.82	127.08
2	ZA	501	NDP	O2N-PN-O1N	2.27	123.45	112.24
2	YN	501	NDP	N3A-C4A-N9A	2.27	130.82	127.08
2	ZD	501	NDP	N3A-C4A-N9A	2.27	130.82	127.08
2	ZX	501	NDP	C5B-C4B-C3B	-2.27	106.69	115.18
2	ZV	501	NDP	O2N-PN-O1N	2.27	123.44	112.24
2	ZC	501	NDP	C5A-C4A-N3A	-2.27	123.80	126.75
2	YX	501	NDP	N3A-C4A-N9A	2.26	130.81	127.08
2	ZR	501	NDP	N3A-C4A-N9A	2.26	130.81	127.08
2	ZG	501	NDP	C5A-C4A-N3A	-2.26	123.80	126.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZF	501	NDP	N3A-C4A-N9A	2.26	130.81	127.08
2	YK	501	NDP	C5A-C4A-N3A	-2.26	123.80	126.75
4	ZH	503	A1JWG	CHC-C4B-C3B	2.26	129.12	125.26
2	YO	501	NDP	O2N-PN-O1N	2.26	123.41	112.24
2	YI	501	NDP	C5A-C4A-N3A	-2.25	123.81	126.75
2	YO	501	NDP	N3A-C4A-N9A	2.25	130.80	127.08
2	ZS	501	NDP	C5A-C4A-N3A	-2.25	123.81	126.75
2	ZY	501	NDP	N3A-C4A-N9A	2.25	130.80	127.08
4	ZV	503	A1JWG	O2D-CGD-O1D	-2.25	119.43	123.84
2	ZL	501	NDP	C3N-C2N-N1N	-2.25	119.88	123.10
2	ZT	501	NDP	N3A-C4A-N9A	2.25	130.79	127.08
4	ZY	503	A1JWG	O2D-CGD-O1D	-2.24	119.45	123.84
2	ZJ	501	NDP	O2N-PN-O1N	2.24	123.33	112.24
2	ZI	501	NDP	C5A-C4A-N3A	-2.24	123.82	126.75
2	ZP	501	NDP	N3A-C4A-N9A	2.24	130.78	127.08
2	ZI	501	NDP	O2N-PN-O1N	2.24	123.33	112.24
2	YM	501	NDP	C5A-C4A-N3A	-2.24	123.83	126.75
4	ZV	503	A1JWG	CHD-C1D-ND	-2.24	121.89	124.26
2	ZU	501	NDP	N3A-C4A-N9A	2.24	130.77	127.08
2	ZL	501	NDP	C5A-C4A-N3A	-2.23	123.83	126.75
2	ZB	501	NDP	N3A-C4A-N9A	2.23	130.76	127.08
4	YR	503	A1JWG	CHC-C4B-C3B	2.23	129.08	125.26
2	YP	501	NDP	C5A-C4A-N3A	-2.23	123.84	126.75
2	ZZ	501	NDP	C5A-C4A-N3A	-2.23	123.84	126.75
2	ZH	501	NDP	N3A-C4A-N9A	2.23	130.76	127.08
2	YZ	501	NDP	C5A-C4A-N3A	-2.23	123.84	126.75
2	ZE	501	NDP	C5A-C4A-N3A	-2.23	123.85	126.75
2	YW	501	NDP	O2N-PN-O1N	2.22	123.24	112.24
2	ZA	501	NDP	N3A-C4A-N9A	2.22	130.75	127.08
2	YS	501	NDP	N3A-C4A-N9A	2.22	130.75	127.08
2	ZY	501	NDP	O2N-PN-O1N	2.22	123.23	112.24
2	ZQ	501	NDP	C5A-C4A-N3A	-2.22	123.85	126.75
2	YK	501	NDP	O2N-PN-O1N	2.22	123.23	112.24
2	YP	501	NDP	O2N-PN-O1N	2.22	123.22	112.24
4	ZR	503	A1JWG	CHD-C1D-ND	-2.22	121.91	124.26
2	YV	501	NDP	O2N-PN-O1N	2.22	123.21	112.24
4	ZD	503	A1JWG	O2D-CGD-O1D	-2.22	119.50	123.84
2	ZJ	501	NDP	C5A-C4A-N3A	-2.22	123.86	126.75
4	ZT	503	A1JWG	CHC-C4B-C3B	2.22	129.05	125.26
2	ZZ	501	NDP	O2N-PN-O1N	2.22	123.19	112.24
2	ZR	501	NDP	C3N-C2N-N1N	-2.22	119.94	123.10
2	ZV	501	NDP	C5A-C4A-N3A	-2.21	123.86	126.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	YS	501	NDP	C3N-C2N-N1N	-2.21	119.94	123.10
2	YO	501	NDP	C5A-C4A-N3A	-2.21	123.86	126.75
2	YJ	501	NDP	N3A-C4A-N9A	2.21	130.72	127.08
2	ZO	501	NDP	C5A-C4A-N3A	-2.21	123.87	126.75
2	YU	501	NDP	C5A-C4A-N3A	-2.21	123.87	126.75
2	YT	501	NDP	O2N-PN-O1N	2.21	123.16	112.24
2	YL	501	NDP	C5A-C4A-N3A	-2.21	123.87	126.75
2	YS	501	NDP	O2N-PN-O1N	2.21	123.15	112.24
2	ZM	501	NDP	C5A-C4A-N3A	-2.21	123.87	126.75
2	ZU	501	NDP	C5A-C4A-N3A	-2.21	123.87	126.75
2	YV	501	NDP	N3A-C4A-N9A	2.20	130.71	127.08
4	YL	503	A1JWG	O2D-CGD-O1D	-2.20	119.53	123.84
2	ZA	501	NDP	C5A-C4A-N3A	-2.20	123.88	126.75
4	ZY	503	A1JWG	C3A-C4A-CHB	-2.20	119.90	122.46
4	YM	503	A1JWG	O2D-CGD-O1D	-2.20	119.54	123.84
2	YW	501	NDP	N3A-C4A-N9A	2.20	130.70	127.08
2	ZQ	501	NDP	C3N-C2N-N1N	-2.20	119.96	123.10
2	YT	501	NDP	C5A-C4A-N3A	-2.20	123.89	126.75
2	ZK	501	NDP	C5A-C4A-N3A	-2.20	123.89	126.75
4	ZZ	503	A1JWG	CHD-C1D-ND	-2.20	121.93	124.26
2	ZY	501	NDP	C5A-C4A-N3A	-2.19	123.89	126.75
2	ZN	501	NDP	C5A-C4A-N3A	-2.19	123.89	126.75
4	ZU	503	A1JWG	CHD-C1D-ND	-2.19	121.94	124.26
4	YL	503	A1JWG	CHD-C1D-ND	-2.18	121.94	124.26
2	ZW	501	NDP	C5A-C4A-N3A	-2.18	123.90	126.75
4	YT	503	A1JWG	CHD-C1D-ND	-2.18	121.95	124.26
2	ZT	501	NDP	C2D-C1D-N1N	-2.18	107.84	113.30
2	YX	501	NDP	C5A-C4A-N3A	-2.18	123.91	126.75
2	YR	501	NDP	N3A-C4A-N9A	2.18	130.67	127.08
4	YX	503	A1JWG	CHD-C1D-ND	-2.18	121.95	124.26
2	YT	501	NDP	C3N-C2N-N1N	-2.18	119.99	123.10
2	ZC	501	NDP	C3N-C2N-N1N	-2.18	119.99	123.10
2	YQ	501	NDP	C5A-C4A-N3A	-2.17	123.92	126.75
2	ZR	501	NDP	C5A-C4A-N3A	-2.17	123.92	126.75
2	YR	501	NDP	C1D-N1N-C2N	-2.17	117.50	121.11
2	ZF	501	NDP	C5A-C4A-N3A	-2.17	123.92	126.75
2	YU	501	NDP	C3N-C2N-N1N	-2.17	120.00	123.10
4	YN	503	A1JWG	CHC-C4B-C3B	2.16	128.96	125.26
2	ZM	501	NDP	O4B-C1B-C2B	-2.16	102.78	106.57
2	YS	501	NDP	C5A-C4A-N3A	-2.16	123.93	126.75
4	YV	503	A1JWG	CHC-C4B-C3B	2.15	128.94	125.26
2	ZT	501	NDP	C5A-C4A-N3A	-2.15	123.94	126.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	YN	501	NDP	C5A-C4A-N3A	-2.15	123.94	126.75
4	ZP	503	A1JWG	CHC-C4B-C3B	2.15	128.94	125.26
4	YJ	503	A1JWG	CHC-C4B-C3B	2.15	128.94	125.26
2	ZB	501	NDP	C5A-C4A-N3A	-2.15	123.95	126.75
4	ZN	503	A1JWG	CHD-C1D-ND	-2.15	121.98	124.26
4	ZJ	503	A1JWG	CHD-C1D-ND	-2.15	121.98	124.26
4	ZB	503	A1JWG	CHD-C1D-ND	-2.15	121.98	124.26
2	ZS	501	NDP	O7N-C7N-C3N	2.14	124.93	120.90
2	ZG	501	NDP	C3N-C2N-N1N	-2.14	120.04	123.10
2	ZD	501	NDP	C2D-C1D-N1N	-2.14	107.94	113.30
2	ZD	501	NDP	C5D-C4D-C3D	-2.14	107.16	115.18
2	ZT	501	NDP	C5D-C4D-C3D	-2.14	107.16	115.18
2	ZK	501	NDP	C3N-C2N-N1N	-2.14	120.05	123.10
2	ZD	501	NDP	C5A-C4A-N3A	-2.14	123.96	126.75
4	ZI	503	A1JWG	CHD-C1D-ND	-2.13	122.00	124.26
2	YW	501	NDP	C5A-C4A-N3A	-2.13	123.97	126.75
4	ZR	503	A1JWG	O2D-CGD-O1D	-2.13	119.67	123.84
4	YP	503	A1JWG	CHD-C1D-ND	-2.13	122.00	124.26
2	YQ	501	NDP	O7N-C7N-C3N	2.13	124.90	120.90
2	YX	501	NDP	C3N-C2N-N1N	-2.12	120.07	123.10
2	ZS	501	NDP	C3N-C2N-N1N	-2.12	120.07	123.10
2	ZG	501	NDP	O7N-C7N-C3N	2.12	124.89	120.90
3	ZD	502	LMG	O1-C7-C8	-2.12	105.79	110.90
2	ZP	501	NDP	C5A-C4A-N3A	-2.12	123.99	126.75
4	YS	503	A1JWG	CHD-C1D-ND	-2.11	122.02	124.26
2	ZH	501	NDP	C2D-C1D-N1N	-2.11	108.01	113.30
2	YU	501	NDP	O7N-C7N-C3N	2.11	124.88	120.90
2	YL	501	NDP	C3N-C2N-N1N	-2.11	120.08	123.10
2	ZE	501	NDP	O4B-C1B-C2B	-2.11	102.86	106.57
2	ZA	501	NDP	O4B-C1B-C2B	-2.11	102.87	106.57
4	ZD	503	A1JWG	CHC-C4B-C3B	2.11	128.86	125.26
2	YJ	501	NDP	O7N-C7N-C3N	2.11	124.86	120.90
2	ZY	501	NDP	O3X-P2B-O2B	-2.11	96.56	105.99
4	YW	503	A1JWG	CHD-C1D-ND	-2.11	122.03	124.26
2	ZU	501	NDP	O3X-P2B-O2B	-2.10	96.56	105.99
2	YJ	501	NDP	C5A-C4A-N3A	-2.10	124.01	126.75
2	YM	501	NDP	O7N-C7N-C3N	2.10	124.86	120.90
2	ZF	501	NDP	C3N-C2N-N1N	-2.10	120.10	123.10
4	ZF	503	A1JWG	CHD-C1D-ND	-2.10	122.03	124.26
2	ZT	501	NDP	O5D-PN-O1N	-2.10	100.87	109.07
4	ZQ	503	A1JWG	CHD-C1D-ND	-2.10	122.04	124.26
2	YJ	501	NDP	O5D-PN-O1N	-2.10	100.87	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	ZM	503	A1JWG	CHD-C1D-ND	-2.09	122.04	124.26
4	YN	503	A1JWG	O2D-CGD-O1D	-2.09	119.75	123.84
2	YJ	501	NDP	C5D-C4D-C3D	-2.09	107.35	115.18
2	ZP	501	NDP	O7N-C7N-C3N	2.09	124.83	120.90
3	ZX	502	LMG	C9-C8-C7	2.09	116.72	111.79
2	YI	501	NDP	O7N-C7N-C3N	2.09	124.83	120.90
2	ZH	501	NDP	O7N-C7N-C3N	2.09	124.83	120.90
2	ZL	501	NDP	O7N-C7N-C3N	2.09	124.83	120.90
2	YY	501	NDP	C5A-C6A-N1A	2.08	121.86	117.39
2	ZD	501	NDP	O7N-C7N-C3N	2.08	124.82	120.90
4	ZH	503	A1JWG	CHD-C1D-ND	-2.08	122.05	124.26
2	YQ	501	NDP	C3N-C2N-N1N	-2.08	120.13	123.10
2	YT	501	NDP	O5D-PN-O1N	-2.08	100.94	109.07
2	ZB	501	NDP	C3N-C2N-N1N	-2.08	120.13	123.10
2	ZH	501	NDP	C5A-C4A-N3A	-2.08	124.04	126.75
4	YZ	503	A1JWG	CHC-C4B-C3B	2.08	128.81	125.26
2	YZ	501	NDP	C5A-C6A-N1A	2.08	121.85	117.39
2	ZZ	501	NDP	O5D-PN-O1N	-2.08	100.95	109.07
2	YK	501	NDP	C5A-C6A-N1A	2.08	121.85	117.39
2	ZO	501	NDP	O7N-C7N-C3N	2.07	124.80	120.90
4	YV	503	A1JWG	CHD-C1D-ND	-2.07	122.06	124.26
2	YZ	501	NDP	O7N-C7N-C3N	2.07	124.80	120.90
2	ZG	501	NDP	C5A-C6A-N1A	2.07	121.84	117.39
4	ZP	503	A1JWG	O2D-CGD-O1D	-2.07	119.79	123.84
2	ZI	501	NDP	C5A-C6A-N1A	2.07	121.83	117.39
2	ZP	501	NDP	C5D-C4D-C3D	-2.07	107.43	115.18
2	ZC	501	NDP	C5A-C6A-N1A	2.07	121.83	117.39
4	YK	503	A1JWG	C3A-C4A-CHB	-2.07	120.05	122.46
2	YZ	501	NDP	C2D-C1D-N1N	-2.07	108.13	113.30
4	YW	503	A1JWG	O2D-CGD-O1D	-2.06	119.80	123.84
2	ZL	501	NDP	C5A-C6A-N1A	2.06	121.82	117.39
2	YO	501	NDP	C5A-C6A-N1A	2.06	121.82	117.39
4	YR	503	A1JWG	CHD-C1D-ND	-2.06	122.07	124.26
2	ZX	501	NDP	O7N-C7N-C3N	2.06	124.78	120.90
4	ZX	503	A1JWG	CHC-C4B-C3B	2.06	128.79	125.26
2	ZX	501	NDP	C5A-C6A-N1A	2.06	121.81	117.39
2	ZP	501	NDP	O5D-PN-O1N	-2.06	101.02	109.07
4	ZL	503	A1JWG	CHC-C4B-C3B	2.06	128.78	125.26
2	ZA	501	NDP	C5A-C6A-N1A	2.06	121.81	117.39
4	YK	503	A1JWG	CHC-C4B-C3B	2.06	128.78	125.26
2	YM	501	NDP	C5A-C6A-N1A	2.05	121.80	117.39
2	ZD	501	NDP	C5A-C6A-N1A	2.05	121.80	117.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZY	501	NDP	C5A-C6A-N1A	2.05	121.80	117.39
2	ZD	501	NDP	O5D-PN-O1N	-2.05	101.05	109.07
2	ZH	501	NDP	C5A-C6A-N1A	2.05	121.79	117.39
2	ZU	501	NDP	C5A-C6A-N1A	2.05	121.79	117.39
2	YR	501	NDP	C5A-C4A-N3A	-2.05	124.08	126.75
2	YT	501	NDP	C5A-C6A-N1A	2.05	121.79	117.39
2	YR	501	NDP	O7N-C7N-C3N	2.05	124.75	120.90
2	ZY	501	NDP	O5D-PN-O1N	-2.05	101.07	109.07
2	YV	501	NDP	C5A-C4A-N3A	-2.05	124.08	126.75
2	ZK	501	NDP	C5A-C6A-N1A	2.05	121.78	117.39
2	YL	501	NDP	O3X-P2B-O2B	-2.05	96.82	105.99
2	ZM	501	NDP	C5A-C6A-N1A	2.04	121.78	117.39
2	YW	501	NDP	C3N-C2N-N1N	-2.04	120.18	123.10
2	YI	501	NDP	C5A-C6A-N1A	2.04	121.78	117.39
2	ZE	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
4	ZW	503	A1JWG	CHC-C4B-C3B	2.04	128.75	125.26
2	ZQ	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
2	YP	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
2	ZS	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
2	ZW	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
2	ZZ	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
2	YP	501	NDP	O3X-P2B-O2B	-2.04	96.86	105.99
2	ZJ	501	NDP	C5A-C6A-N1A	2.04	121.77	117.39
2	ZM	501	NDP	C4B-O4B-C1B	-2.04	104.98	109.47
2	YR	501	NDP	C5D-C4D-C3D	-2.04	107.55	115.18
2	ZT	501	NDP	C5A-C6A-N1A	2.04	121.76	117.39
2	ZO	501	NDP	C5A-C6A-N1A	2.03	121.76	117.39
2	YU	501	NDP	C5A-C6A-N1A	2.03	121.76	117.39
2	ZN	501	NDP	C5A-C6A-N1A	2.03	121.76	117.39
2	ZV	501	NDP	C5A-C6A-N1A	2.03	121.76	117.39
2	ZR	501	NDP	C5A-C6A-N1A	2.03	121.75	117.39
2	ZT	501	NDP	O7N-C7N-C3N	2.03	124.72	120.90
2	YS	501	NDP	C5A-C6A-N1A	2.03	121.75	117.39
4	ZM	503	A1JWG	O2D-CGD-O1D	-2.03	119.87	123.84
2	YL	501	NDP	C5A-C6A-N1A	2.03	121.75	117.39
2	YW	501	NDP	C5A-C6A-N1A	2.03	121.75	117.39
2	YJ	501	NDP	C5A-C6A-N1A	2.03	121.75	117.39
2	YN	501	NDP	C5A-C6A-N1A	2.03	121.75	117.39
2	YI	501	NDP	C3N-C2N-N1N	-2.03	120.20	123.10
2	YW	501	NDP	O5D-PN-O1N	-2.03	101.14	109.07
2	ZK	501	NDP	O7N-C7N-C3N	2.03	124.72	120.90
2	ZV	501	NDP	C3N-C2N-N1N	-2.03	120.20	123.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	ZF	501	NDP	C5A-C6A-N1A	2.03	121.74	117.39
2	YX	501	NDP	C5A-C6A-N1A	2.03	121.74	117.39
2	ZU	501	NDP	O4B-C1B-C2B	-2.03	103.02	106.57
2	YQ	501	NDP	C5A-C6A-N1A	2.03	121.74	117.39
2	YS	501	NDP	O5D-PN-O1N	-2.02	101.16	109.07
2	YN	501	NDP	O7N-C7N-C3N	2.02	124.71	120.90
2	ZP	501	NDP	C5A-C6A-N1A	2.02	121.74	117.39
4	YM	503	A1JWG	CHC-C4B-C3B	2.02	128.72	125.26
2	YY	501	NDP	O7N-C7N-C3N	2.02	124.71	120.90
3	YY	502	LMG	O1-C7-C8	-2.02	106.02	110.90
4	ZC	503	A1JWG	CHC-C4B-C3B	2.02	128.71	125.26
2	ZB	501	NDP	C5A-C6A-N1A	2.02	121.72	117.39
3	ZT	502	LMG	C1-C2-C3	-2.02	105.80	110.00
2	YV	501	NDP	C5A-C6A-N1A	2.01	121.72	117.39
2	ZQ	501	NDP	O7N-C7N-C3N	2.01	124.69	120.90
2	ZE	501	NDP	C4B-O4B-C1B	-2.01	105.03	109.47
4	ZI	503	A1JWG	CHC-C4B-C3B	2.01	128.70	125.26
2	ZA	501	NDP	C4B-O4B-C1B	-2.01	105.03	109.47
2	YR	501	NDP	C5A-C6A-N1A	2.01	121.71	117.39
2	YS	501	NDP	O7N-C7N-C3N	2.01	124.68	120.90
4	ZE	503	A1JWG	CHC-C4B-C3B	2.00	128.68	125.26

All (131) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	YI	503	A1JWG	NC
4	YI	503	A1JWG	NA
4	YI	503	A1JWG	NB
4	YJ	503	A1JWG	NC
4	YJ	503	A1JWG	NA
4	YJ	503	A1JWG	NB
4	YK	503	A1JWG	NC
4	YK	503	A1JWG	NA
4	YK	503	A1JWG	NB
4	YL	503	A1JWG	NC
4	YL	503	A1JWG	NA
4	YL	503	A1JWG	NB
4	YM	503	A1JWG	NC
4	YM	503	A1JWG	NA
4	YM	503	A1JWG	NB
4	YN	503	A1JWG	NC
4	YN	503	A1JWG	NA

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Mol	Chain	Res	Type	Atom
4	YN	503	A1JWG	NB
4	YO	503	A1JWG	NC
4	YO	503	A1JWG	NA
4	YO	503	A1JWG	NB
4	YP	503	A1JWG	NC
4	YP	503	A1JWG	NA
4	YP	503	A1JWG	NB
4	YQ	503	A1JWG	NC
4	YQ	503	A1JWG	NA
4	YQ	503	A1JWG	NB
4	YR	503	A1JWG	NC
4	YR	503	A1JWG	NA
4	YR	503	A1JWG	NB
4	YS	503	A1JWG	NC
4	YS	503	A1JWG	NA
4	YS	503	A1JWG	NB
4	YT	503	A1JWG	NC
4	YT	503	A1JWG	NA
4	YT	503	A1JWG	NB
4	YU	503	A1JWG	NC
4	YU	503	A1JWG	NA
4	YU	503	A1JWG	NB
4	YV	503	A1JWG	NC
4	YV	503	A1JWG	NA
4	YV	503	A1JWG	NB
4	YW	503	A1JWG	NC
4	YW	503	A1JWG	NA
4	YW	503	A1JWG	NB
4	YX	503	A1JWG	NC
4	YX	503	A1JWG	NA
4	YX	503	A1JWG	NB
4	YY	503	A1JWG	NC
4	YY	503	A1JWG	NA
4	YY	503	A1JWG	NB
4	YZ	503	A1JWG	NC
4	YZ	503	A1JWG	NA
4	YZ	503	A1JWG	NB
4	ZA	503	A1JWG	NC
4	ZA	503	A1JWG	NA
4	ZA	503	A1JWG	NB
4	ZB	503	A1JWG	NC
4	ZB	503	A1JWG	NB

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Mol	Chain	Res	Type	Atom
4	ZC	503	A1JWG	NC
4	ZC	503	A1JWG	NA
4	ZC	503	A1JWG	NB
4	ZD	503	A1JWG	NC
4	ZD	503	A1JWG	NA
4	ZD	503	A1JWG	NB
4	ZE	503	A1JWG	NC
4	ZE	503	A1JWG	NA
4	ZE	503	A1JWG	NB
4	ZF	503	A1JWG	NC
4	ZF	503	A1JWG	NA
4	ZF	503	A1JWG	NB
4	ZG	503	A1JWG	NC
4	ZG	503	A1JWG	NA
4	ZG	503	A1JWG	NB
4	ZH	503	A1JWG	NC
4	ZH	503	A1JWG	NA
4	ZH	503	A1JWG	NB
4	ZI	503	A1JWG	NC
4	ZI	503	A1JWG	NA
4	ZI	503	A1JWG	NB
4	ZJ	503	A1JWG	NC
4	ZJ	503	A1JWG	NA
4	ZJ	503	A1JWG	NB
4	ZK	503	A1JWG	NC
4	ZK	503	A1JWG	NA
4	ZK	503	A1JWG	NB
4	ZL	503	A1JWG	NC
4	ZL	503	A1JWG	NA
4	ZL	503	A1JWG	NB
4	ZM	503	A1JWG	NC
4	ZM	503	A1JWG	NA
4	ZM	503	A1JWG	NB
4	ZN	503	A1JWG	NC
4	ZN	503	A1JWG	NA
4	ZN	503	A1JWG	NB
4	ZO	503	A1JWG	NC
4	ZO	503	A1JWG	NA
4	ZO	503	A1JWG	NB
4	ZP	503	A1JWG	NC
4	ZP	503	A1JWG	NA
4	ZP	503	A1JWG	NB

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Mol	Chain	Res	Type	Atom
4	ZQ	503	A1JWG	NC
4	ZQ	503	A1JWG	NA
4	ZQ	503	A1JWG	NB
4	ZR	503	A1JWG	NC
4	ZR	503	A1JWG	NA
4	ZR	503	A1JWG	NB
4	ZS	503	A1JWG	NC
4	ZS	503	A1JWG	NA
4	ZS	503	A1JWG	NB
4	ZT	503	A1JWG	NC
4	ZT	503	A1JWG	NA
4	ZT	503	A1JWG	NB
4	ZU	503	A1JWG	NC
4	ZU	503	A1JWG	NA
4	ZU	503	A1JWG	NB
4	ZV	503	A1JWG	NC
4	ZV	503	A1JWG	NA
4	ZV	503	A1JWG	NB
4	ZW	503	A1JWG	NC
4	ZW	503	A1JWG	NA
4	ZW	503	A1JWG	NB
4	ZX	503	A1JWG	NC
4	ZX	503	A1JWG	NA
4	ZX	503	A1JWG	NB
4	ZY	503	A1JWG	NC
4	ZY	503	A1JWG	NA
4	ZY	503	A1JWG	NB
4	ZZ	503	A1JWG	NC
4	ZZ	503	A1JWG	NA
4	ZZ	503	A1JWG	NB

All (1002) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	YI	501	NDP	C5B-O5B-PA-O1A
2	YI	501	NDP	C2B-O2B-P2B-O1X
2	YI	501	NDP	C2B-O2B-P2B-O2X
2	YI	501	NDP	C5D-O5D-PN-O1N
2	YI	501	NDP	C5D-O5D-PN-O2N
2	YJ	501	NDP	C5B-O5B-PA-O1A
2	YJ	501	NDP	C5D-O5D-PN-O3
2	YJ	501	NDP	C5D-O5D-PN-O2N

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Mol	Chain	Res	Type	Atoms
2	YJ	501	NDP	C2N-C3N-C7N-N7N
2	YK	501	NDP	C5B-O5B-PA-O2A
2	YK	501	NDP	C5B-O5B-PA-O3
2	YK	501	NDP	O4B-C4B-C5B-O5B
2	YK	501	NDP	C3B-C4B-C5B-O5B
2	YK	501	NDP	C5D-O5D-PN-O3
2	YL	501	NDP	C5D-O5D-PN-O2N
2	YM	501	NDP	C5B-O5B-PA-O1A
2	YM	501	NDP	C5D-O5D-PN-O1N
2	YM	501	NDP	C5D-O5D-PN-O2N
2	YN	501	NDP	C5B-O5B-PA-O1A
2	YN	501	NDP	O4D-C4D-C5D-O5D
2	YN	501	NDP	O4D-C1D-N1N-C6N
2	YN	501	NDP	C2N-C3N-C7N-N7N
2	YO	501	NDP	C5B-O5B-PA-O2A
2	YO	501	NDP	C5B-O5B-PA-O3
2	YO	501	NDP	PN-O3-PA-O5B
2	YO	501	NDP	O4B-C4B-C5B-O5B
2	YO	501	NDP	C3B-C4B-C5B-O5B
2	YO	501	NDP	C5D-O5D-PN-O3
2	YP	501	NDP	C5B-O5B-PA-O2A
2	YP	501	NDP	C5B-O5B-PA-O3
2	YP	501	NDP	C3B-C4B-C5B-O5B
2	YP	501	NDP	C5D-O5D-PN-O3
2	YP	501	NDP	C5D-O5D-PN-O1N
2	YQ	501	NDP	C5B-O5B-PA-O1A
2	YQ	501	NDP	O4B-C4B-C5B-O5B
2	YQ	501	NDP	C5D-O5D-PN-O3
2	YQ	501	NDP	C5D-O5D-PN-O2N
2	YR	501	NDP	C5D-O5D-PN-O1N
2	YR	501	NDP	C5D-O5D-PN-O2N
2	YR	501	NDP	C2N-C3N-C7N-N7N
2	YS	501	NDP	C5B-O5B-PA-O2A
2	YS	501	NDP	C2B-O2B-P2B-O3X
2	YS	501	NDP	C5D-O5D-PN-O1N
2	YT	501	NDP	C5B-O5B-PA-O1A
2	YT	501	NDP	C5B-O5B-PA-O2A
2	YT	501	NDP	PN-O3-PA-O5B
2	YT	501	NDP	C5D-O5D-PN-O1N
2	YT	501	NDP	C5D-O5D-PN-O2N
2	YU	501	NDP	C5B-O5B-PA-O1A
2	YU	501	NDP	C5D-O5D-PN-O3

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Mol	Chain	Res	Type	Atoms
2	YU	501	NDP	C5D-O5D-PN-O2N
2	YV	501	NDP	C5D-O5D-PN-O3
2	YV	501	NDP	O4D-C4D-C5D-O5D
2	YV	501	NDP	C2N-C3N-C7N-N7N
2	YW	501	NDP	C5B-O5B-PA-O1A
2	YW	501	NDP	C5B-O5B-PA-O2A
2	YW	501	NDP	PN-O3-PA-O5B
2	YW	501	NDP	C5D-O5D-PN-O1N
2	YX	501	NDP	C5D-O5D-PN-O2N
2	YY	501	NDP	C5B-O5B-PA-O1A
2	YY	501	NDP	C5D-O5D-PN-O3
2	YY	501	NDP	C5D-O5D-PN-O1N
2	YY	501	NDP	C5D-O5D-PN-O2N
2	YZ	501	NDP	C5B-O5B-PA-O1A
2	YZ	501	NDP	C5D-O5D-PN-O3
2	ZA	501	NDP	C5B-O5B-PA-O2A
2	ZA	501	NDP	C5B-O5B-PA-O3
2	ZA	501	NDP	PN-O3-PA-O5B
2	ZA	501	NDP	O4B-C4B-C5B-O5B
2	ZA	501	NDP	C3B-C4B-C5B-O5B
2	ZA	501	NDP	C5D-O5D-PN-O3
2	ZB	501	NDP	C5D-O5D-PN-O1N
2	ZB	501	NDP	C5D-O5D-PN-O2N
2	ZC	501	NDP	C5B-O5B-PA-O1A
2	ZC	501	NDP	C5D-O5D-PN-O3
2	ZC	501	NDP	C5D-O5D-PN-O2N
2	ZD	501	NDP	C5B-O5B-PA-O1A
2	ZD	501	NDP	PN-O3-PA-O5B
2	ZD	501	NDP	C5D-O5D-PN-O3
2	ZD	501	NDP	C5D-O5D-PN-O2N
2	ZD	501	NDP	C2N-C3N-C7N-N7N
2	ZE	501	NDP	C5B-O5B-PA-O1A
2	ZE	501	NDP	C5D-O5D-PN-O2N
2	ZF	501	NDP	C5D-O5D-PN-O2N
2	ZG	501	NDP	C5B-O5B-PA-O1A
2	ZG	501	NDP	O4B-C4B-C5B-O5B
2	ZG	501	NDP	C5D-O5D-PN-O3
2	ZG	501	NDP	C5D-O5D-PN-O2N
2	ZH	501	NDP	C5B-O5B-PA-O3
2	ZH	501	NDP	PN-O3-PA-O5B
2	ZH	501	NDP	C3B-C4B-C5B-O5B
2	ZH	501	NDP	C5D-O5D-PN-O1N

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Mol	Chain	Res	Type	Atoms
2	ZH	501	NDP	C2N-C3N-C7N-N7N
2	ZI	501	NDP	C5B-O5B-PA-O2A
2	ZI	501	NDP	C5B-O5B-PA-O3
2	ZI	501	NDP	O4B-C4B-C5B-O5B
2	ZI	501	NDP	C3B-C4B-C5B-O5B
2	ZI	501	NDP	C5D-O5D-PN-O3
2	ZJ	501	NDP	C5B-O5B-PA-O2A
2	ZJ	501	NDP	PN-O3-PA-O5B
2	ZJ	501	NDP	C3B-C4B-C5B-O5B
2	ZJ	501	NDP	C5D-O5D-PN-O3
2	ZK	501	NDP	C5B-O5B-PA-O1A
2	ZK	501	NDP	O4B-C4B-C5B-O5B
2	ZK	501	NDP	C5D-O5D-PN-O3
2	ZL	501	NDP	C5B-O5B-PA-O1A
2	ZL	501	NDP	C5B-O5B-PA-O2A
2	ZL	501	NDP	C5D-O5D-PN-O3
2	ZM	501	NDP	C5B-O5B-PA-O1A
2	ZM	501	NDP	C5D-O5D-PN-O2N
2	ZN	501	NDP	C5D-O5D-PN-O2N
2	ZO	501	NDP	C5B-O5B-PA-O1A
2	ZO	501	NDP	C5D-O5D-PN-O3
2	ZO	501	NDP	C5D-O5D-PN-O2N
2	ZP	501	NDP	C5B-O5B-PA-O1A
2	ZP	501	NDP	C5D-O5D-PN-O3
2	ZP	501	NDP	C5D-O5D-PN-O2N
2	ZP	501	NDP	C2N-C3N-C7N-N7N
2	ZQ	501	NDP	C5B-O5B-PA-O1A
2	ZQ	501	NDP	C5D-O5D-PN-O2N
2	ZR	501	NDP	C2B-O2B-P2B-O1X
2	ZR	501	NDP	C5D-O5D-PN-O2N
2	ZS	501	NDP	C5B-O5B-PA-O1A
2	ZS	501	NDP	C5B-O5B-PA-O2A
2	ZS	501	NDP	C3B-C4B-C5B-O5B
2	ZS	501	NDP	C5D-O5D-PN-O3
2	ZT	501	NDP	C5B-O5B-PA-O1A
2	ZT	501	NDP	PN-O3-PA-O5B
2	ZT	501	NDP	C2B-O2B-P2B-O2X
2	ZT	501	NDP	C5D-O5D-PN-O3
2	ZT	501	NDP	C5D-O5D-PN-O2N
2	ZT	501	NDP	C2N-C3N-C7N-N7N
2	ZU	501	NDP	C5B-O5B-PA-O2A
2	ZU	501	NDP	C5B-O5B-PA-O3

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Mol	Chain	Res	Type	Atoms
2	ZU	501	NDP	PN-O3-PA-O5B
2	ZU	501	NDP	O4B-C4B-C5B-O5B
2	ZU	501	NDP	C3B-C4B-C5B-O5B
2	ZU	501	NDP	C5D-O5D-PN-O3
2	ZV	501	NDP	C5B-O5B-PA-O2A
2	ZV	501	NDP	C5B-O5B-PA-O3
2	ZV	501	NDP	PN-O3-PA-O5B
2	ZV	501	NDP	O4B-C4B-C5B-O5B
2	ZV	501	NDP	C3B-C4B-C5B-O5B
2	ZV	501	NDP	C2B-O2B-P2B-O3X
2	ZV	501	NDP	C5D-O5D-PN-O3
2	ZW	501	NDP	C5B-O5B-PA-O1A
2	ZW	501	NDP	C5D-O5D-PN-O3
2	ZX	501	NDP	C5B-O5B-PA-O1A
2	ZX	501	NDP	O4B-C4B-C5B-O5B
2	ZX	501	NDP	C5D-O5D-PN-O3
2	ZY	501	NDP	C5B-O5B-PA-O2A
2	ZY	501	NDP	C5D-O5D-PN-O1N
2	ZY	501	NDP	O4D-C4D-C5D-O5D
2	ZZ	501	NDP	C5B-O5B-PA-O1A
2	ZZ	501	NDP	C5B-O5B-PA-O2A
2	ZZ	501	NDP	PN-O3-PA-O5B
2	ZZ	501	NDP	C5D-O5D-PN-O1N
2	ZZ	501	NDP	C5D-O5D-PN-O2N
3	YI	502	LMG	O9-C10-O7-C8
3	YI	502	LMG	C11-C10-O7-C8
3	YK	502	LMG	C11-C10-O7-C8
3	YM	502	LMG	C11-C10-O7-C8
3	YO	502	LMG	C11-C10-O7-C8
3	YR	502	LMG	O9-C10-O7-C8
3	YR	502	LMG	C11-C10-O7-C8
3	YS	502	LMG	O9-C10-O7-C8
3	YS	502	LMG	C11-C10-O7-C8
3	YT	502	LMG	C11-C10-O7-C8
3	YU	502	LMG	O9-C10-O7-C8
3	YU	502	LMG	C11-C10-O7-C8
3	YV	502	LMG	C2-C1-O1-C7
3	YV	502	LMG	O6-C1-O1-C7
3	YV	502	LMG	C11-C10-O7-C8
3	YW	502	LMG	C2-C1-O1-C7
3	YW	502	LMG	O6-C1-O1-C7
3	YW	502	LMG	C11-C10-O7-C8

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Mol	Chain	Res	Type	Atoms
3	YX	502	LMG	O9-C10-O7-C8
3	YX	502	LMG	C11-C10-O7-C8
3	YY	502	LMG	C11-C10-O7-C8
3	ZA	502	LMG	O9-C10-O7-C8
3	ZA	502	LMG	C11-C10-O7-C8
3	ZB	502	LMG	C11-C10-O7-C8
3	ZE	502	LMG	C11-C10-O7-C8
3	ZG	502	LMG	C2-C1-O1-C7
3	ZG	502	LMG	O6-C1-O1-C7
3	ZG	502	LMG	O9-C10-O7-C8
3	ZG	502	LMG	C11-C10-O7-C8
3	ZH	502	LMG	C11-C10-O7-C8
3	ZK	502	LMG	C11-C10-O7-C8
3	ZL	502	LMG	O9-C10-O7-C8
3	ZL	502	LMG	C11-C10-O7-C8
3	ZM	502	LMG	C2-C1-O1-C7
3	ZM	502	LMG	O6-C1-O1-C7
3	ZM	502	LMG	O9-C10-O7-C8
3	ZM	502	LMG	C11-C10-O7-C8
3	ZN	502	LMG	C2-C1-O1-C7
3	ZN	502	LMG	O6-C1-O1-C7
3	ZP	502	LMG	C11-C10-O7-C8
3	ZR	502	LMG	O9-C10-O7-C8
3	ZR	502	LMG	C11-C10-O7-C8
3	ZS	502	LMG	C11-C10-O7-C8
3	ZT	502	LMG	O9-C10-O7-C8
3	ZT	502	LMG	C11-C10-O7-C8
3	ZU	502	LMG	O9-C10-O7-C8
3	ZU	502	LMG	C11-C10-O7-C8
3	ZW	502	LMG	O9-C10-O7-C8
3	ZW	502	LMG	C11-C10-O7-C8
3	ZX	502	LMG	C2-C1-O1-C7
3	ZX	502	LMG	O6-C1-O1-C7
3	ZX	502	LMG	C8-C7-O1-C1
3	ZX	502	LMG	C11-C10-O7-C8
3	ZY	502	LMG	O9-C10-O7-C8
3	ZY	502	LMG	C11-C10-O7-C8
3	ZZ	502	LMG	O9-C10-O7-C8
3	ZZ	502	LMG	C11-C10-O7-C8
4	YI	503	A1JWG	CHA-CBD-CGD-O2D
4	YJ	503	A1JWG	C1A-C2A-CAA-CBA
4	YJ	503	A1JWG	C2B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
4	YJ	503	A1JWG	C4B-C3B-CAB-CBB
4	YM	503	A1JWG	CAD-CBD-CGD-O2D
4	YN	503	A1JWG	C1A-C2A-CAA-CBA
4	YN	503	A1JWG	C3A-C2A-CAA-CBA
4	YN	503	A1JWG	C2B-C3B-CAB-CBB
4	YN	503	A1JWG	C4B-C3B-CAB-CBB
4	YP	503	A1JWG	CHA-CBD-CGD-O1D
4	YQ	503	A1JWG	C2B-C3B-CAB-CBB
4	YQ	503	A1JWG	CHA-CBD-CGD-O2D
4	YR	503	A1JWG	C1A-C2A-CAA-CBA
4	YR	503	A1JWG	C3A-C2A-CAA-CBA
4	YR	503	A1JWG	C2B-C3B-CAB-CBB
4	YR	503	A1JWG	C4B-C3B-CAB-CBB
4	YS	503	A1JWG	CBD-CGD-O2D-CED
4	YT	503	A1JWG	CHA-CBD-CGD-O2D
4	YU	503	A1JWG	C1A-C2A-CAA-CBA
4	YU	503	A1JWG	C3A-C2A-CAA-CBA
4	YU	503	A1JWG	C2B-C3B-CAB-CBB
4	YV	503	A1JWG	C2B-C3B-CAB-CBB
4	YV	503	A1JWG	C4B-C3B-CAB-CBB
4	YW	503	A1JWG	CBD-CGD-O2D-CED
4	YX	503	A1JWG	CHA-CBD-CGD-O1D
4	ZD	503	A1JWG	C1A-C2A-CAA-CBA
4	ZD	503	A1JWG	C3A-C2A-CAA-CBA
4	ZD	503	A1JWG	C2B-C3B-CAB-CBB
4	ZD	503	A1JWG	C4B-C3B-CAB-CBB
4	ZE	503	A1JWG	C2B-C3B-CAB-CBB
4	ZG	503	A1JWG	C2B-C3B-CAB-CBB
4	ZH	503	A1JWG	C1A-C2A-CAA-CBA
4	ZH	503	A1JWG	C3A-C2A-CAA-CBA
4	ZH	503	A1JWG	C2B-C3B-CAB-CBB
4	ZH	503	A1JWG	C4B-C3B-CAB-CBB
4	ZI	503	A1JWG	CHA-CBD-CGD-O2D
4	ZJ	503	A1JWG	CHA-CBD-CGD-O1D
4	ZK	503	A1JWG	C2B-C3B-CAB-CBB
4	ZL	503	A1JWG	C1A-C2A-CAA-CBA
4	ZM	503	A1JWG	CHA-CBD-CGD-O2D
4	ZN	503	A1JWG	CHA-CBD-CGD-O1D
4	ZO	503	A1JWG	C2B-C3B-CAB-CBB
4	ZP	503	A1JWG	C1A-C2A-CAA-CBA
4	ZP	503	A1JWG	C2B-C3B-CAB-CBB
4	ZP	503	A1JWG	C4B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
4	ZQ	503	A1JWG	CBD-CGD-O2D-CED
4	ZS	503	A1JWG	C2B-C3B-CAB-CBB
4	ZT	503	A1JWG	C1A-C2A-CAA-CBA
4	ZT	503	A1JWG	C2B-C3B-CAB-CBB
4	ZT	503	A1JWG	C4B-C3B-CAB-CBB
4	ZU	503	A1JWG	CHA-CBD-CGD-O2D
4	ZW	503	A1JWG	CHA-CBD-CGD-O1D
4	ZX	503	A1JWG	CHA-CBD-CGD-O1D
4	ZY	503	A1JWG	C2B-C3B-CAB-CBB
4	ZY	503	A1JWG	C4B-C3B-CAB-CBB
3	YK	502	LMG	O9-C10-O7-C8
3	YM	502	LMG	O9-C10-O7-C8
3	YY	502	LMG	O9-C10-O7-C8
3	ZE	502	LMG	O9-C10-O7-C8
3	ZS	502	LMG	O9-C10-O7-C8
3	YO	502	LMG	O9-C10-O7-C8
3	YT	502	LMG	O9-C10-O7-C8
3	YV	502	LMG	O9-C10-O7-C8
3	YW	502	LMG	O9-C10-O7-C8
3	ZH	502	LMG	O9-C10-O7-C8
3	ZK	502	LMG	O9-C10-O7-C8
3	ZP	502	LMG	O9-C10-O7-C8
3	ZX	502	LMG	O9-C10-O7-C8
4	YW	503	A1JWG	O1D-CGD-O2D-CED
4	ZQ	503	A1JWG	O1D-CGD-O2D-CED
3	YW	502	LMG	C29-C28-O8-C9
4	YS	503	A1JWG	O1D-CGD-O2D-CED
3	YW	502	LMG	O10-C28-O8-C9
4	YY	503	A1JWG	CBD-CGD-O2D-CED
4	ZK	503	A1JWG	CBD-CGD-O2D-CED
4	ZO	503	A1JWG	CBD-CGD-O2D-CED
3	YN	502	LMG	O10-C28-O8-C9
3	YR	502	LMG	O10-C28-O8-C9
3	ZH	502	LMG	O10-C28-O8-C9
3	ZU	502	LMG	O10-C28-O8-C9
3	YS	502	LMG	O6-C5-C6-O5
3	ZB	502	LMG	O9-C10-O7-C8
3	YO	502	LMG	O10-C28-O8-C9
3	YT	502	LMG	O10-C28-O8-C9
3	ZV	502	LMG	O10-C28-O8-C9
3	ZZ	502	LMG	O10-C28-O8-C9
4	YI	503	A1JWG	C2A-CAA-CBA-CGA

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Mol	Chain	Res	Type	Atoms
4	YK	503	A1JWG	C2A-CAA-CBA-CGA
4	YQ	503	A1JWG	C2A-CAA-CBA-CGA
4	YS	503	A1JWG	C2A-CAA-CBA-CGA
4	YT	503	A1JWG	C2A-CAA-CBA-CGA
4	YV	503	A1JWG	C2A-CAA-CBA-CGA
4	YY	503	A1JWG	C2A-CAA-CBA-CGA
4	ZA	503	A1JWG	C2A-CAA-CBA-CGA
4	ZB	503	A1JWG	C2A-CAA-CBA-CGA
4	ZC	503	A1JWG	C2A-CAA-CBA-CGA
4	ZE	503	A1JWG	C2A-CAA-CBA-CGA
4	ZF	503	A1JWG	C2A-CAA-CBA-CGA
4	ZG	503	A1JWG	C2A-CAA-CBA-CGA
4	ZJ	503	A1JWG	C2A-CAA-CBA-CGA
4	ZN	503	A1JWG	C2A-CAA-CBA-CGA
4	ZO	503	A1JWG	C2A-CAA-CBA-CGA
4	ZS	503	A1JWG	C2A-CAA-CBA-CGA
4	ZY	503	A1JWG	C2A-CAA-CBA-CGA
3	ZY	502	LMG	O10-C28-O8-C9
3	YN	502	LMG	C29-C28-O8-C9
3	YR	502	LMG	C29-C28-O8-C9
3	YT	502	LMG	C29-C28-O8-C9
3	ZH	502	LMG	C29-C28-O8-C9
3	ZV	502	LMG	C29-C28-O8-C9
3	ZY	502	LMG	C29-C28-O8-C9
3	ZZ	502	LMG	C29-C28-O8-C9
3	ZO	502	LMG	C11-C10-O7-C8
3	YX	502	LMG	O10-C28-O8-C9
3	ZJ	502	LMG	O10-C28-O8-C9
3	YO	502	LMG	C29-C28-O8-C9
3	ZU	502	LMG	C29-C28-O8-C9
2	YV	501	NDP	O4D-C1D-N1N-C6N
4	YM	503	A1JWG	CBD-CGD-O2D-CED
4	YU	503	A1JWG	CBD-CGD-O2D-CED
3	YM	502	LMG	O10-C28-O8-C9
3	YP	502	LMG	O10-C28-O8-C9
3	ZC	502	LMG	O10-C28-O8-C9
3	ZJ	502	LMG	C29-C28-O8-C9
3	ZP	502	LMG	C29-C28-O8-C9
3	ZP	502	LMG	O6-C5-C6-O5
4	YL	503	A1JWG	CBD-CGD-O2D-CED
4	ZS	503	A1JWG	CBD-CGD-O2D-CED
3	YI	502	LMG	O10-C28-O8-C9

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Mol	Chain	Res	Type	Atoms
3	YX	502	LMG	C29-C28-O8-C9
2	YP	501	NDP	O4B-C4B-C5B-O5B
2	YS	501	NDP	O4B-C4B-C5B-O5B
2	YS	501	NDP	C3B-C4B-C5B-O5B
2	YS	501	NDP	O4D-C4D-C5D-O5D
2	YT	501	NDP	O4B-C4B-C5B-O5B
2	YT	501	NDP	C3B-C4B-C5B-O5B
2	YU	501	NDP	O4B-C4B-C5B-O5B
2	YU	501	NDP	C3B-C4B-C5B-O5B
2	YW	501	NDP	O4B-C4B-C5B-O5B
2	YW	501	NDP	C3B-C4B-C5B-O5B
2	YW	501	NDP	O4D-C4D-C5D-O5D
2	YW	501	NDP	C3D-C4D-C5D-O5D
2	YZ	501	NDP	O4B-C4B-C5B-O5B
2	YZ	501	NDP	C3B-C4B-C5B-O5B
2	ZC	501	NDP	O4B-C4B-C5B-O5B
2	ZC	501	NDP	C3B-C4B-C5B-O5B
2	ZG	501	NDP	C3B-C4B-C5B-O5B
2	ZJ	501	NDP	O4B-C4B-C5B-O5B
2	ZL	501	NDP	O4B-C4B-C5B-O5B
2	ZL	501	NDP	C3B-C4B-C5B-O5B
2	ZO	501	NDP	O4B-C4B-C5B-O5B
2	ZO	501	NDP	C3B-C4B-C5B-O5B
2	ZQ	501	NDP	O4B-C4B-C5B-O5B
2	ZQ	501	NDP	C3B-C4B-C5B-O5B
2	ZS	501	NDP	O4B-C4B-C5B-O5B
2	ZW	501	NDP	O4B-C4B-C5B-O5B
2	ZW	501	NDP	C3B-C4B-C5B-O5B
2	ZX	501	NDP	C3B-C4B-C5B-O5B
2	ZY	501	NDP	O4B-C4B-C5B-O5B
2	ZY	501	NDP	C3B-C4B-C5B-O5B
2	ZY	501	NDP	C3D-C4D-C5D-O5D
2	ZZ	501	NDP	O4B-C4B-C5B-O5B
2	ZZ	501	NDP	C3B-C4B-C5B-O5B
3	ZW	502	LMG	O10-C28-O8-C9
3	YZ	502	LMG	O10-C28-O8-C9
3	YP	502	LMG	C29-C28-O8-C9
3	YZ	502	LMG	C29-C28-O8-C9
4	YO	503	A1JWG	C2A-CAA-CBA-CGA
4	YP	503	A1JWG	C2A-CAA-CBA-CGA
4	YX	503	A1JWG	C2A-CAA-CBA-CGA
4	ZK	503	A1JWG	C2A-CAA-CBA-CGA

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Mol	Chain	Res	Type	Atoms
4	ZV	503	A1JWG	C2A-CAA-CBA-CGA
3	YR	502	LMG	O6-C1-O1-C7
3	YN	502	LMG	C11-C10-O7-C8
3	YM	502	LMG	C29-C28-O8-C9
3	ZC	502	LMG	C29-C28-O8-C9
3	YS	502	LMG	C4-C5-C6-O5
3	ZT	502	LMG	C29-C28-O8-C9
4	YI	503	A1JWG	CBD-CGD-O2D-CED
4	ZW	503	A1JWG	O1D-CGD-O2D-CED
3	YI	502	LMG	C29-C28-O8-C9
4	ZZ	503	A1JWG	CBD-CGD-O2D-CED
2	YI	501	NDP	O4D-C1D-N1N-C6N
2	YU	501	NDP	O4D-C1D-N1N-C6N
2	ZK	501	NDP	O4D-C1D-N1N-C6N
2	ZL	501	NDP	O4D-C1D-N1N-C6N
2	ZS	501	NDP	O4D-C1D-N1N-C6N
3	ZW	502	LMG	C29-C28-O8-C9
3	YR	502	LMG	C2-C1-O1-C7
3	ZT	502	LMG	C2-C1-O1-C7
3	YJ	502	LMG	C11-C10-O7-C8
2	YN	501	NDP	C3D-C4D-C5D-O5D
2	YQ	501	NDP	C3B-C4B-C5B-O5B
2	YS	501	NDP	C3D-C4D-C5D-O5D
2	YV	501	NDP	C3D-C4D-C5D-O5D
2	ZH	501	NDP	O4B-C4B-C5B-O5B
2	ZH	501	NDP	O4D-C4D-C5D-O5D
2	ZH	501	NDP	C3D-C4D-C5D-O5D
2	ZK	501	NDP	C3B-C4B-C5B-O5B
2	ZM	501	NDP	O4B-C4B-C5B-O5B
3	ZP	502	LMG	O10-C28-O8-C9
4	ZU	503	A1JWG	CBD-CGD-O2D-CED
2	ZG	501	NDP	O4D-C1D-N1N-C6N
4	YM	503	A1JWG	C2A-CAA-CBA-CGA
4	YW	503	A1JWG	C2A-CAA-CBA-CGA
4	ZQ	503	A1JWG	C2A-CAA-CBA-CGA
3	ZT	502	LMG	O6-C1-O1-C7
3	ZP	502	LMG	C4-C5-C6-O5
3	YV	502	LMG	C4-C5-C6-O5
3	ZT	502	LMG	O10-C28-O8-C9
3	ZT	502	LMG	O6-C5-C6-O5
4	YU	503	A1JWG	C2A-CAA-CBA-CGA
4	ZU	503	A1JWG	C2A-CAA-CBA-CGA

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Mol	Chain	Res	Type	Atoms
2	YM	501	NDP	O4D-C1D-N1N-C6N
2	YY	501	NDP	O4D-C1D-N1N-C6N
2	ZC	501	NDP	O4D-C1D-N1N-C6N
2	ZO	501	NDP	O4D-C1D-N1N-C6N
4	YN	503	A1JWG	CBD-CGD-O2D-CED
3	ZV	502	LMG	O7-C8-C9-O8
3	ZO	502	LMG	O9-C10-O7-C8
4	YI	503	A1JWG	C3A-C2A-CAA-CBA
4	YJ	503	A1JWG	C3A-C2A-CAA-CBA
4	YL	503	A1JWG	C3A-C2A-CAA-CBA
4	YM	503	A1JWG	C3A-C2A-CAA-CBA
4	YP	503	A1JWG	C3A-C2A-CAA-CBA
4	YQ	503	A1JWG	C3A-C2A-CAA-CBA
4	YS	503	A1JWG	C3A-C2A-CAA-CBA
4	YT	503	A1JWG	C3A-C2A-CAA-CBA
4	YV	503	A1JWG	C3A-C2A-CAA-CBA
4	YX	503	A1JWG	C3A-C2A-CAA-CBA
4	YY	503	A1JWG	C3A-C2A-CAA-CBA
4	YZ	503	A1JWG	C3A-C2A-CAA-CBA
4	ZB	503	A1JWG	C3A-C2A-CAA-CBA
4	ZC	503	A1JWG	C3A-C2A-CAA-CBA
4	ZF	503	A1JWG	C3A-C2A-CAA-CBA
4	ZG	503	A1JWG	C3A-C2A-CAA-CBA
4	ZJ	503	A1JWG	C3A-C2A-CAA-CBA
4	ZK	503	A1JWG	C3A-C2A-CAA-CBA
4	ZL	503	A1JWG	C3A-C2A-CAA-CBA
4	ZN	503	A1JWG	C3A-C2A-CAA-CBA
4	ZO	503	A1JWG	C3A-C2A-CAA-CBA
4	ZP	503	A1JWG	C3A-C2A-CAA-CBA
4	ZQ	503	A1JWG	C3A-C2A-CAA-CBA
4	ZR	503	A1JWG	C3A-C2A-CAA-CBA
4	ZS	503	A1JWG	C3A-C2A-CAA-CBA
4	ZT	503	A1JWG	C3A-C2A-CAA-CBA
4	ZU	503	A1JWG	C3A-C2A-CAA-CBA
4	ZW	503	A1JWG	C3A-C2A-CAA-CBA
4	ZX	503	A1JWG	C3A-C2A-CAA-CBA
4	ZZ	503	A1JWG	C3A-C2A-CAA-CBA
3	ZQ	502	LMG	O6-C5-C6-O5
2	YQ	501	NDP	O4D-C1D-N1N-C6N
4	ZD	503	A1JWG	CBD-CGD-O2D-CED
4	ZR	503	A1JWG	C2A-CAA-CBA-CGA
2	YV	501	NDP	C3B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
2	ZM	501	NDP	C3B-C4B-C5B-O5B
2	YK	501	NDP	O4D-C1D-N1N-C6N
2	ZI	501	NDP	O4D-C1D-N1N-C6N
2	ZW	501	NDP	O4D-C1D-N1N-C6N
4	YI	503	A1JWG	C2B-C3B-CAB-CBB
4	YK	503	A1JWG	C2B-C3B-CAB-CBB
4	YO	503	A1JWG	C2B-C3B-CAB-CBB
4	YS	503	A1JWG	C2B-C3B-CAB-CBB
4	YW	503	A1JWG	C2B-C3B-CAB-CBB
4	YY	503	A1JWG	C2B-C3B-CAB-CBB
4	ZA	503	A1JWG	C2B-C3B-CAB-CBB
4	ZC	503	A1JWG	C2B-C3B-CAB-CBB
4	ZI	503	A1JWG	C2B-C3B-CAB-CBB
4	ZJ	503	A1JWG	C2B-C3B-CAB-CBB
4	ZL	503	A1JWG	C2B-C3B-CAB-CBB
4	ZM	503	A1JWG	C2B-C3B-CAB-CBB
4	ZN	503	A1JWG	C2B-C3B-CAB-CBB
4	ZQ	503	A1JWG	C2B-C3B-CAB-CBB
4	ZR	503	A1JWG	C2B-C3B-CAB-CBB
4	ZU	503	A1JWG	C2B-C3B-CAB-CBB
4	ZV	503	A1JWG	C2B-C3B-CAB-CBB
4	ZZ	503	A1JWG	C2B-C3B-CAB-CBB
3	YN	502	LMG	O9-C10-O7-C8
4	YI	503	A1JWG	C4B-C3B-CAB-CBB
4	YK	503	A1JWG	C4B-C3B-CAB-CBB
4	YO	503	A1JWG	C4B-C3B-CAB-CBB
4	YQ	503	A1JWG	C4B-C3B-CAB-CBB
4	YS	503	A1JWG	C4B-C3B-CAB-CBB
4	YU	503	A1JWG	C4B-C3B-CAB-CBB
4	YW	503	A1JWG	C4B-C3B-CAB-CBB
4	YY	503	A1JWG	C4B-C3B-CAB-CBB
4	ZA	503	A1JWG	C4B-C3B-CAB-CBB
4	ZC	503	A1JWG	C4B-C3B-CAB-CBB
4	ZE	503	A1JWG	C4B-C3B-CAB-CBB
4	ZG	503	A1JWG	C4B-C3B-CAB-CBB
4	ZI	503	A1JWG	C4B-C3B-CAB-CBB
4	ZJ	503	A1JWG	C4B-C3B-CAB-CBB
4	ZK	503	A1JWG	C4B-C3B-CAB-CBB
4	ZM	503	A1JWG	C4B-C3B-CAB-CBB
4	ZO	503	A1JWG	C4B-C3B-CAB-CBB
4	ZQ	503	A1JWG	C4B-C3B-CAB-CBB
4	ZR	503	A1JWG	C4B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
4	ZS	503	A1JWG	C4B-C3B-CAB-CBB
4	ZU	503	A1JWG	C4B-C3B-CAB-CBB
4	ZV	503	A1JWG	C4B-C3B-CAB-CBB
3	ZA	502	LMG	O6-C5-C6-O5
2	ZK	501	NDP	O4D-C4D-C5D-O5D
4	ZZ	503	A1JWG	C2A-CAA-CBA-CGA
3	ZG	502	LMG	O6-C5-C6-O5
4	YI	503	A1JWG	C1A-C2A-CAA-CBA
4	YL	503	A1JWG	C1A-C2A-CAA-CBA
4	YM	503	A1JWG	C1A-C2A-CAA-CBA
4	YP	503	A1JWG	C1A-C2A-CAA-CBA
4	YQ	503	A1JWG	C1A-C2A-CAA-CBA
4	YS	503	A1JWG	C1A-C2A-CAA-CBA
4	YT	503	A1JWG	C1A-C2A-CAA-CBA
4	YV	503	A1JWG	C1A-C2A-CAA-CBA
4	YX	503	A1JWG	C1A-C2A-CAA-CBA
4	YY	503	A1JWG	C1A-C2A-CAA-CBA
4	YZ	503	A1JWG	C1A-C2A-CAA-CBA
4	ZB	503	A1JWG	C1A-C2A-CAA-CBA
4	ZC	503	A1JWG	C1A-C2A-CAA-CBA
4	ZF	503	A1JWG	C1A-C2A-CAA-CBA
4	ZG	503	A1JWG	C1A-C2A-CAA-CBA
4	ZJ	503	A1JWG	C1A-C2A-CAA-CBA
4	ZK	503	A1JWG	C1A-C2A-CAA-CBA
4	ZN	503	A1JWG	C1A-C2A-CAA-CBA
4	ZO	503	A1JWG	C1A-C2A-CAA-CBA
4	ZQ	503	A1JWG	C1A-C2A-CAA-CBA
4	ZR	503	A1JWG	C1A-C2A-CAA-CBA
4	ZS	503	A1JWG	C1A-C2A-CAA-CBA
4	ZU	503	A1JWG	C1A-C2A-CAA-CBA
4	ZW	503	A1JWG	C1A-C2A-CAA-CBA
4	ZX	503	A1JWG	C1A-C2A-CAA-CBA
4	ZZ	503	A1JWG	C1A-C2A-CAA-CBA
2	YO	501	NDP	O4D-C1D-N1N-C6N
2	ZA	501	NDP	O4D-C1D-N1N-C6N
4	YR	503	A1JWG	CBD-CGD-O2D-CED
3	ZV	502	LMG	C7-C8-C9-O8
3	YW	502	LMG	O6-C5-C6-O5
3	ZD	502	LMG	C29-C28-O8-C9
3	ZC	502	LMG	O6-C5-C6-O5
3	YJ	502	LMG	C29-C28-O8-C9
2	ZE	501	NDP	O4D-C1D-N1N-C6N

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Mol	Chain	Res	Type	Atoms
2	YY	501	NDP	O4B-C4B-C5B-O5B
4	ZL	503	A1JWG	O1D-CGD-O2D-CED
3	YQ	502	LMG	O6-C5-C6-O5
3	YV	502	LMG	O6-C5-C6-O5
3	YJ	502	LMG	O9-C10-O7-C8
3	ZX	502	LMG	O6-C5-C6-O5
3	ZF	502	LMG	O6-C5-C6-O5
3	ZV	502	LMG	C11-C10-O7-C8
3	ZZ	502	LMG	O7-C8-C9-O8
2	ZM	501	NDP	O4D-C1D-N1N-C6N
2	ZN	501	NDP	O4D-C1D-N1N-C6N
2	ZV	501	NDP	O4D-C1D-N1N-C6N
2	YV	501	NDP	PA-O3-PN-O1N
2	ZH	501	NDP	C2D-C1D-N1N-C6N
2	YM	501	NDP	O4B-C4B-C5B-O5B
2	YQ	501	NDP	O4D-C4D-C5D-O5D
2	YZ	501	NDP	O4D-C4D-C5D-O5D
2	ZE	501	NDP	O4B-C4B-C5B-O5B
2	ZK	501	NDP	C3D-C4D-C5D-O5D
2	ZX	501	NDP	O4D-C4D-C5D-O5D
3	ZM	502	LMG	O6-C5-C6-O5
4	YL	503	A1JWG	C2A-CAA-CBA-CGA
3	YW	502	LMG	C4-C5-C6-O5
2	YL	501	NDP	O4D-C1D-N1N-C6N
2	YP	501	NDP	O4D-C1D-N1N-C6N
2	YX	501	NDP	O4D-C1D-N1N-C6N
2	ZB	501	NDP	O4D-C1D-N1N-C6N
2	ZJ	501	NDP	O4D-C1D-N1N-C6N
2	ZQ	501	NDP	O4D-C1D-N1N-C6N
2	ZU	501	NDP	O4D-C1D-N1N-C6N
2	ZX	501	NDP	O4D-C1D-N1N-C6N
3	ZT	502	LMG	O1-C7-C8-C9
3	ZT	502	LMG	C7-C8-C9-O8
3	ZZ	502	LMG	C7-C8-C9-O8
3	ZM	502	LMG	C4-C5-C6-O5
3	ZO	502	LMG	O7-C8-C9-O8
3	ZW	502	LMG	C4-C5-C6-O5
2	YJ	501	NDP	PN-O3-PA-O5B
2	YK	501	NDP	PN-O3-PA-O5B
2	YN	501	NDP	PN-O3-PA-O5B
2	YP	501	NDP	PN-O3-PA-O5B
2	YS	501	NDP	PN-O3-PA-O5B

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Mol	Chain	Res	Type	Atoms
2	ZI	501	NDP	PN-O3-PA-O5B
2	ZL	501	NDP	PN-O3-PA-O5B
2	ZP	501	NDP	PN-O3-PA-O5B
2	ZY	501	NDP	PN-O3-PA-O5B
4	ZC	503	A1JWG	O1D-CGD-O2D-CED
2	YQ	501	NDP	C3D-C4D-C5D-O5D
2	YV	501	NDP	O4B-C4B-C5B-O5B
2	YY	501	NDP	C3B-C4B-C5B-O5B
2	YZ	501	NDP	C3D-C4D-C5D-O5D
2	ZS	501	NDP	O4D-C4D-C5D-O5D
2	YK	501	NDP	C2B-O2B-P2B-O1X
2	YO	501	NDP	C2B-O2B-P2B-O1X
2	YV	501	NDP	C2B-O2B-P2B-O1X
2	ZH	501	NDP	C2B-O2B-P2B-O1X
4	YJ	503	A1JWG	CAD-CBD-CGD-O2D
4	YP	503	A1JWG	C2B-C3B-CAB-CBB
4	ZB	503	A1JWG	C2B-C3B-CAB-CBB
4	ZF	503	A1JWG	C2B-C3B-CAB-CBB
4	ZR	503	A1JWG	CAD-CBD-CGD-O2D
4	ZT	503	A1JWG	CAD-CBD-CGD-O2D
4	ZW	503	A1JWG	C2B-C3B-CAB-CBB
4	ZX	503	A1JWG	C2B-C3B-CAB-CBB
4	ZZ	503	A1JWG	CAD-CBD-CGD-O2D
3	YT	502	LMG	C4-C5-C6-O5
3	YT	502	LMG	O6-C5-C6-O5
3	YT	502	LMG	C7-C8-C9-O8
3	YV	502	LMG	C7-C8-C9-O8
3	ZF	502	LMG	C7-C8-C9-O8
4	ZL	503	A1JWG	C4B-C3B-CAB-CBB
4	ZZ	503	A1JWG	C4B-C3B-CAB-CBB
2	ZD	501	NDP	C2D-C1D-N1N-C6N
4	YQ	503	A1JWG	CHA-CBD-CGD-O1D
4	YT	503	A1JWG	CHA-CBD-CGD-O1D
4	YY	503	A1JWG	CHA-CBD-CGD-O2D
4	YZ	503	A1JWG	CHA-CBD-CGD-O1D
4	ZB	503	A1JWG	CHA-CBD-CGD-O1D
4	ZC	503	A1JWG	CHA-CBD-CGD-O1D
4	ZD	503	A1JWG	CHA-CBD-CGD-O2D
4	ZF	503	A1JWG	CHA-CBD-CGD-O1D
4	ZG	503	A1JWG	CHA-CBD-CGD-O2D
4	ZI	503	A1JWG	CHA-CBD-CGD-O1D
4	ZK	503	A1JWG	CHA-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
4	ZM	503	A1JWG	CHA-CBD-CGD-O1D
4	ZO	503	A1JWG	CHA-CBD-CGD-O2D
4	ZS	503	A1JWG	CHA-CBD-CGD-O2D
4	ZU	503	A1JWG	CHA-CBD-CGD-O1D
2	YW	501	NDP	O4D-C1D-N1N-C6N
3	YJ	502	LMG	O1-C7-C8-O7
3	YN	502	LMG	O1-C7-C8-O7
3	ZF	502	LMG	O7-C8-C9-O8
3	ZT	502	LMG	O1-C7-C8-O7
3	ZT	502	LMG	O7-C8-C9-O8
3	ZY	502	LMG	O1-C7-C8-O7
2	YJ	501	NDP	C5B-O5B-PA-O3
2	YL	501	NDP	C5D-O5D-PN-O3
2	YN	501	NDP	C5B-O5B-PA-O3
2	YQ	501	NDP	C5B-O5B-PA-O3
2	YR	501	NDP	C5D-O5D-PN-O3
2	YS	501	NDP	C5B-O5B-PA-O3
2	YS	501	NDP	C5D-O5D-PN-O3
2	YT	501	NDP	C5D-O5D-PN-O3
2	YU	501	NDP	C5B-O5B-PA-O3
2	YW	501	NDP	C2B-O2B-P2B-O2X
2	YW	501	NDP	C2B-O2B-P2B-O3X
2	YW	501	NDP	C5D-O5D-PN-O3
2	YX	501	NDP	C5D-O5D-PN-O3
2	YZ	501	NDP	C5B-O5B-PA-O3
2	ZC	501	NDP	C5B-O5B-PA-O3
2	ZE	501	NDP	C5D-O5D-PN-O3
2	ZF	501	NDP	C2B-O2B-P2B-O3X
2	ZF	501	NDP	C5D-O5D-PN-O3
2	ZG	501	NDP	C5B-O5B-PA-O3
2	ZJ	501	NDP	C5B-O5B-PA-O3
2	ZK	501	NDP	C5B-O5B-PA-O3
2	ZM	501	NDP	C5B-O5B-PA-O3
2	ZM	501	NDP	C5D-O5D-PN-O3
2	ZN	501	NDP	C5D-O5D-PN-O3
2	ZO	501	NDP	C5B-O5B-PA-O3
2	ZQ	501	NDP	C5B-O5B-PA-O3
2	ZQ	501	NDP	C5D-O5D-PN-O3
2	ZR	501	NDP	C2B-O2B-P2B-O2X
2	ZR	501	NDP	C5D-O5D-PN-O3
2	ZW	501	NDP	C5B-O5B-PA-O3
2	ZX	501	NDP	C5B-O5B-PA-O3

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Mol	Chain	Res	Type	Atoms
2	ZY	501	NDP	C5B-O5B-PA-O3
2	ZY	501	NDP	C5D-O5D-PN-O3
2	ZZ	501	NDP	C5D-O5D-PN-O3
2	YI	501	NDP	O4B-C4B-C5B-O5B
2	ZS	501	NDP	C3D-C4D-C5D-O5D
2	ZX	501	NDP	C3D-C4D-C5D-O5D
2	YK	501	NDP	PA-O3-PN-O2N
2	YV	501	NDP	PA-O3-PN-O2N
2	ZA	501	NDP	PA-O3-PN-O2N
2	ZU	501	NDP	PA-O3-PN-O2N
3	ZN	502	LMG	C4-C5-C6-O5
4	YO	503	A1JWG	C1A-C2A-CAA-CBA
2	YR	501	NDP	O4D-C1D-N1N-C6N
2	YZ	501	NDP	O4D-C1D-N1N-C6N
2	ZF	501	NDP	O4D-C1D-N1N-C6N
2	ZR	501	NDP	O4D-C1D-N1N-C6N
2	ZY	501	NDP	O4D-C1D-N1N-C6N
2	ZT	501	NDP	C2D-C1D-N1N-C6N
2	YK	501	NDP	C5D-O5D-PN-O1N
2	YL	501	NDP	C5D-O5D-PN-O1N
2	YN	501	NDP	C5B-O5B-PA-O2A
2	YO	501	NDP	C5D-O5D-PN-O1N
2	YQ	501	NDP	C5D-O5D-PN-O1N
2	YS	501	NDP	C5D-O5D-PN-O2N
2	YU	501	NDP	C5B-O5B-PA-O2A
2	YU	501	NDP	C5D-O5D-PN-O1N
2	YV	501	NDP	C5D-O5D-PN-O1N
2	YW	501	NDP	C5D-O5D-PN-O2N
2	YX	501	NDP	C5D-O5D-PN-O1N
2	YZ	501	NDP	C5B-O5B-PA-O2A
2	YZ	501	NDP	C5D-O5D-PN-O1N
2	ZA	501	NDP	C5D-O5D-PN-O1N
2	ZC	501	NDP	C5B-O5B-PA-O2A
2	ZC	501	NDP	C5D-O5D-PN-O1N
2	ZE	501	NDP	C5D-O5D-PN-O1N
2	ZF	501	NDP	C5D-O5D-PN-O1N
2	ZG	501	NDP	C5B-O5B-PA-O2A
2	ZG	501	NDP	C5D-O5D-PN-O1N
2	ZH	501	NDP	C5B-O5B-PA-O2A
2	ZI	501	NDP	C5D-O5D-PN-O1N
2	ZJ	501	NDP	C5D-O5D-PN-O1N
2	ZK	501	NDP	C5D-O5D-PN-O1N

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Mol	Chain	Res	Type	Atoms
2	ZL	501	NDP	C5D-O5D-PN-O1N
2	ZM	501	NDP	C5D-O5D-PN-O1N
2	ZN	501	NDP	C5D-O5D-PN-O1N
2	ZO	501	NDP	C5B-O5B-PA-O2A
2	ZO	501	NDP	C5D-O5D-PN-O1N
2	ZQ	501	NDP	C5B-O5B-PA-O2A
2	ZQ	501	NDP	C5D-O5D-PN-O1N
2	ZR	501	NDP	C5D-O5D-PN-O1N
2	ZS	501	NDP	C5D-O5D-PN-O1N
2	ZU	501	NDP	C5D-O5D-PN-O1N
2	ZV	501	NDP	C5D-O5D-PN-O1N
2	ZW	501	NDP	C5B-O5B-PA-O2A
2	ZW	501	NDP	C5D-O5D-PN-O1N
2	ZW	501	NDP	C5D-O5D-PN-O2N
2	ZX	501	NDP	C5B-O5B-PA-O2A
2	ZX	501	NDP	C5D-O5D-PN-O1N
2	ZY	501	NDP	C5D-O5D-PN-O2N
4	ZW	503	A1JWG	C2A-CAA-CBA-CGA
4	YJ	503	A1JWG	CAD-CBD-CGD-O1D
4	YM	503	A1JWG	CAD-CBD-CGD-O1D
3	ZV	502	LMG	C4-C5-C6-O5
3	ZD	502	LMG	C4-C5-C6-O5
2	ZE	501	NDP	C3B-C4B-C5B-O5B
2	ZW	501	NDP	O4D-C4D-C5D-O5D
2	YS	501	NDP	O4D-C1D-N1N-C6N
3	ZW	502	LMG	O6-C5-C6-O5
2	YT	501	NDP	C2D-C1D-N1N-C6N
2	ZH	501	NDP	C2D-C1D-N1N-C2N
3	ZZ	502	LMG	C4-C5-C6-O5
2	YN	501	NDP	C2N-C3N-C7N-O7N
2	ZH	501	NDP	C2N-C3N-C7N-O7N
3	YN	502	LMG	O1-C7-C8-C9
3	YO	502	LMG	O1-C7-C8-C9
3	ZY	502	LMG	O1-C7-C8-C9
3	YO	502	LMG	O1-C7-C8-O7
3	YT	502	LMG	O7-C8-C9-O8
3	YV	502	LMG	O7-C8-C9-O8
3	ZT	502	LMG	C8-C7-O1-C1
3	ZE	502	LMG	O6-C5-C6-O5
3	ZN	502	LMG	O6-C5-C6-O5
4	YZ	503	A1JWG	O1D-CGD-O2D-CED
3	ZJ	502	LMG	C4-C5-C6-O5

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Mol	Chain	Res	Type	Atoms
2	ZH	501	NDP	O4D-C1D-N1N-C6N
3	ZV	502	LMG	O9-C10-O7-C8
2	YJ	501	NDP	C2D-C1D-N1N-C6N
3	ZV	502	LMG	O6-C5-C6-O5
3	ZD	502	LMG	O10-C28-O8-C9
4	ZL	503	A1JWG	C2A-CAA-CBA-CGA
4	ZX	503	A1JWG	C2A-CAA-CBA-CGA
2	YJ	501	NDP	O4D-C1D-N1N-C6N
2	YT	501	NDP	O4D-C1D-N1N-C6N
2	ZT	501	NDP	O4D-C1D-N1N-C6N
2	ZW	501	NDP	C3D-C4D-C5D-O5D
2	ZD	501	NDP	O4D-C1D-N1N-C6N
2	ZZ	501	NDP	O4D-C1D-N1N-C6N
2	YK	501	NDP	PA-O3-PN-O1N
2	YO	501	NDP	PA-O3-PN-O2N
2	ZA	501	NDP	PA-O3-PN-O1N
2	ZI	501	NDP	PA-O3-PN-O2N
2	ZL	501	NDP	PA-O3-PN-O1N
2	ZL	501	NDP	PA-O3-PN-O2N
2	ZU	501	NDP	PA-O3-PN-O1N
2	ZV	501	NDP	PA-O3-PN-O2N
3	ZZ	502	LMG	O6-C5-C6-O5
3	YX	502	LMG	C4-C5-C6-O5
4	ZM	503	A1JWG	CBD-CGD-O2D-CED
4	YX	503	A1JWG	C2B-C3B-CAB-CBB
3	YJ	502	LMG	O10-C28-O8-C9
2	ZP	501	NDP	C2D-C1D-N1N-C6N
2	ZZ	501	NDP	C2D-C1D-N1N-C6N
4	ZM	503	A1JWG	C2A-CAA-CBA-CGA
2	YJ	501	NDP	O4D-C4D-C5D-O5D
2	YM	501	NDP	C3B-C4B-C5B-O5B
2	YX	501	NDP	O4B-C4B-C5B-O5B
2	ZF	501	NDP	O4B-C4B-C5B-O5B
2	ZJ	501	NDP	C3D-C4D-C5D-O5D
2	ZR	501	NDP	O4B-C4B-C5B-O5B
4	ZV	503	A1JWG	CAA-CBA-CGA-O1A
3	YP	502	LMG	C4-C5-C6-O5
3	YZ	502	LMG	O6-C1-O1-C7
2	ZP	501	NDP	O4D-C1D-N1N-C6N
4	ZN	503	A1JWG	C4B-C3B-CAB-CBB
4	ZW	503	A1JWG	C4B-C3B-CAB-CBB
4	ZP	503	A1JWG	CBD-CGD-O2D-CED

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Mol	Chain	Res	Type	Atoms
4	YV	503	A1JWG	CAA-CBA-CGA-O1A
3	ZJ	502	LMG	O6-C5-C6-O5
4	ZI	503	A1JWG	C2A-CAA-CBA-CGA
4	YO	503	A1JWG	C3A-C2A-CAA-CBA
4	YO	503	A1JWG	CAA-CBA-CGA-O2A
4	ZE	503	A1JWG	CAA-CBA-CGA-O1A
4	ZE	503	A1JWG	CAA-CBA-CGA-O2A
4	YI	503	A1JWG	CAA-CBA-CGA-O2A
4	YX	503	A1JWG	CAA-CBA-CGA-O1A
4	ZA	503	A1JWG	CAA-CBA-CGA-O1A
4	ZR	503	A1JWG	CBD-CGD-O2D-CED
2	YP	501	NDP	C3D-C4D-C5D-O5D
2	ZD	501	NDP	O4B-C4B-C5B-O5B
2	ZP	501	NDP	O4B-C4B-C5B-O5B
4	YK	503	A1JWG	CAA-CBA-CGA-O1A
4	YQ	503	A1JWG	CAA-CBA-CGA-O1A
4	ZJ	503	A1JWG	CAA-CBA-CGA-O1A
4	YN	503	A1JWG	CAA-CBA-CGA-O2A
4	YO	503	A1JWG	CAA-CBA-CGA-O1A
4	ZC	503	A1JWG	CAA-CBA-CGA-O1A
4	ZA	503	A1JWG	C1A-C2A-CAA-CBA
4	YP	503	A1JWG	CAA-CBA-CGA-O1A
4	YY	503	A1JWG	CAA-CBA-CGA-O2A
4	ZK	503	A1JWG	CAA-CBA-CGA-O2A
4	ZV	503	A1JWG	CAA-CBA-CGA-O2A
2	YJ	501	NDP	C2D-C1D-N1N-C2N
2	ZD	501	NDP	C2D-C1D-N1N-C2N
2	ZT	501	NDP	C2D-C1D-N1N-C2N
4	YK	503	A1JWG	CAA-CBA-CGA-O2A
4	ZM	503	A1JWG	CAA-CBA-CGA-O2A
4	ZU	503	A1JWG	CAA-CBA-CGA-O1A
4	YL	503	A1JWG	CAA-CBA-CGA-O2A
4	ZA	503	A1JWG	CAA-CBA-CGA-O2A
4	YI	503	A1JWG	CAA-CBA-CGA-O1A
4	YR	503	A1JWG	CAA-CBA-CGA-O2A
4	YW	503	A1JWG	CAA-CBA-CGA-O2A
4	ZR	503	A1JWG	CAA-CBA-CGA-O2A
4	ZY	503	A1JWG	CAA-CBA-CGA-O2A
3	YX	502	LMG	O6-C5-C6-O5
3	ZQ	502	LMG	C4-C5-C6-O5
2	ZN	501	NDP	O4B-C4B-C5B-O5B
2	ZT	501	NDP	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
4	YJ	503	A1JWG	CAA-CBA-CGA-O2A
4	YT	503	A1JWG	CAA-CBA-CGA-O2A
4	ZB	503	A1JWG	CAA-CBA-CGA-O1A
4	ZI	503	A1JWG	CAA-CBA-CGA-O2A
4	YS	503	A1JWG	CAA-CBA-CGA-O2A
4	ZO	503	A1JWG	CAA-CBA-CGA-O2A
4	ZU	503	A1JWG	CAA-CBA-CGA-O2A
4	ZQ	503	A1JWG	CAA-CBA-CGA-O2A
4	ZF	503	A1JWG	CAA-CBA-CGA-O2A
4	ZG	503	A1JWG	CAA-CBA-CGA-O2A
4	ZJ	503	A1JWG	CAA-CBA-CGA-O2A
3	ZU	502	LMG	C4-C5-C6-O5
2	YO	501	NDP	PA-O3-PN-O1N
2	YW	501	NDP	PA-O3-PN-O2N
2	ZS	501	NDP	PA-O3-PN-O2N
3	YP	502	LMG	O6-C5-C6-O5
4	YP	503	A1JWG	CAA-CBA-CGA-O2A
4	YV	503	A1JWG	CAA-CBA-CGA-O2A
4	ZN	503	A1JWG	CAA-CBA-CGA-O2A
4	ZP	503	A1JWG	CAA-CBA-CGA-O2A
4	ZT	503	A1JWG	CAA-CBA-CGA-O2A
4	ZZ	503	A1JWG	CAA-CBA-CGA-O2A
3	ZU	502	LMG	O6-C5-C6-O5
3	YJ	502	LMG	O1-C7-C8-C9
4	YX	503	A1JWG	CAA-CBA-CGA-O2A
2	ZQ	501	NDP	O4D-C4D-C5D-O5D
4	ZD	503	A1JWG	CAA-CBA-CGA-O2A
4	ZB	503	A1JWG	CAA-CBA-CGA-O2A
4	ZS	503	A1JWG	CAA-CBA-CGA-O2A
4	YM	503	A1JWG	CAA-CBA-CGA-O1A
4	YM	503	A1JWG	CAA-CBA-CGA-O2A
4	ZC	503	A1JWG	CAA-CBA-CGA-O2A
4	ZN	503	A1JWG	CAA-CBA-CGA-O1A
3	YW	502	LMG	O1-C7-C8-O7
4	ZF	503	A1JWG	CAA-CBA-CGA-O1A
2	YI	501	NDP	C3B-C4B-C5B-O5B
2	YU	501	NDP	C3D-C4D-C5D-O5D
2	ZV	501	NDP	C3D-C4D-C5D-O5D
4	YQ	503	A1JWG	CAA-CBA-CGA-O2A
4	YT	503	A1JWG	CAA-CBA-CGA-O1A
4	YU	503	A1JWG	CAA-CBA-CGA-O2A
2	YT	501	NDP	C2D-C1D-N1N-C2N

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Mol	Chain	Res	Type	Atoms
3	YJ	502	LMG	C4-C5-C6-O5
4	ZH	503	A1JWG	CAA-CBA-CGA-O2A
4	YK	503	A1JWG	CAD-CBD-CGD-O2D
4	YL	503	A1JWG	CAD-CBD-CGD-O2D
4	YN	503	A1JWG	CAD-CBD-CGD-O2D
4	YO	503	A1JWG	CAD-CBD-CGD-O2D
4	ZA	503	A1JWG	CAD-CBD-CGD-O2D
4	ZE	503	A1JWG	CAD-CBD-CGD-O2D
4	ZP	503	A1JWG	CAD-CBD-CGD-O2D
4	ZV	503	A1JWG	CAD-CBD-CGD-O2D
4	ZY	503	A1JWG	CAD-CBD-CGD-O2D
2	ZH	501	NDP	O4D-C1D-N1N-C2N
4	ZL	503	A1JWG	CAA-CBA-CGA-O2A
4	ZP	503	A1JWG	CAA-CBA-CGA-O1A
4	ZX	503	A1JWG	CAA-CBA-CGA-O1A
3	YR	502	LMG	C7-C8-C9-O8
3	ZP	502	LMG	C7-C8-C9-O8
4	YL	503	A1JWG	CAA-CBA-CGA-O1A
4	ZR	503	A1JWG	CAA-CBA-CGA-O1A
4	ZY	503	A1JWG	CAA-CBA-CGA-O1A
4	YN	503	A1JWG	CAA-CBA-CGA-O1A
4	YS	503	A1JWG	CAA-CBA-CGA-O1A
4	YU	503	A1JWG	CAA-CBA-CGA-O1A
4	YZ	503	A1JWG	CAA-CBA-CGA-O1A
4	ZK	503	A1JWG	CAA-CBA-CGA-O1A
4	ZL	503	A1JWG	CAA-CBA-CGA-O1A
4	ZS	503	A1JWG	CAA-CBA-CGA-O1A
4	ZT	503	A1JWG	CAA-CBA-CGA-O1A
4	ZW	503	A1JWG	CAA-CBA-CGA-O1A
2	YP	501	NDP	O4D-C4D-C5D-O5D
2	YU	501	NDP	O4D-C4D-C5D-O5D
2	ZB	501	NDP	O4B-C4B-C5B-O5B
2	ZJ	501	NDP	O4D-C4D-C5D-O5D
2	ZZ	501	NDP	O4D-C4D-C5D-O5D
4	ZM	503	A1JWG	CAA-CBA-CGA-O1A
4	YS	503	A1JWG	CHA-CBD-CGD-O2D
4	YU	503	A1JWG	CHA-CBD-CGD-O2D
4	YW	503	A1JWG	CHA-CBD-CGD-O2D
4	ZB	503	A1JWG	CHA-CBD-CGD-O2D
4	ZC	503	A1JWG	CHA-CBD-CGD-O2D
4	ZF	503	A1JWG	CHA-CBD-CGD-O2D
4	ZL	503	A1JWG	CHA-CBD-CGD-O1D

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Mol	Chain	Res	Type	Atoms
4	ZN	503	A1JWG	CHA-CBD-CGD-O2D
4	ZQ	503	A1JWG	CHA-CBD-CGD-O2D
4	YY	503	A1JWG	CAA-CBA-CGA-O1A
4	ZG	503	A1JWG	CAA-CBA-CGA-O1A
4	ZO	503	A1JWG	CAA-CBA-CGA-O1A
4	ZZ	503	A1JWG	CAA-CBA-CGA-O1A
4	ZQ	503	A1JWG	CAA-CBA-CGA-O1A
3	YQ	502	LMG	C11-C10-O7-C8
3	ZR	502	LMG	O7-C8-C9-O8
3	ZM	502	LMG	C29-C28-O8-C9
4	ZD	503	A1JWG	C2A-CAA-CBA-CGA
4	ZD	503	A1JWG	CAA-CBA-CGA-O1A
2	YI	501	NDP	C5B-O5B-PA-O3
2	YI	501	NDP	C5D-O5D-PN-O3
2	YK	501	NDP	C2B-O2B-P2B-O3X
2	YM	501	NDP	C5D-O5D-PN-O3
2	YT	501	NDP	C5B-O5B-PA-O3
2	YV	501	NDP	C2B-O2B-P2B-O3X
2	YW	501	NDP	C5B-O5B-PA-O3
2	YY	501	NDP	C5B-O5B-PA-O3
2	YY	501	NDP	C2B-O2B-P2B-O2X
2	ZB	501	NDP	C5D-O5D-PN-O3
2	ZD	501	NDP	C5B-O5B-PA-O3
2	ZE	501	NDP	C5B-O5B-PA-O3
2	ZI	501	NDP	C2B-O2B-P2B-O2X
2	ZL	501	NDP	C5B-O5B-PA-O3
2	ZP	501	NDP	C5B-O5B-PA-O3
2	ZS	501	NDP	C5B-O5B-PA-O3
2	ZU	501	NDP	C2B-O2B-P2B-O3X
2	ZY	501	NDP	C2B-O2B-P2B-O3X
2	ZZ	501	NDP	C5B-O5B-PA-O3
4	YW	503	A1JWG	CAA-CBA-CGA-O1A
4	YR	503	A1JWG	CAA-CBA-CGA-O1A
2	ZP	501	NDP	C2D-C1D-N1N-C2N
3	ZY	502	LMG	O6-C5-C6-O5
2	YJ	501	NDP	O4B-C4B-C5B-O5B
2	YL	501	NDP	O4B-C4B-C5B-O5B
2	YT	501	NDP	O4D-C4D-C5D-O5D
2	ZT	501	NDP	O4D-C4D-C5D-O5D
2	YP	501	NDP	PA-O3-PN-O2N
2	YQ	501	NDP	PA-O3-PN-O1N
2	YU	501	NDP	PA-O3-PN-O1N

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Mol	Chain	Res	Type	Atoms
2	ZJ	501	NDP	PA-O3-PN-O2N
2	ZS	501	NDP	PA-O3-PN-O1N
2	ZY	501	NDP	PA-O3-PN-O2N
2	ZZ	501	NDP	PA-O3-PN-O2N
4	YJ	503	A1JWG	CAA-CBA-CGA-O1A
4	YK	503	A1JWG	C1A-C2A-CAA-CBA
4	ZE	503	A1JWG	C1A-C2A-CAA-CBA
3	YW	502	LMG	O1-C7-C8-C9
4	ZI	503	A1JWG	CAA-CBA-CGA-O1A
3	YZ	502	LMG	C2-C1-O1-C7
2	YJ	501	NDP	C5B-O5B-PA-O2A
2	YQ	501	NDP	C5B-O5B-PA-O2A
2	YV	501	NDP	C5B-O5B-PA-O1A
2	ZK	501	NDP	C5B-O5B-PA-O2A
2	ZM	501	NDP	C5B-O5B-PA-O2A
2	YN	501	NDP	O4B-C4B-C5B-O5B
2	YR	501	NDP	O4B-C4B-C5B-O5B
2	ZD	501	NDP	O4D-C4D-C5D-O5D
2	ZM	501	NDP	O4D-C4D-C5D-O5D
2	ZP	501	NDP	O4D-C4D-C5D-O5D
4	ZX	503	A1JWG	O1D-CGD-O2D-CED
3	ZZ	502	LMG	C7-C8-O7-C10
4	YN	503	A1JWG	CAD-CBD-CGD-O1D
4	ZL	503	A1JWG	CAD-CBD-CGD-O1D
4	ZP	503	A1JWG	CAD-CBD-CGD-O1D
4	ZT	503	A1JWG	CAD-CBD-CGD-O1D
4	ZA	503	A1JWG	C3A-C2A-CAA-CBA
4	ZW	503	A1JWG	CAA-CBA-CGA-O2A

There are no ring outliers.

70 monomers are involved in 65 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	YT	501	NDP	3	0
4	YK	503	A1JWG	1	0
2	YY	501	NDP	1	0
4	ZZ	503	A1JWG	1	0
4	YZ	503	A1JWG	1	0
2	ZU	501	NDP	1	0
4	ZR	503	A1JWG	1	0
2	YS	501	NDP	3	0
3	ZQ	502	LMG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	ZN	503	A1JWG	1	0
3	ZA	502	LMG	1	0
3	YS	502	LMG	1	0
4	YT	503	A1JWG	1	0
4	ZL	503	A1JWG	1	0
2	YL	501	NDP	1	0
2	ZX	501	NDP	2	0
3	ZY	502	LMG	1	0
4	YV	503	A1JWG	1	0
2	ZZ	501	NDP	3	0
2	ZK	501	NDP	1	0
3	ZX	502	LMG	1	0
4	ZF	503	A1JWG	1	0
4	ZB	503	A1JWG	1	0
4	ZX	503	A1JWG	1	0
4	ZJ	503	A1JWG	1	0
2	YM	501	NDP	2	0
2	YQ	501	NDP	1	0
3	YK	502	LMG	1	0
2	YK	501	NDP	1	0
2	ZH	501	NDP	2	0
3	ZM	502	LMG	1	0
2	ZY	501	NDP	2	0
4	YS	503	A1JWG	1	0
4	ZU	503	A1JWG	1	0
2	ZB	501	NDP	1	0
4	ZI	503	A1JWG	1	0
3	ZE	502	LMG	1	0
2	YR	501	NDP	1	0
4	ZC	503	A1JWG	1	0
2	YN	501	NDP	2	0
2	ZI	501	NDP	1	0
2	ZL	501	NDP	4	0
2	YP	501	NDP	1	0
4	YX	503	A1JWG	1	0
4	ZH	503	A1JWG	1	0
2	ZS	501	NDP	1	0
2	ZW	501	NDP	1	0
4	YU	503	A1JWG	1	0
2	ZR	501	NDP	1	0
4	YP	503	A1JWG	1	0
2	ZG	501	NDP	1	0

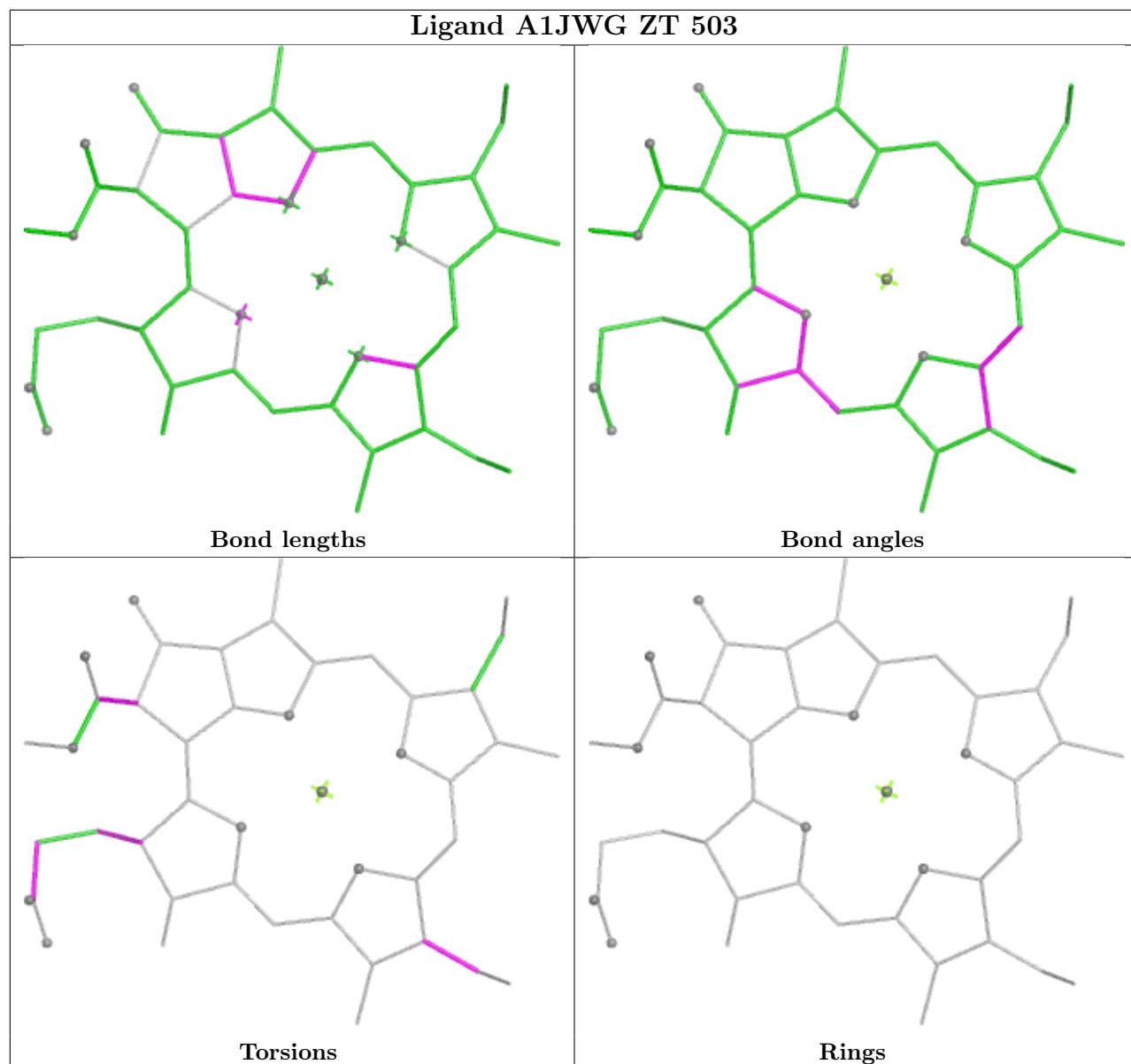
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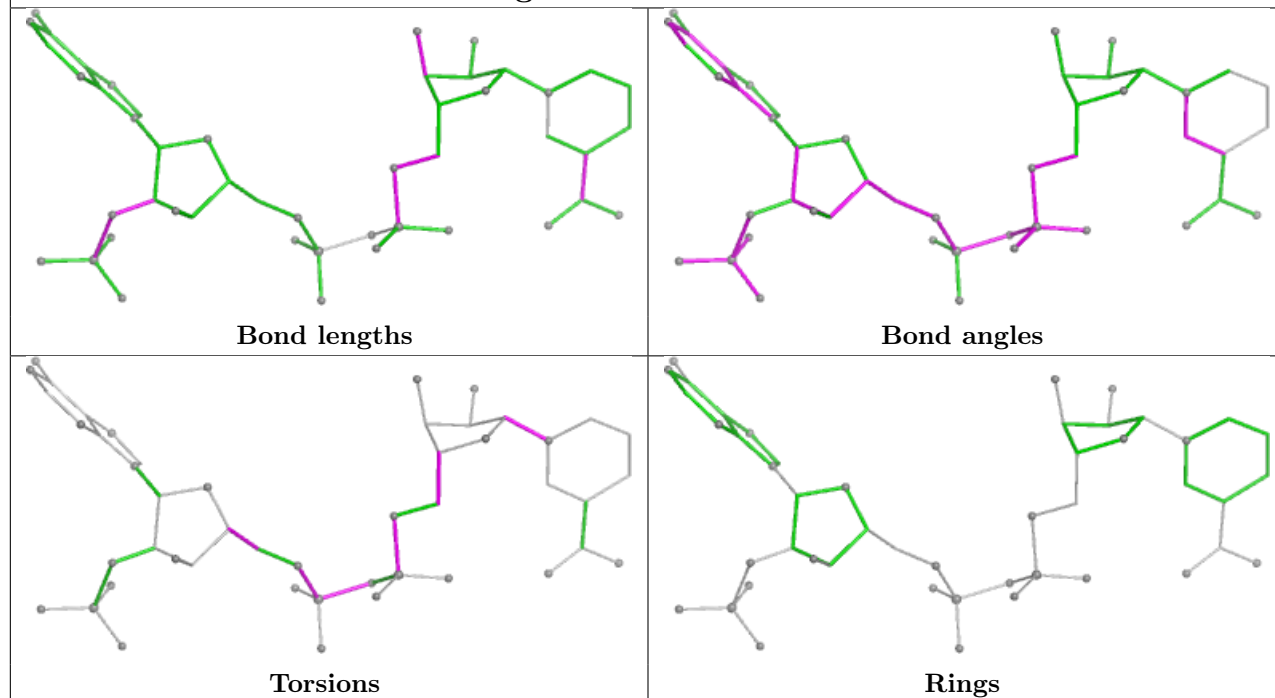
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	ZM	501	NDP	1	0
4	YL	503	A1JWG	1	0
2	YW	501	NDP	2	0
4	ZM	503	A1JWG	1	0
4	YM	503	A1JWG	1	0
2	YZ	501	NDP	2	0
3	YO	502	LMG	1	0
3	YW	502	LMG	1	0
2	ZF	501	NDP	1	0
2	ZJ	501	NDP	1	0
2	YV	501	NDP	1	0
2	YX	501	NDP	1	0
2	YU	501	NDP	2	0
2	ZC	501	NDP	2	0
2	ZN	501	NDP	1	0
3	ZI	502	LMG	1	0
2	ZT	501	NDP	1	0
2	ZD	501	NDP	1	0
2	YI	501	NDP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

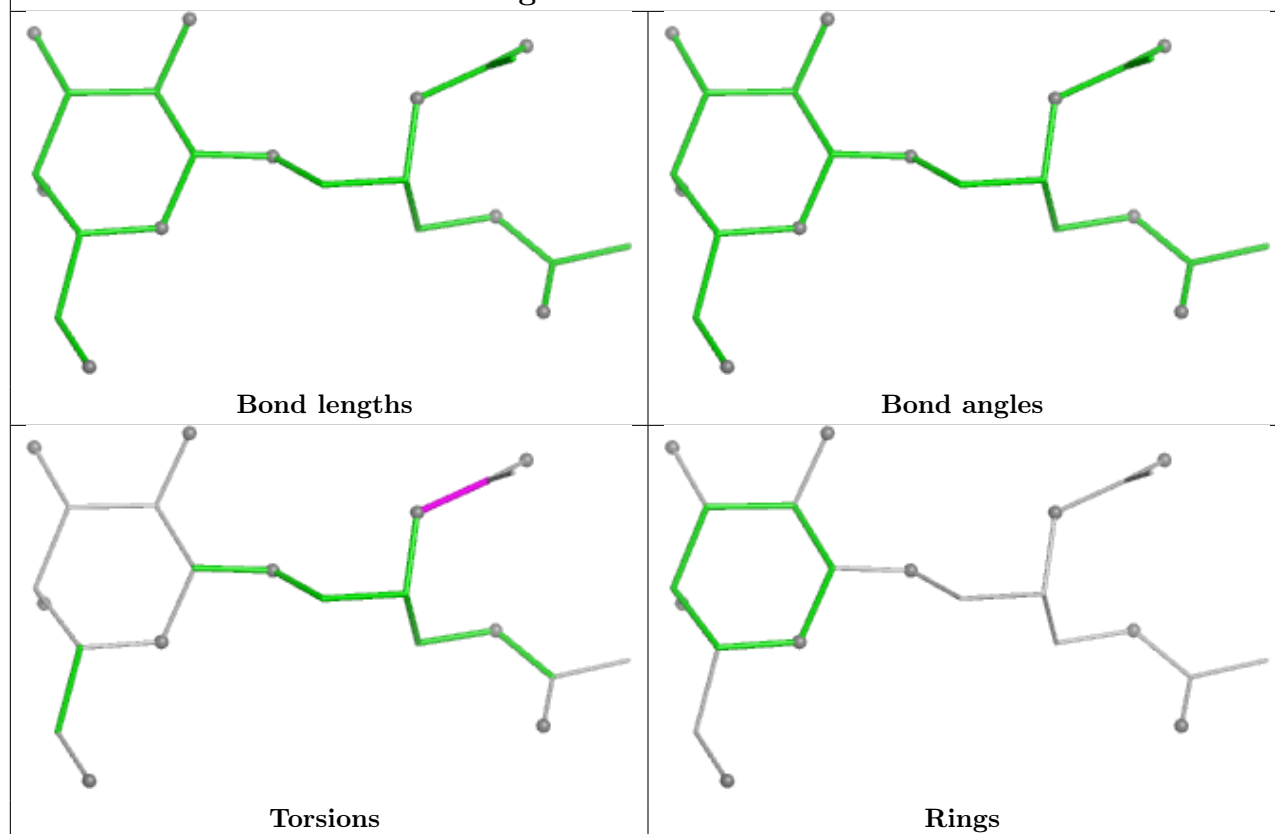
Ligand A1JWG ZT 503



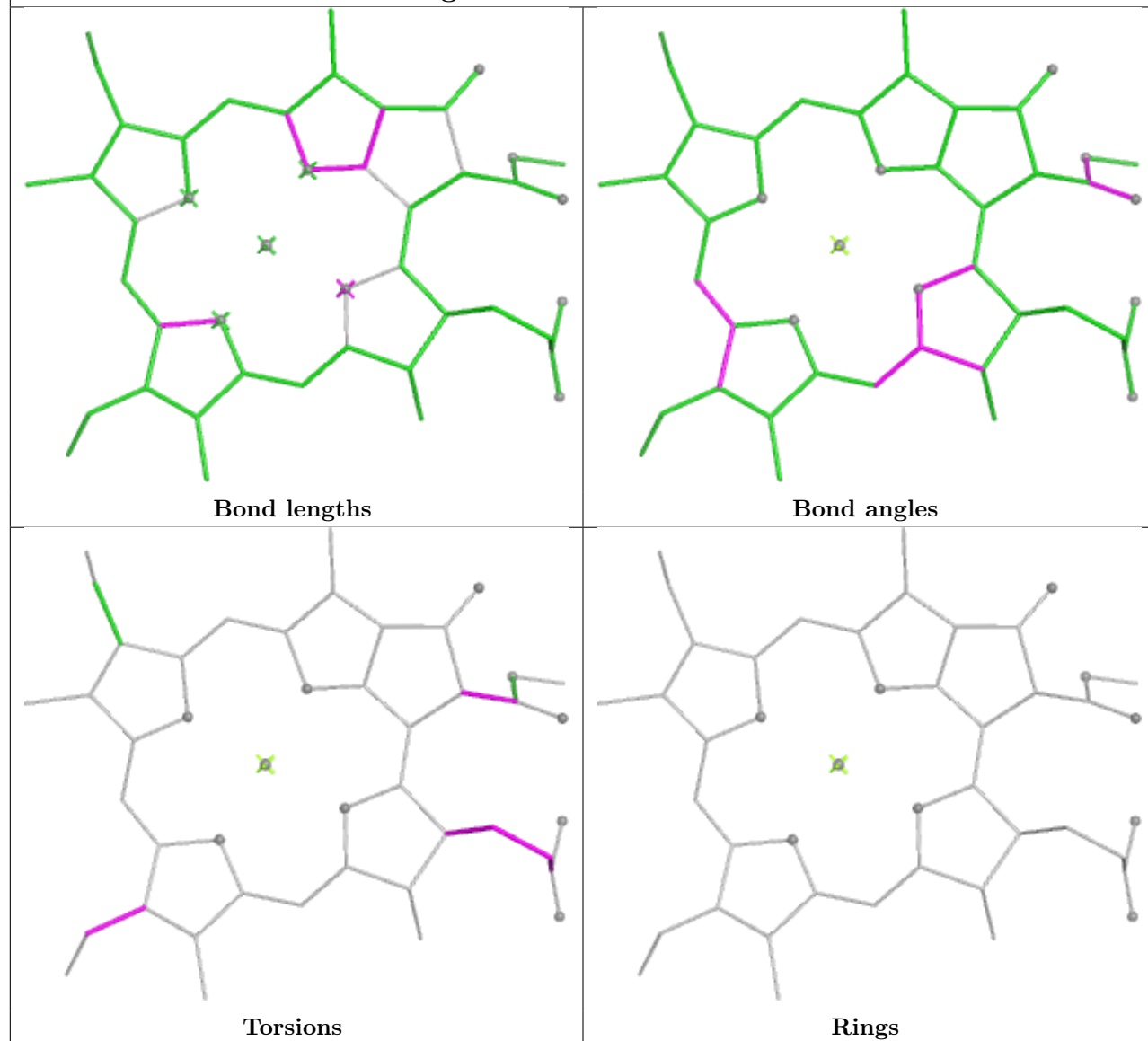
Ligand NDP YT 501

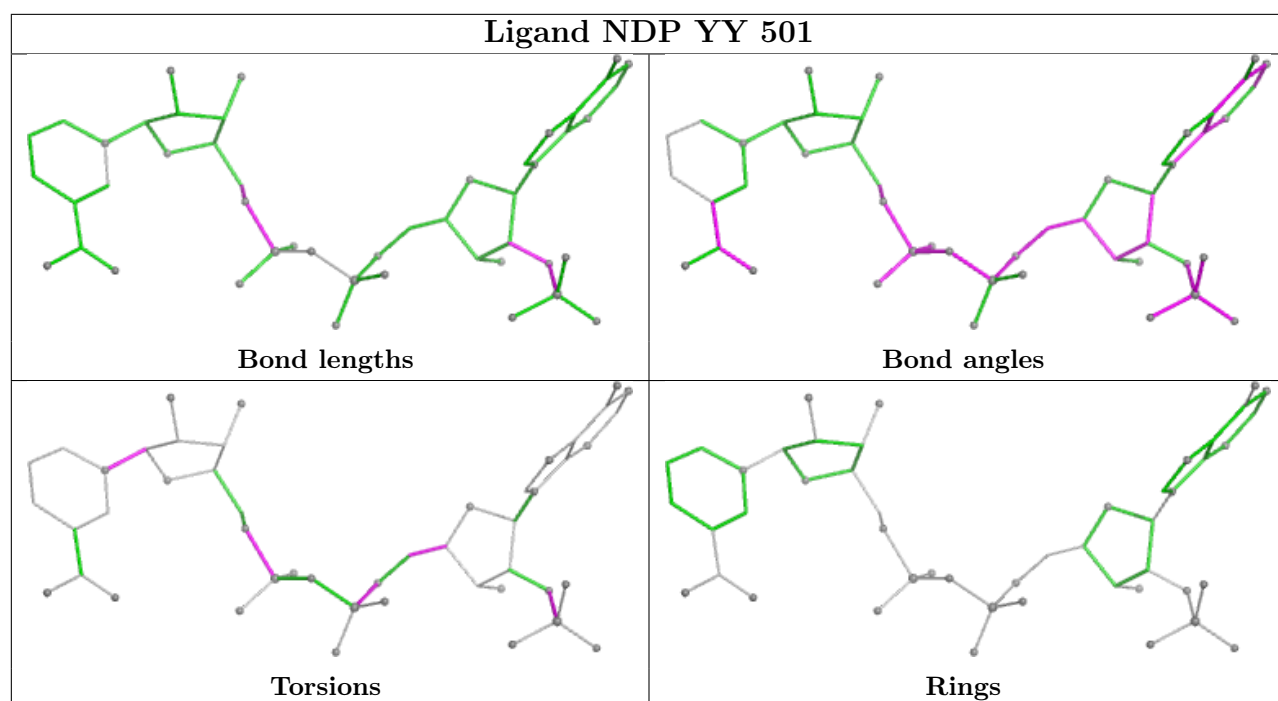


Ligand LMG ZS 502

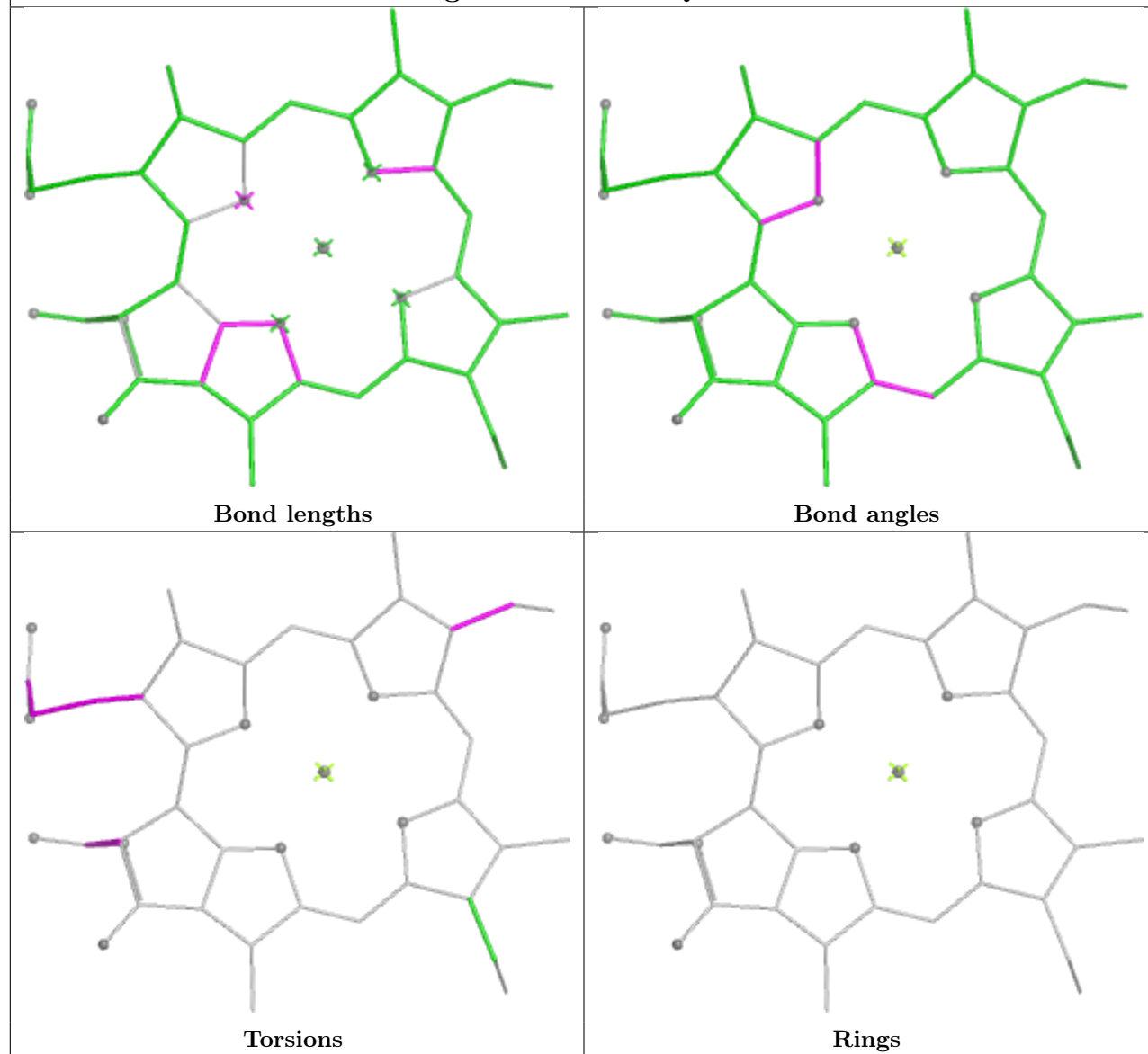


Ligand A1JWG YK 503

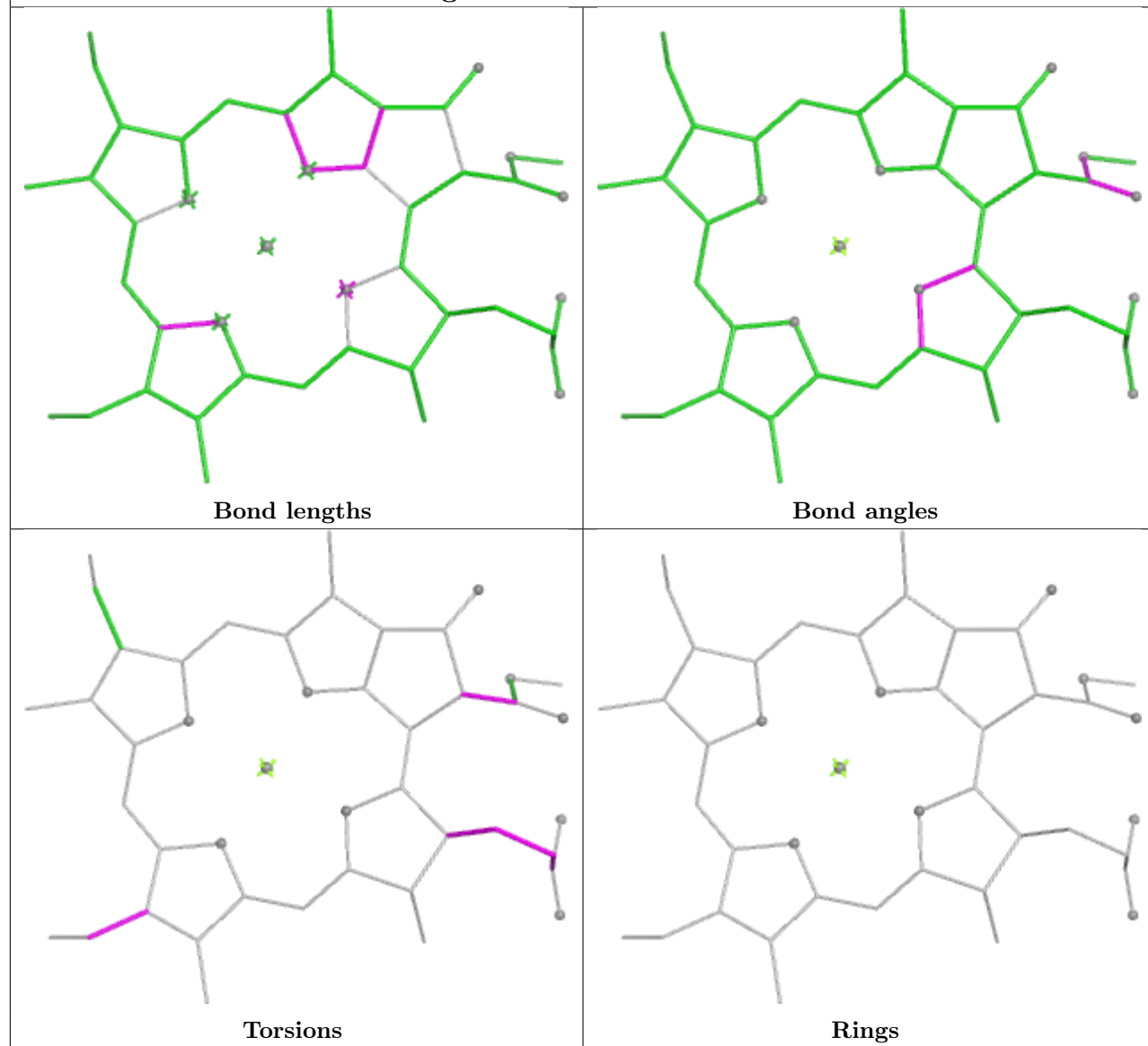




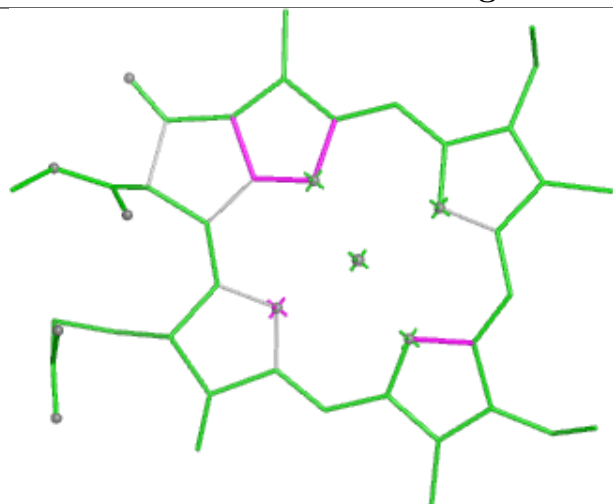
Ligand A1JWG ZQ 503



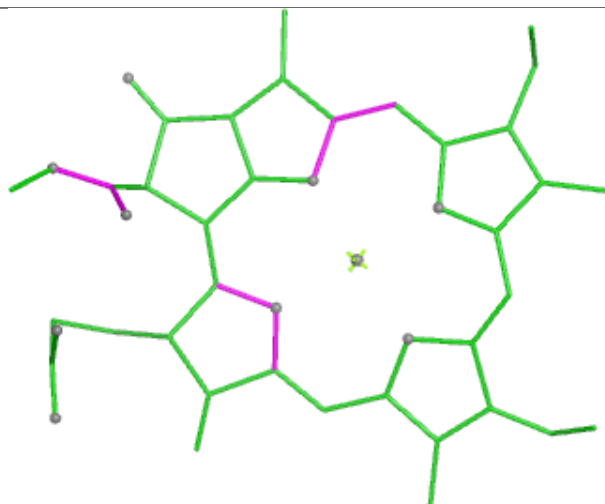
Ligand A1JWG YO 503



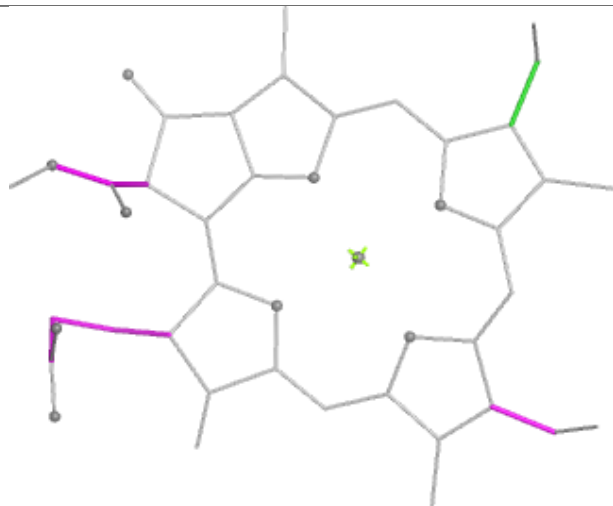
Ligand A1JWG ZZ 503



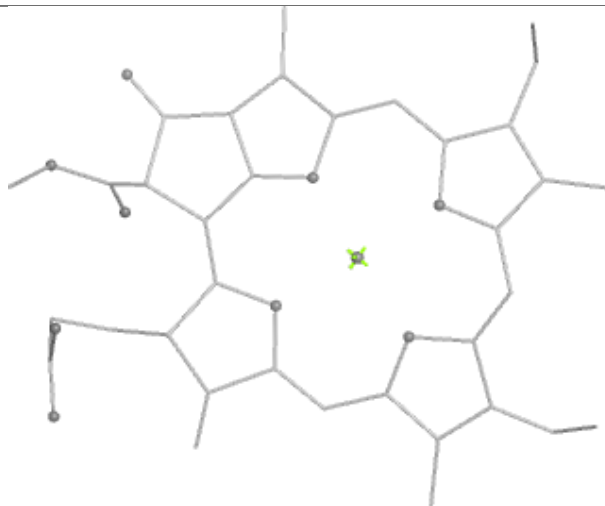
Bond lengths



Bond angles

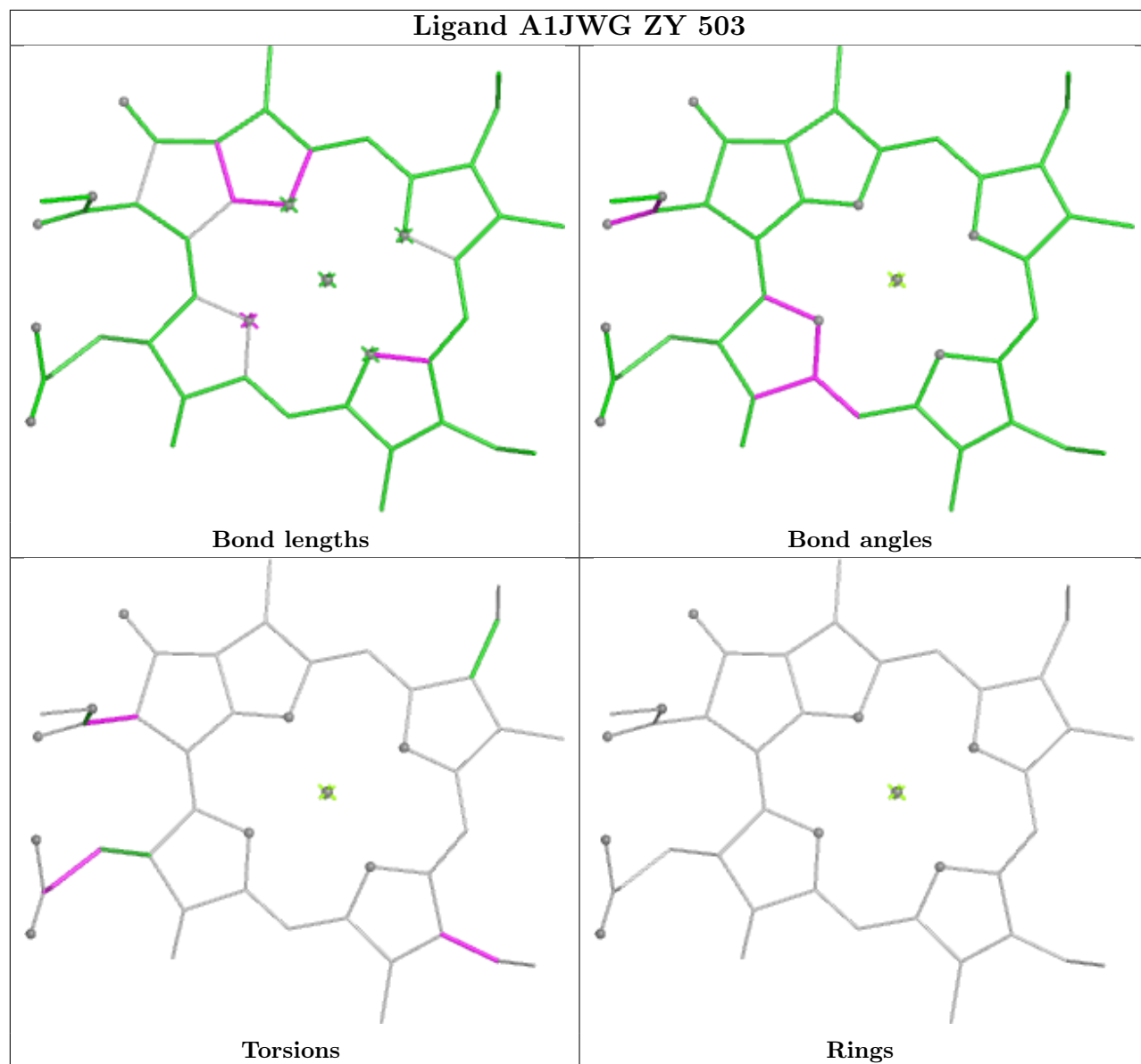


Torsions

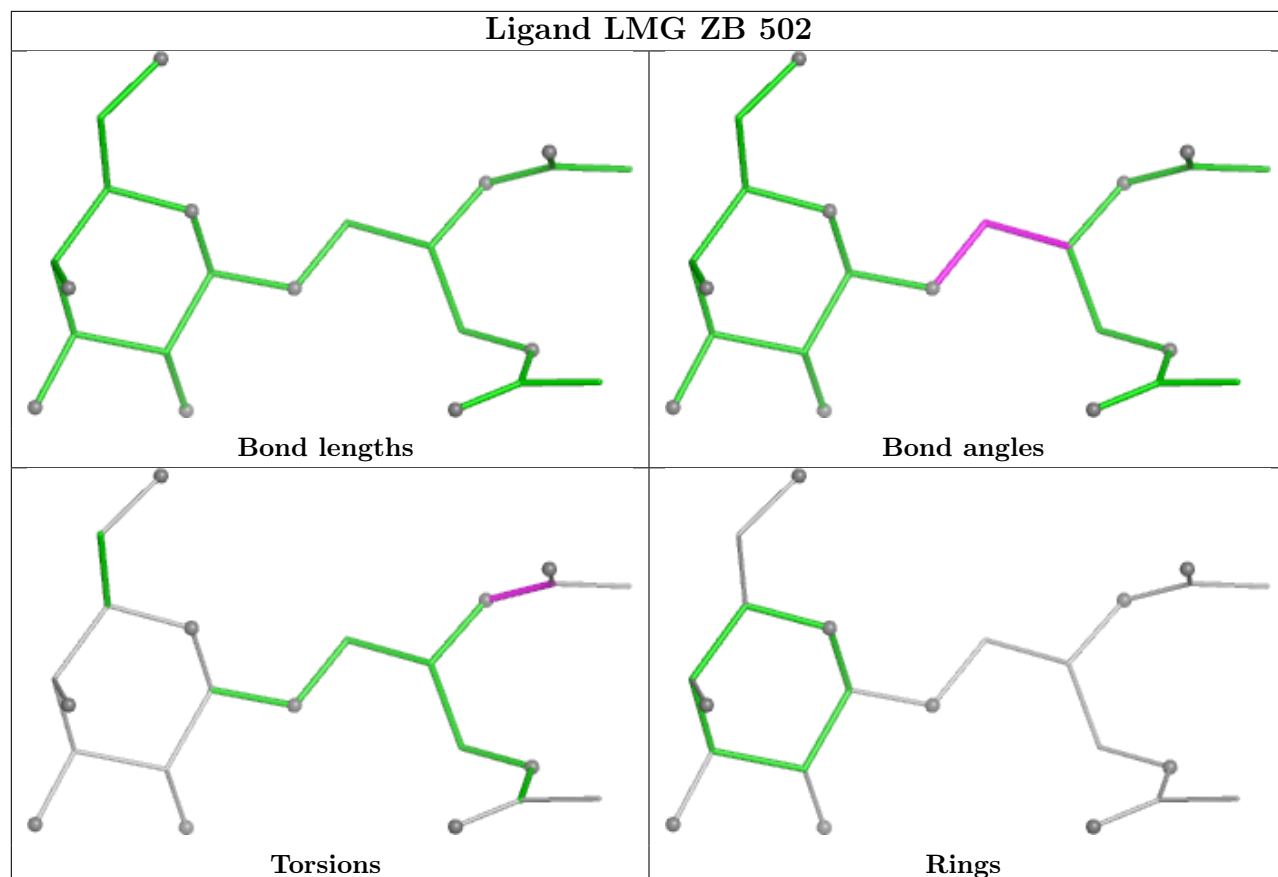


Rings

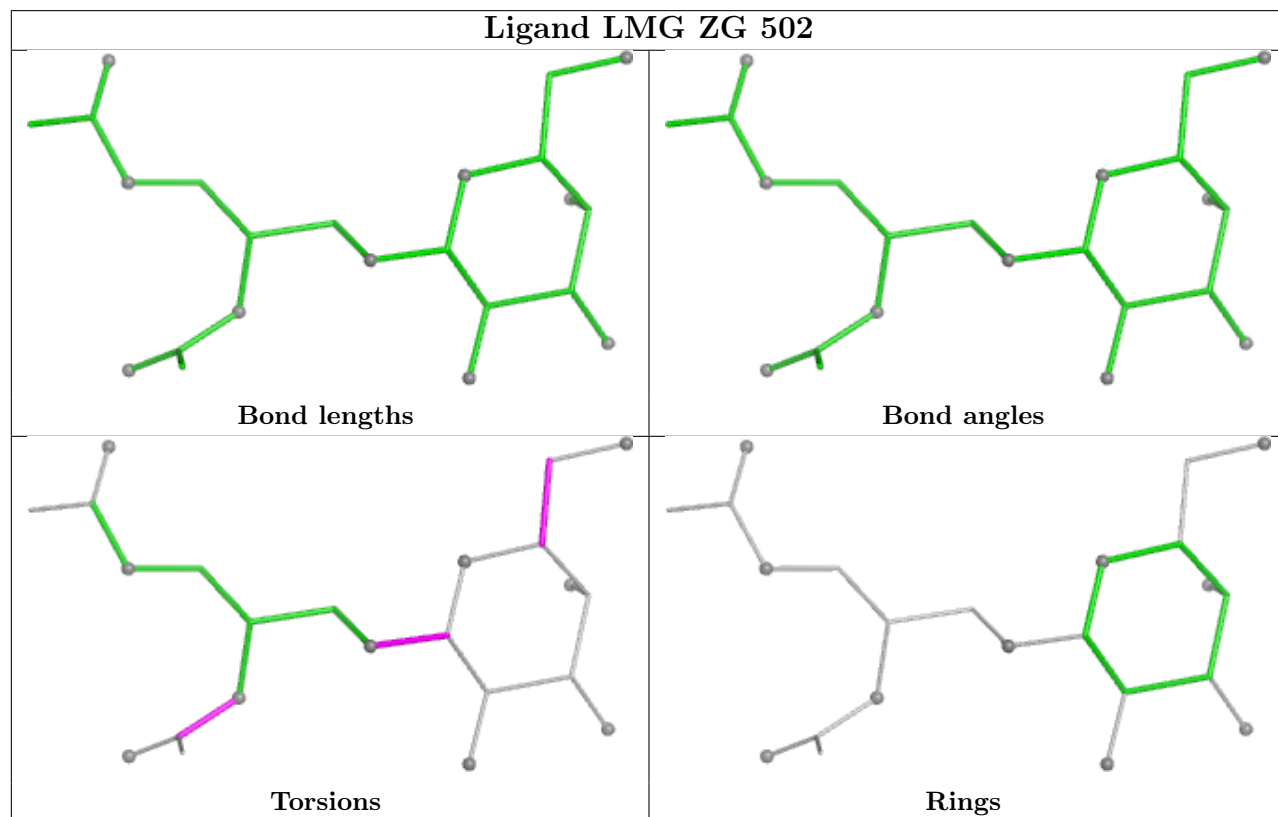
Ligand A1JWG ZY 503

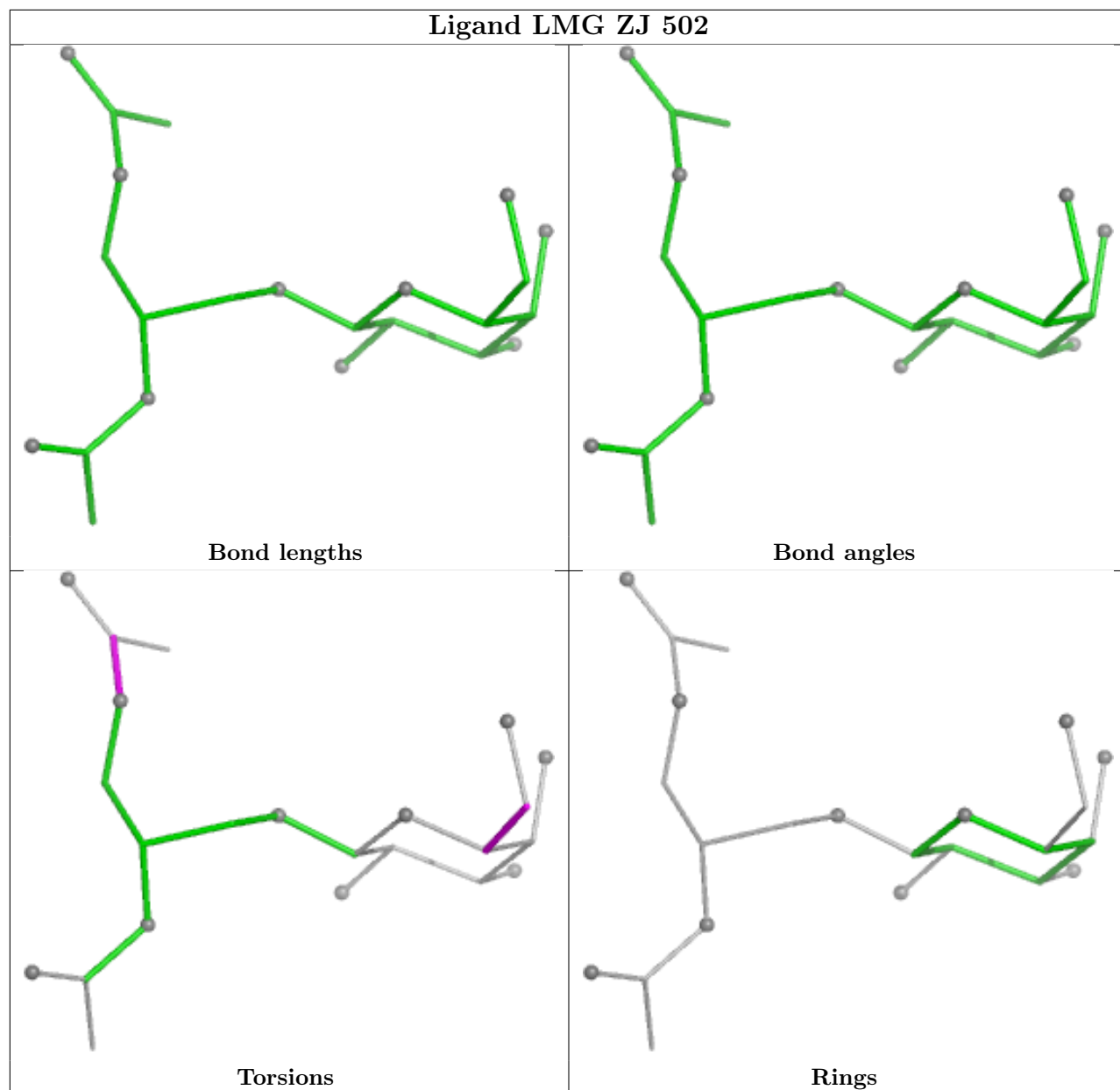


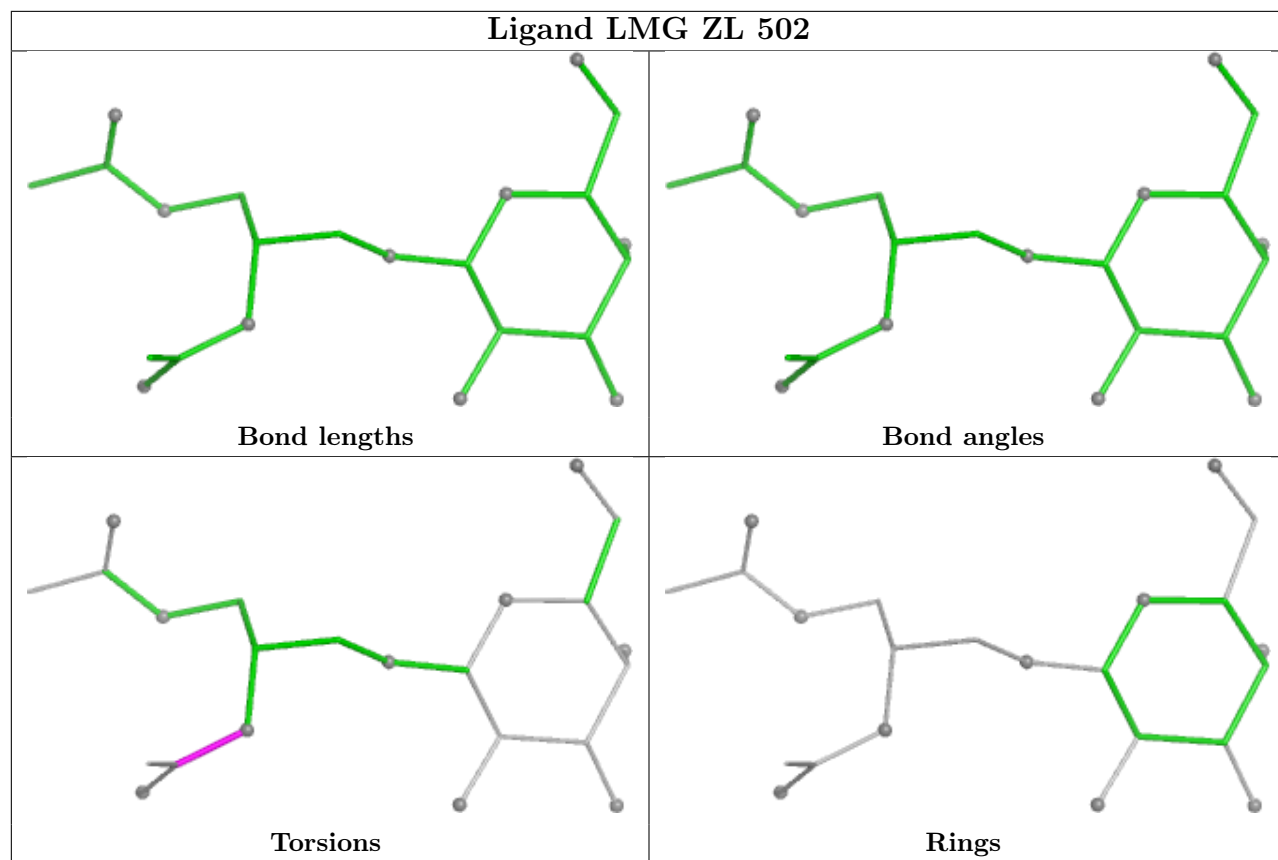
Ligand LMG ZB 502



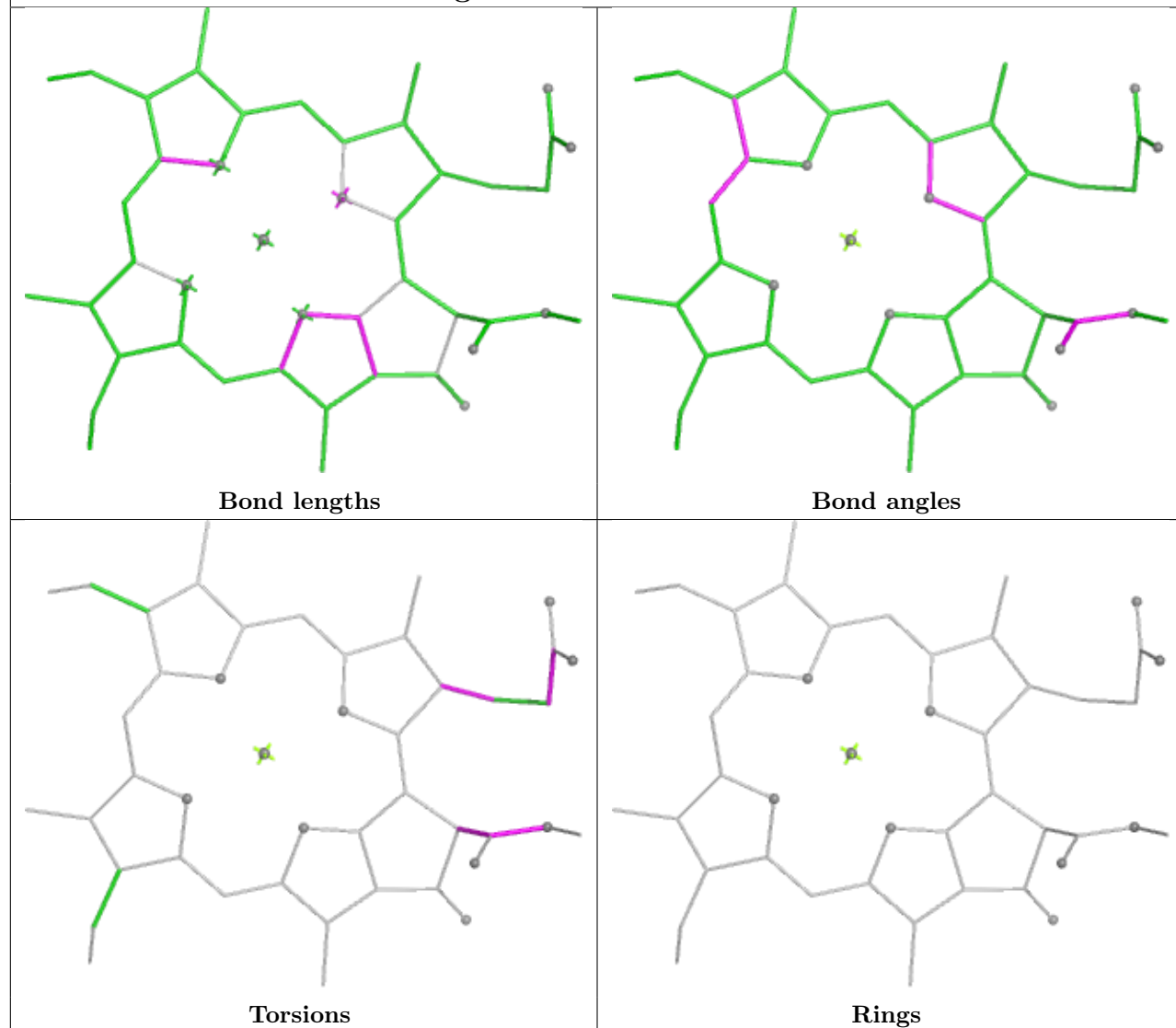
Ligand LMG ZG 502

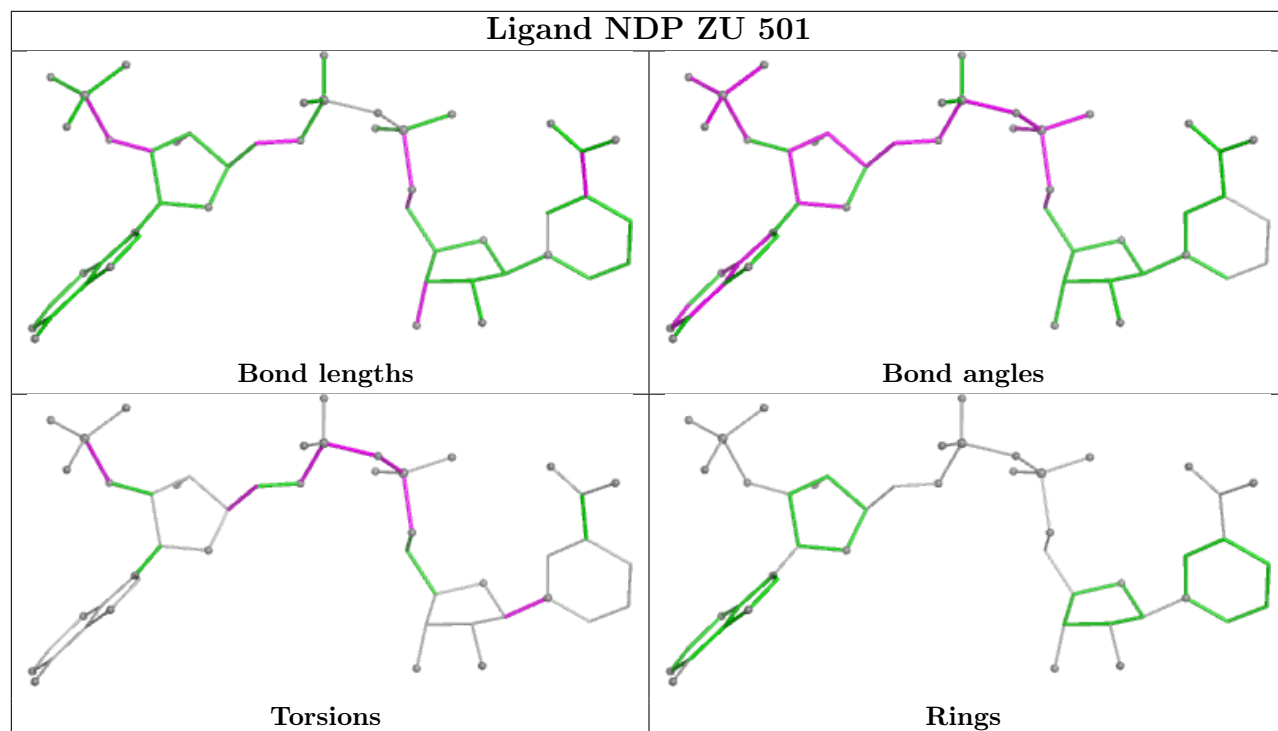


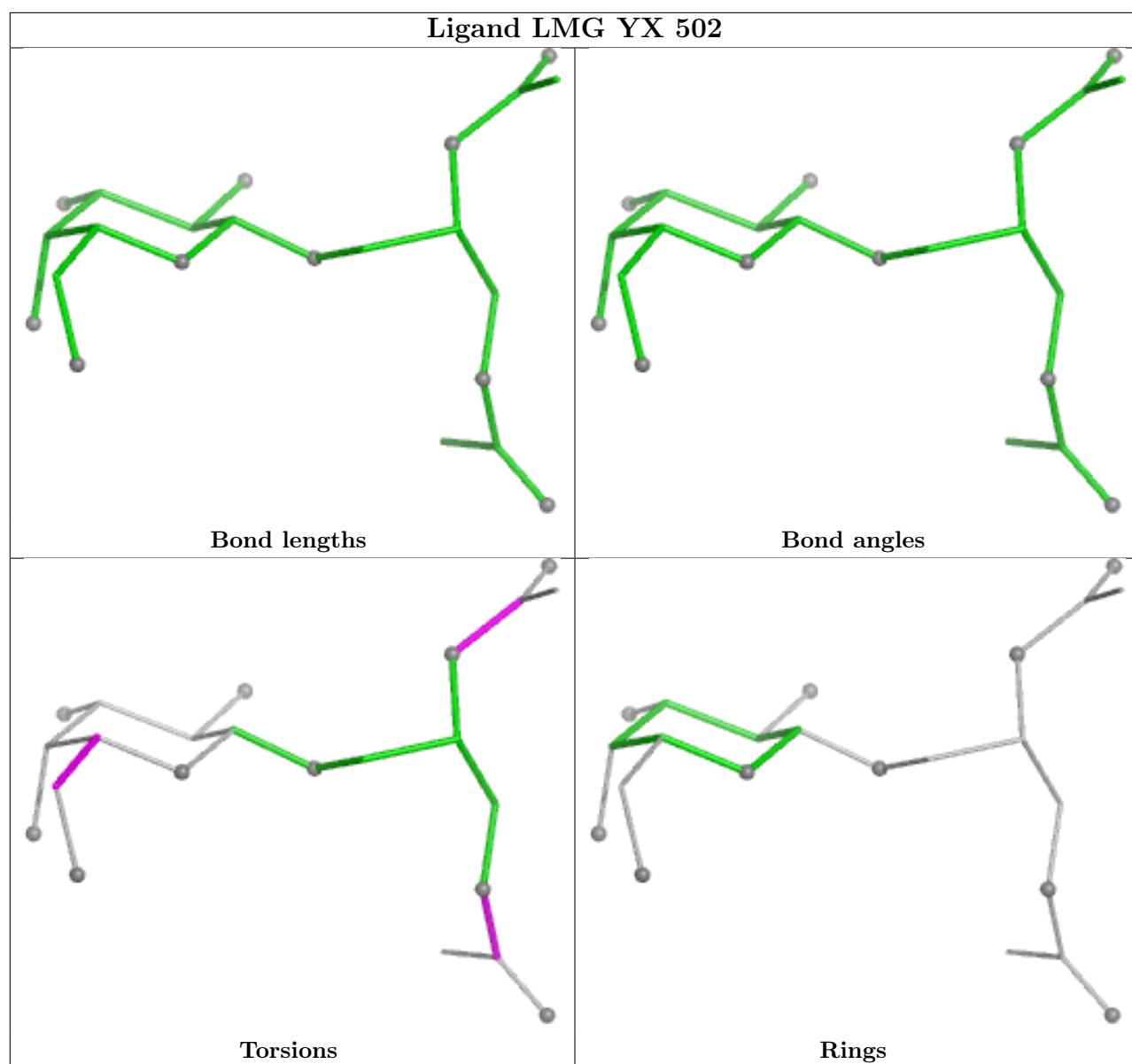




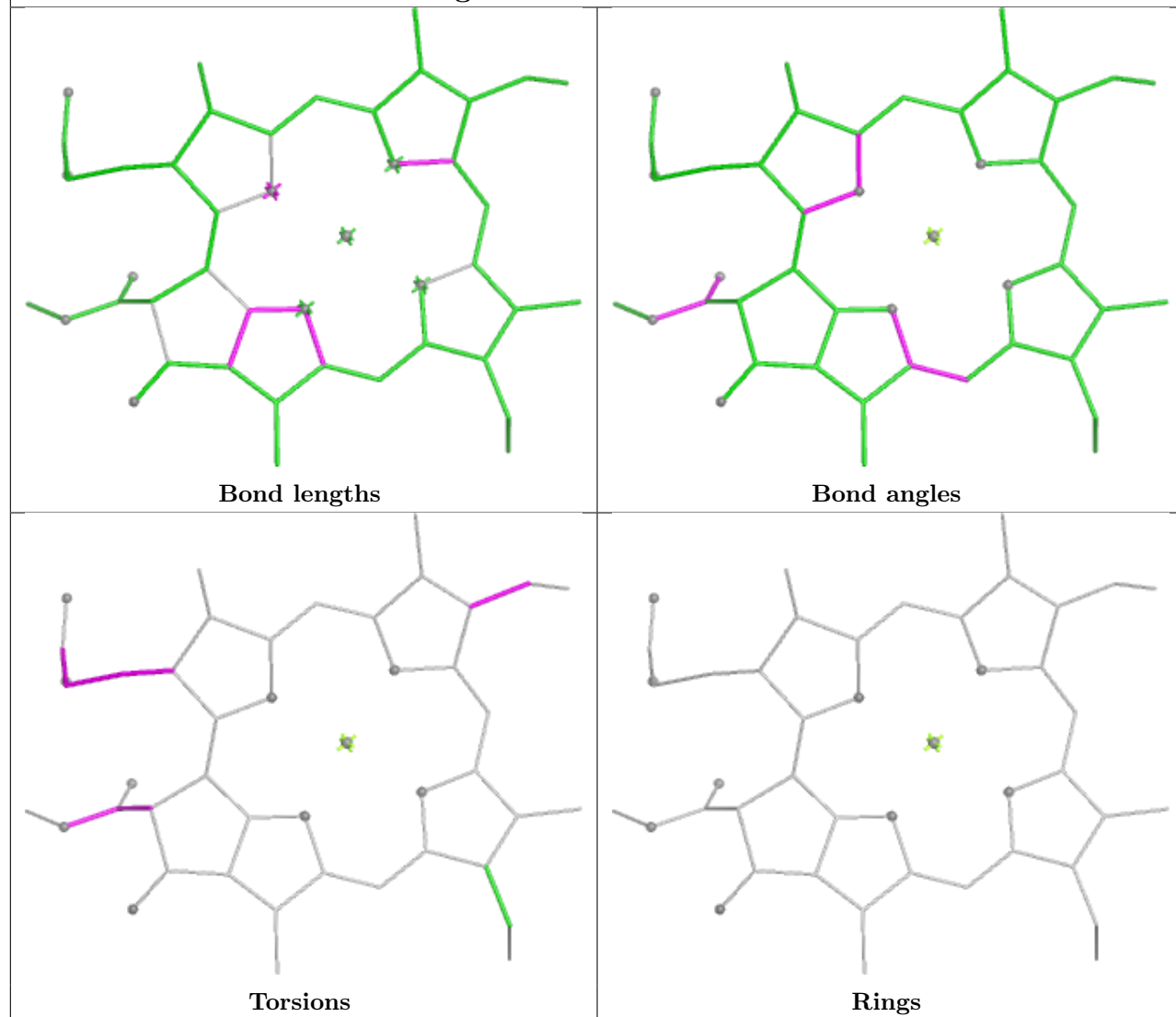
Ligand A1JWG YZ 503

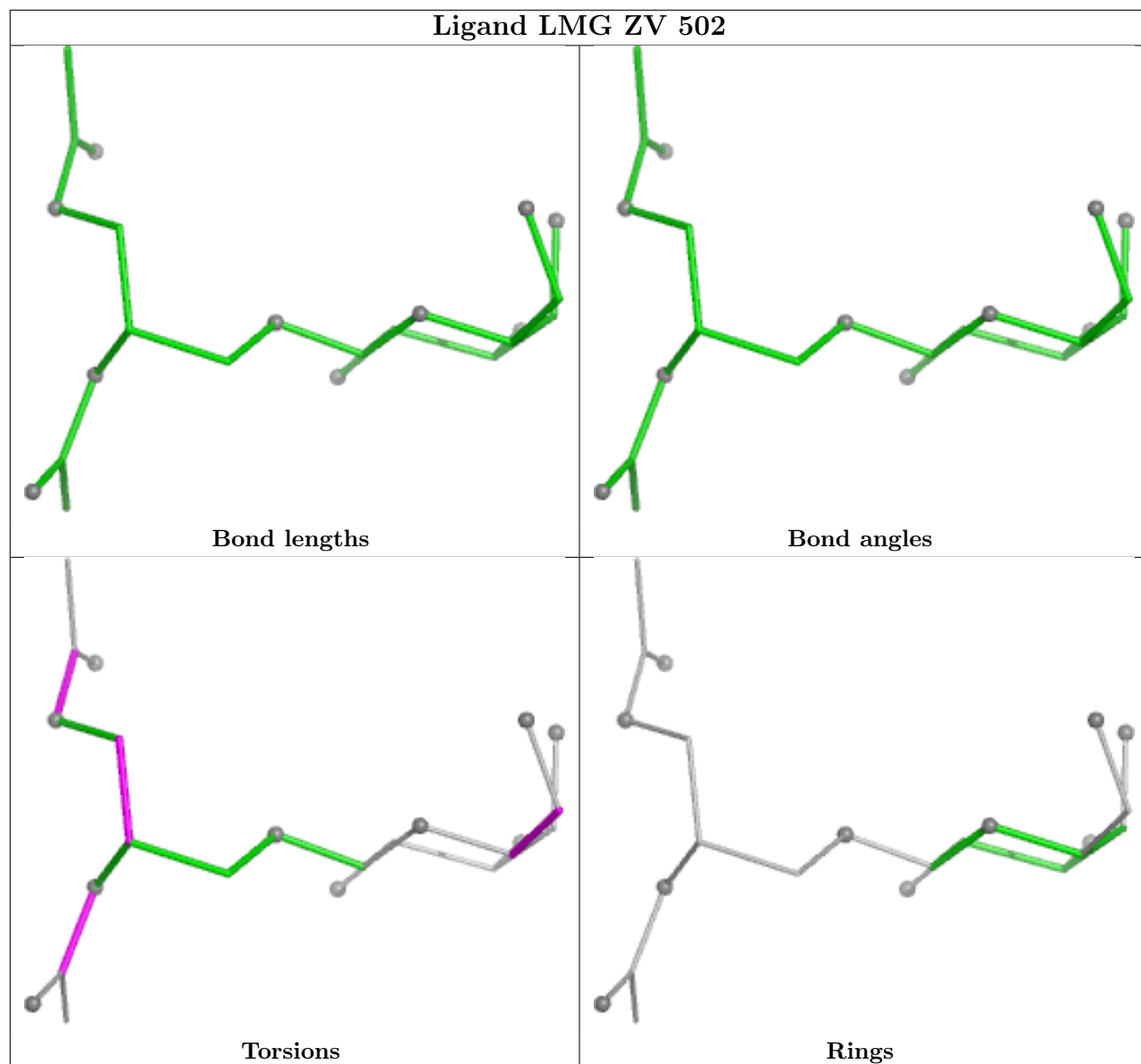




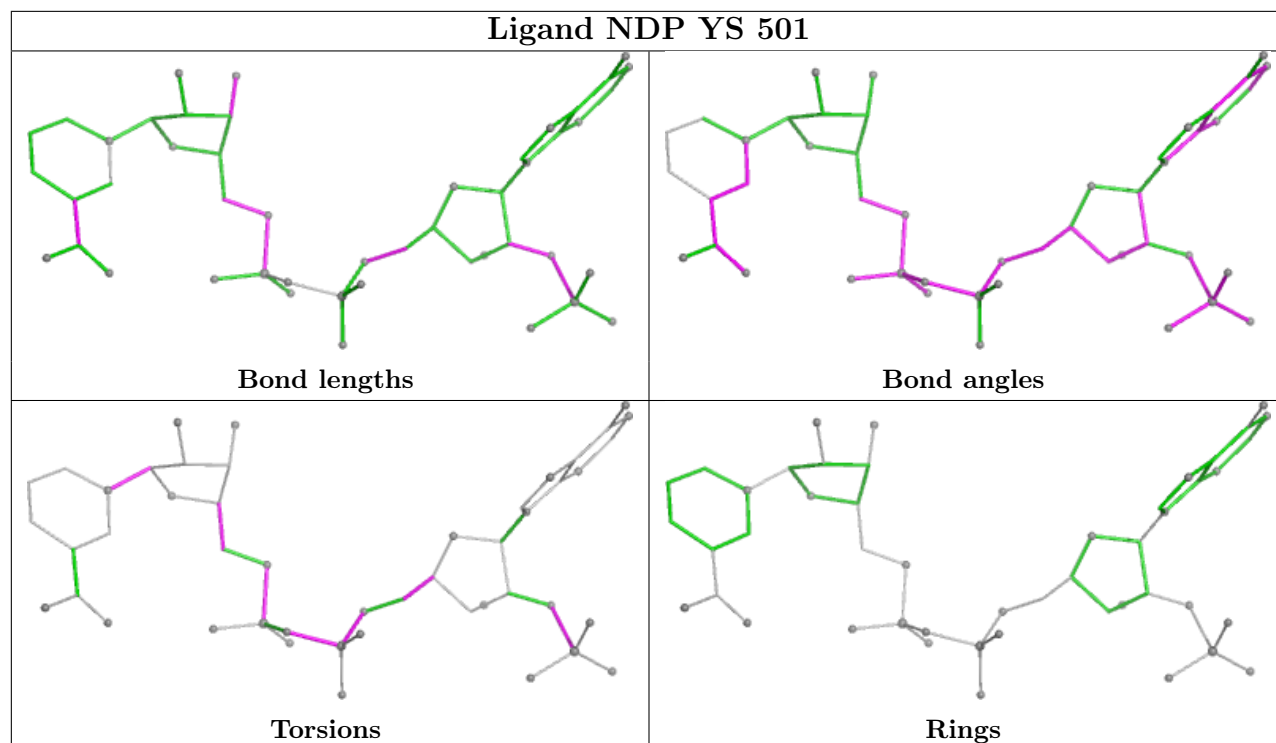


Ligand A1JWG ZR 503

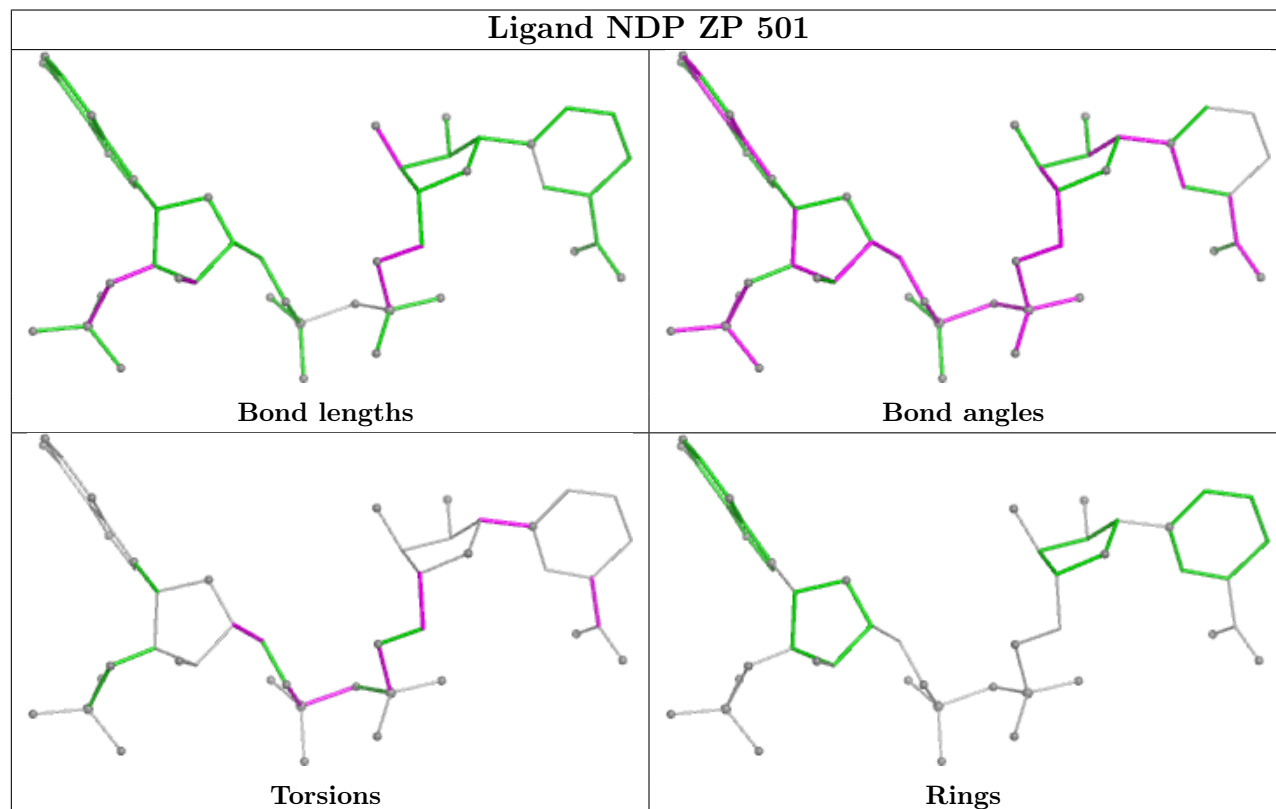




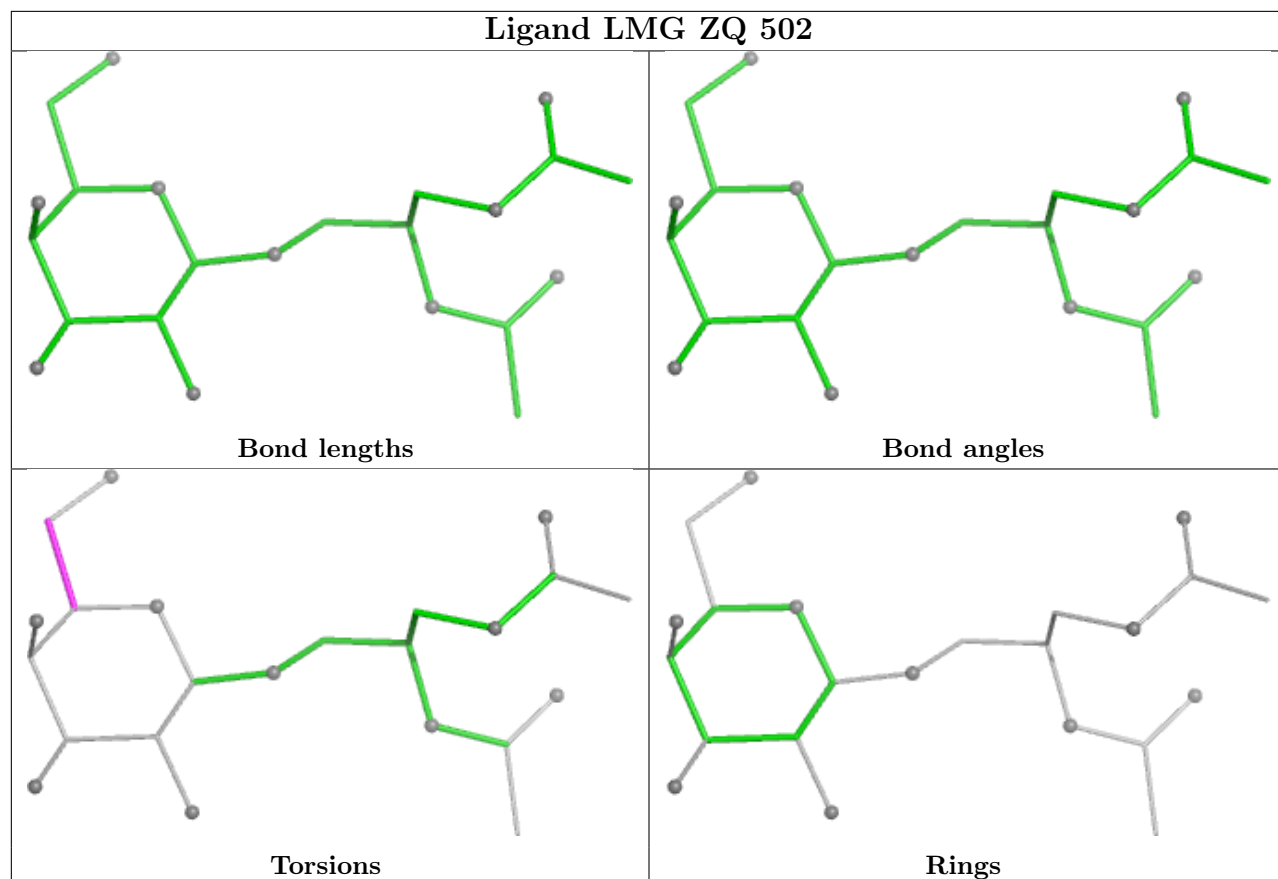
Ligand NDP YS 501



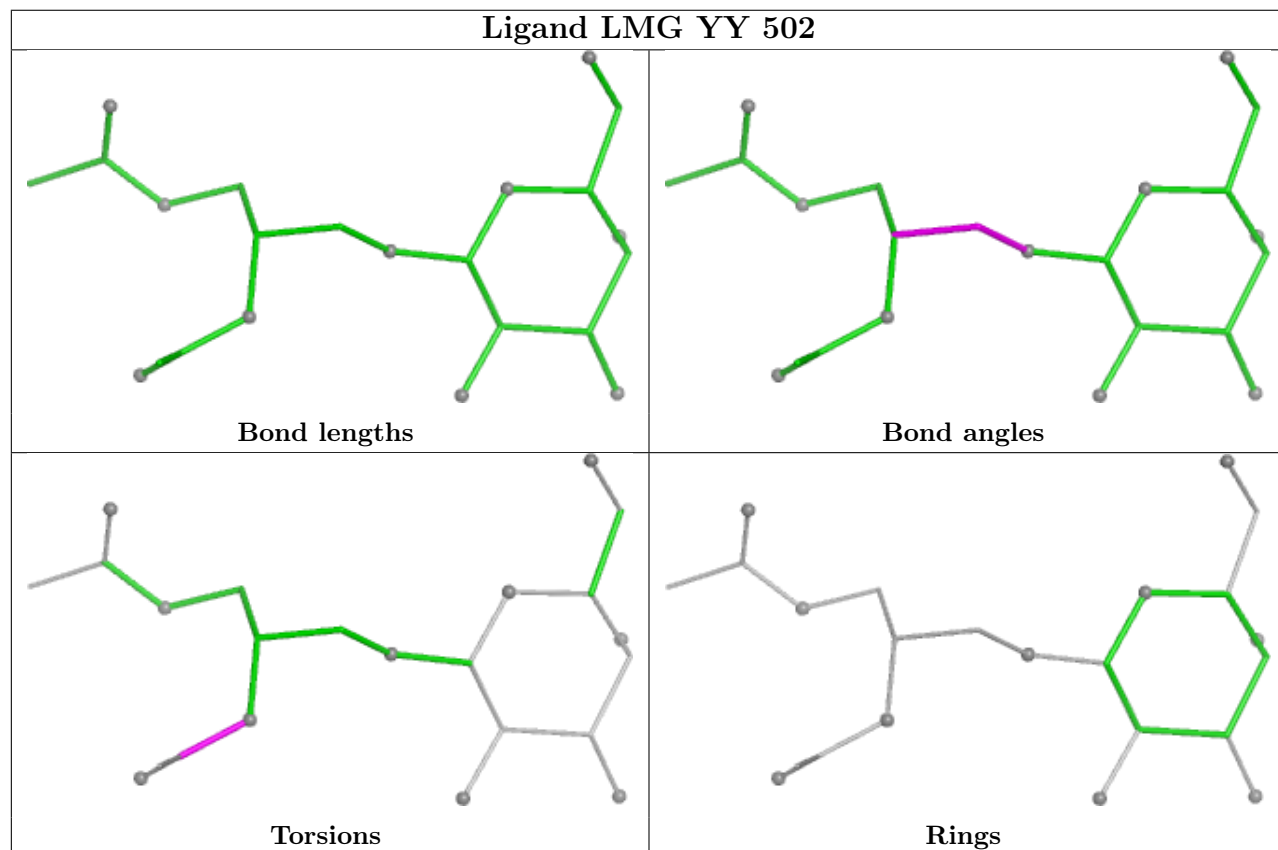
Ligand NDP ZP 501

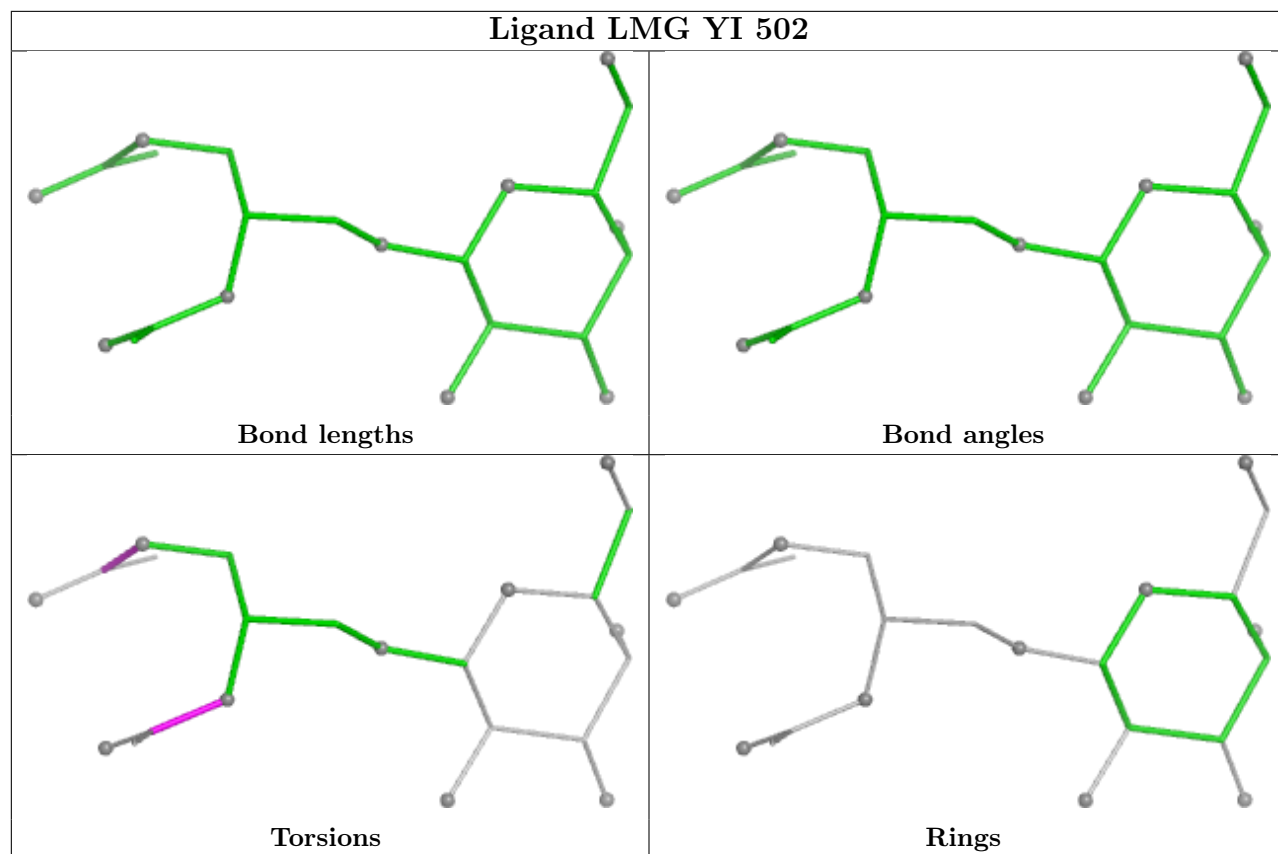


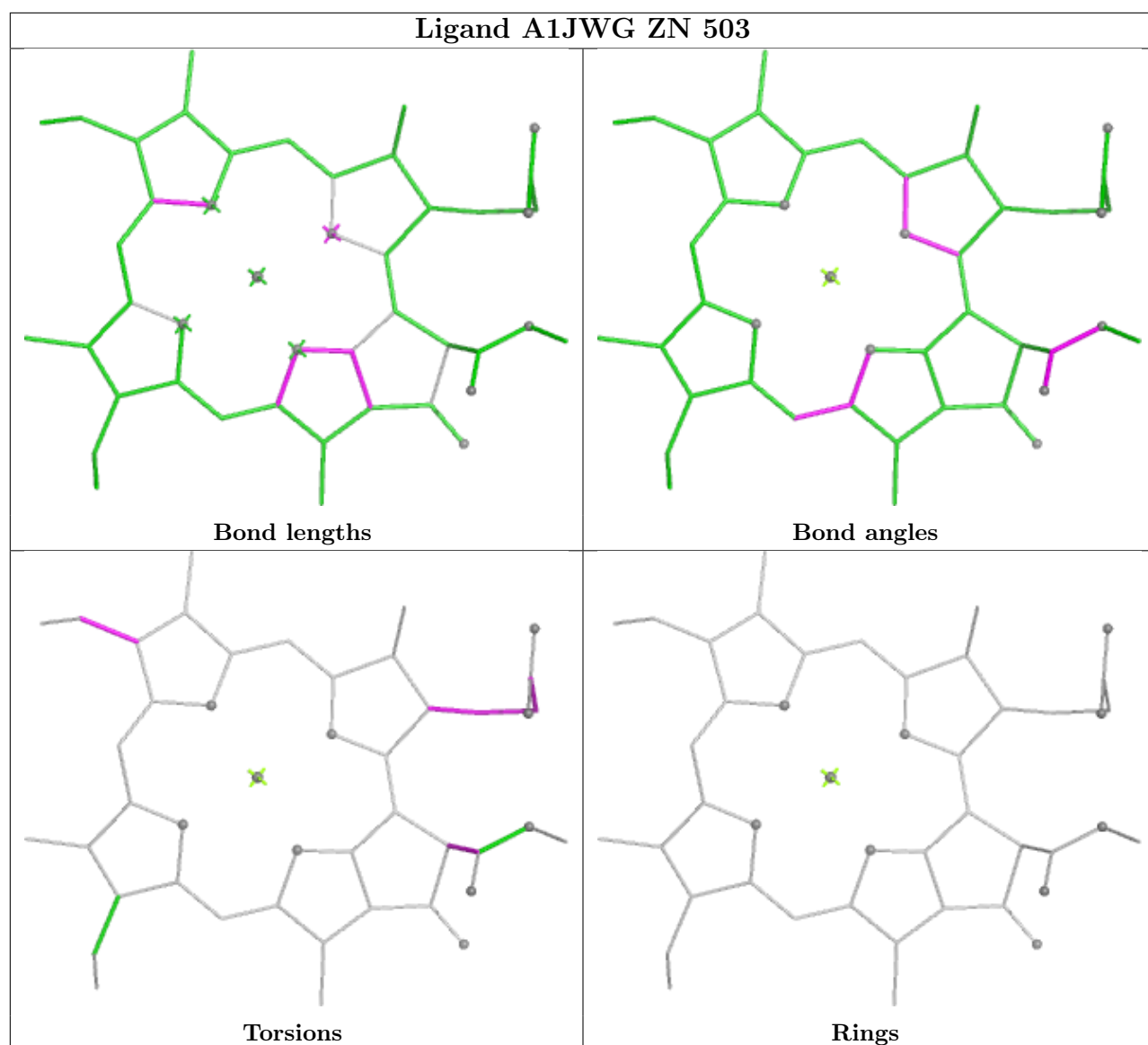
Ligand LMG ZQ 502

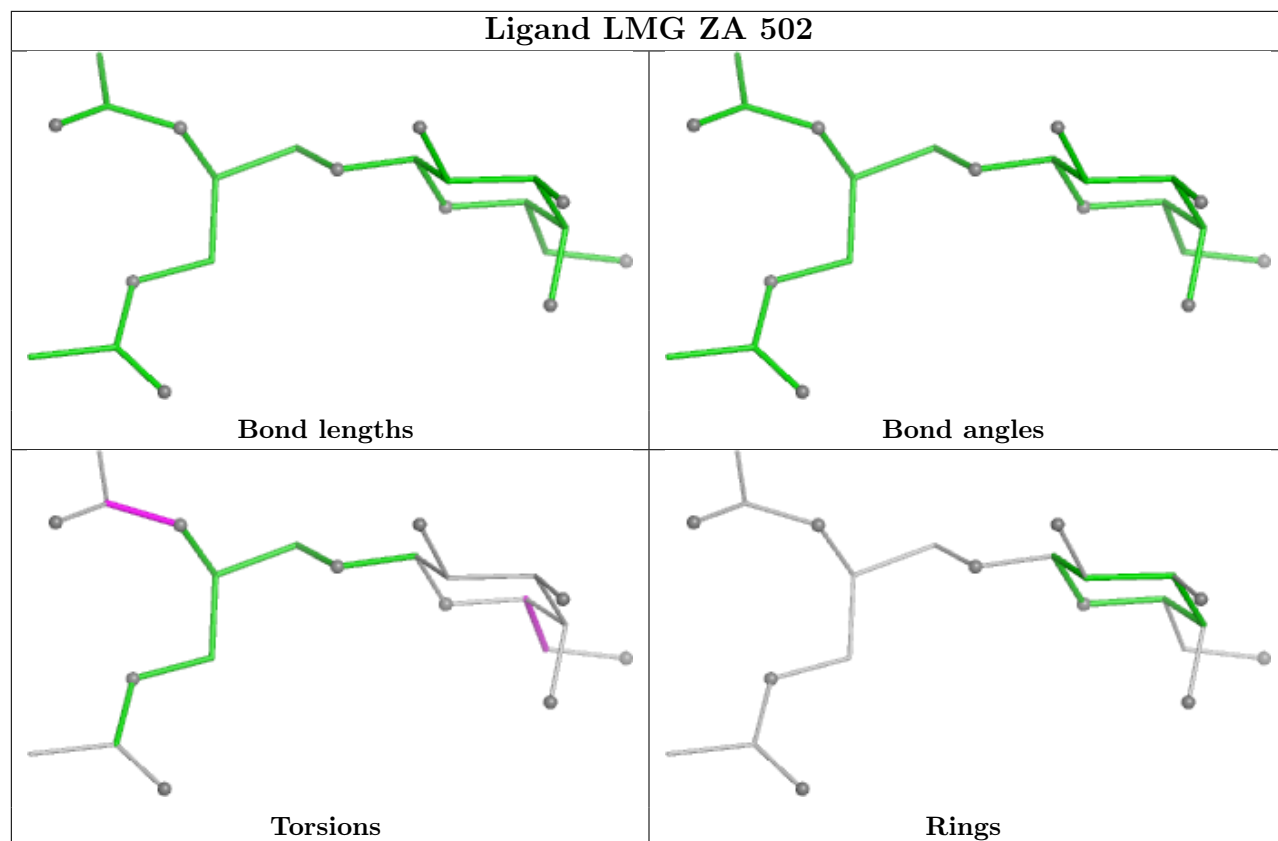


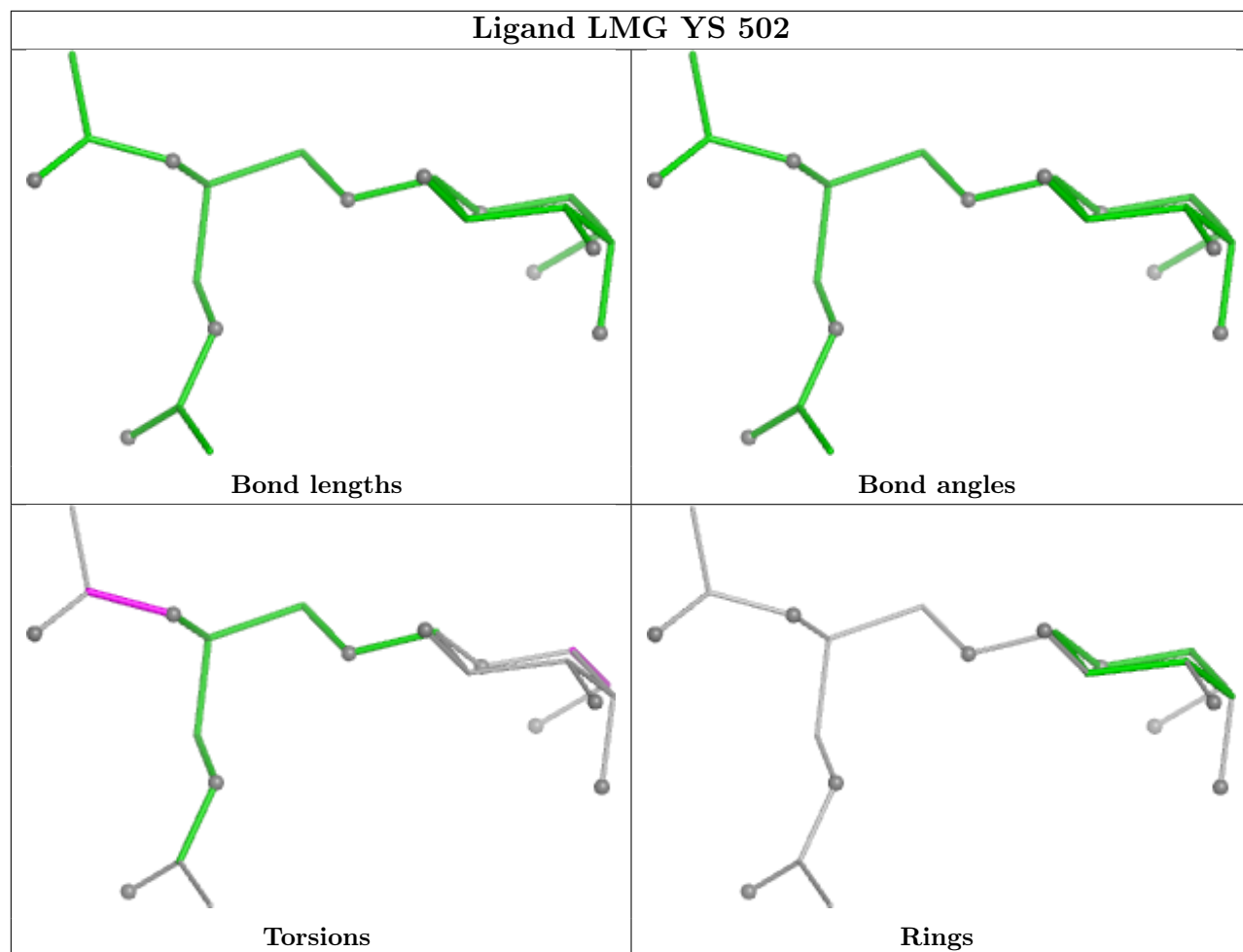
Ligand LMG YY 502



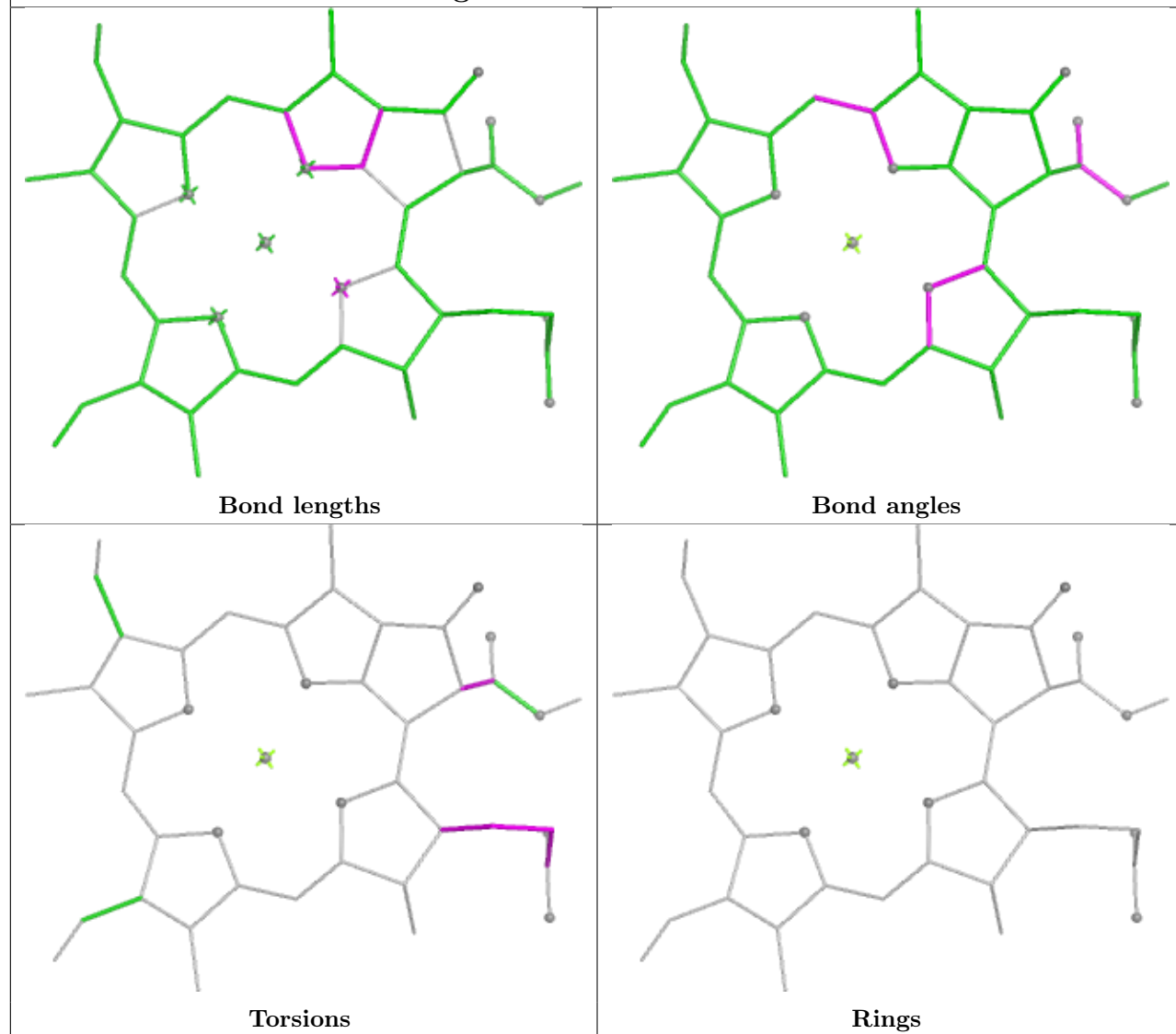




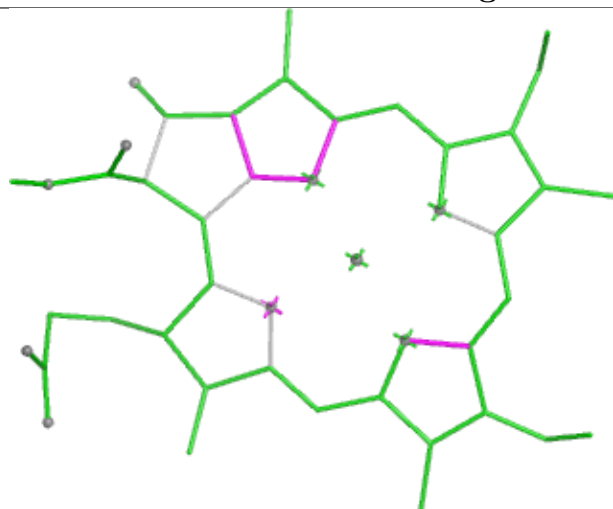




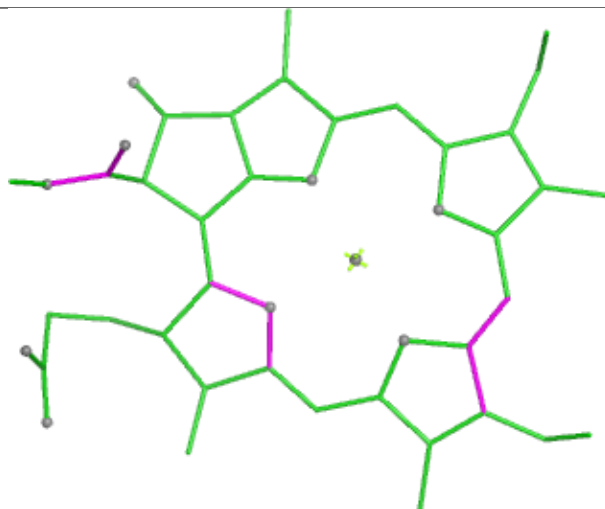
Ligand A1JWG YT 503



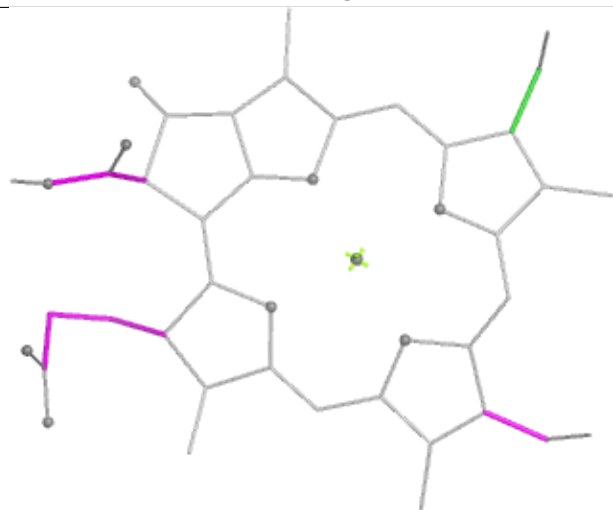
Ligand A1JWG ZL 503



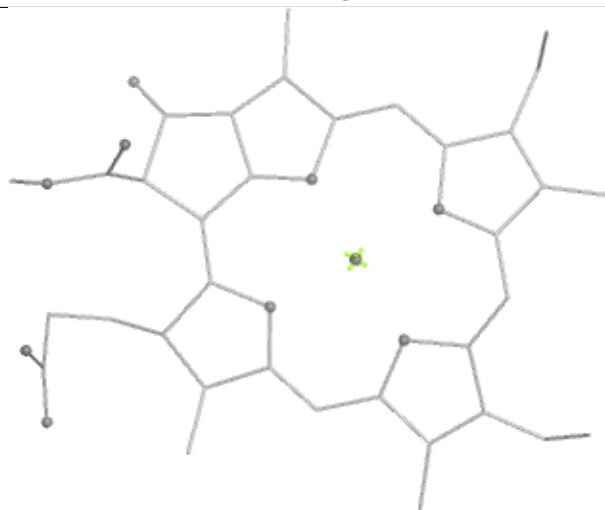
Bond lengths



Bond angles

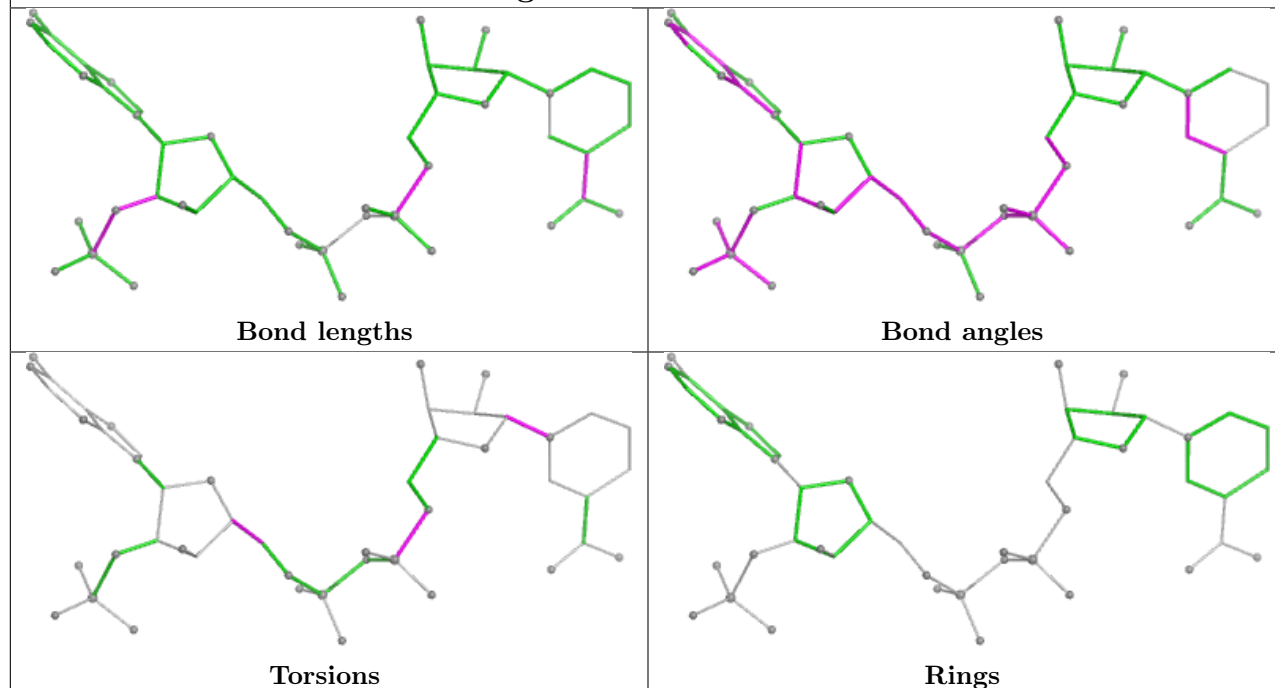


Torsions

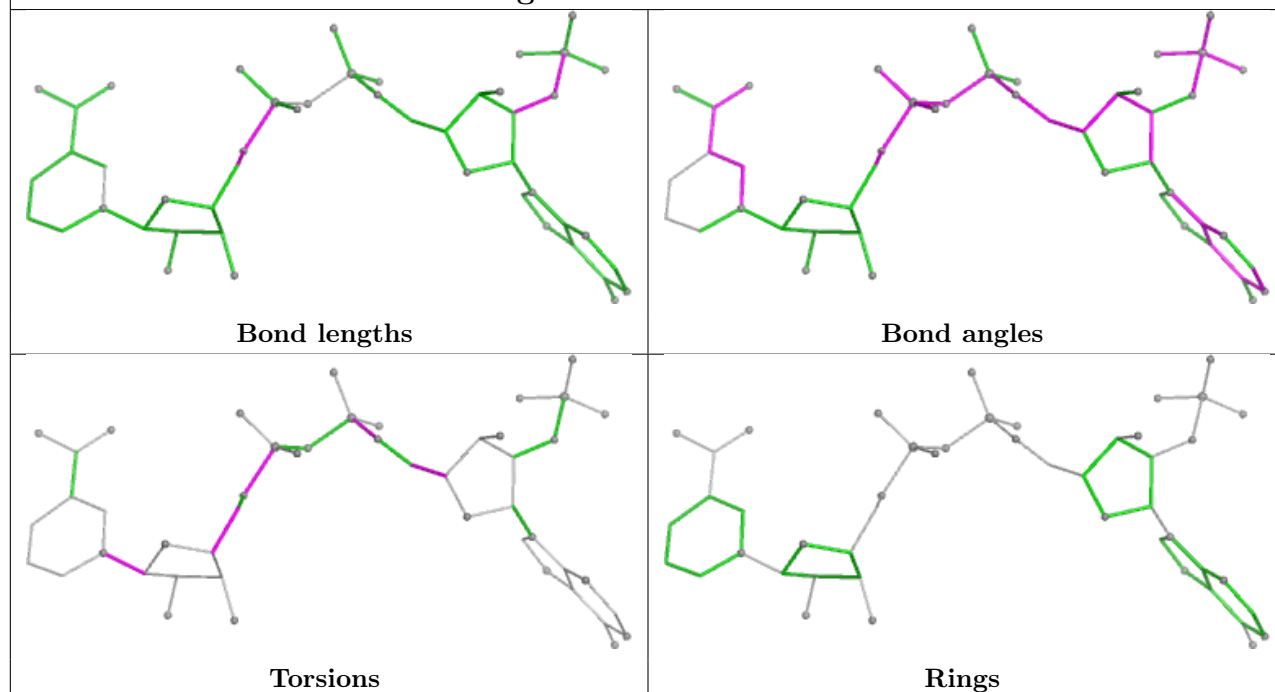


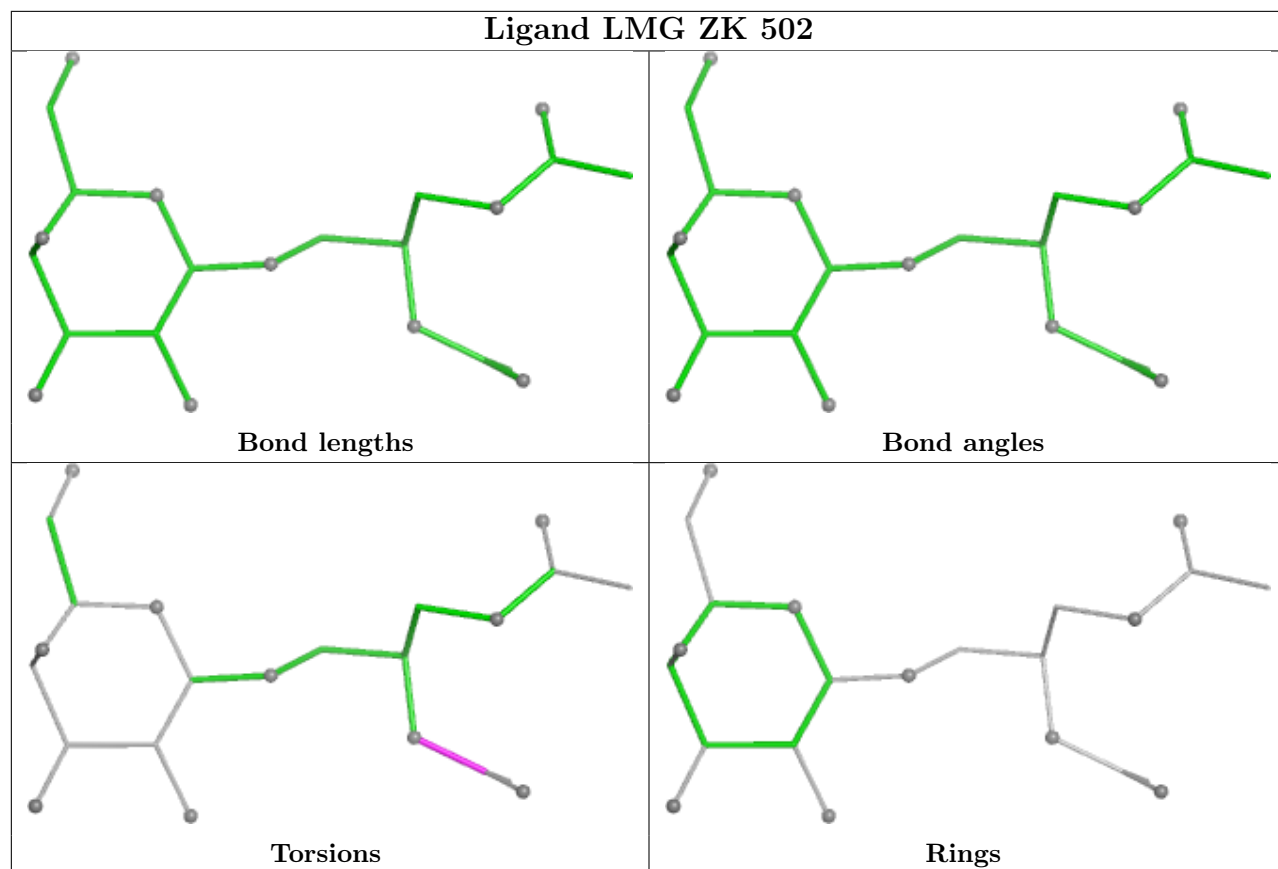
Rings

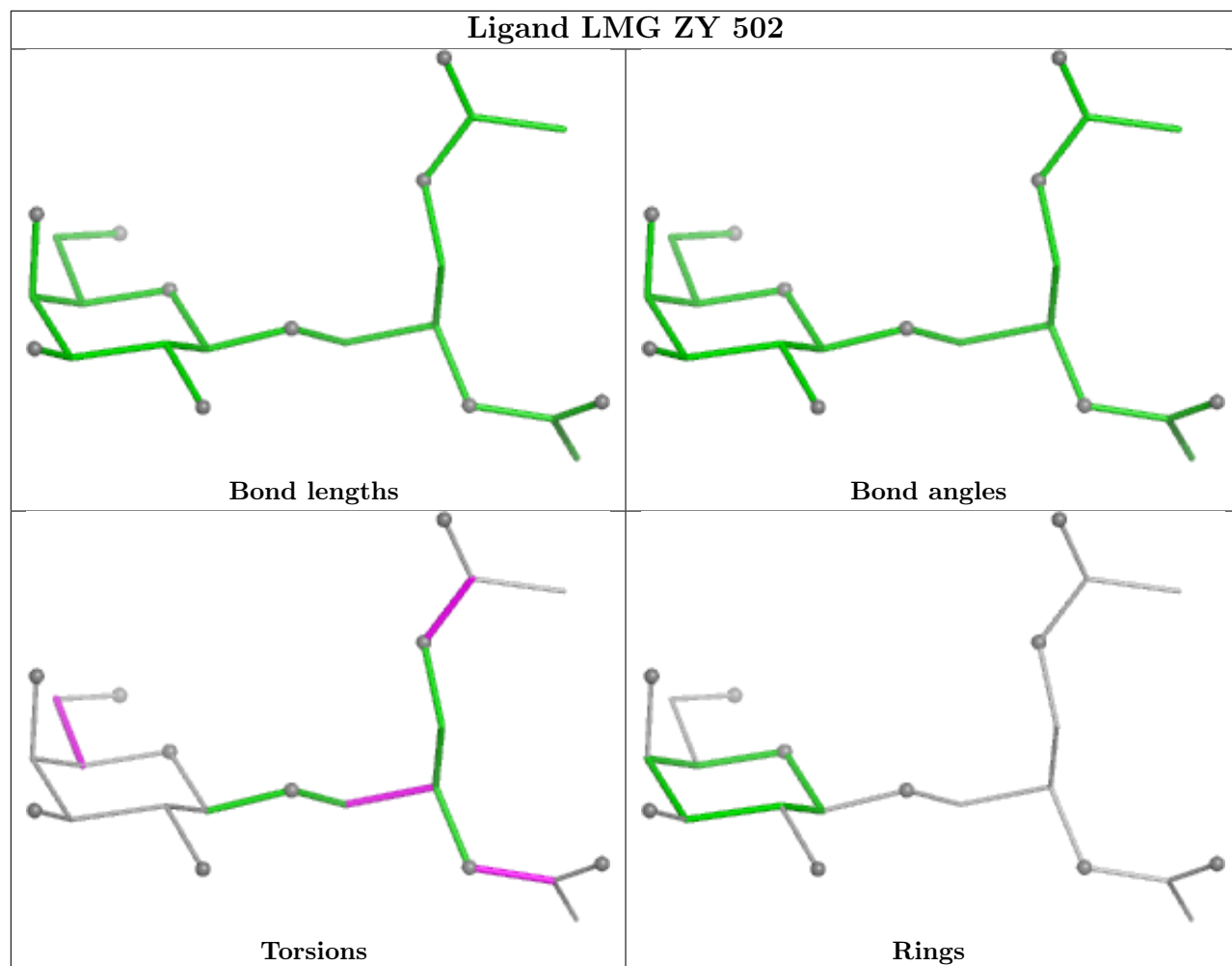
Ligand NDP YL 501

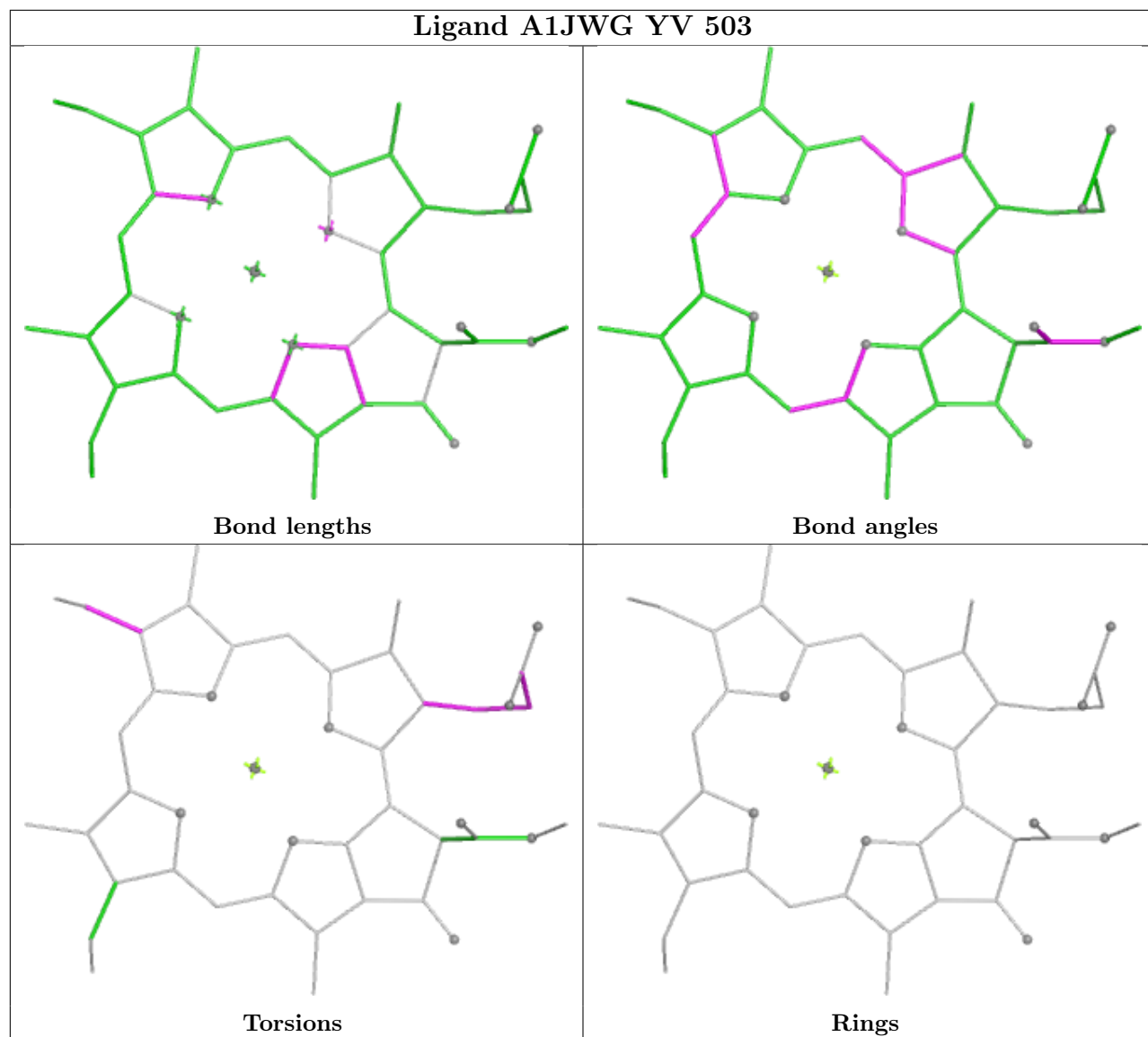


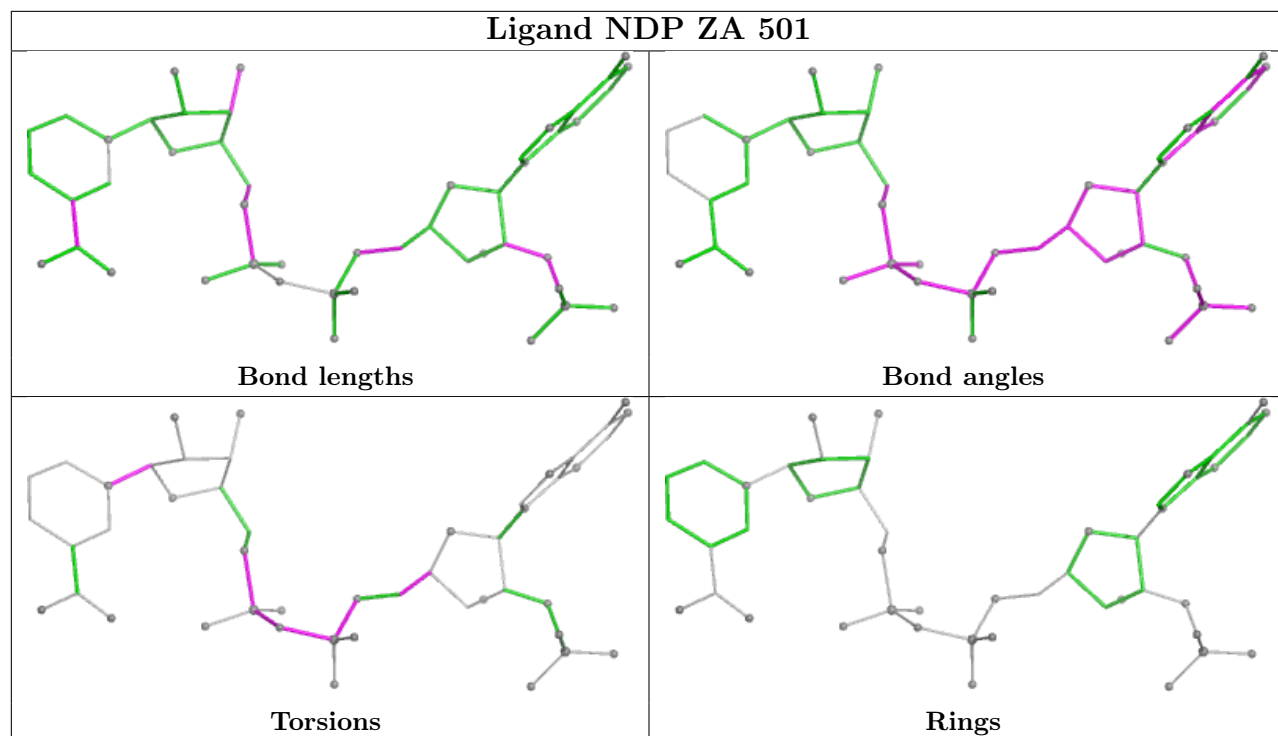
Ligand NDP ZX 501

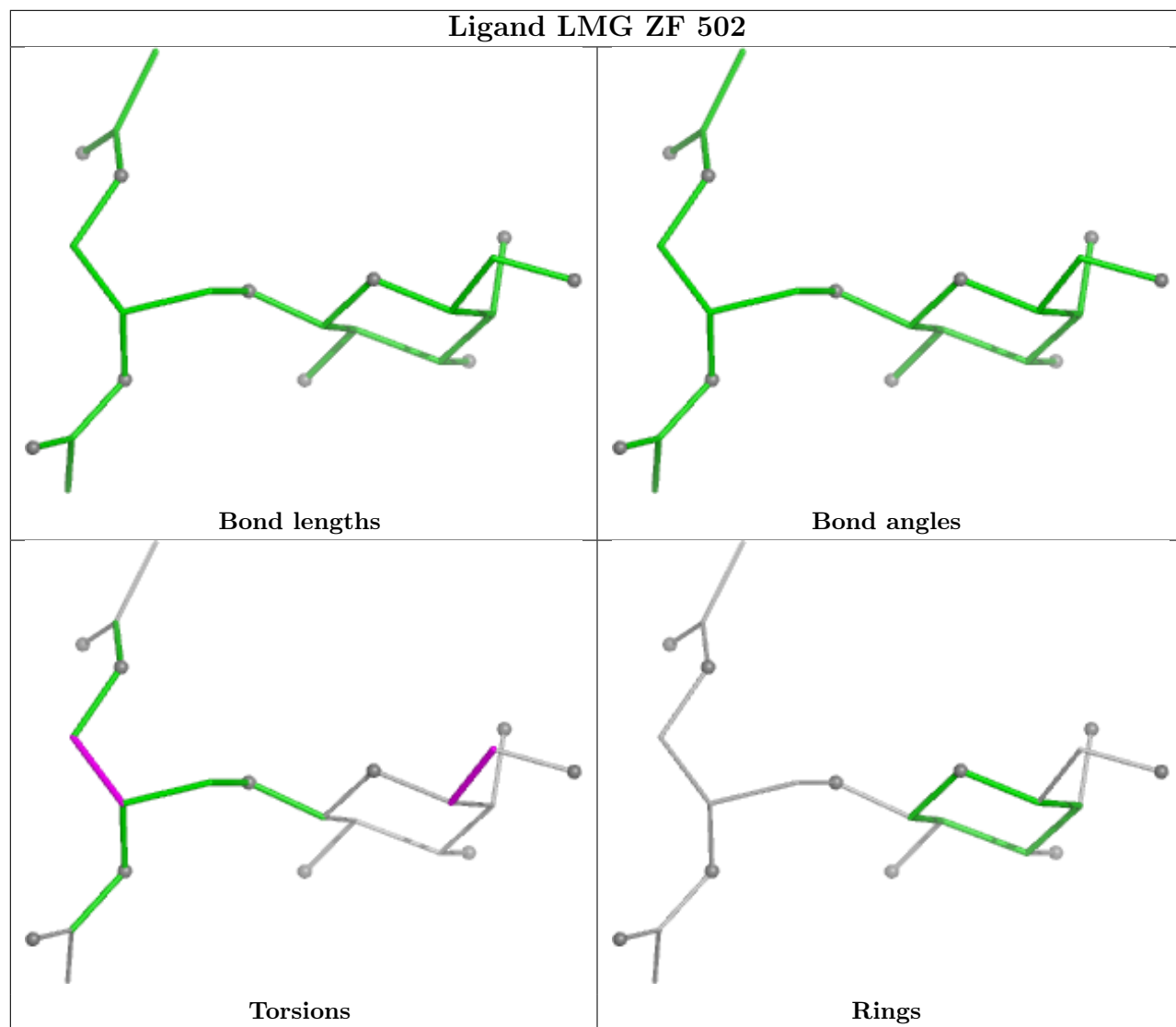




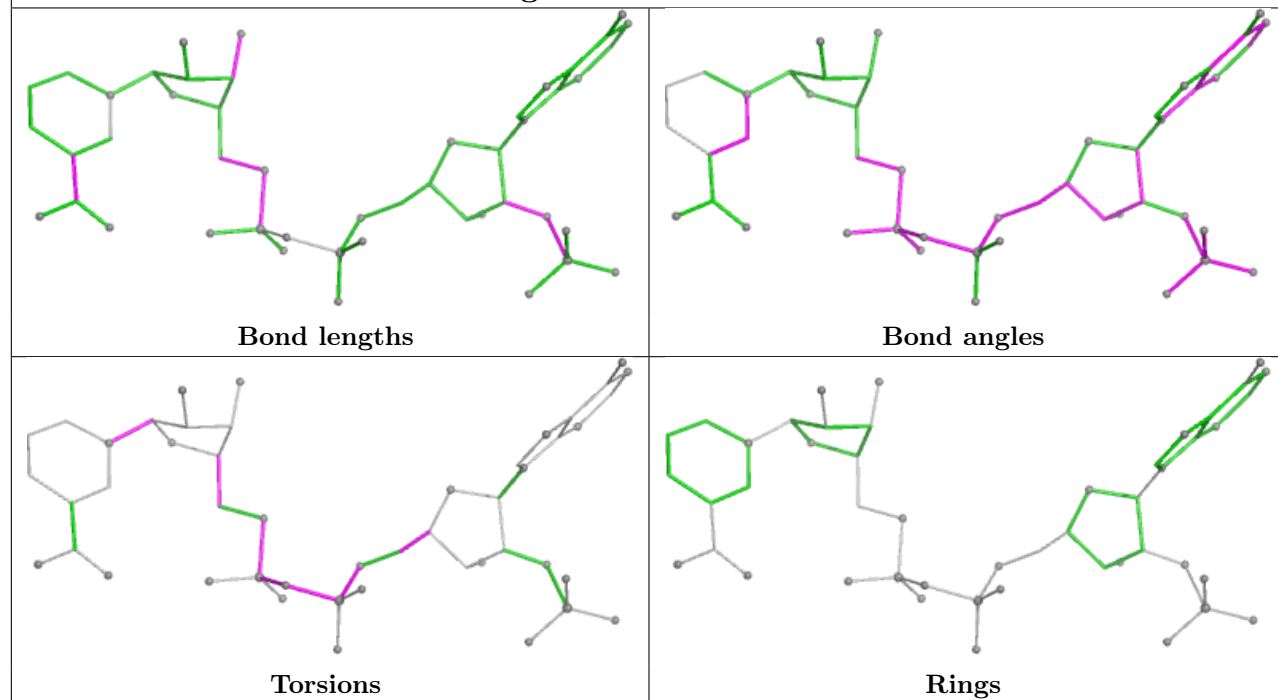




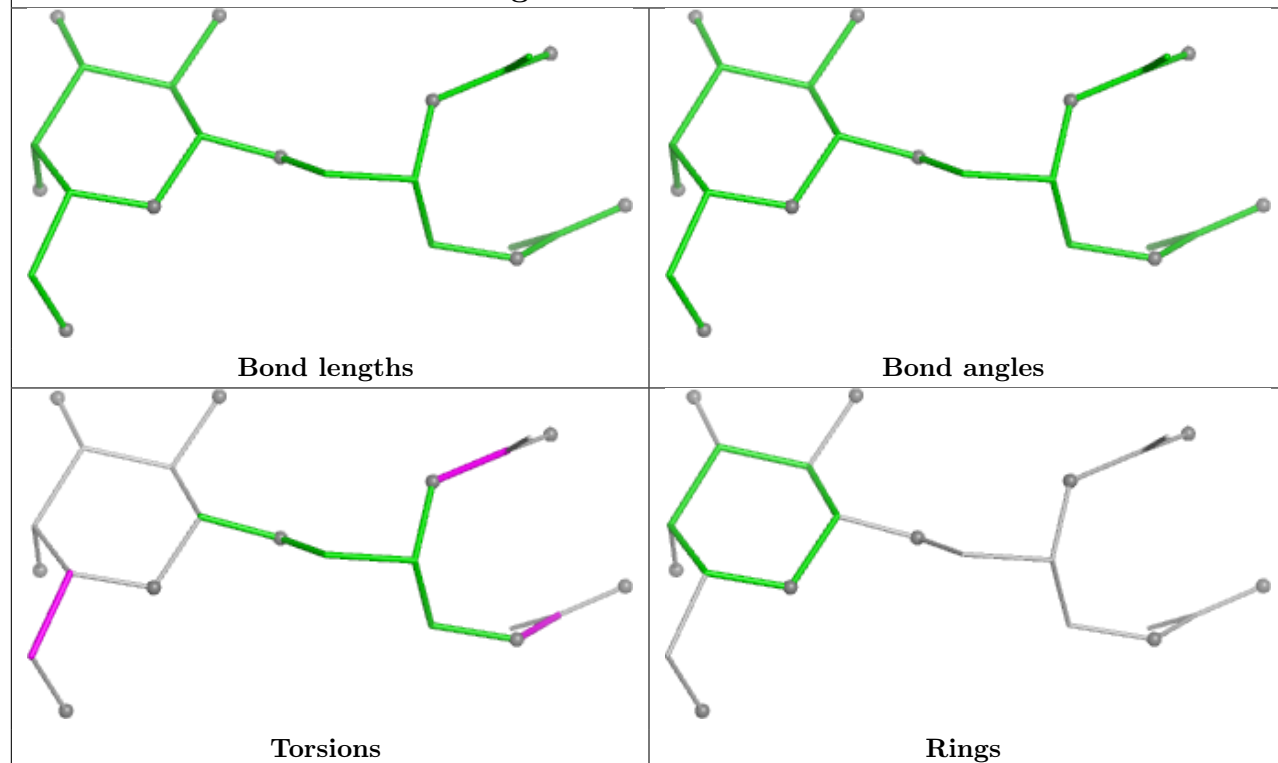


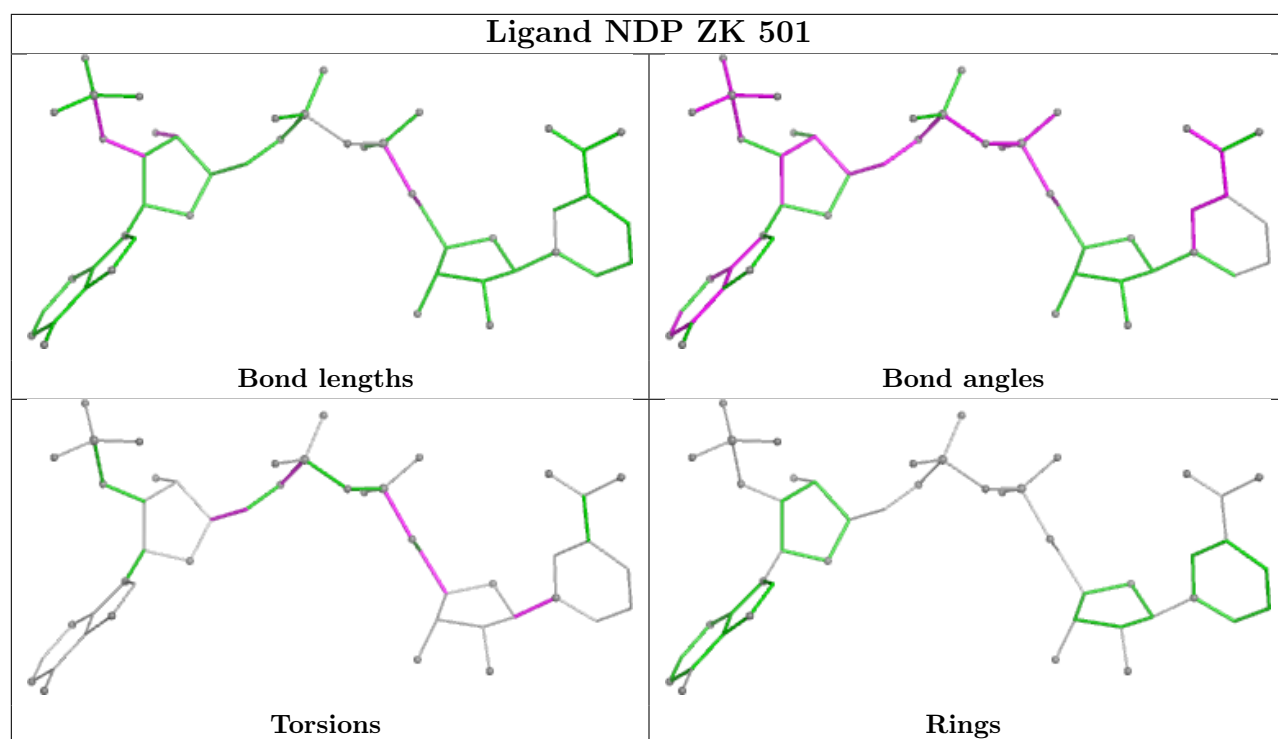


Ligand NDP ZZ 501

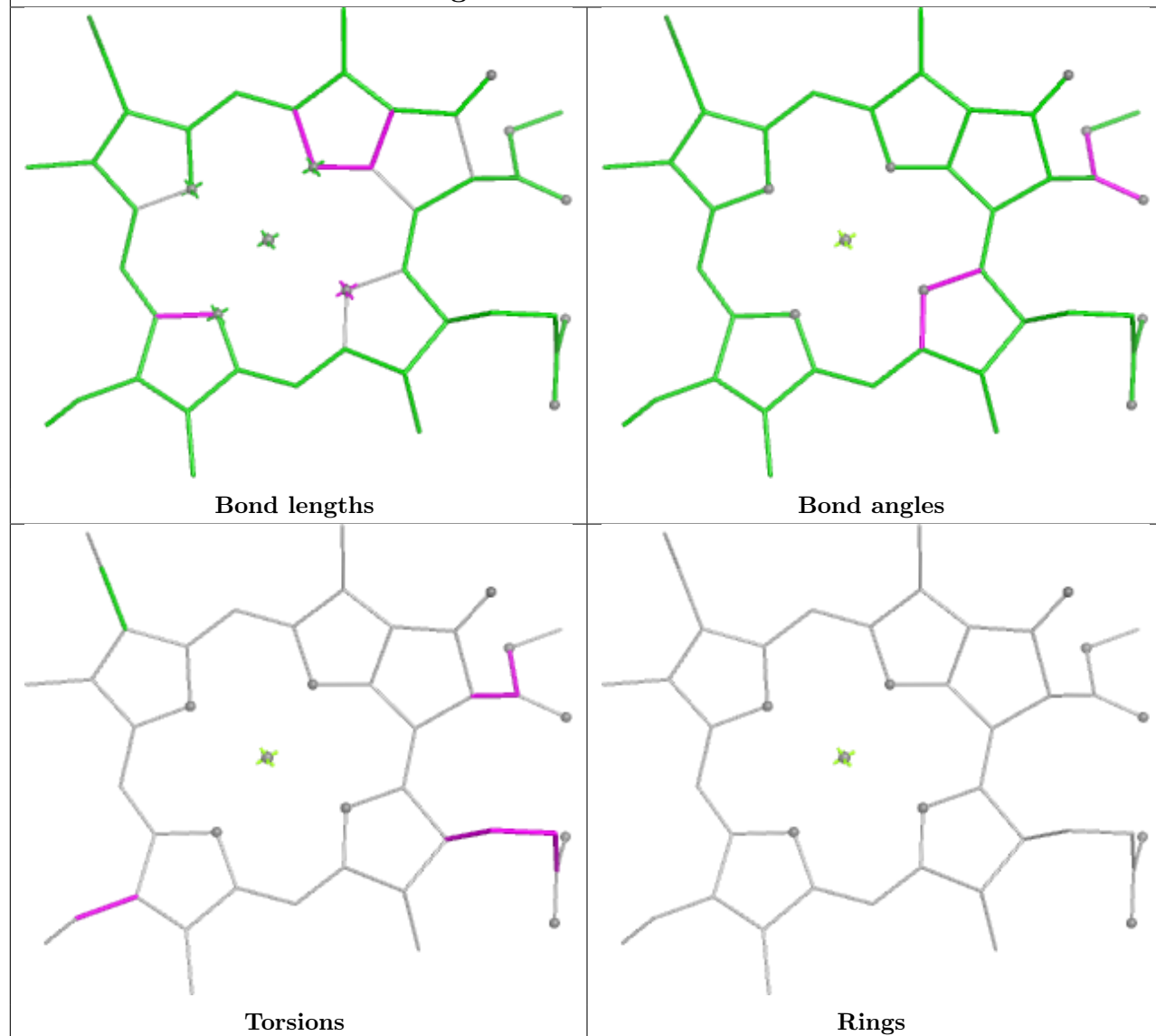


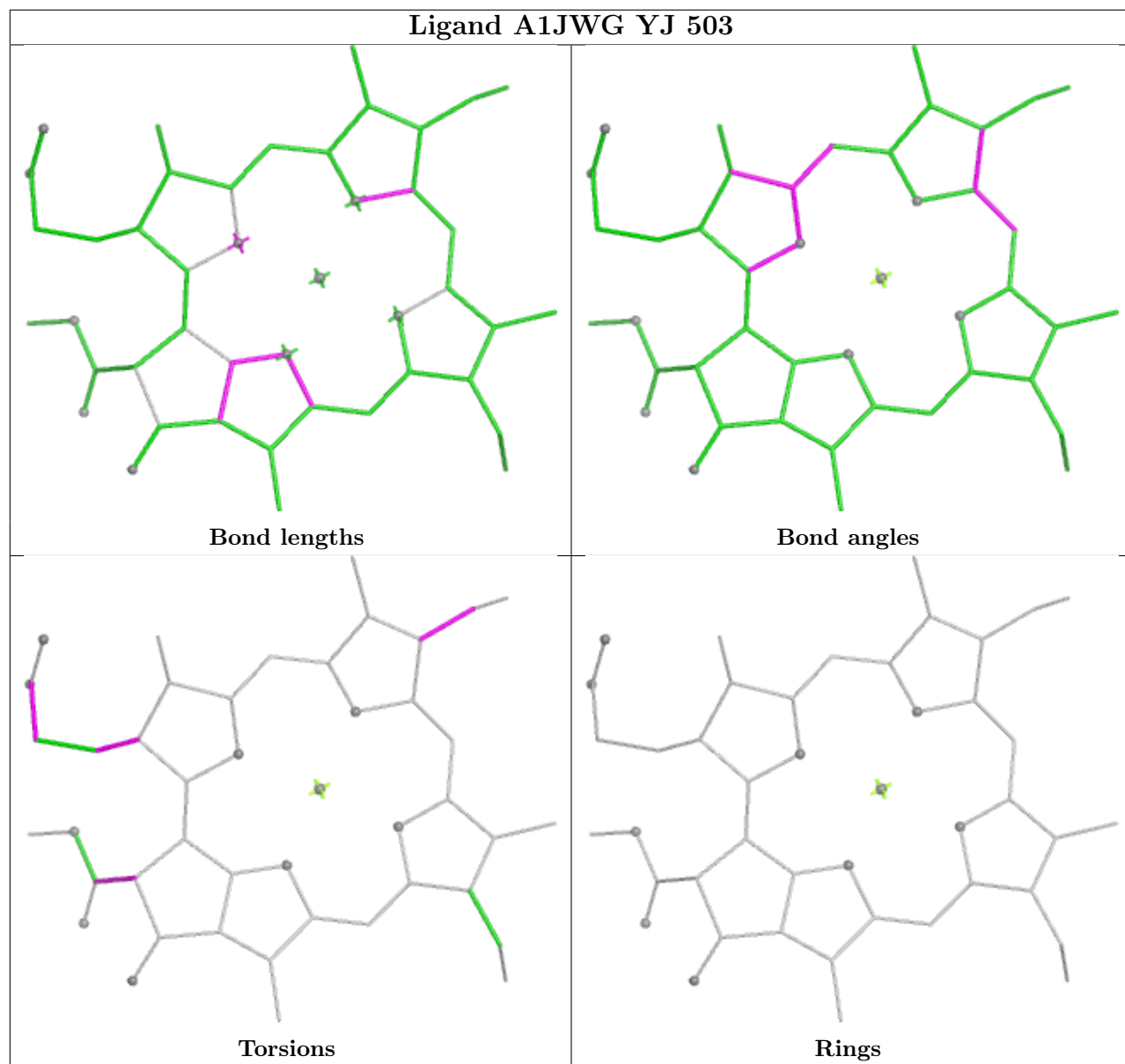
Ligand LMG ZW 502

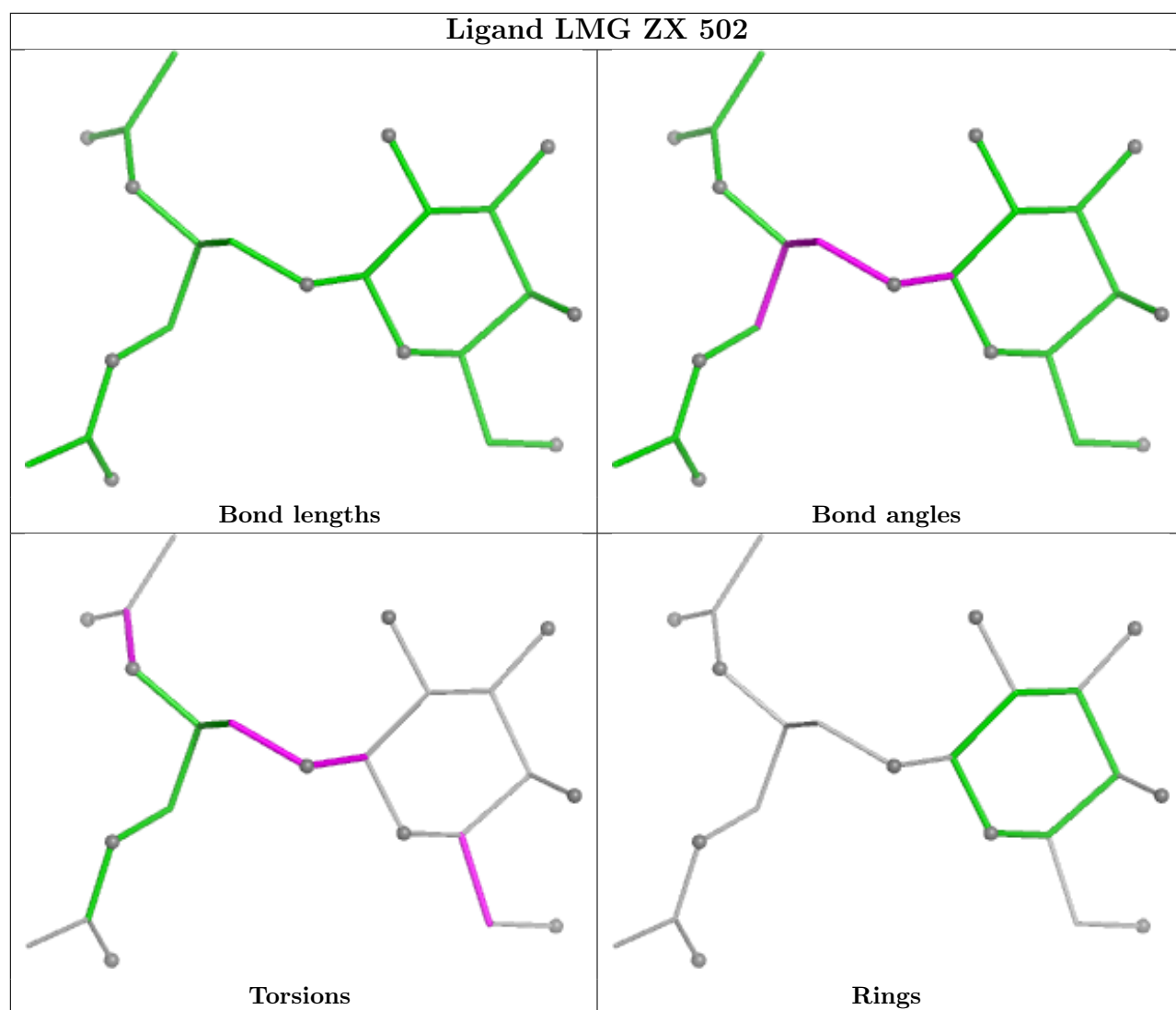




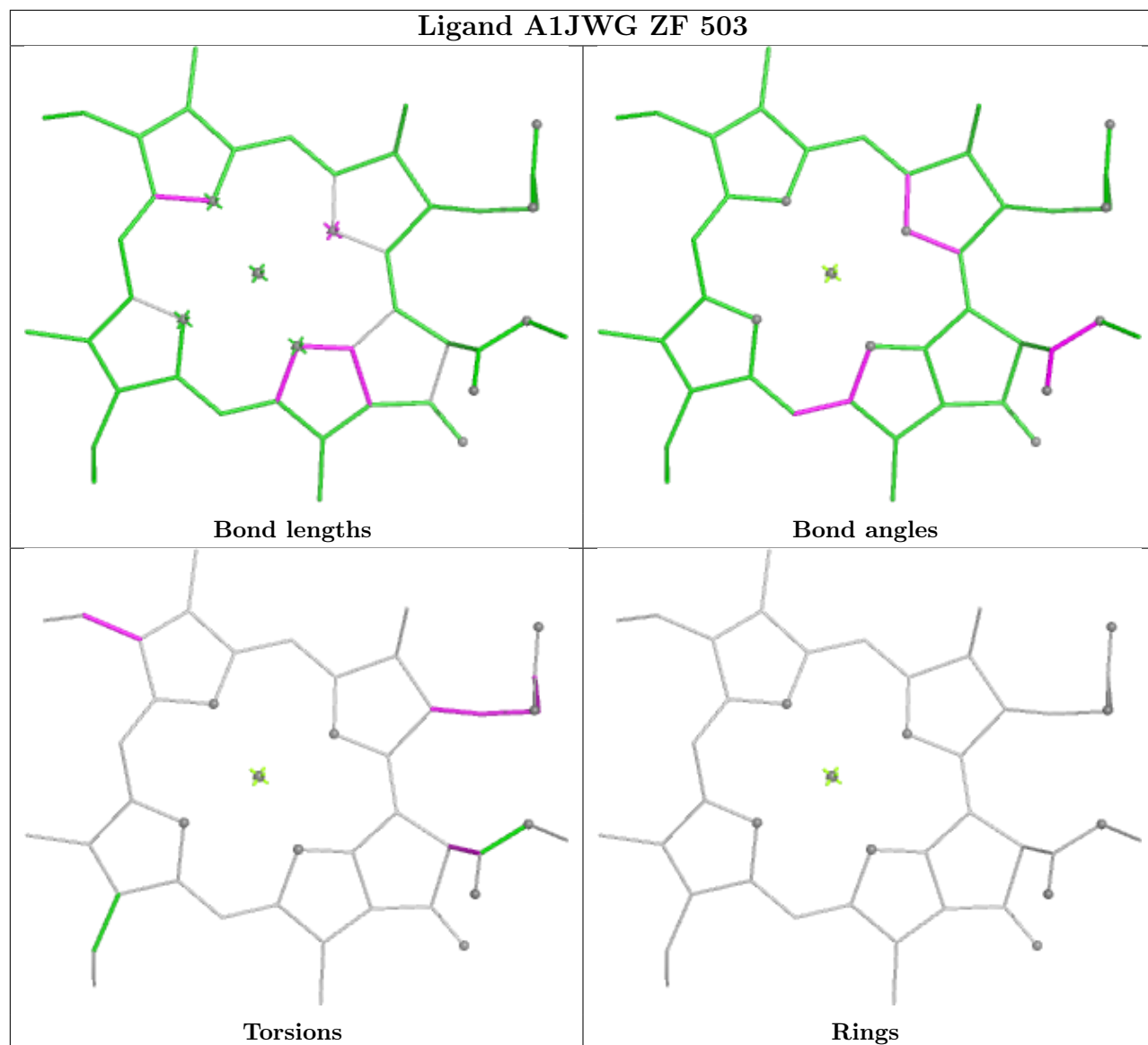
Ligand A1JWG YI 503



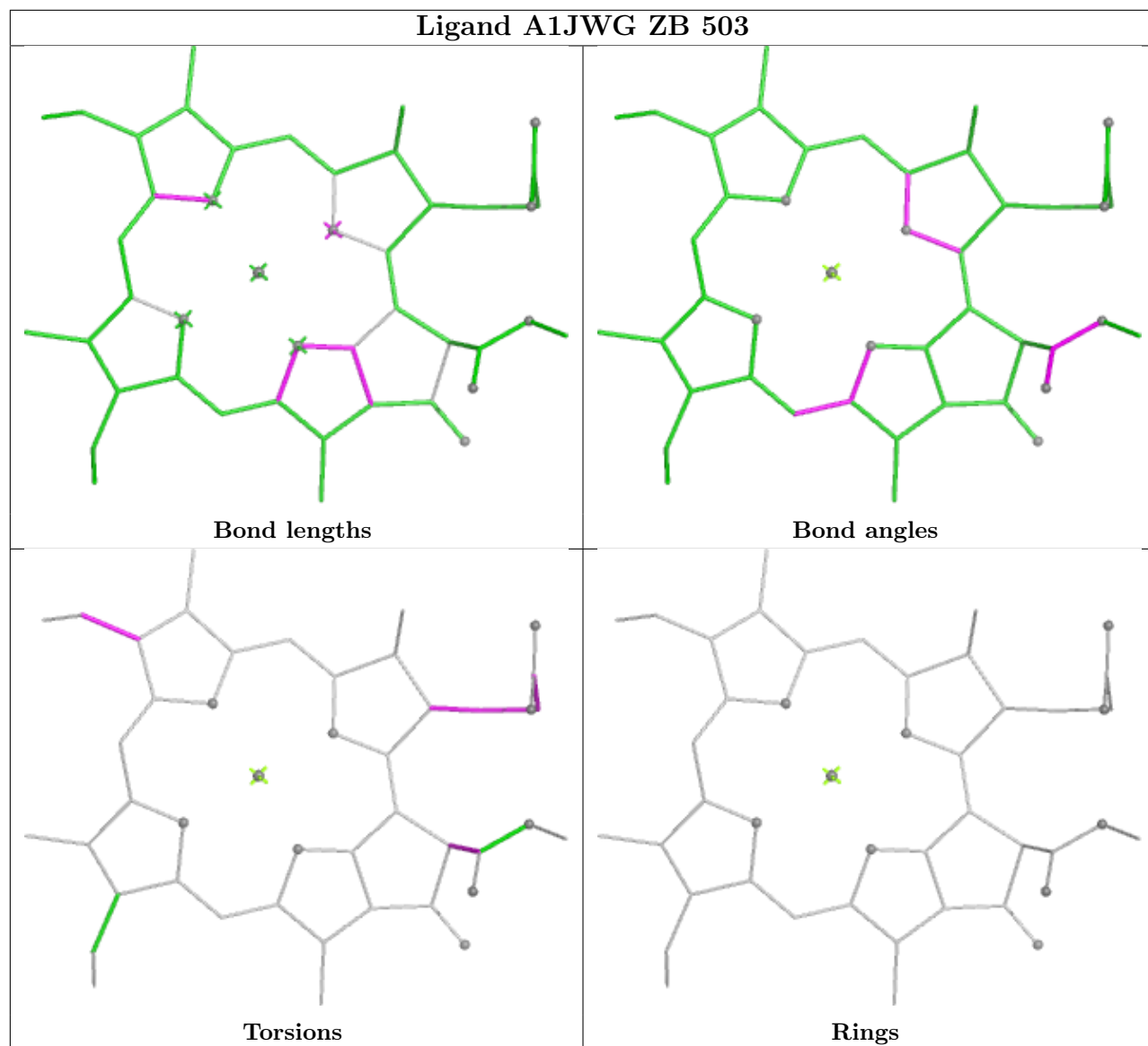


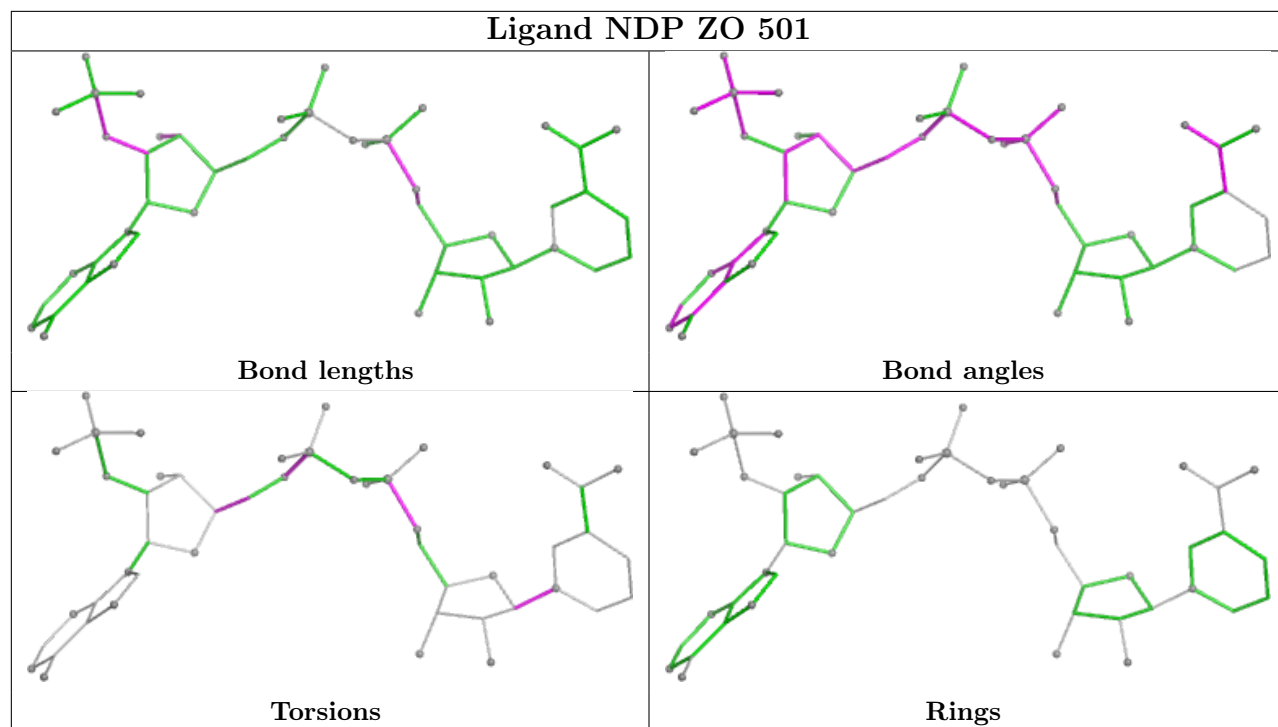


Ligand A1JWG ZF 503

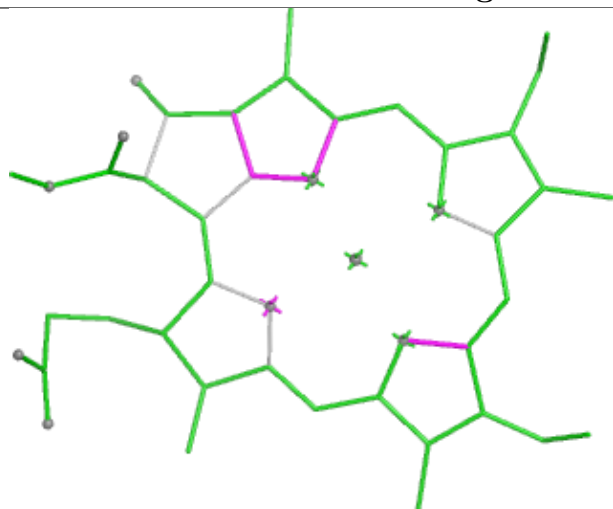


Ligand A1JWG ZB 503

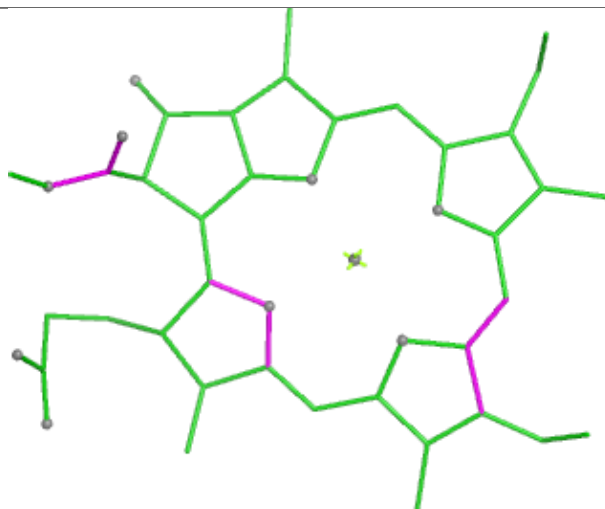




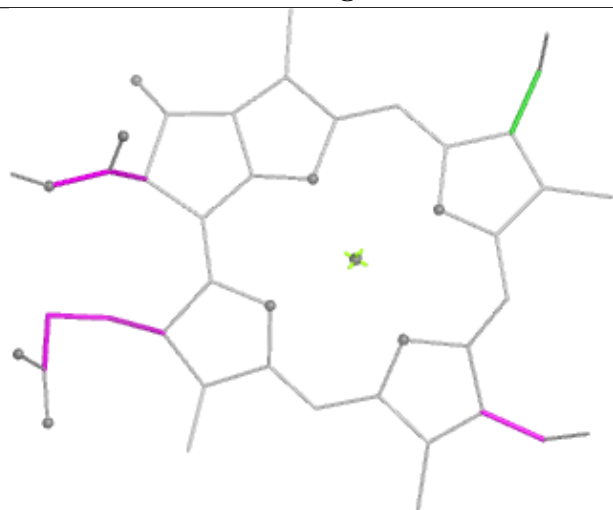
Ligand A1JWG ZX 503



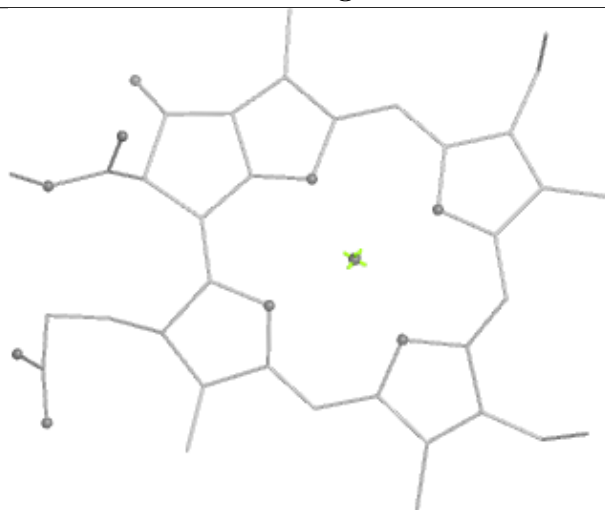
Bond lengths



Bond angles

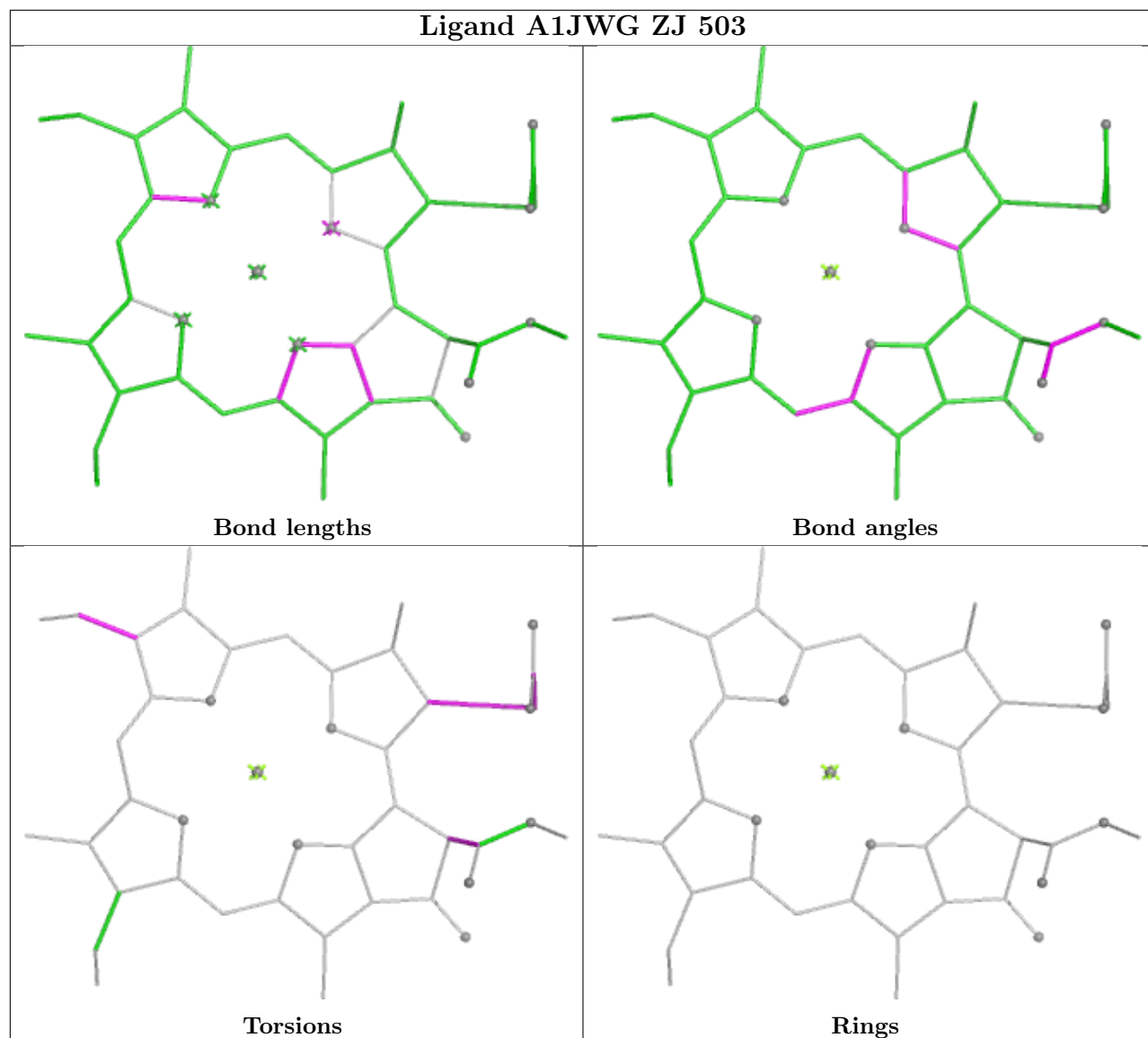


Torsions

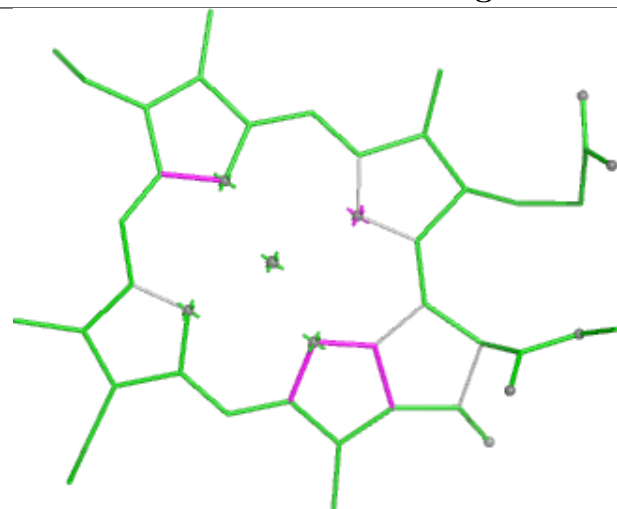


Rings

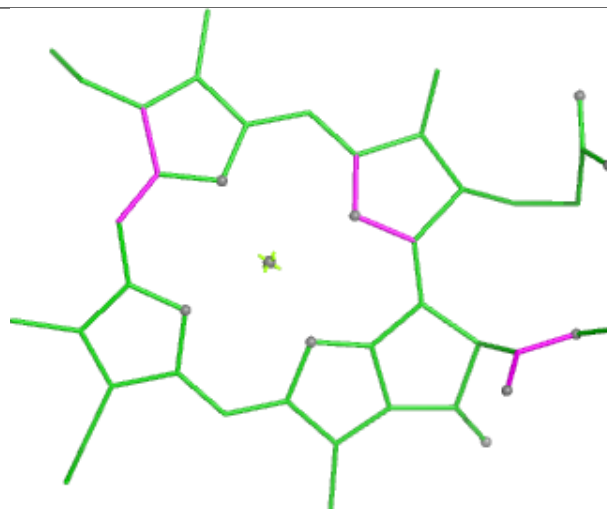
Ligand A1JWG ZJ 503



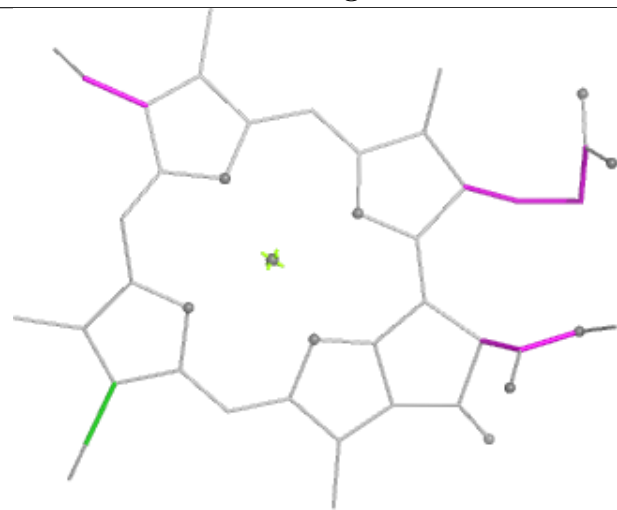
Ligand A1JWG ZW 503



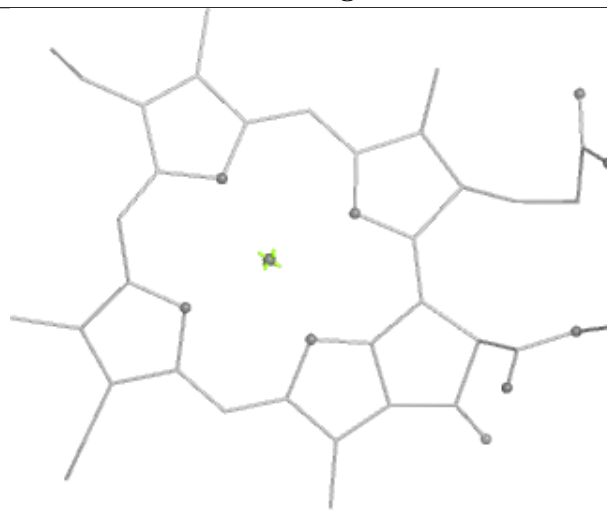
Bond lengths



Bond angles

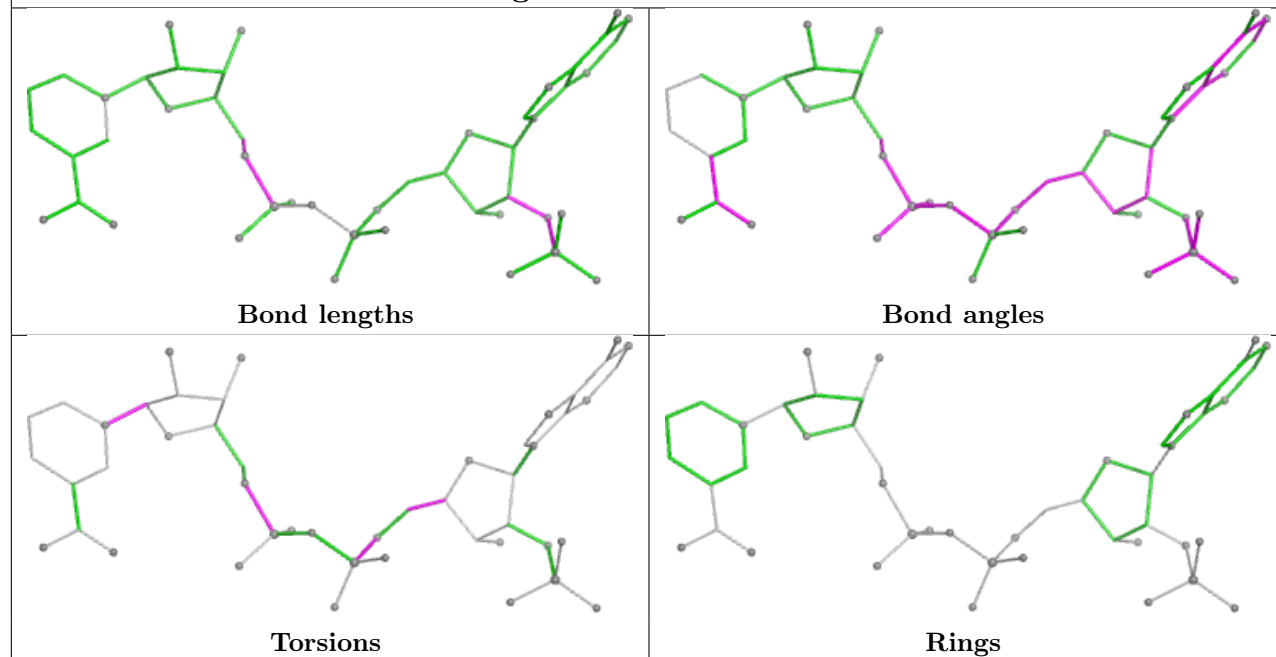


Torsions

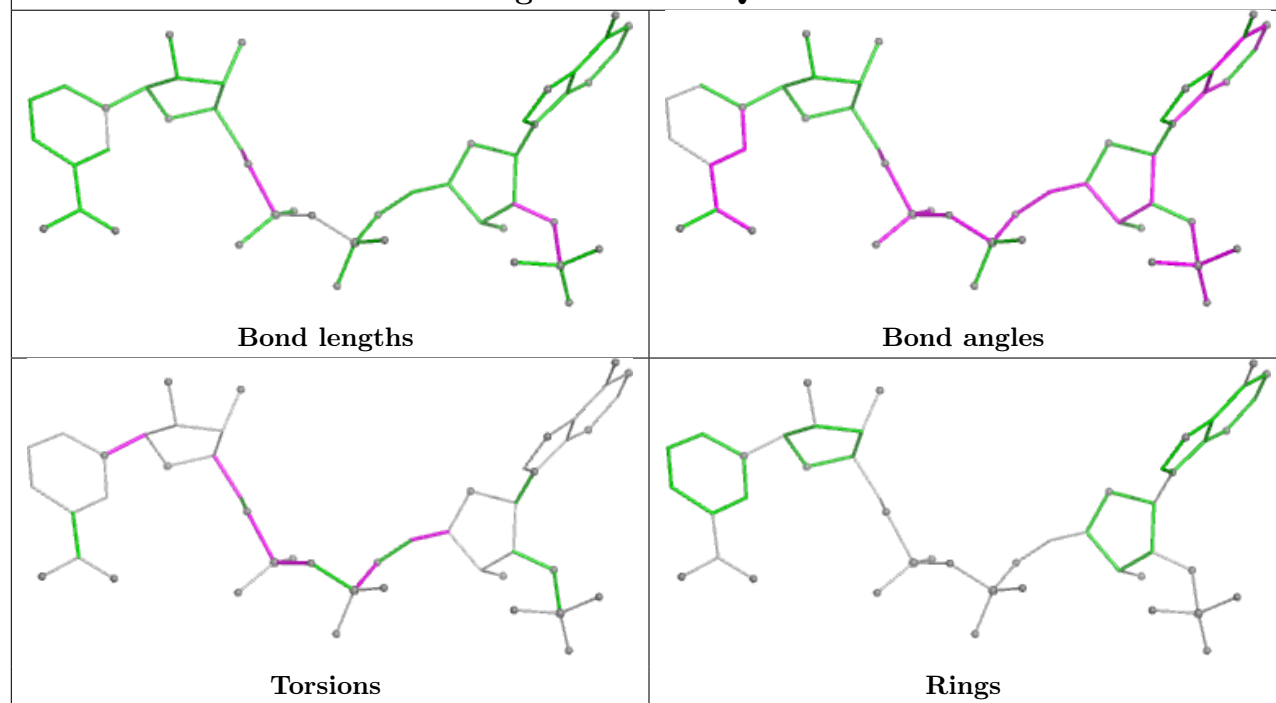


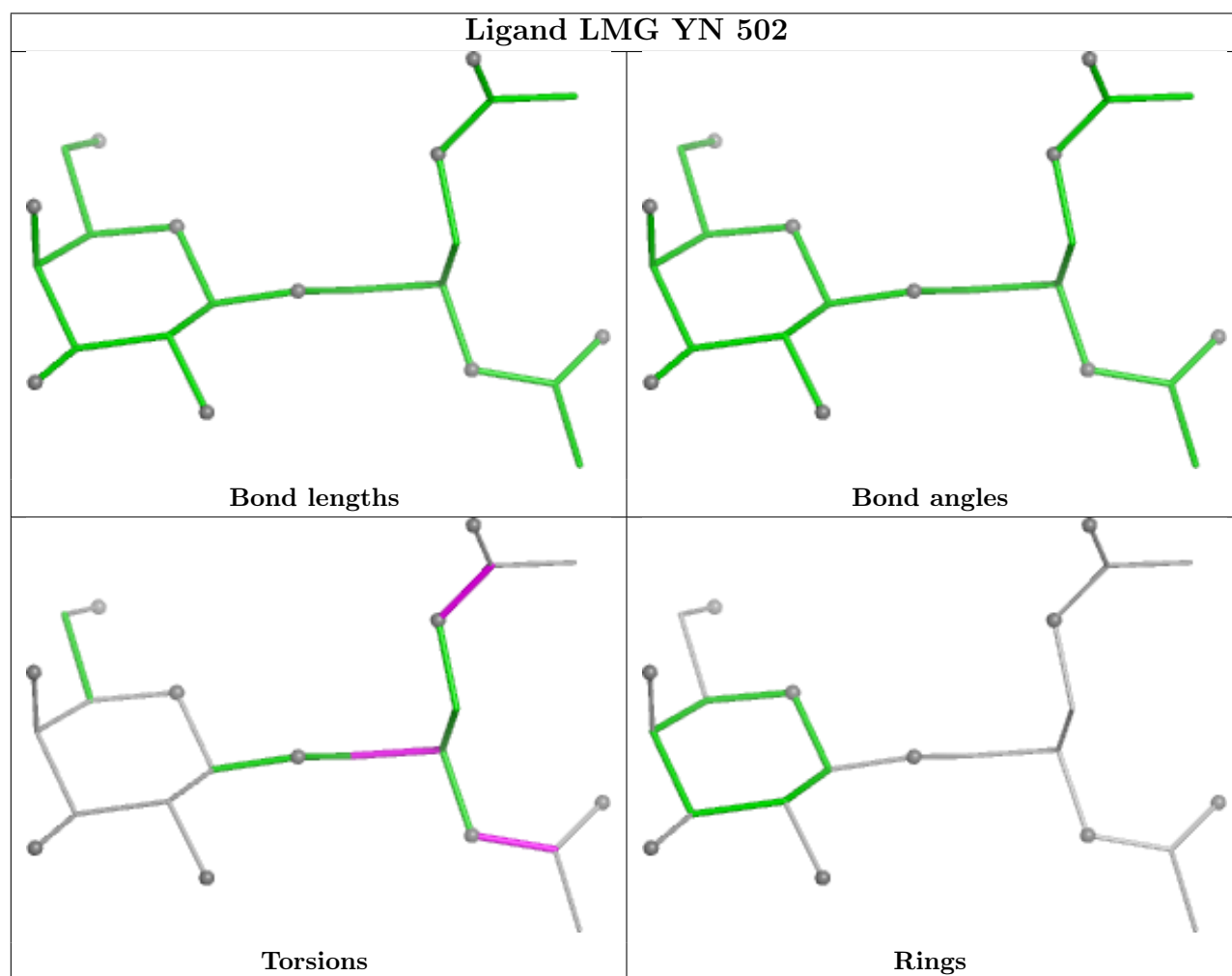
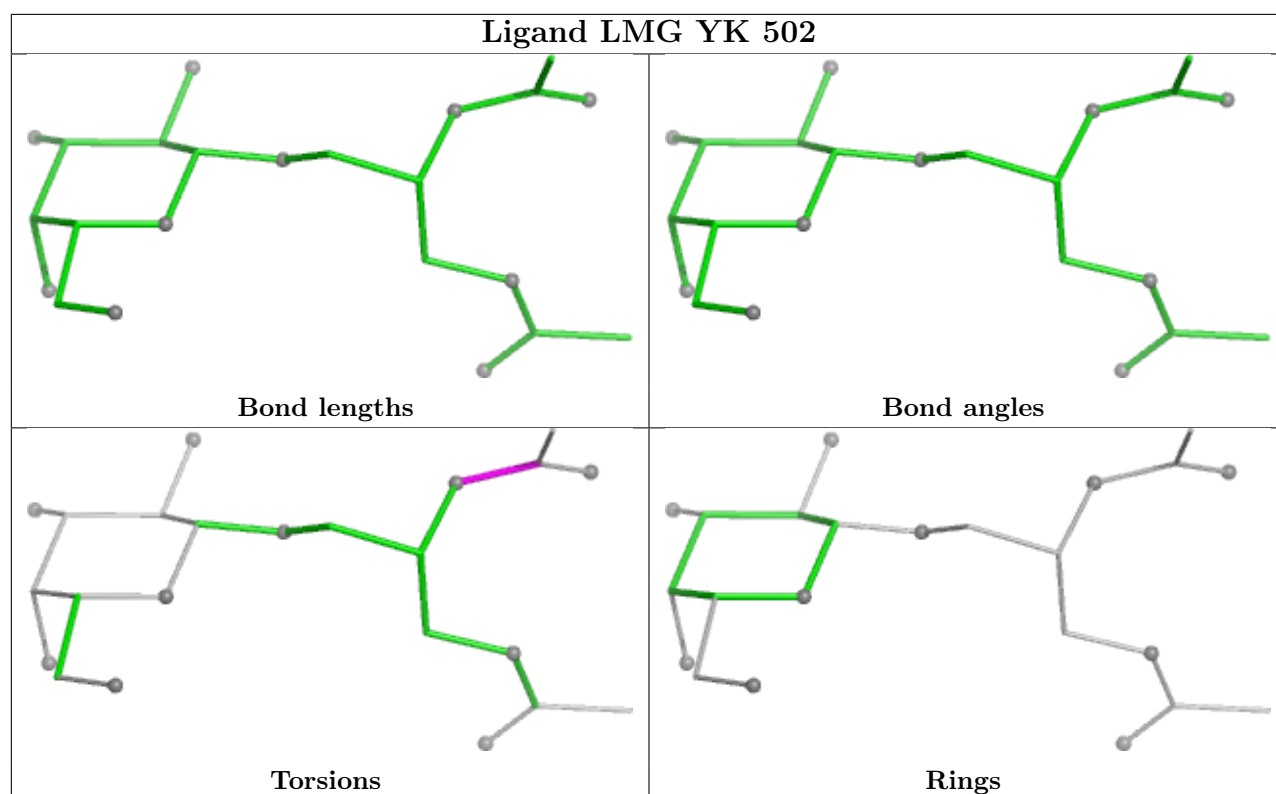
Rings

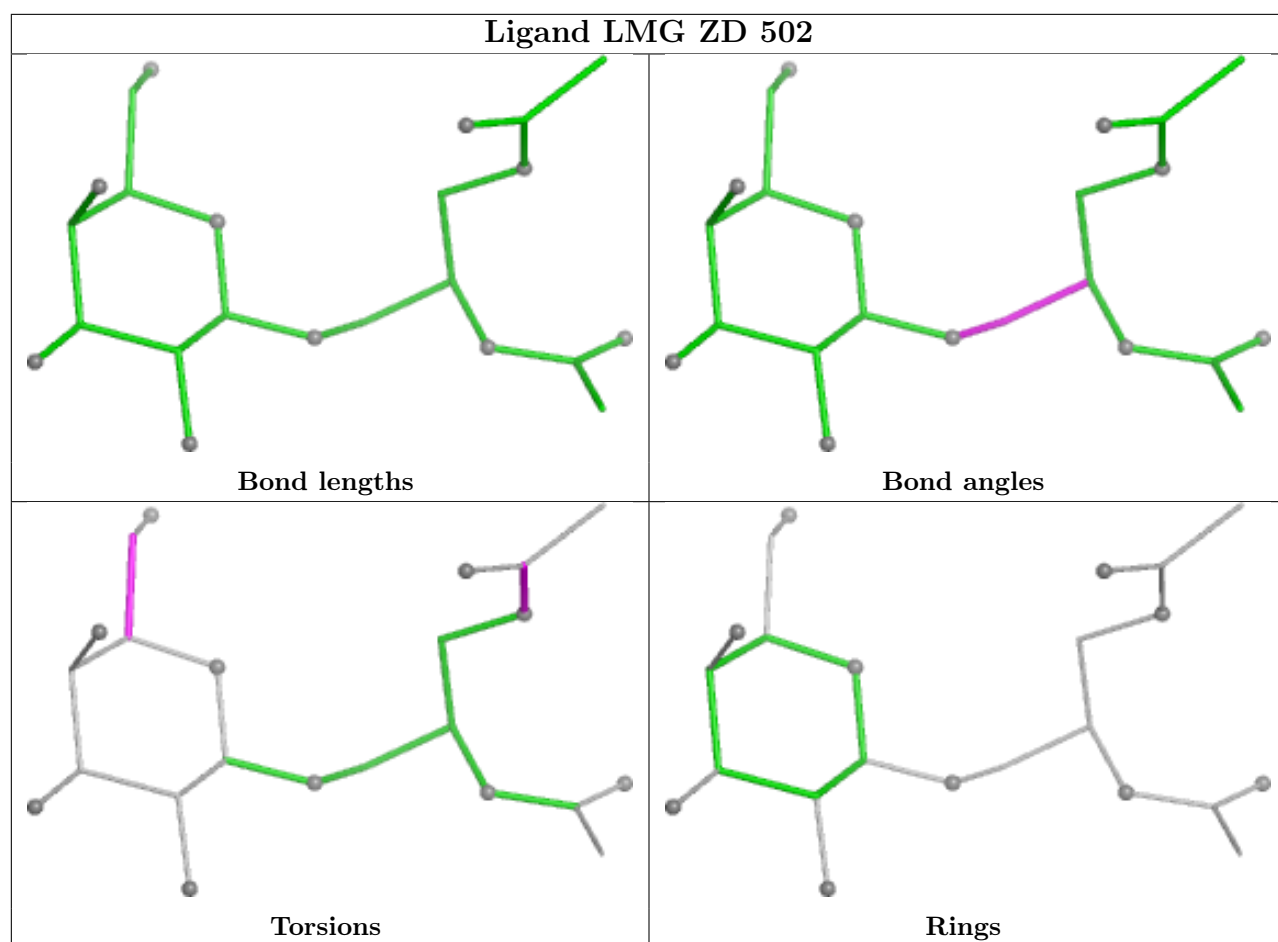
Ligand NDP YM 501



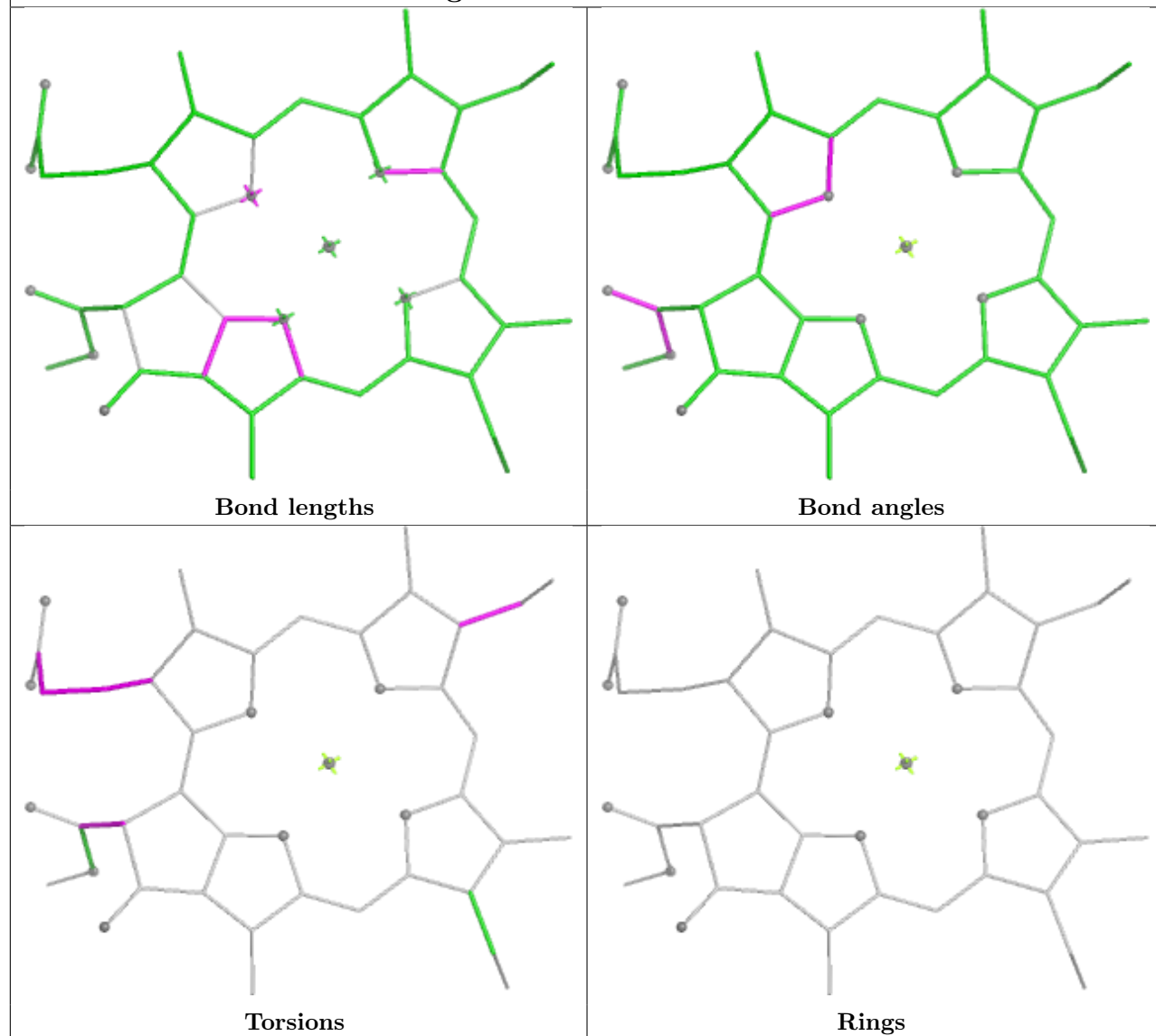
Ligand NDP YQ 501

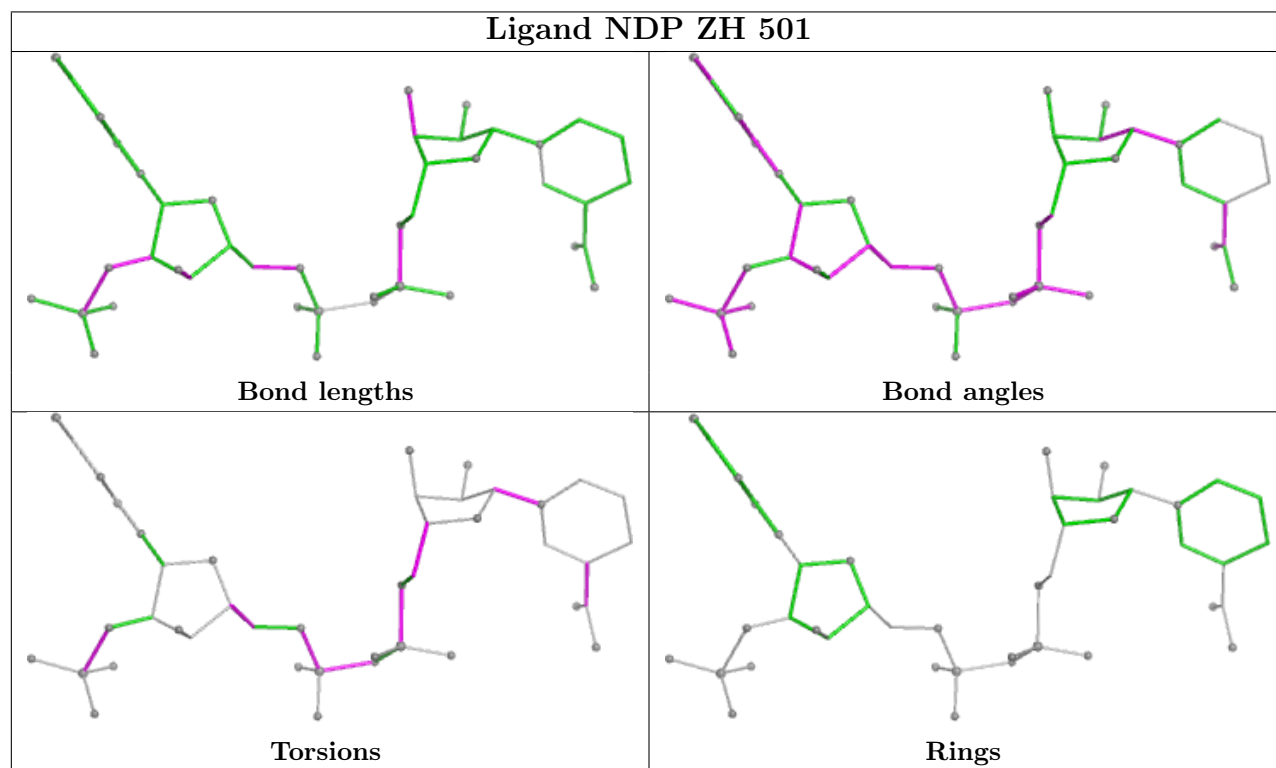
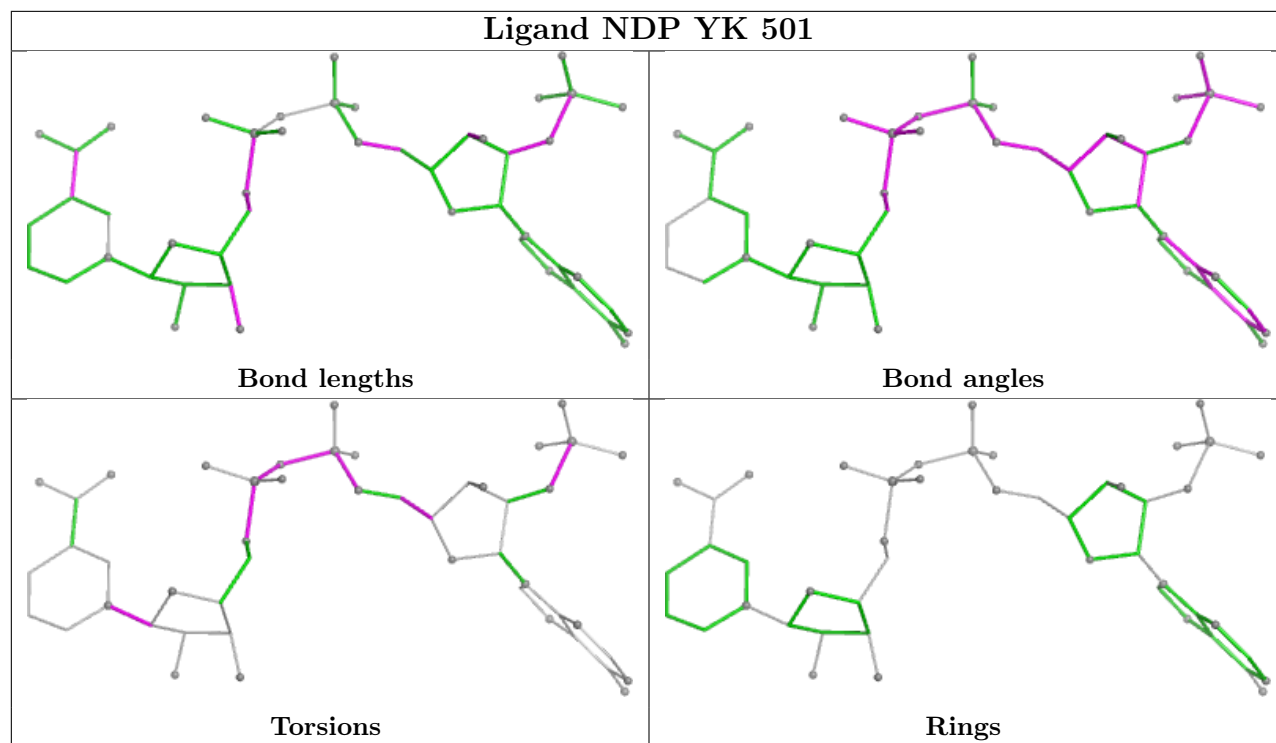


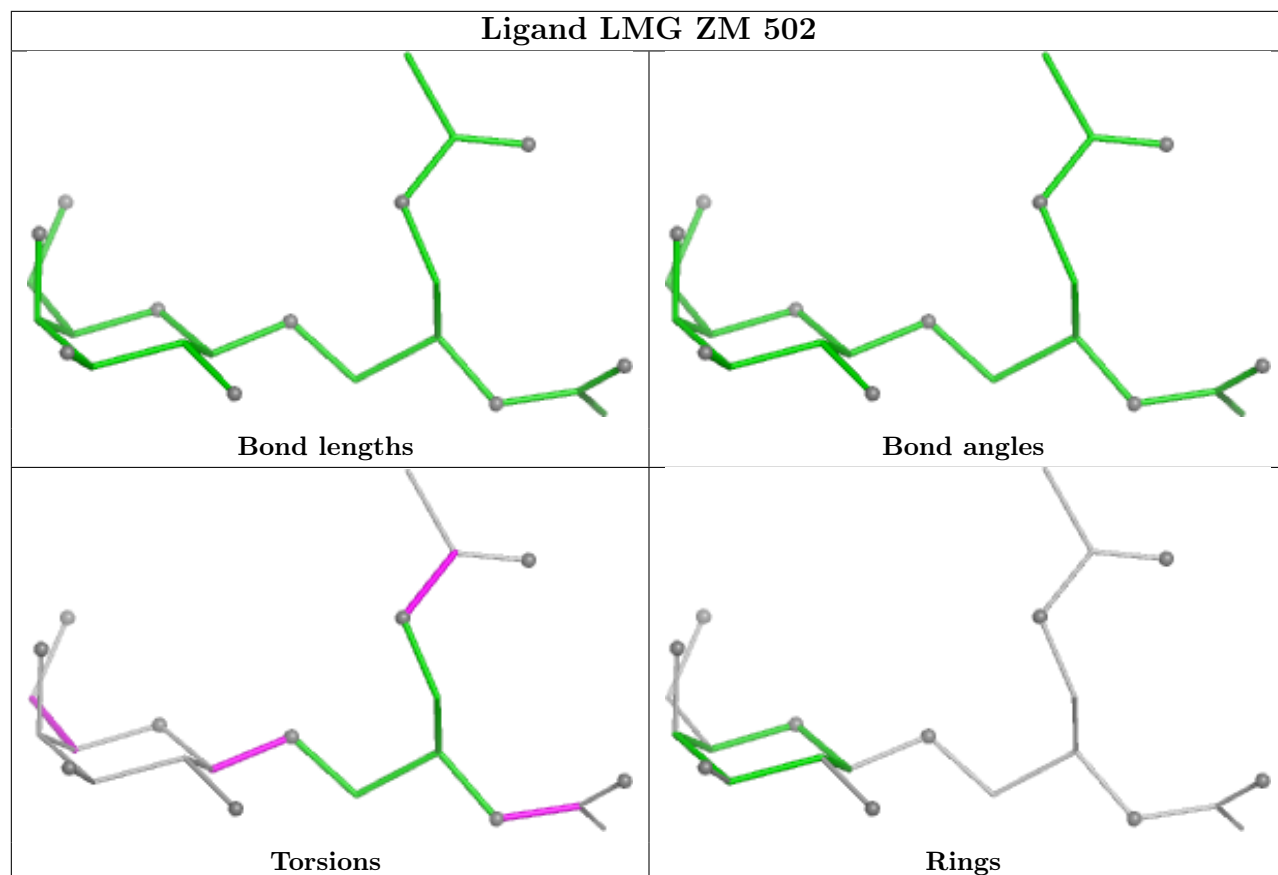


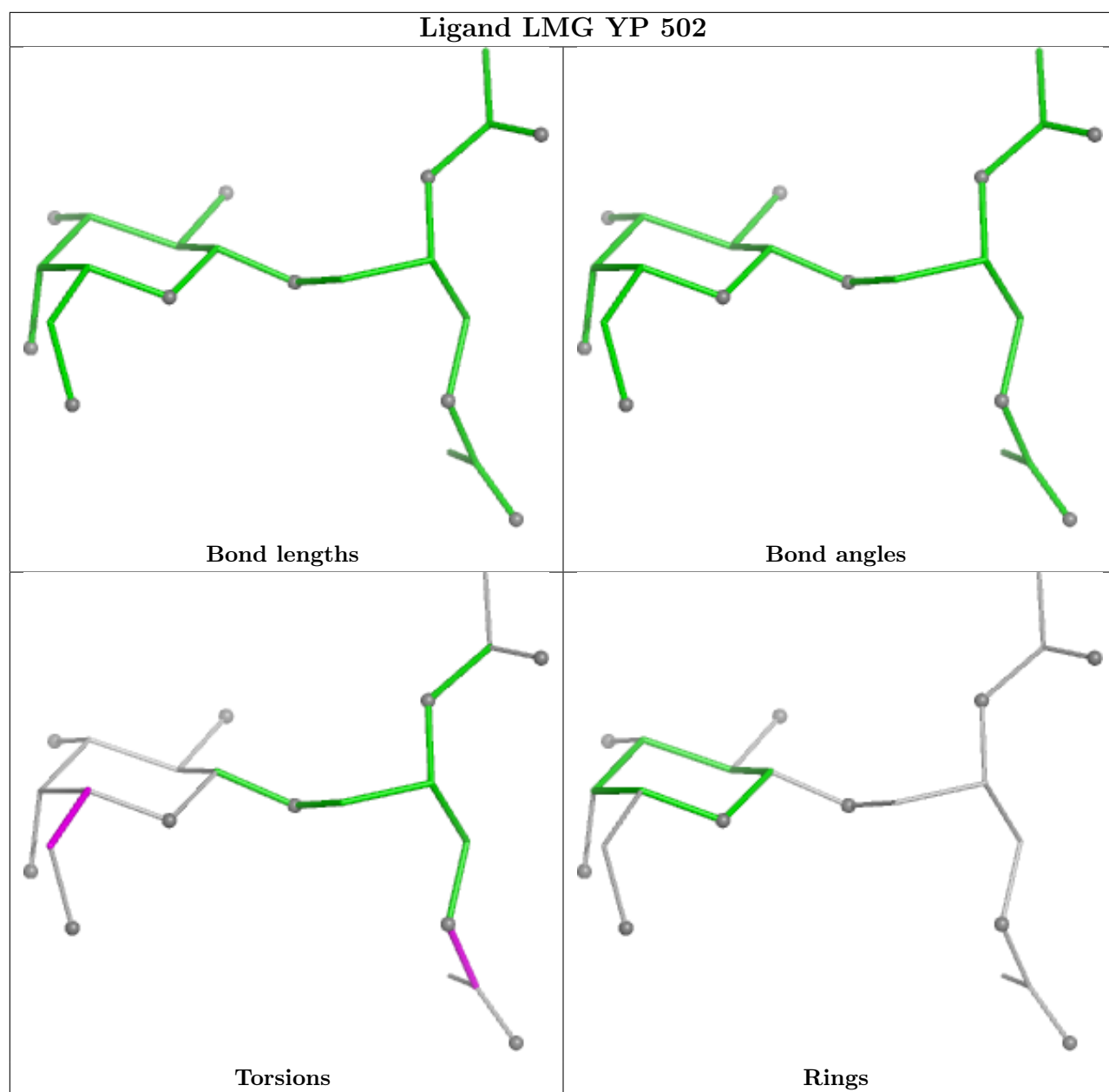


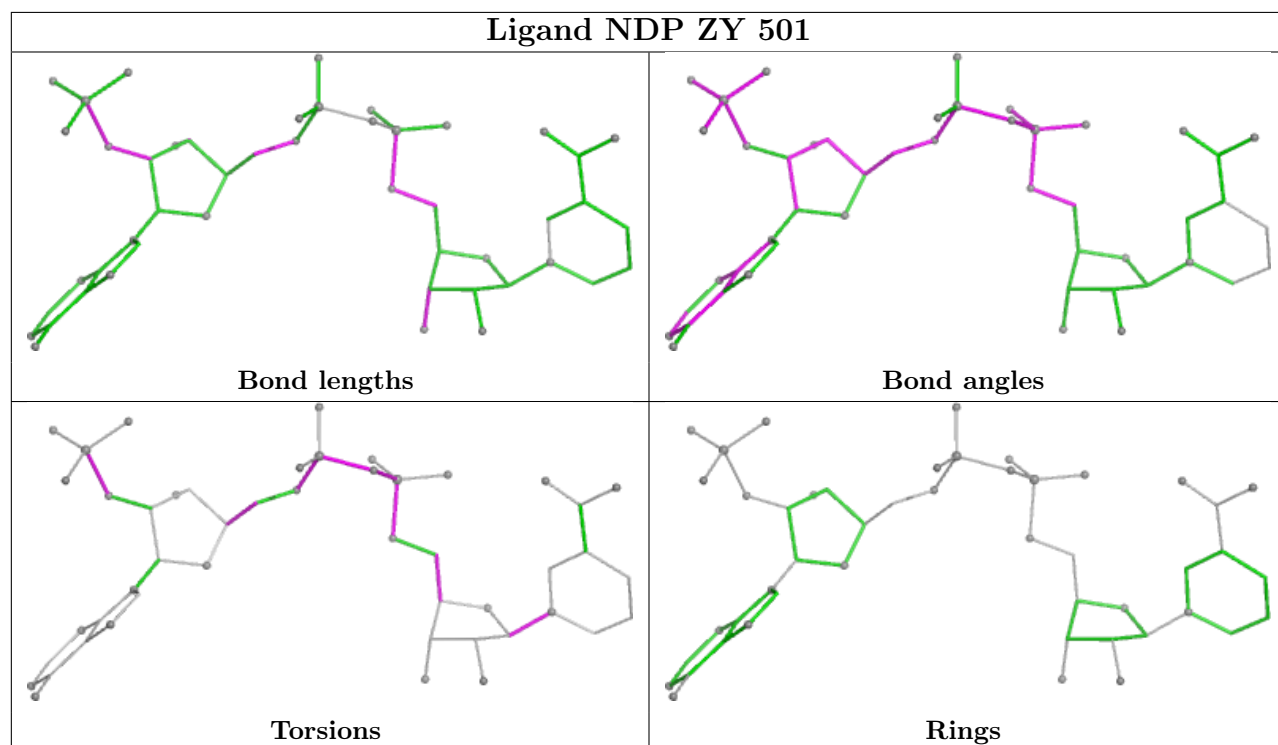
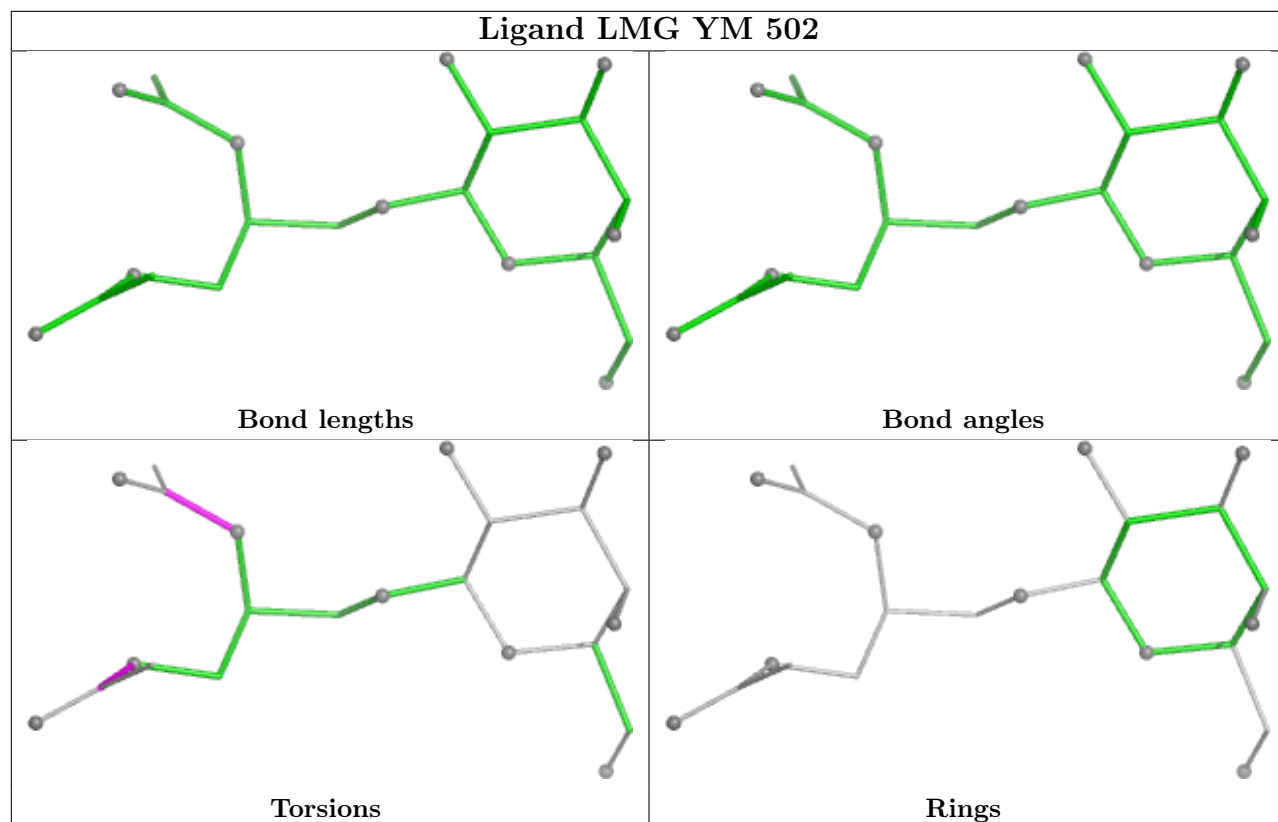
Ligand A1JWG ZG 503

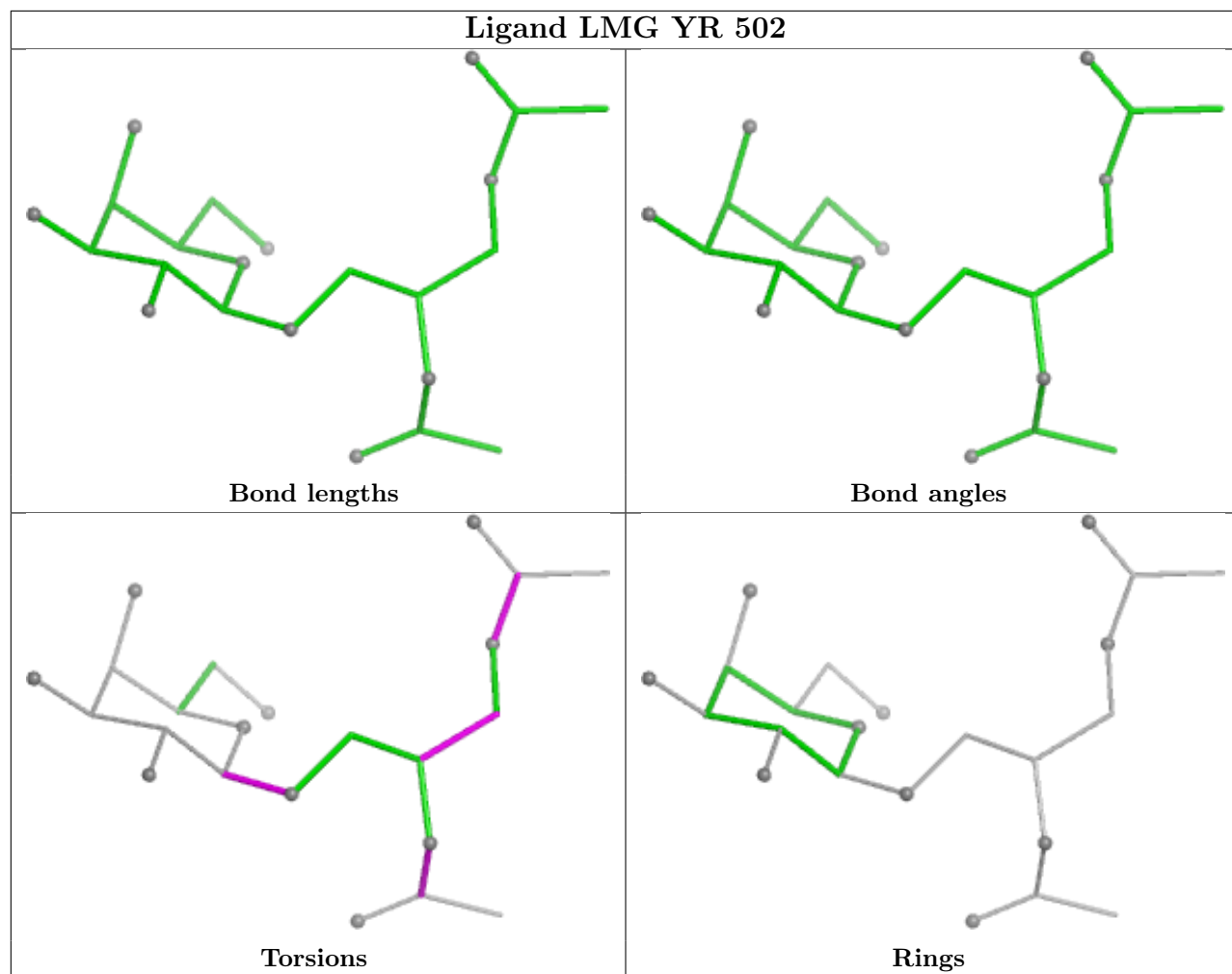


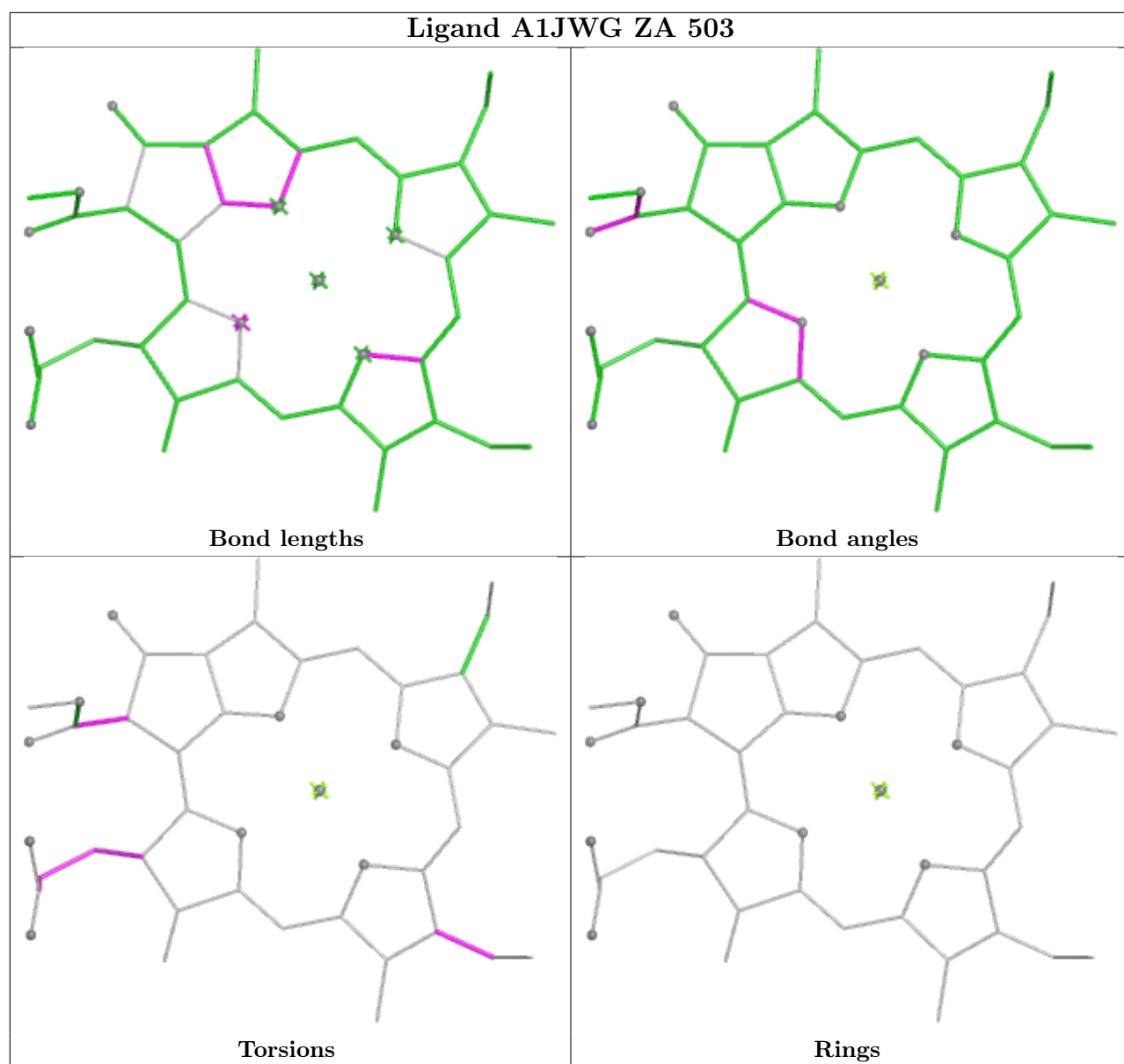




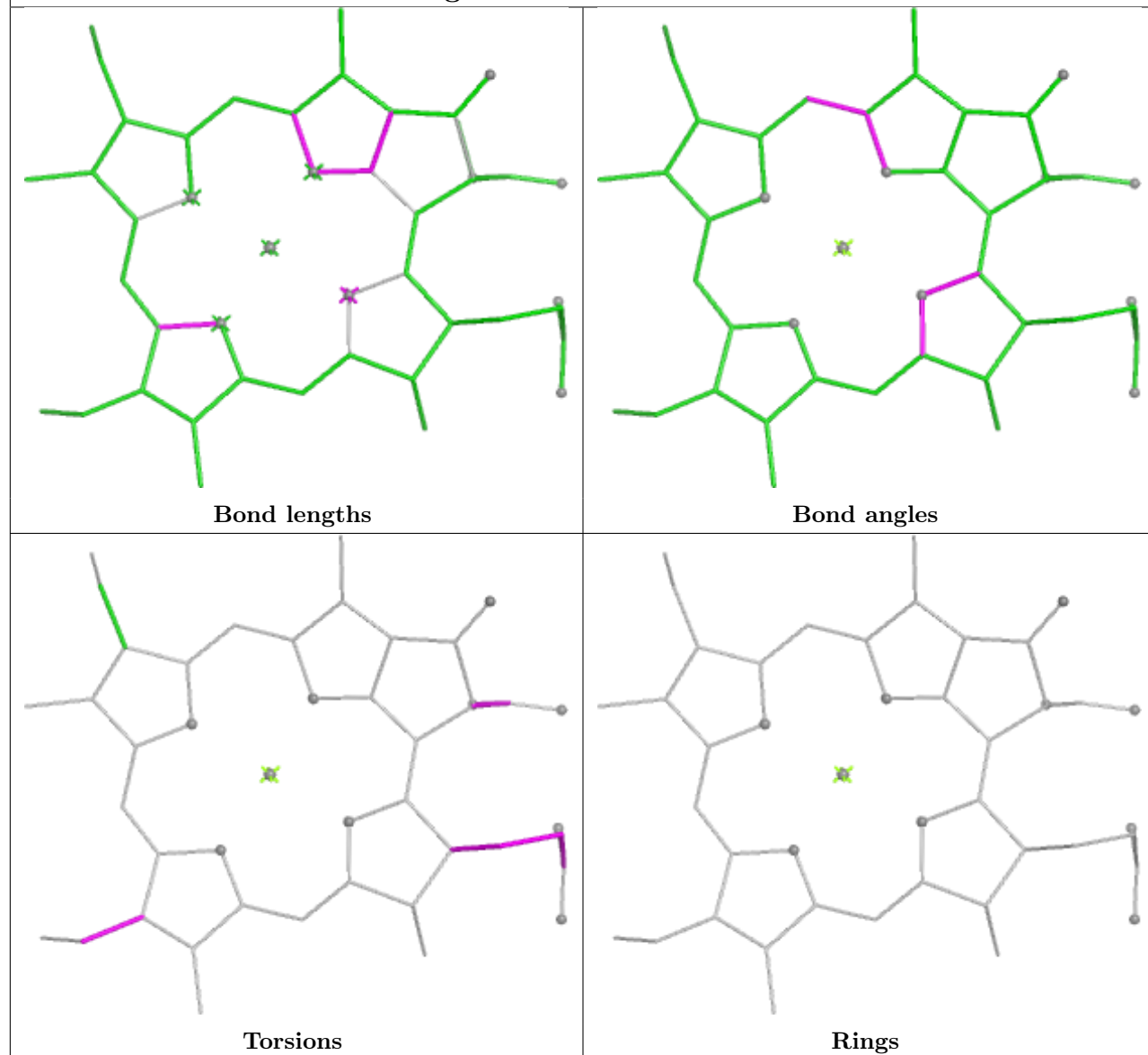


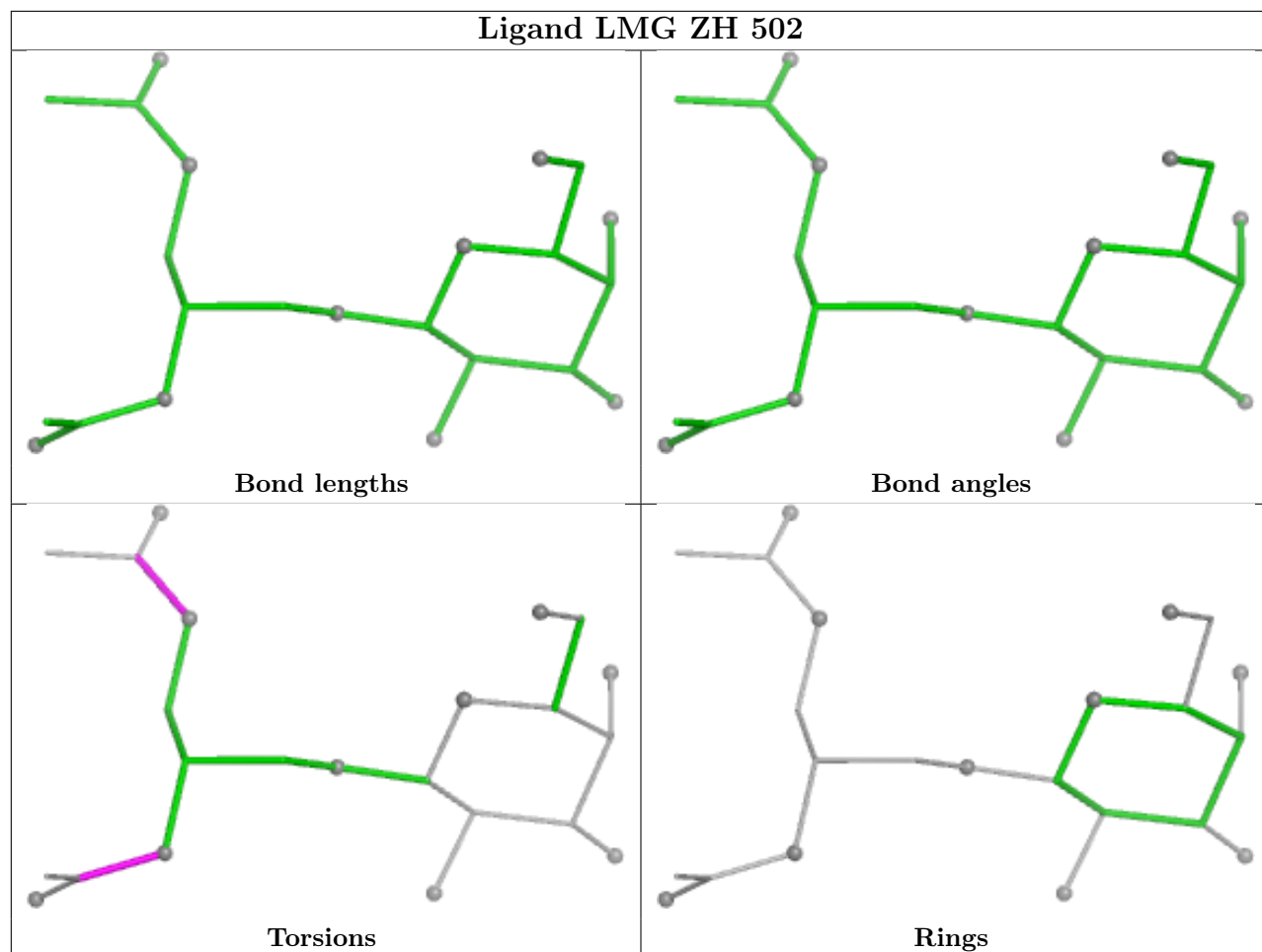




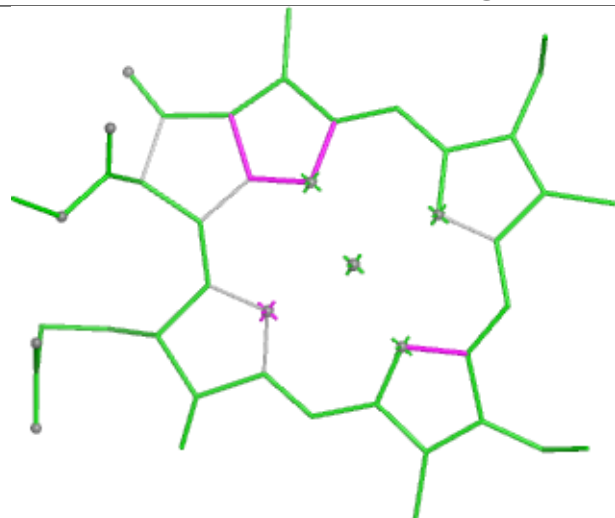


Ligand A1JWG YS 503

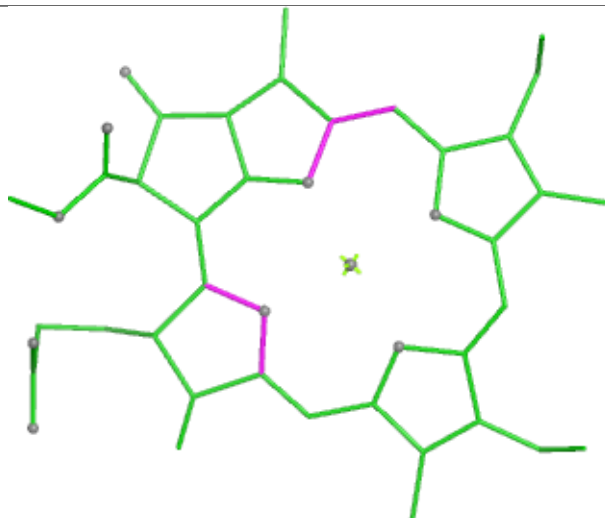




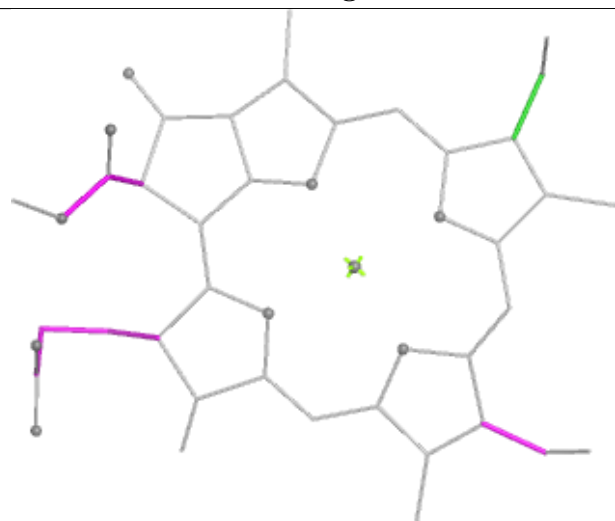
Ligand A1JWG ZU 503



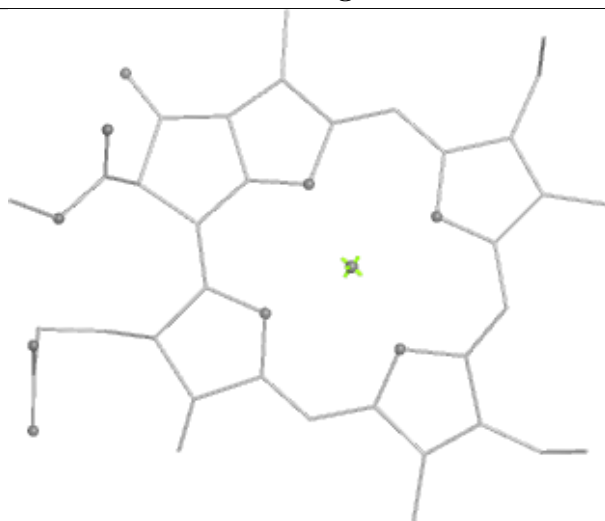
Bond lengths



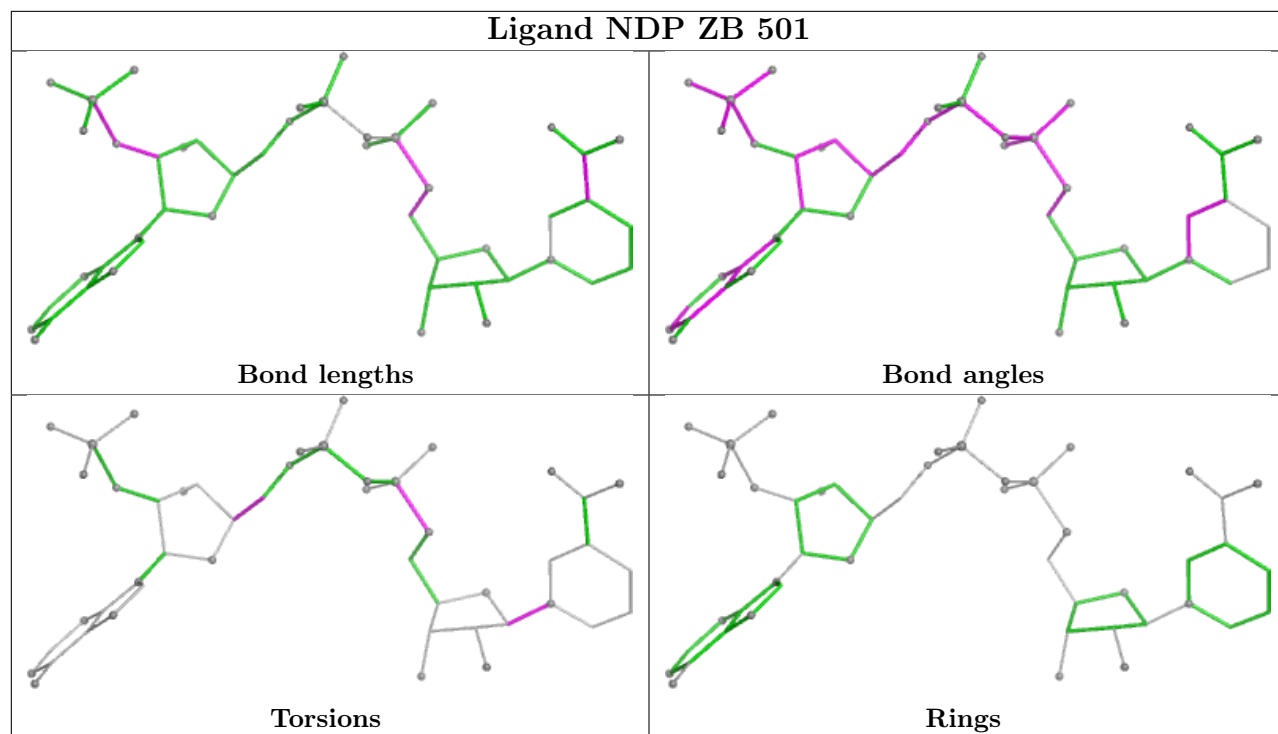
Bond angles



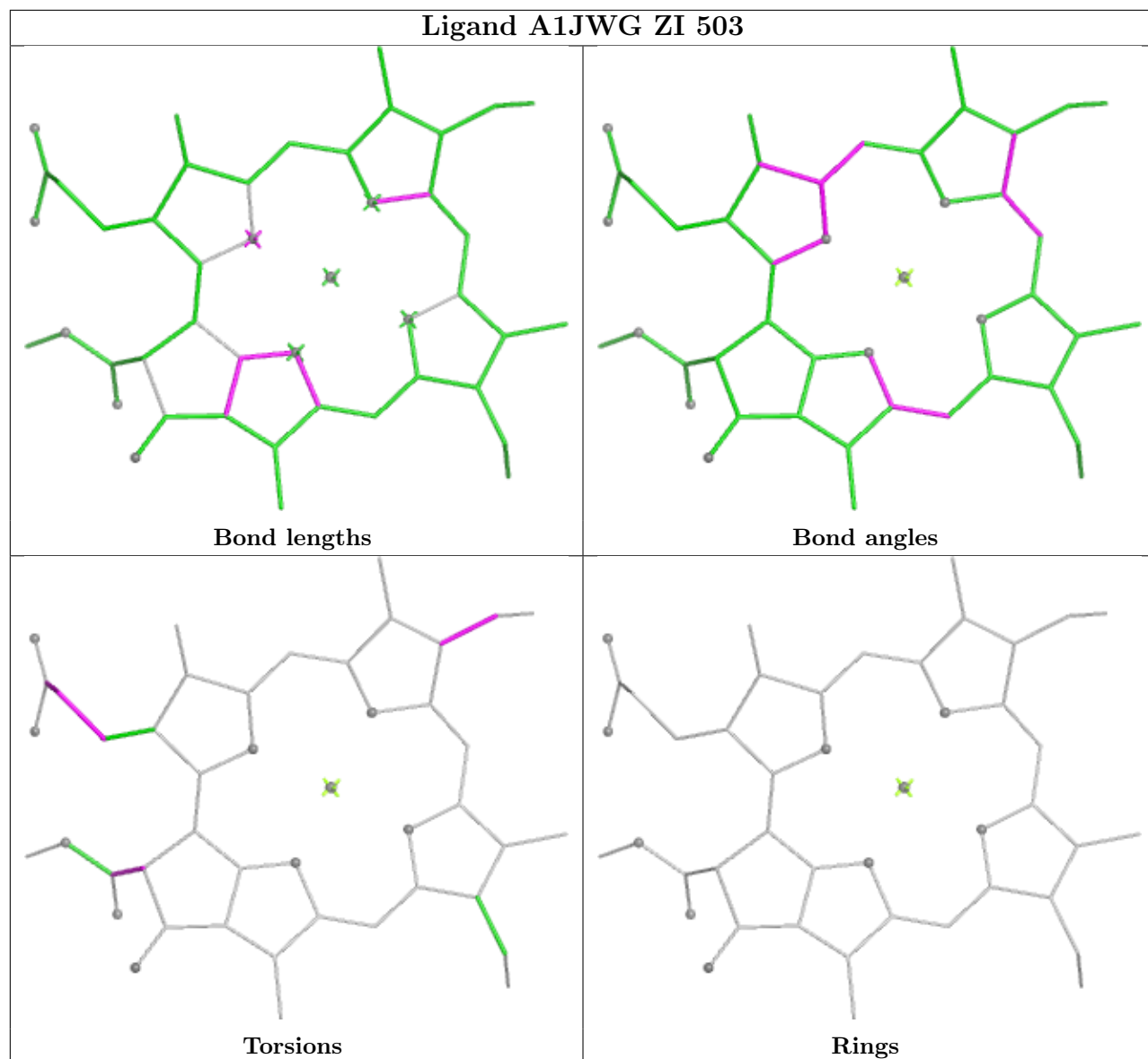
Torsions

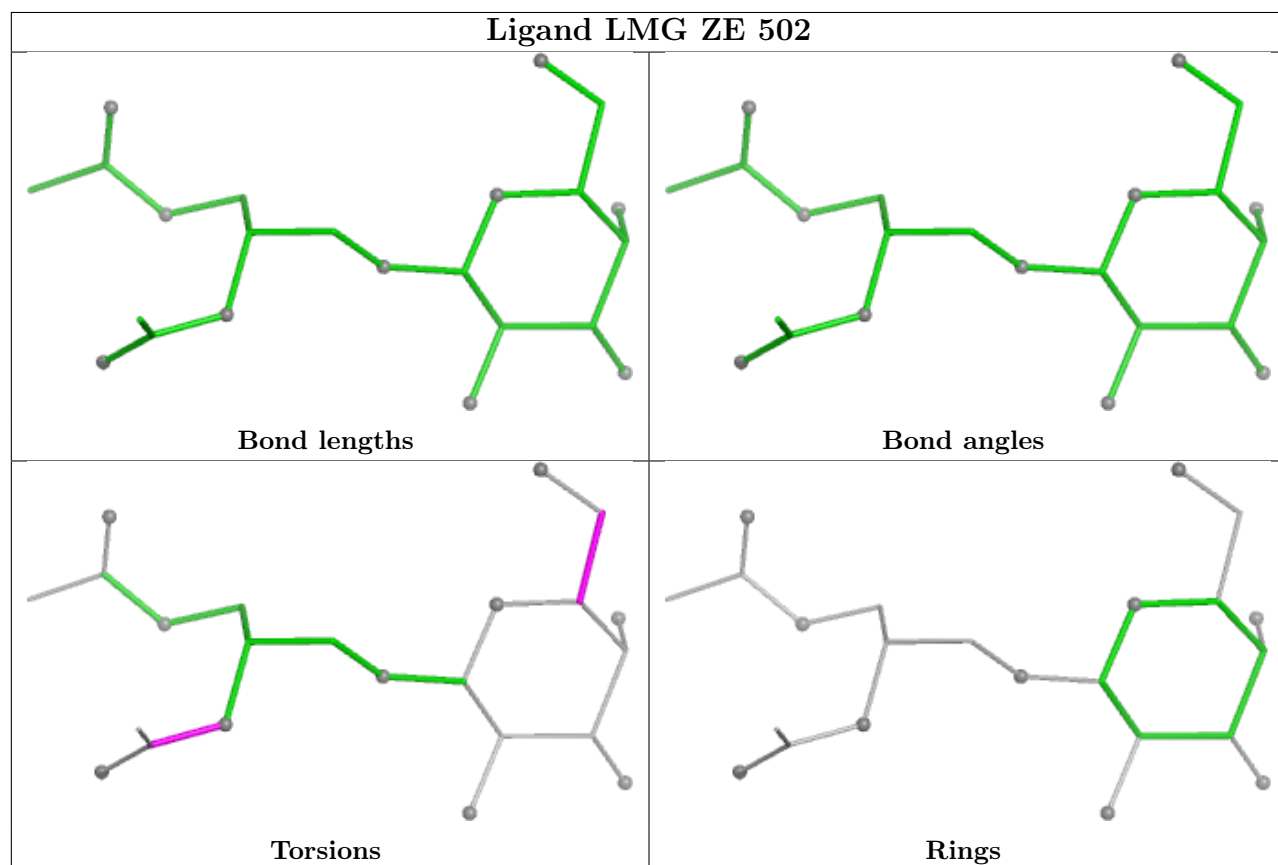
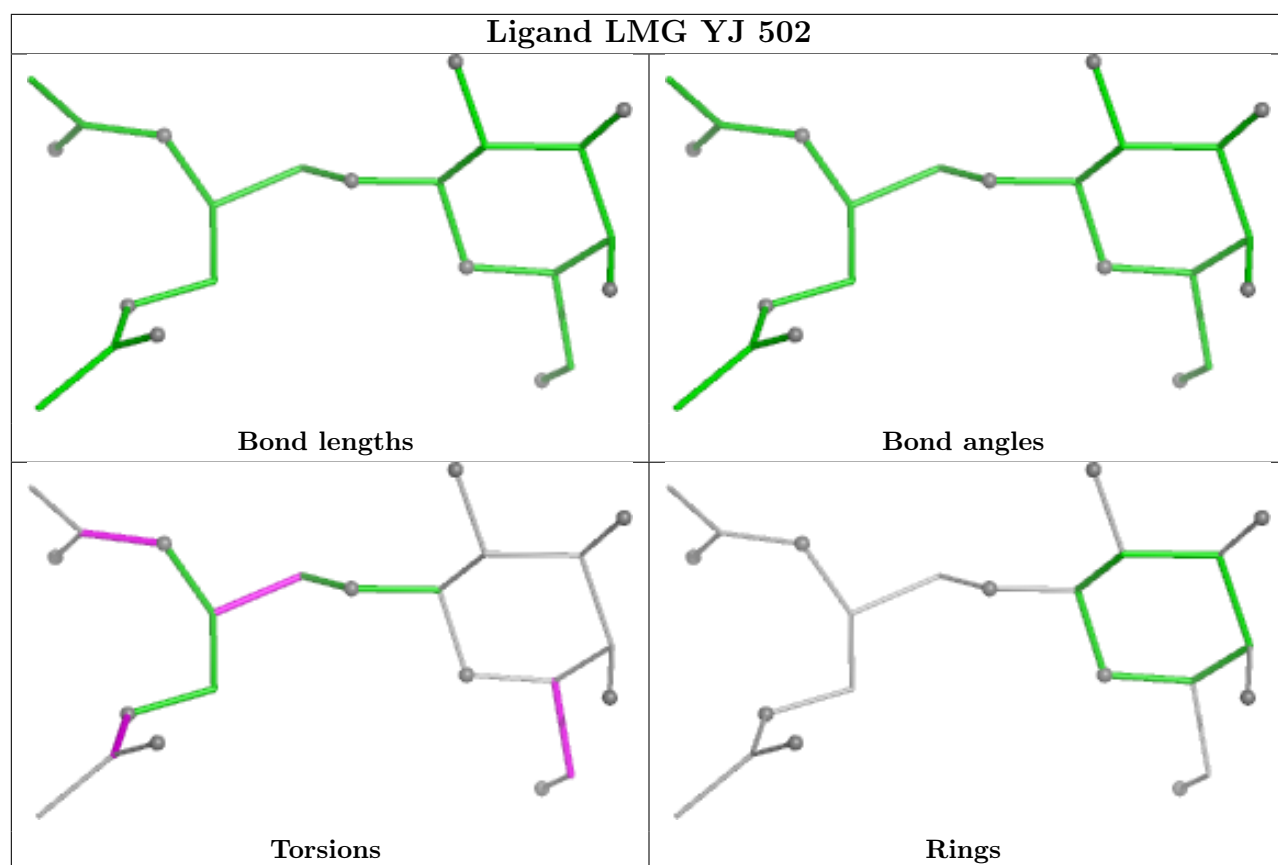


Rings

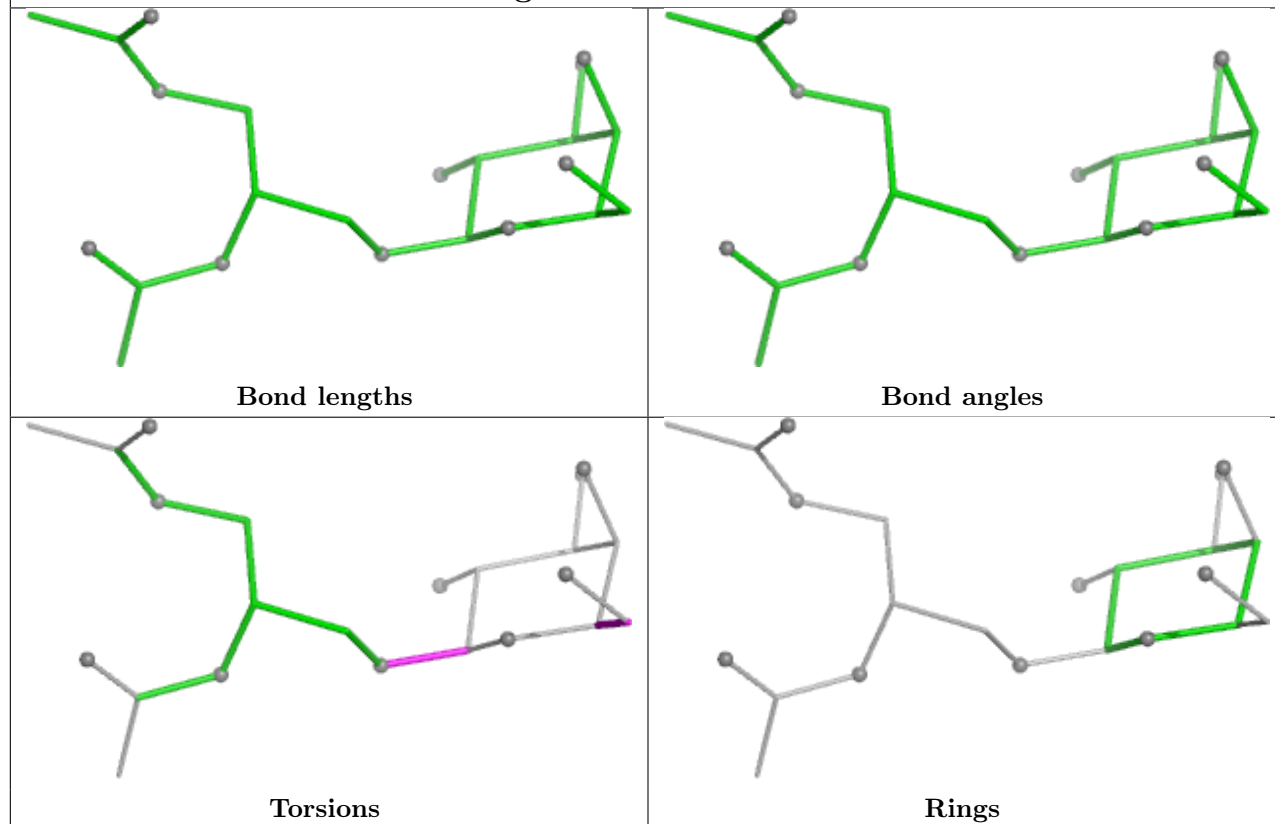


Ligand A1JWG ZI 503

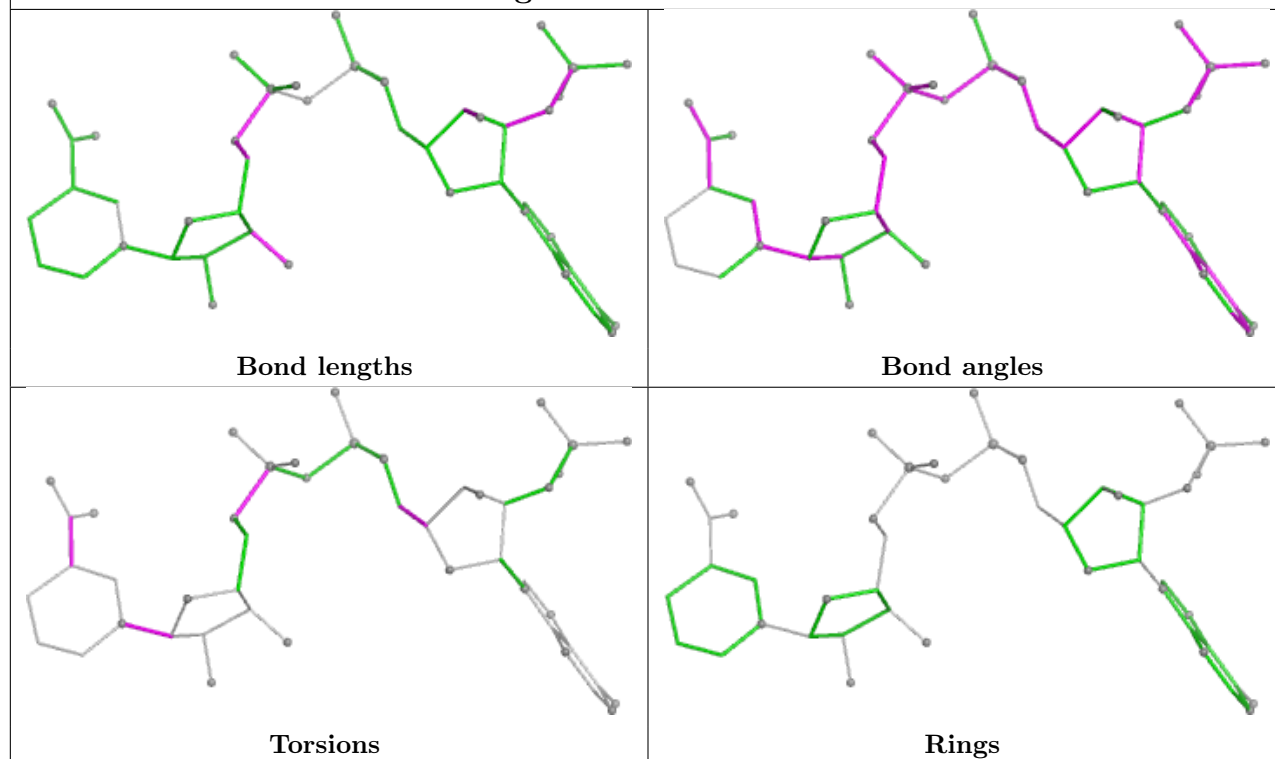




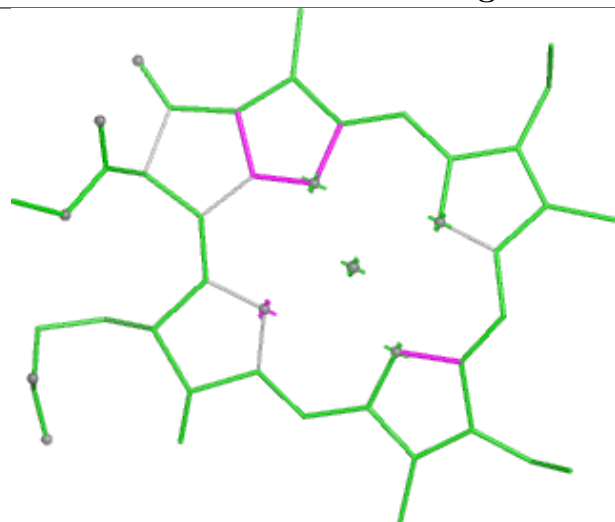
Ligand LMG ZN 502



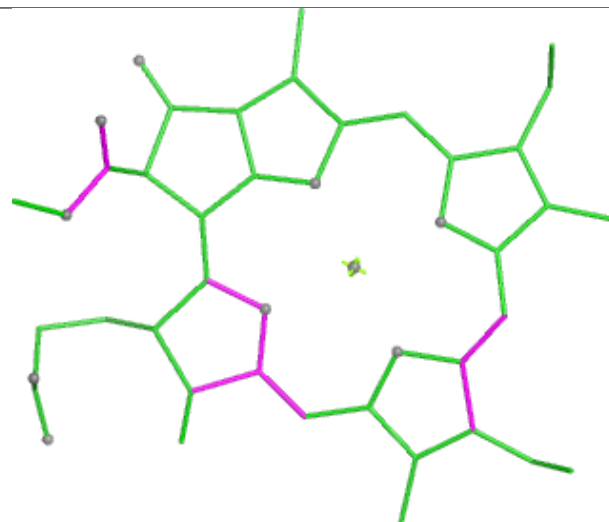
Ligand NDP YR 501



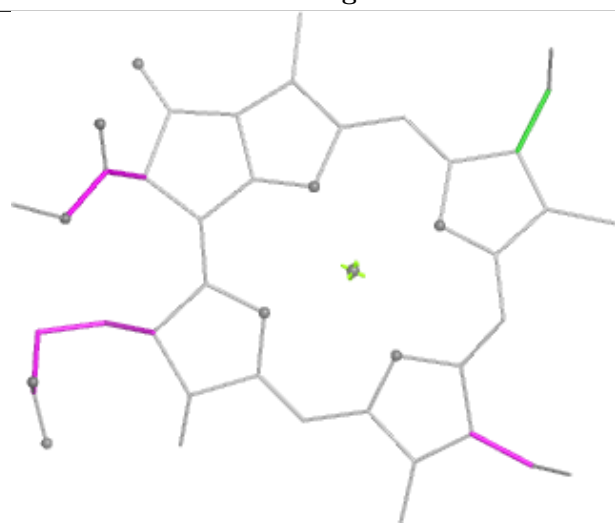
Ligand A1JWG ZD 503



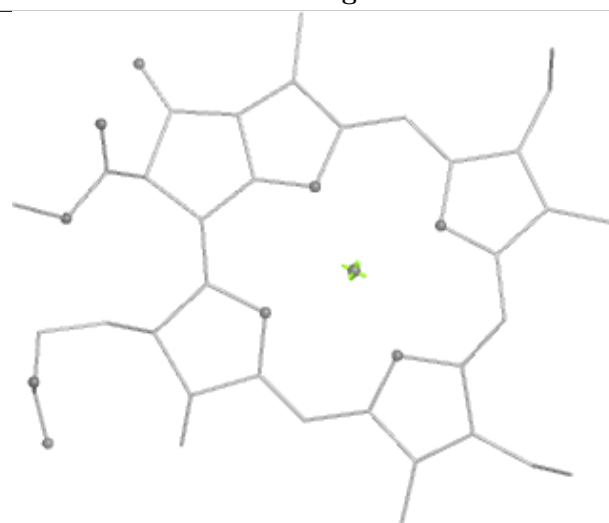
Bond lengths



Bond angles

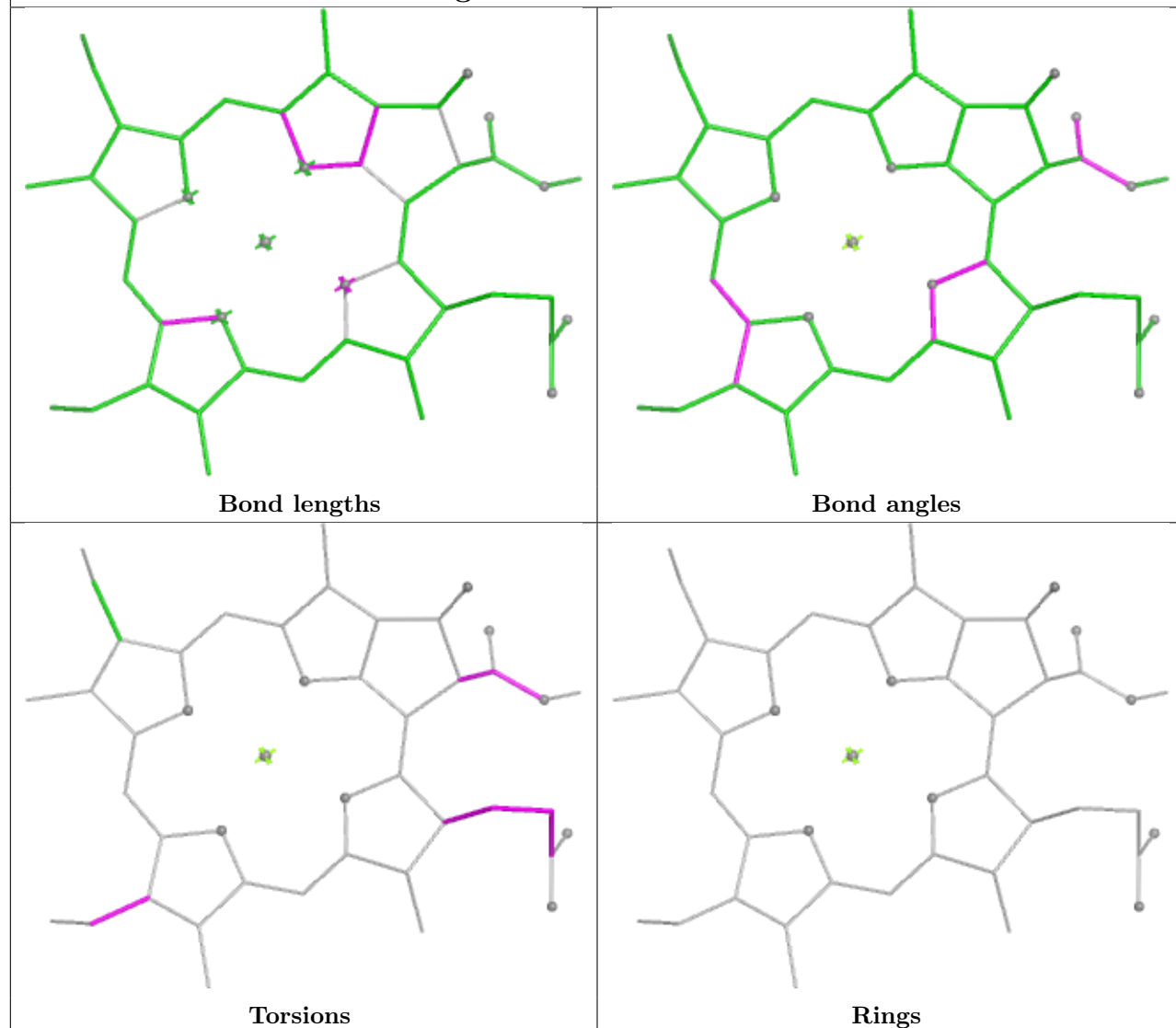


Torsions

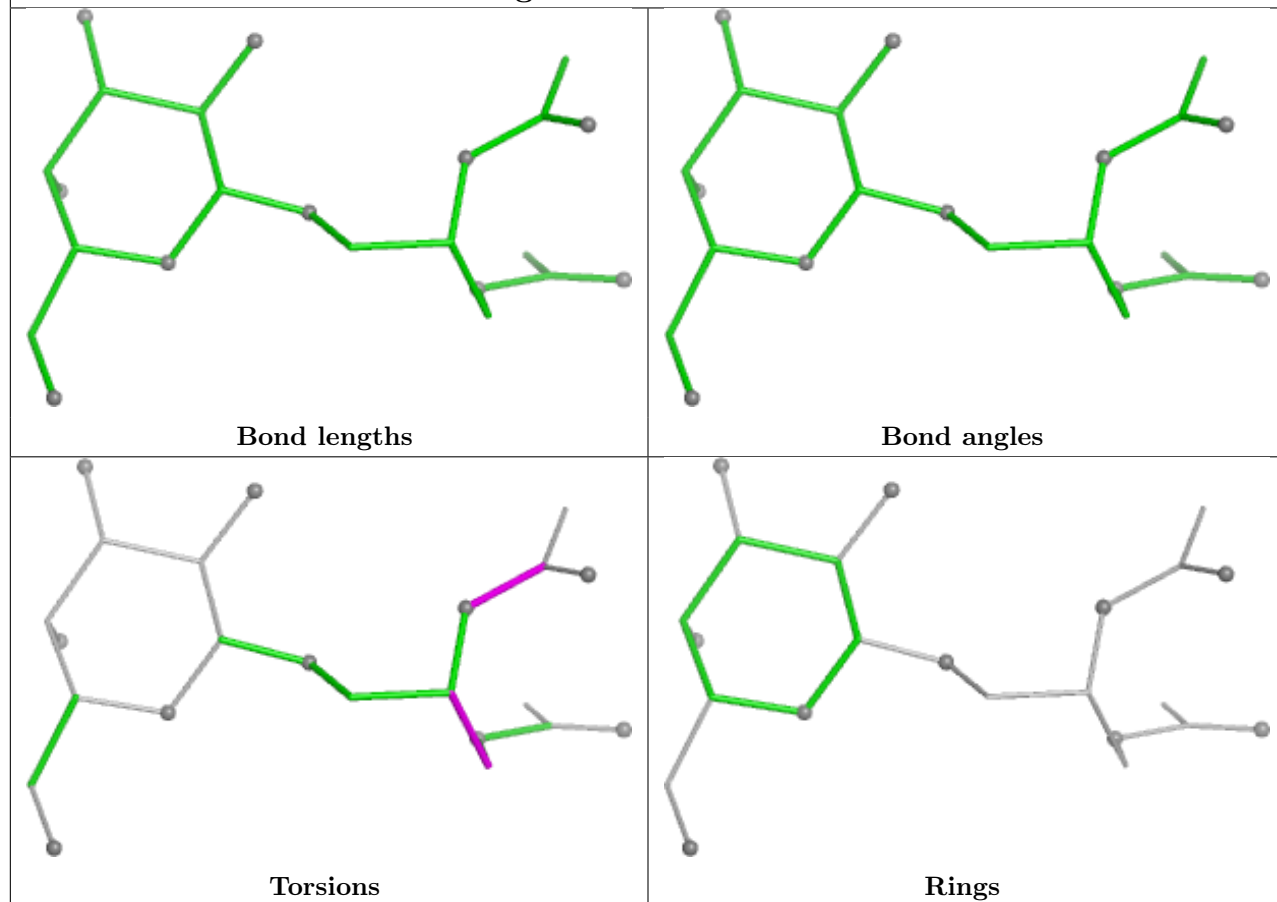


Rings

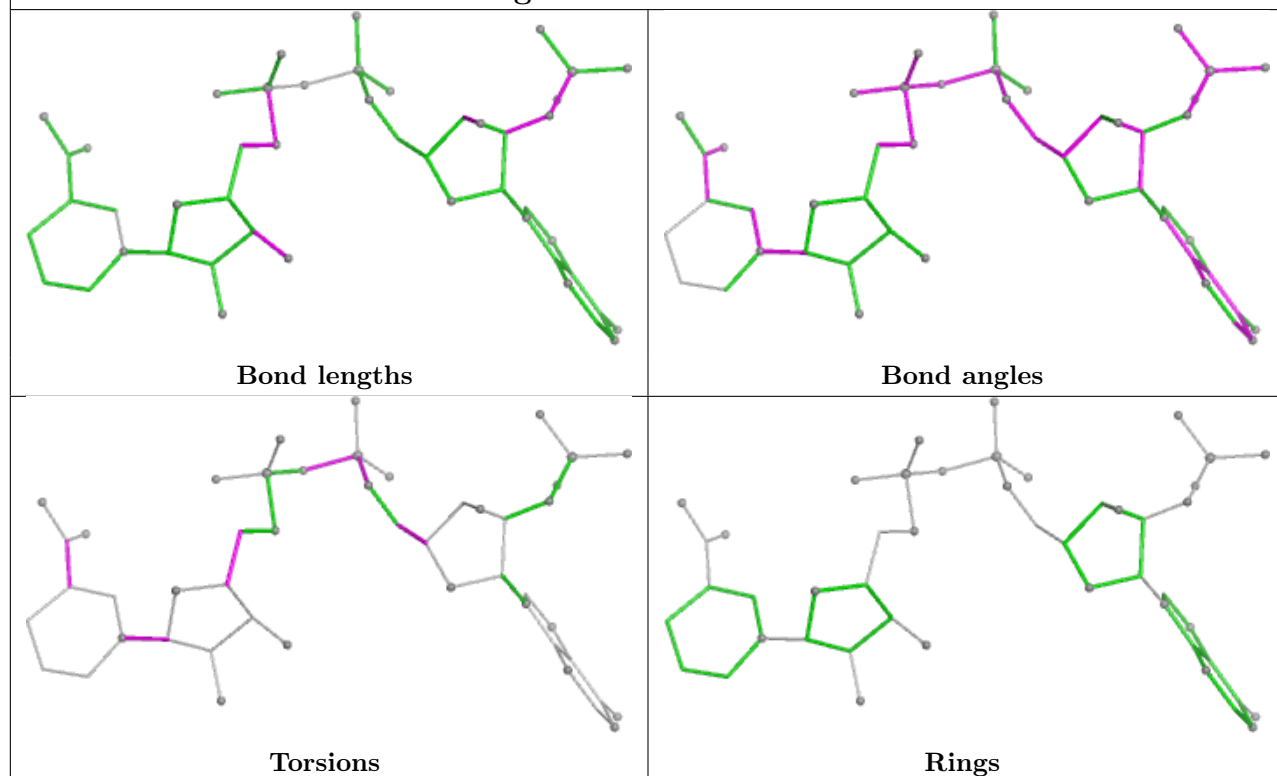
Ligand A1JWG ZC 503



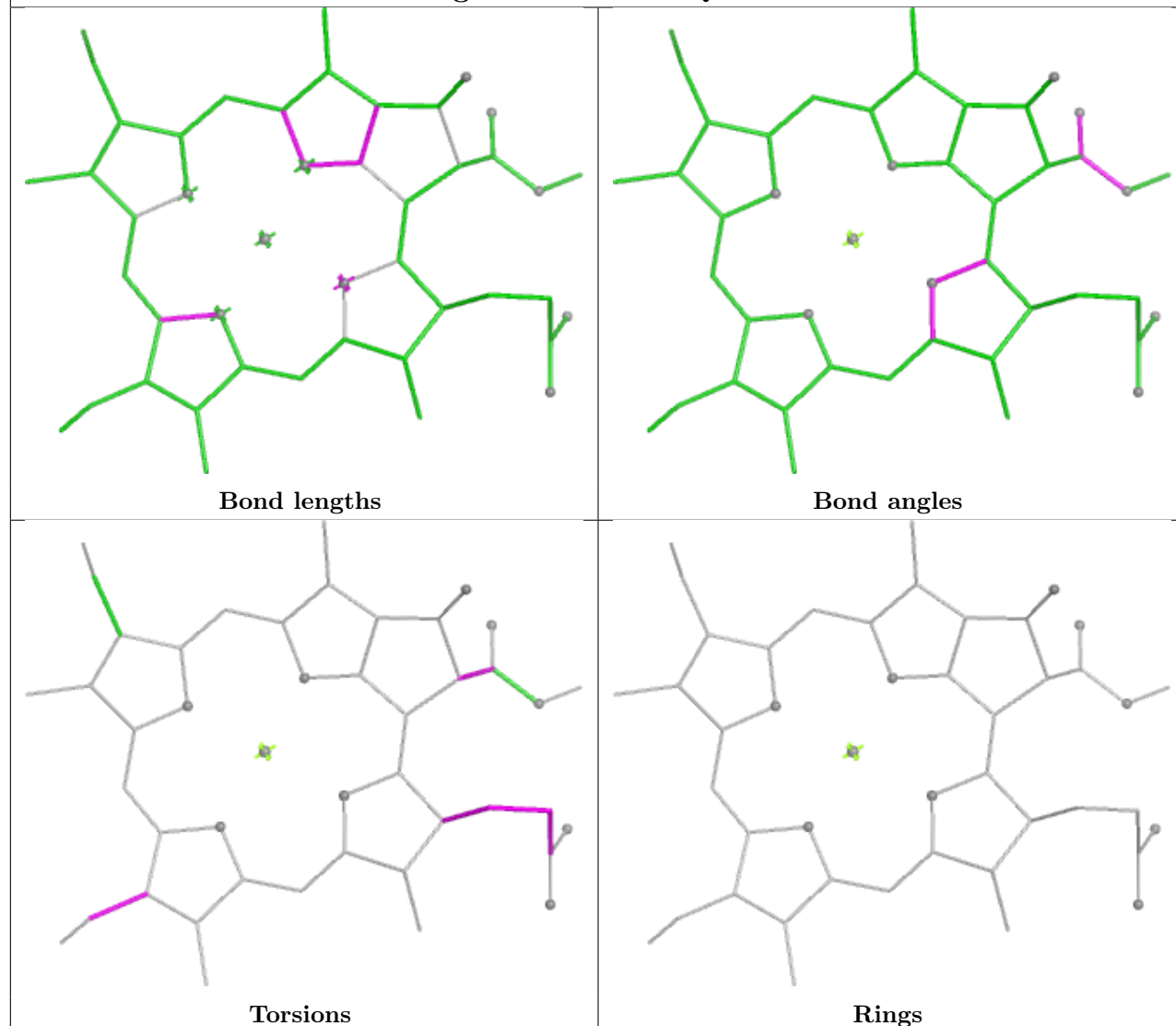
Ligand LMG ZO 502

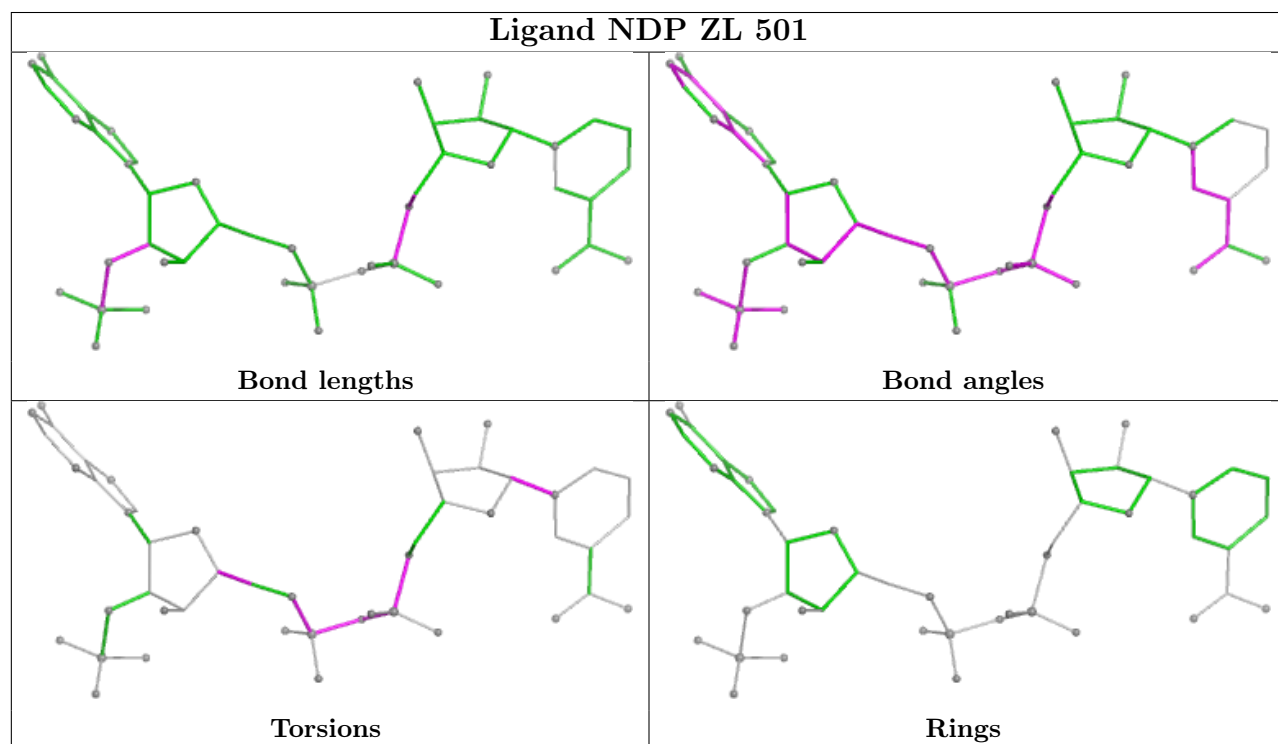
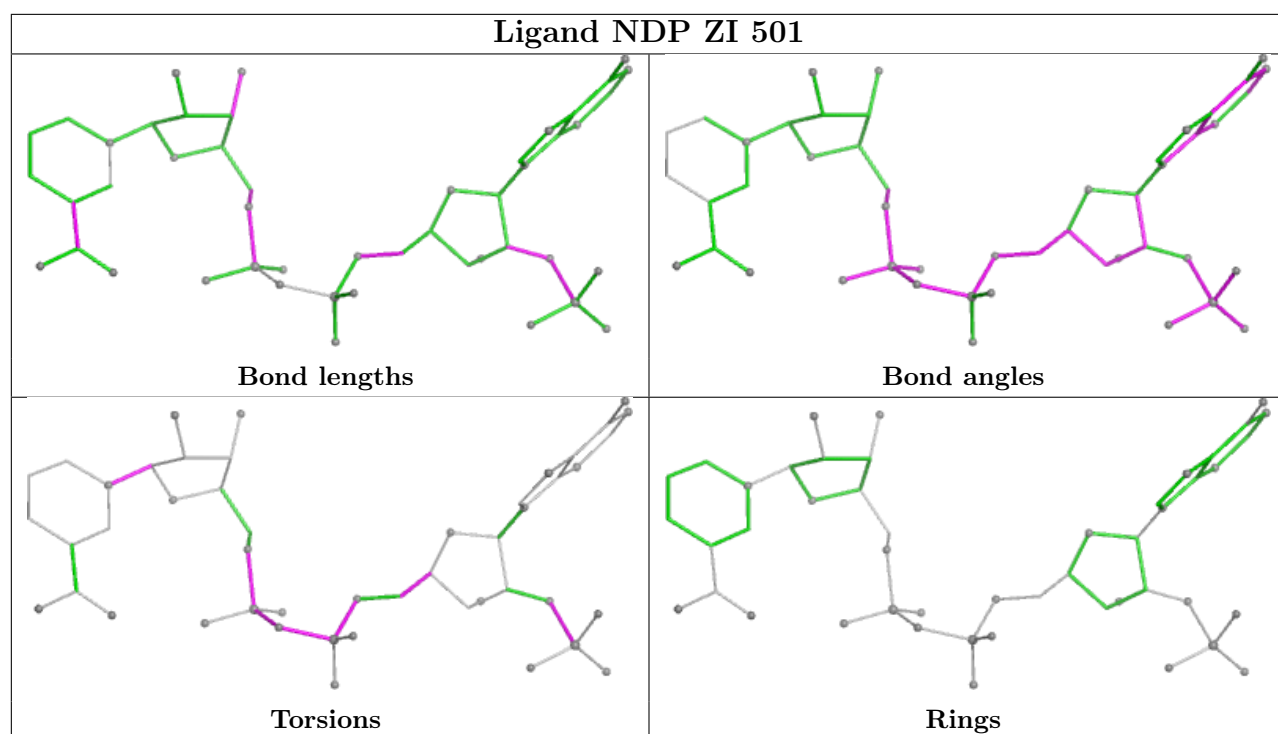


Ligand NDP YN 501

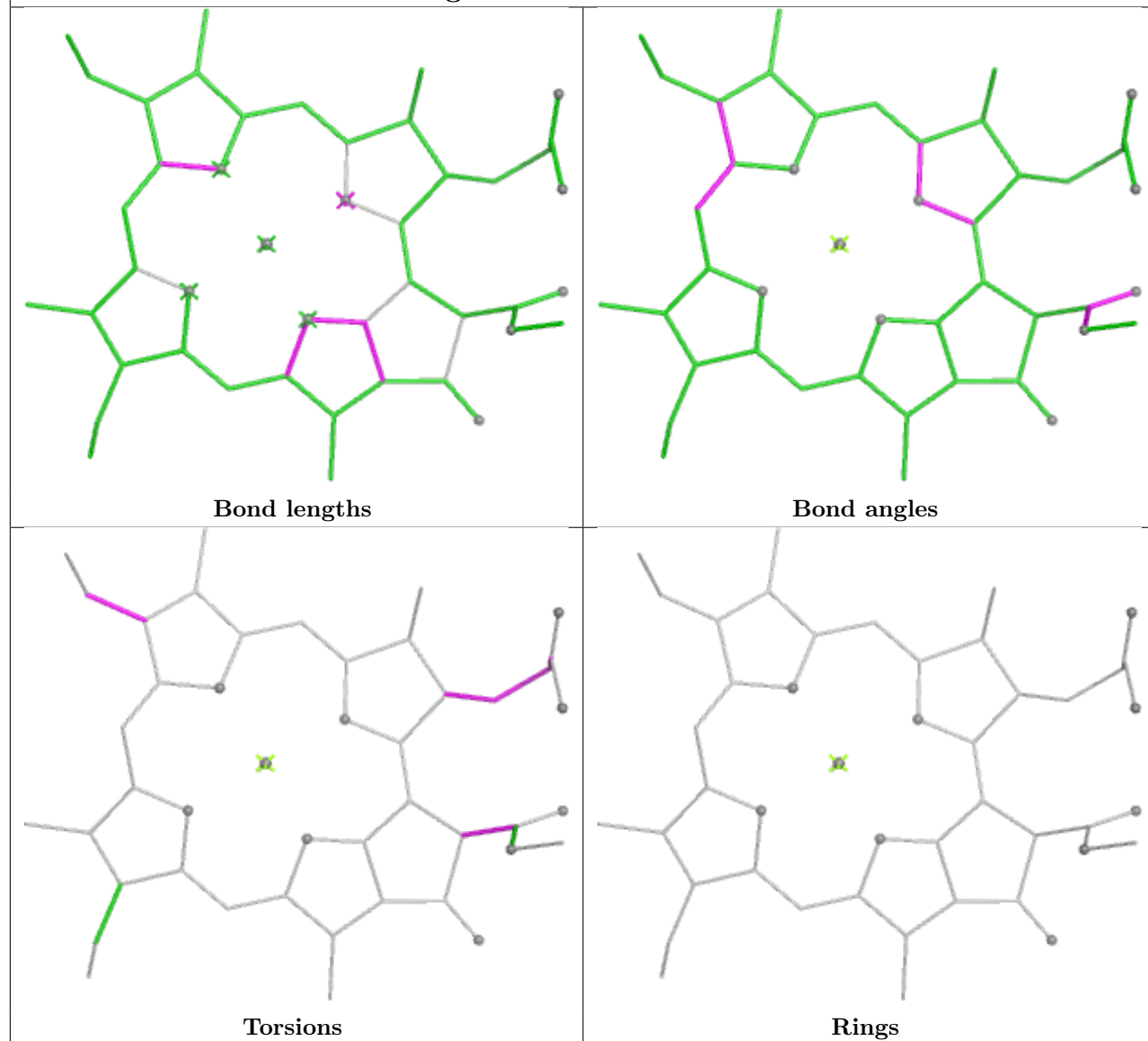


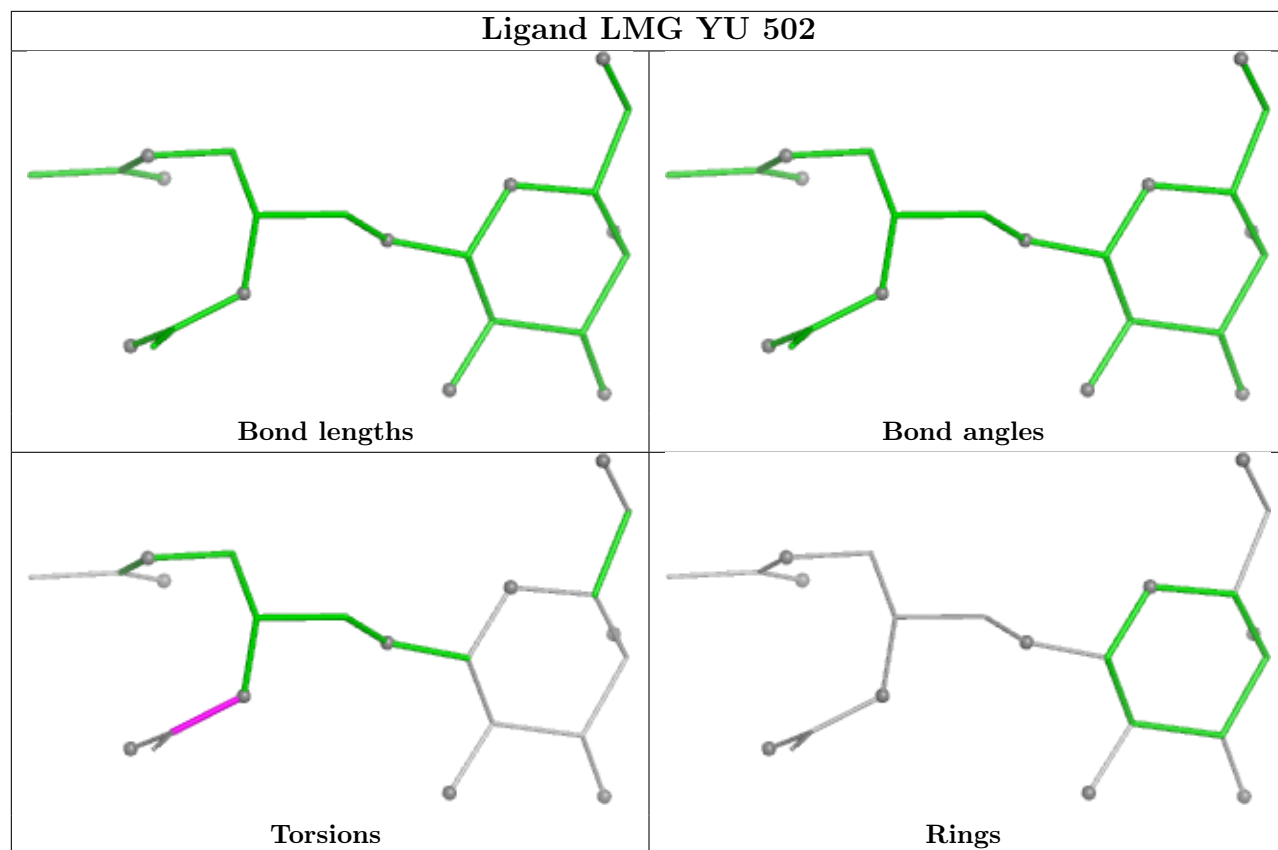
Ligand A1JWG YQ 503



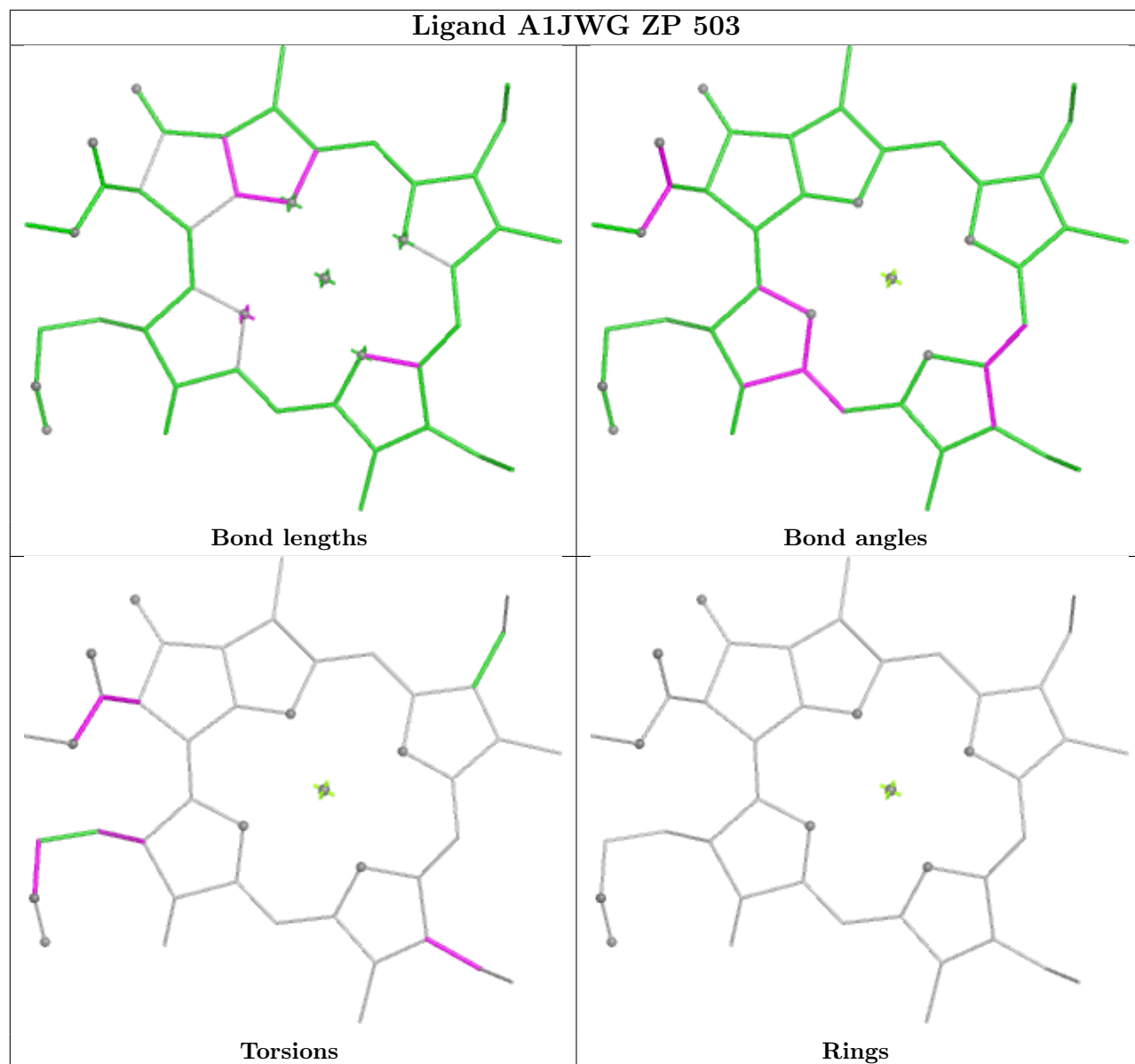


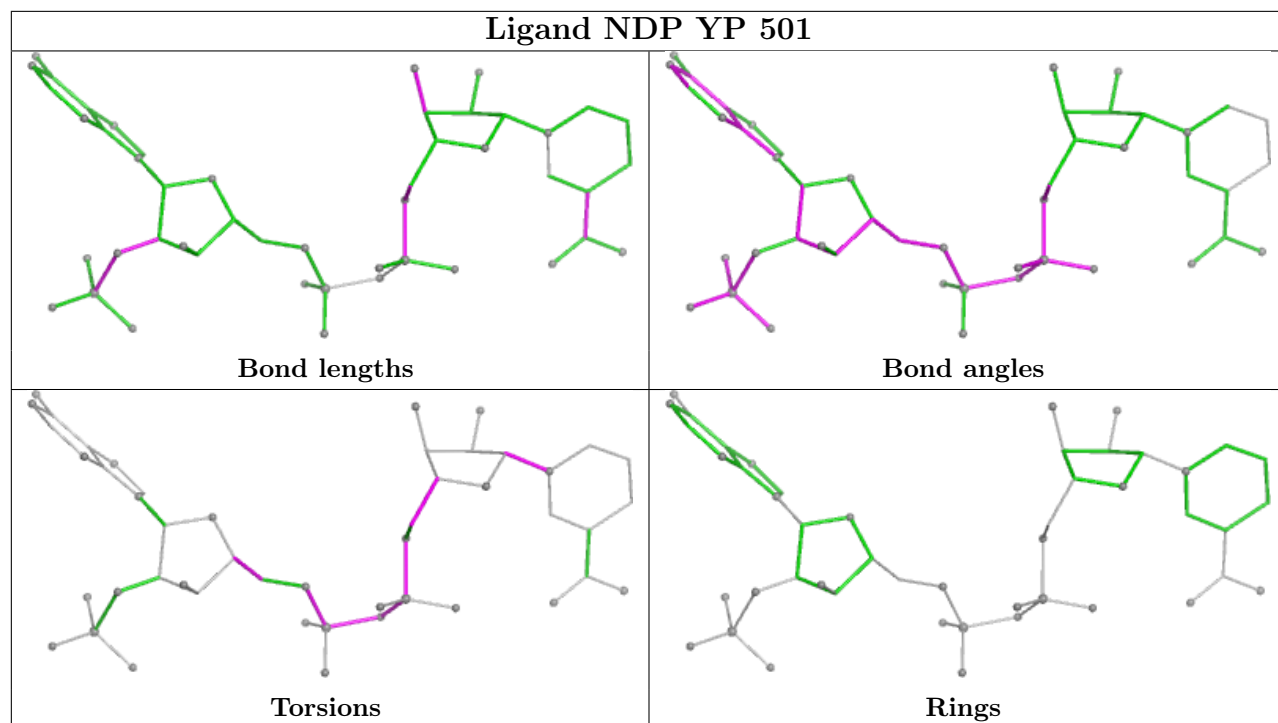
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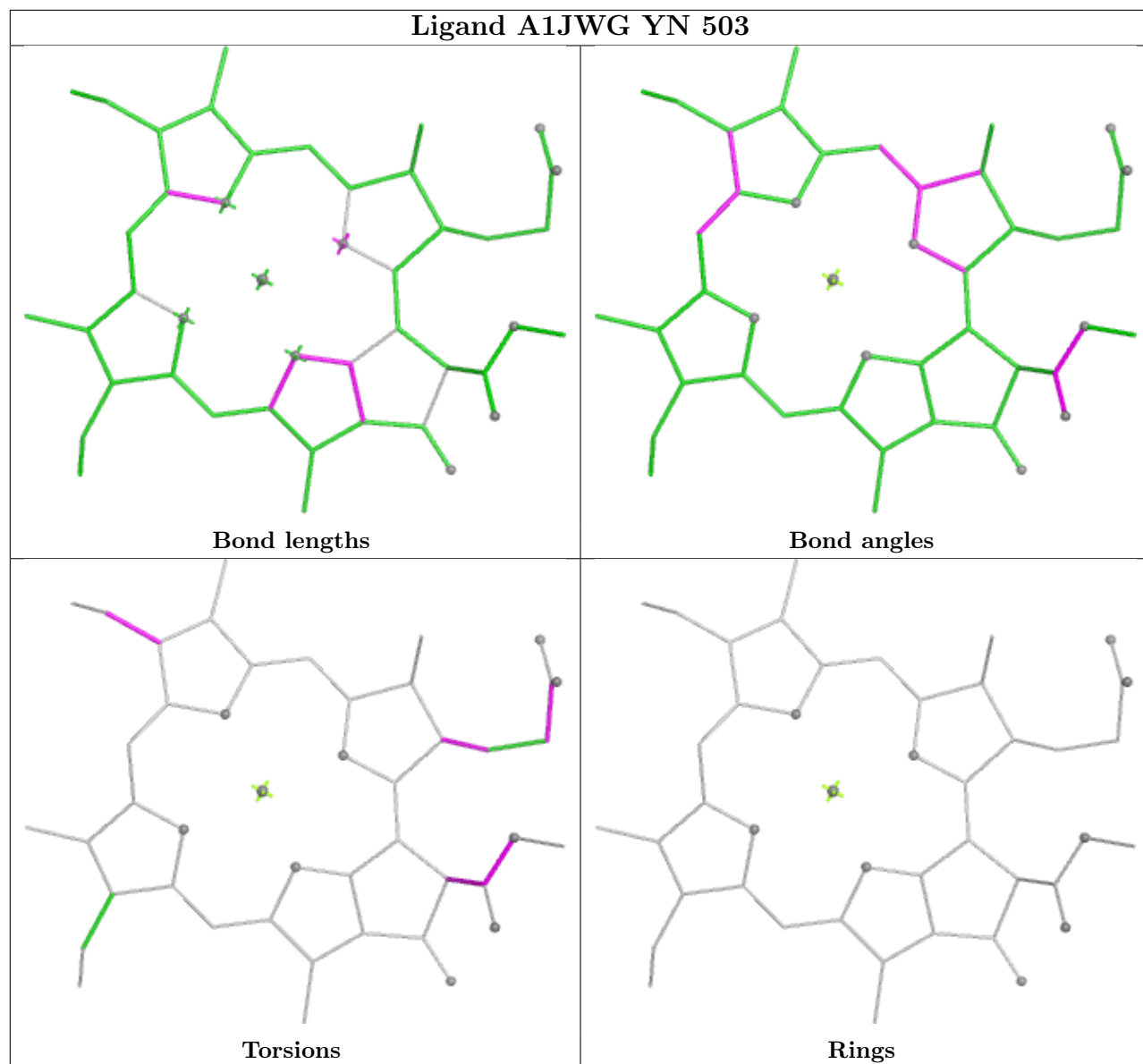


Ligand A1JWG ZP 503

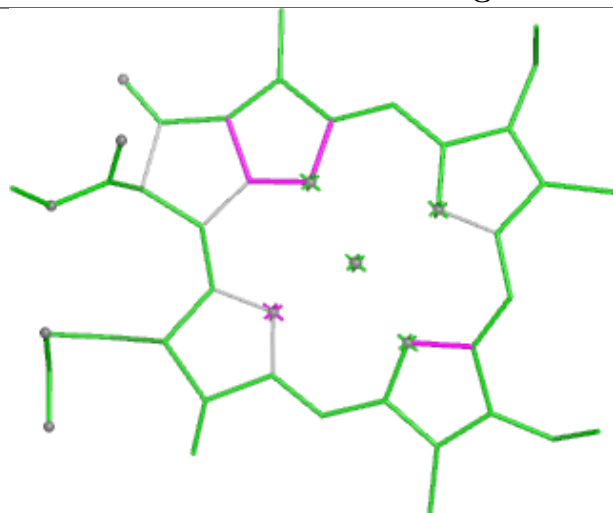




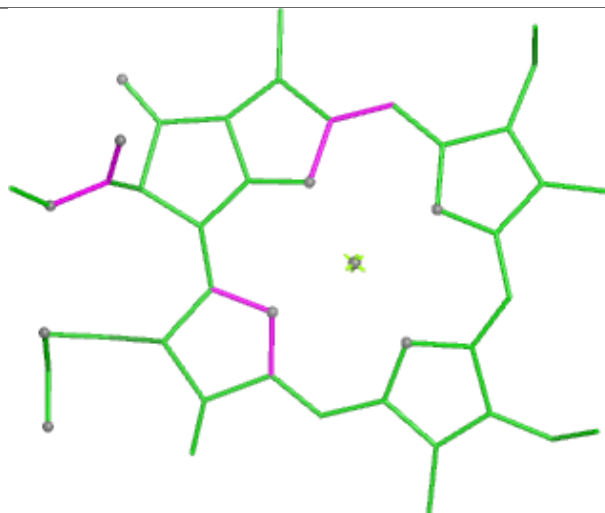
Ligand A1JWG YN 503



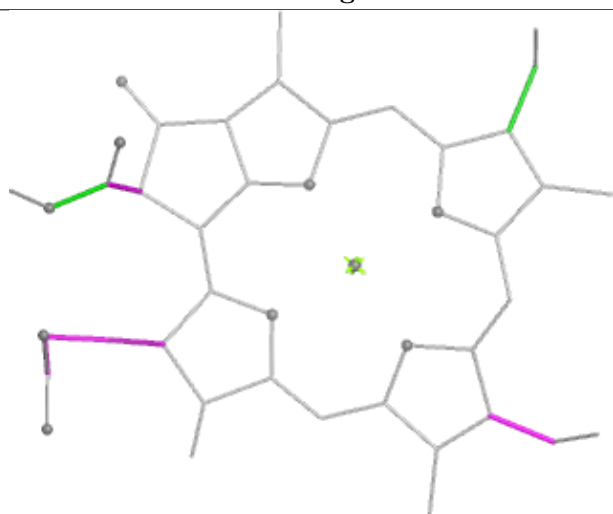
Ligand A1JWG YX 503



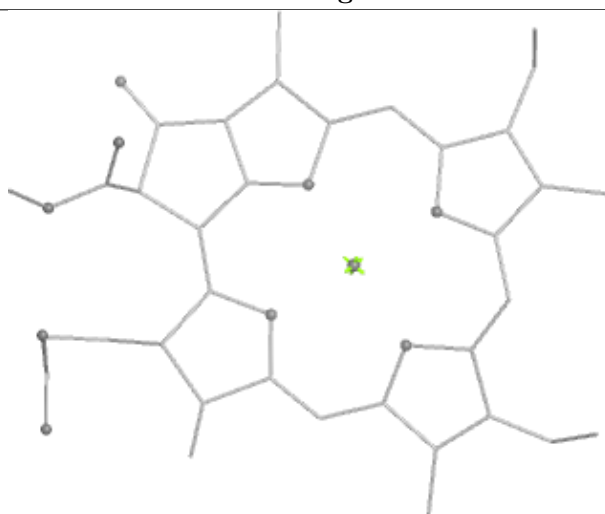
Bond lengths



Bond angles

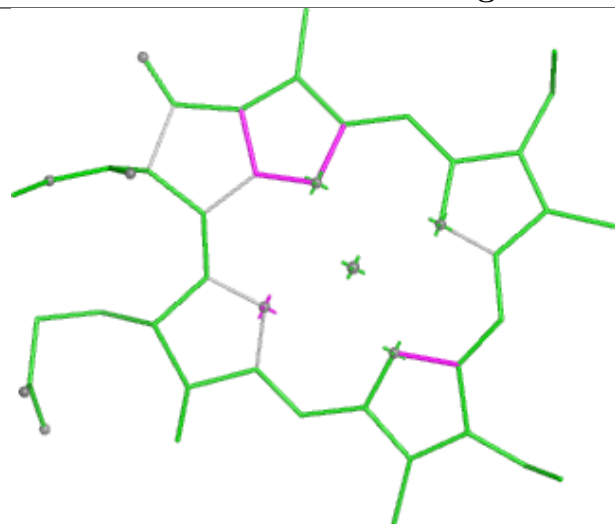


Torsions

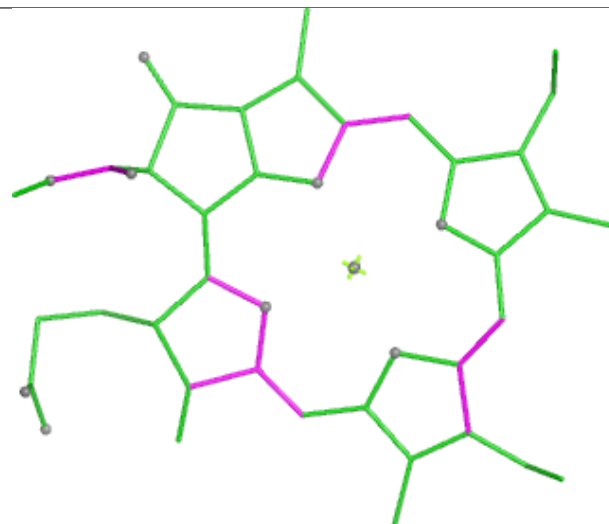


Rings

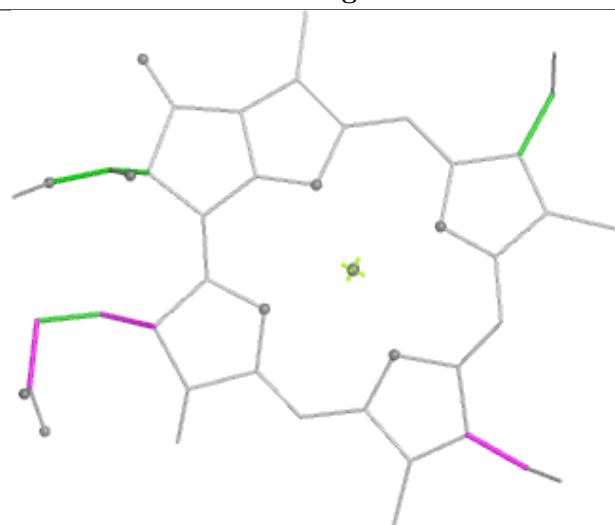
Ligand A1JWG ZH 503



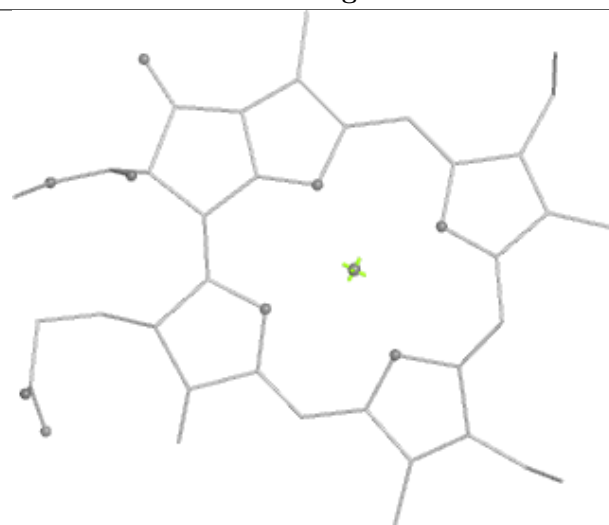
Bond lengths



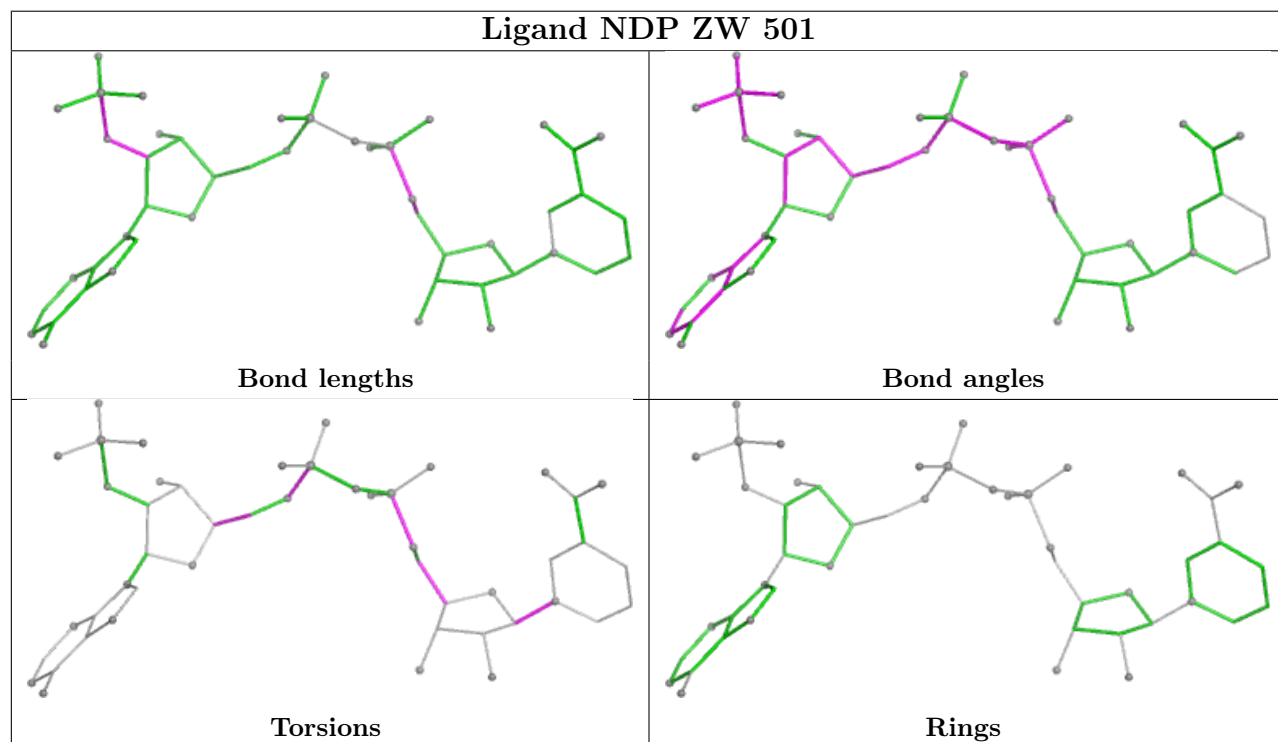
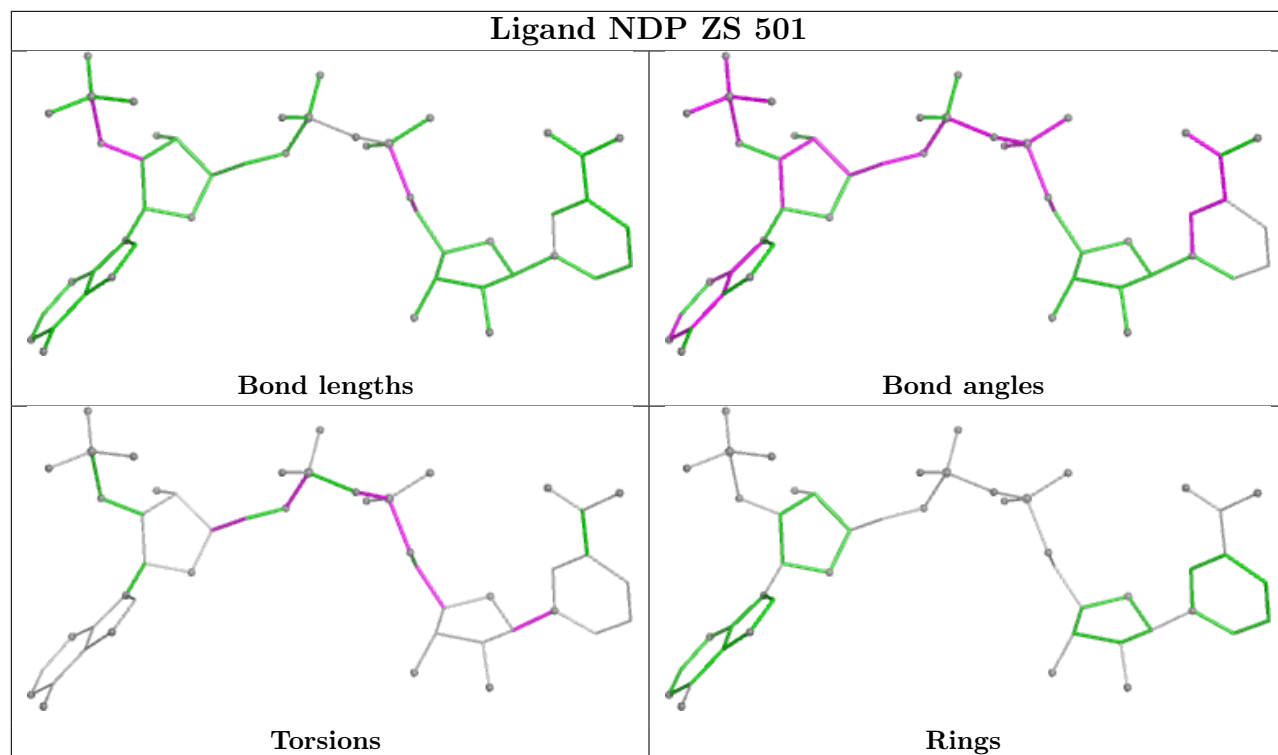
Bond angles



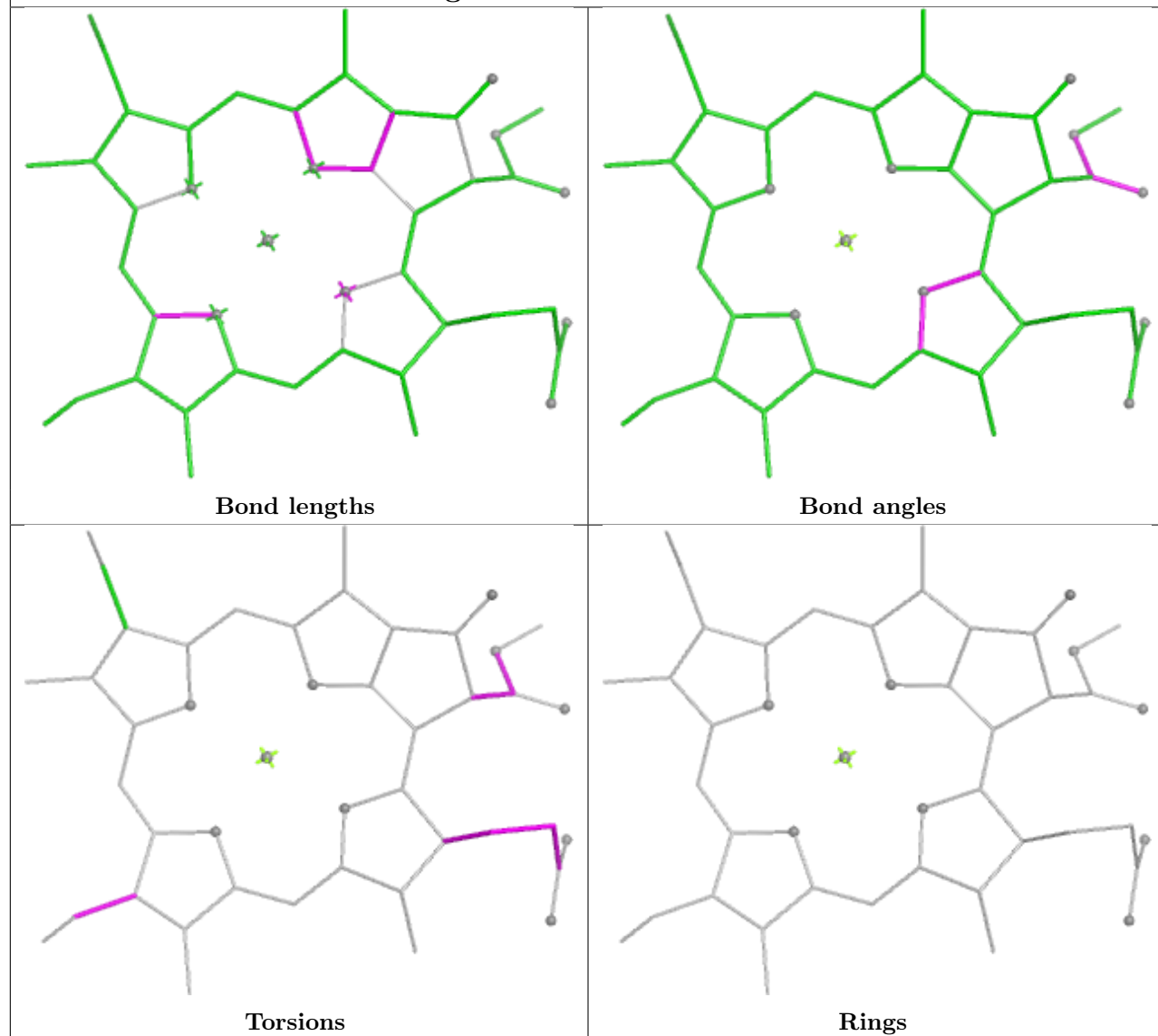
Torsions

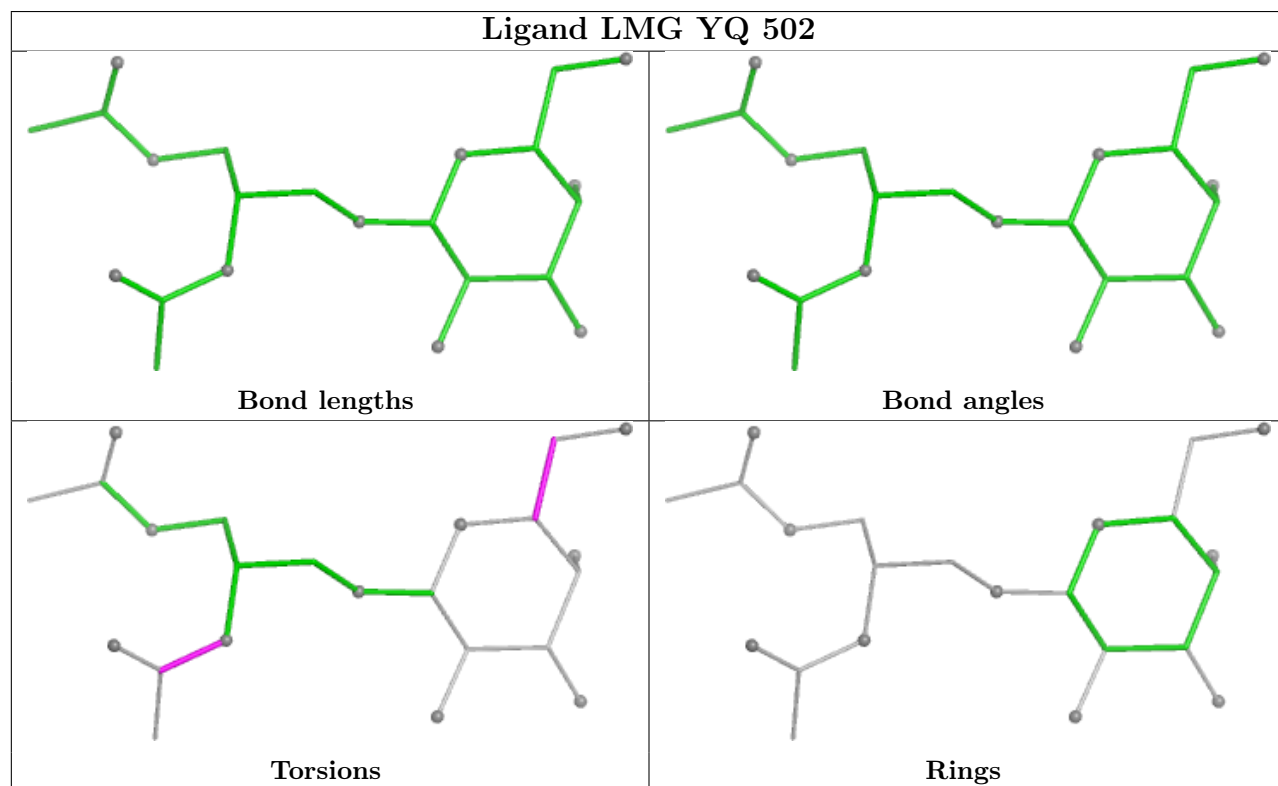
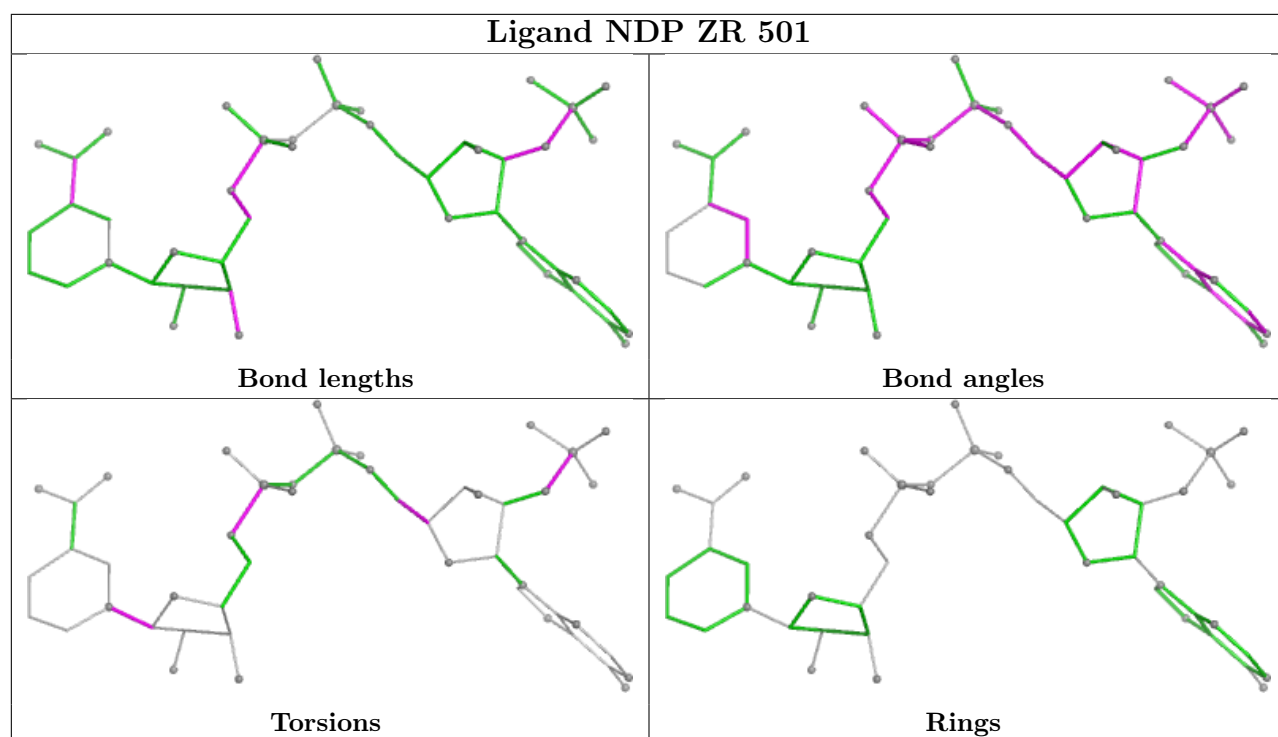


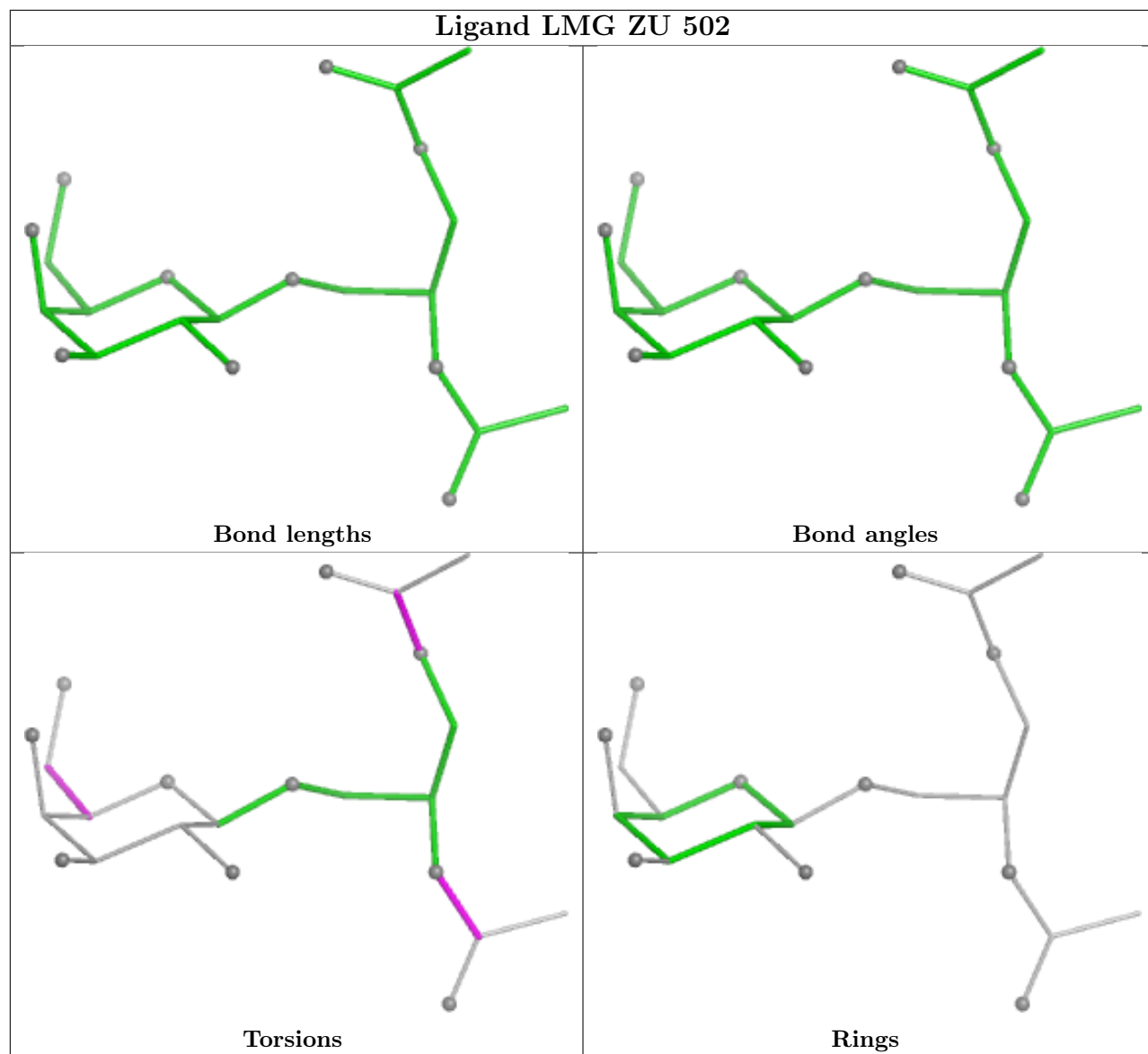
Rings



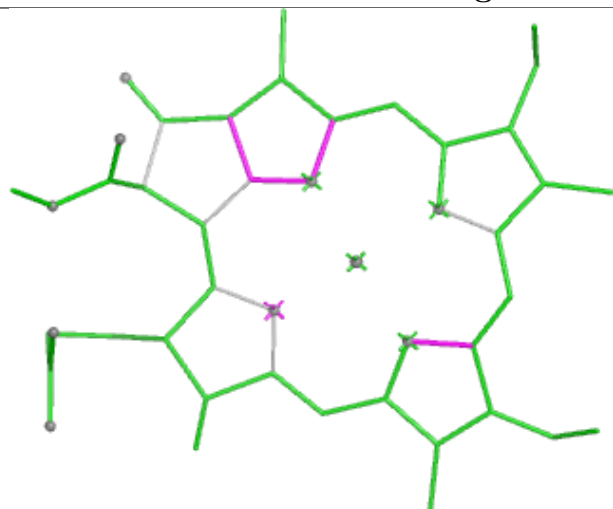
Ligand A1JWG YU 503



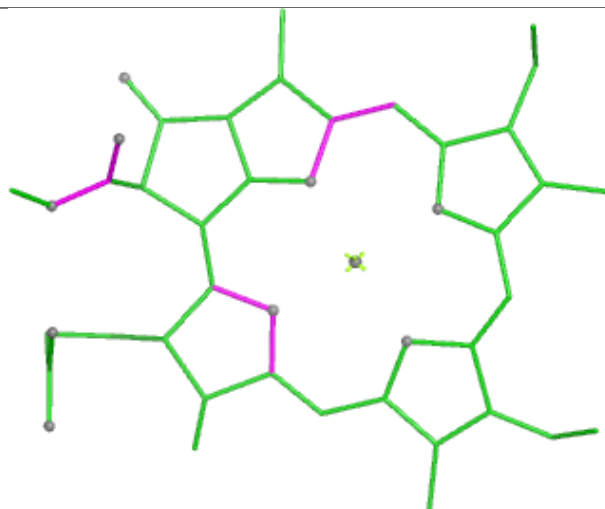




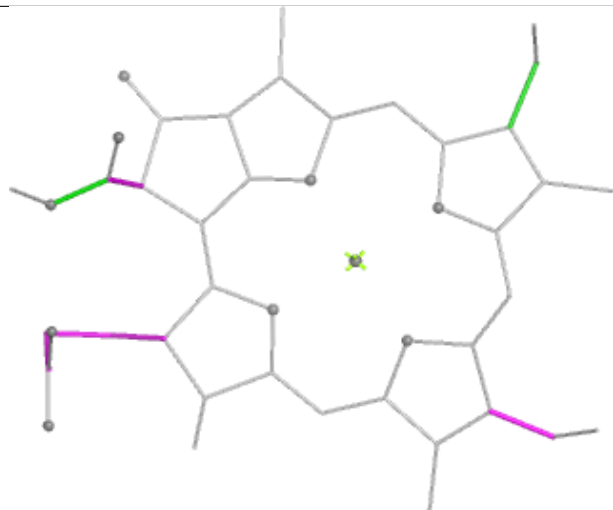
Ligand A1JWG YP 503



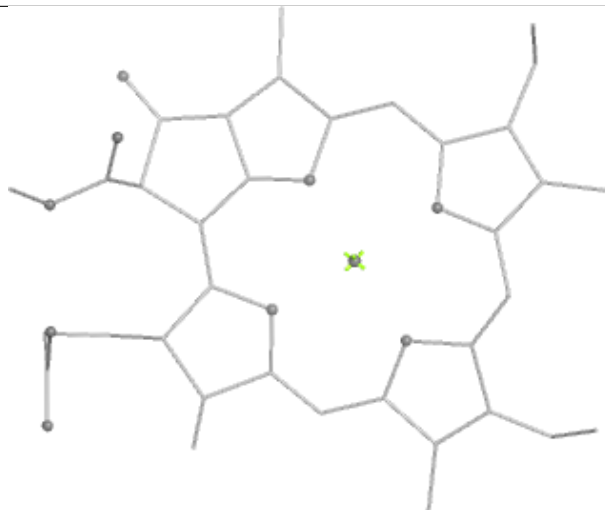
Bond lengths



Bond angles

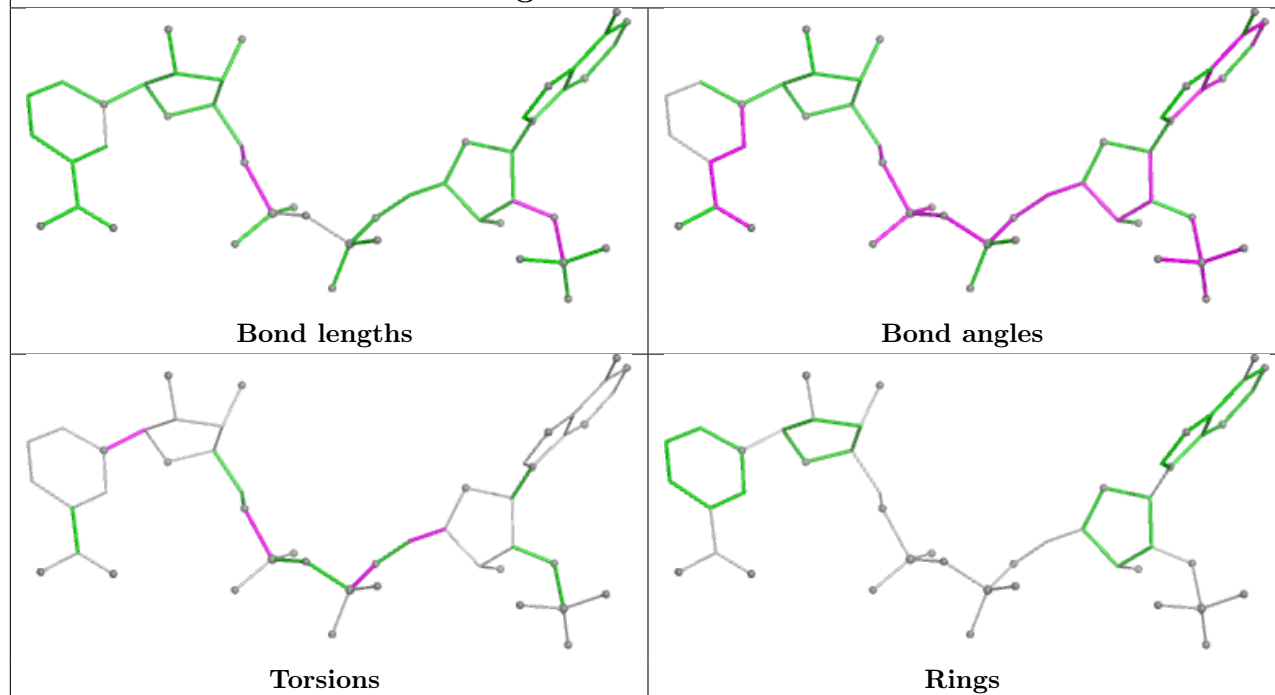


Torsions

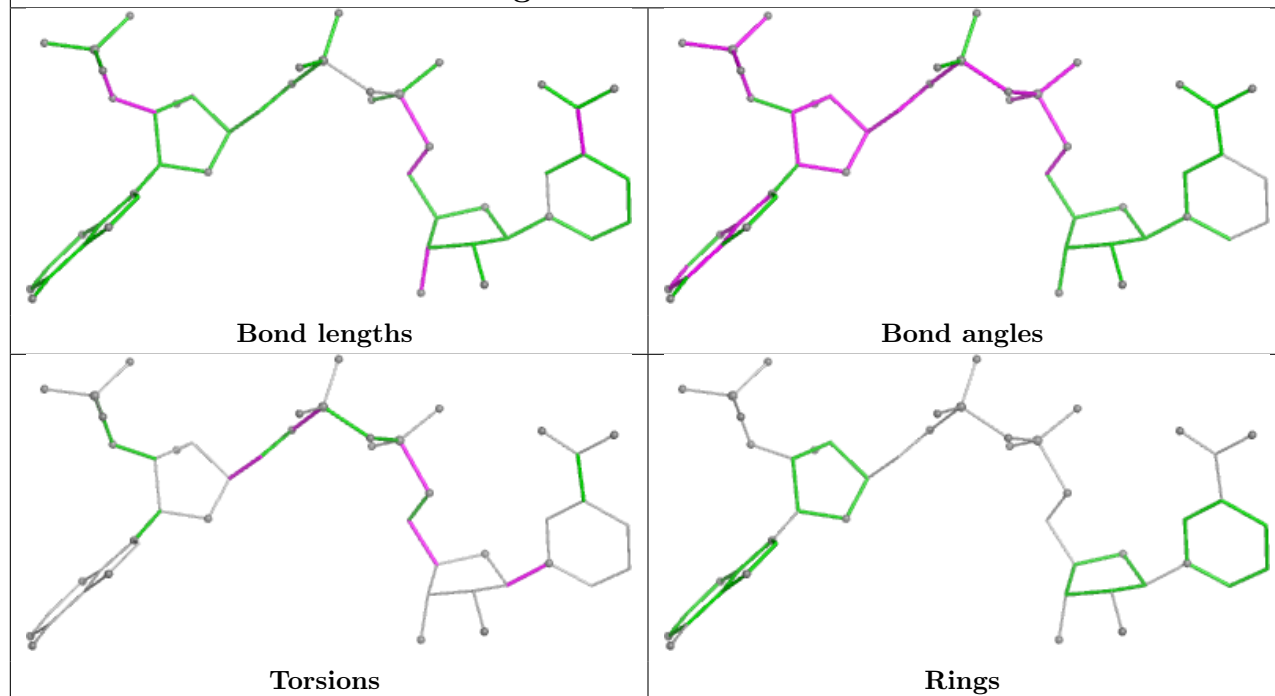


Rings

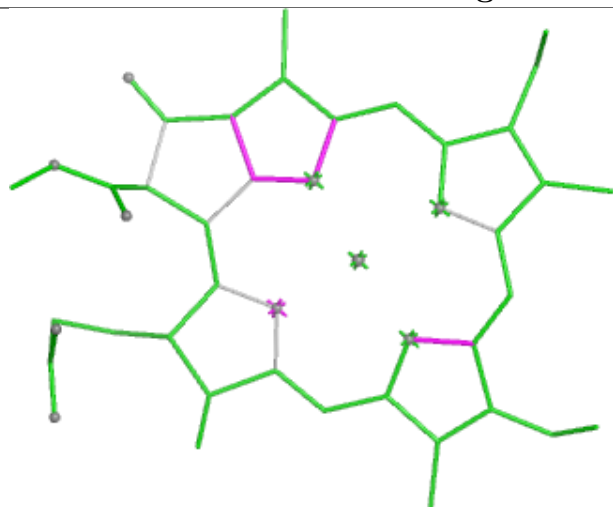
Ligand NDP ZG 501



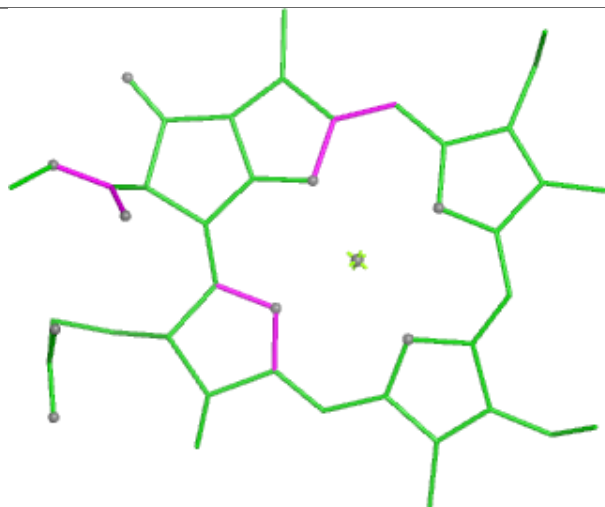
Ligand NDP ZM 501



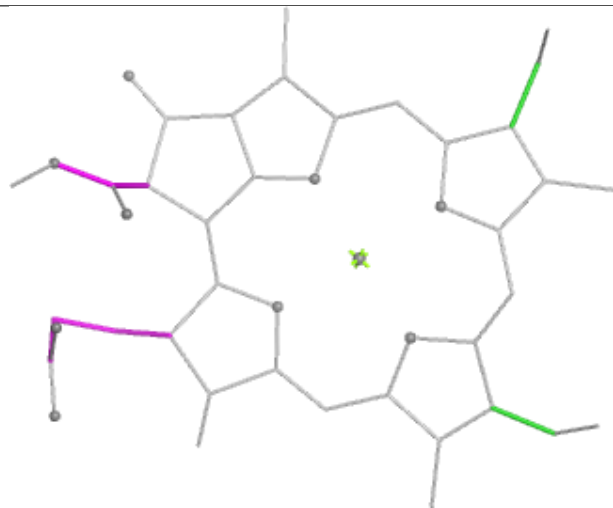
Ligand A1JWG YL 503



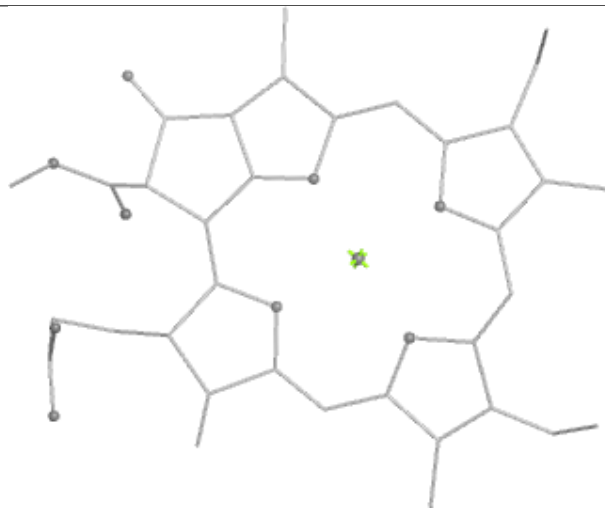
Bond lengths



Bond angles

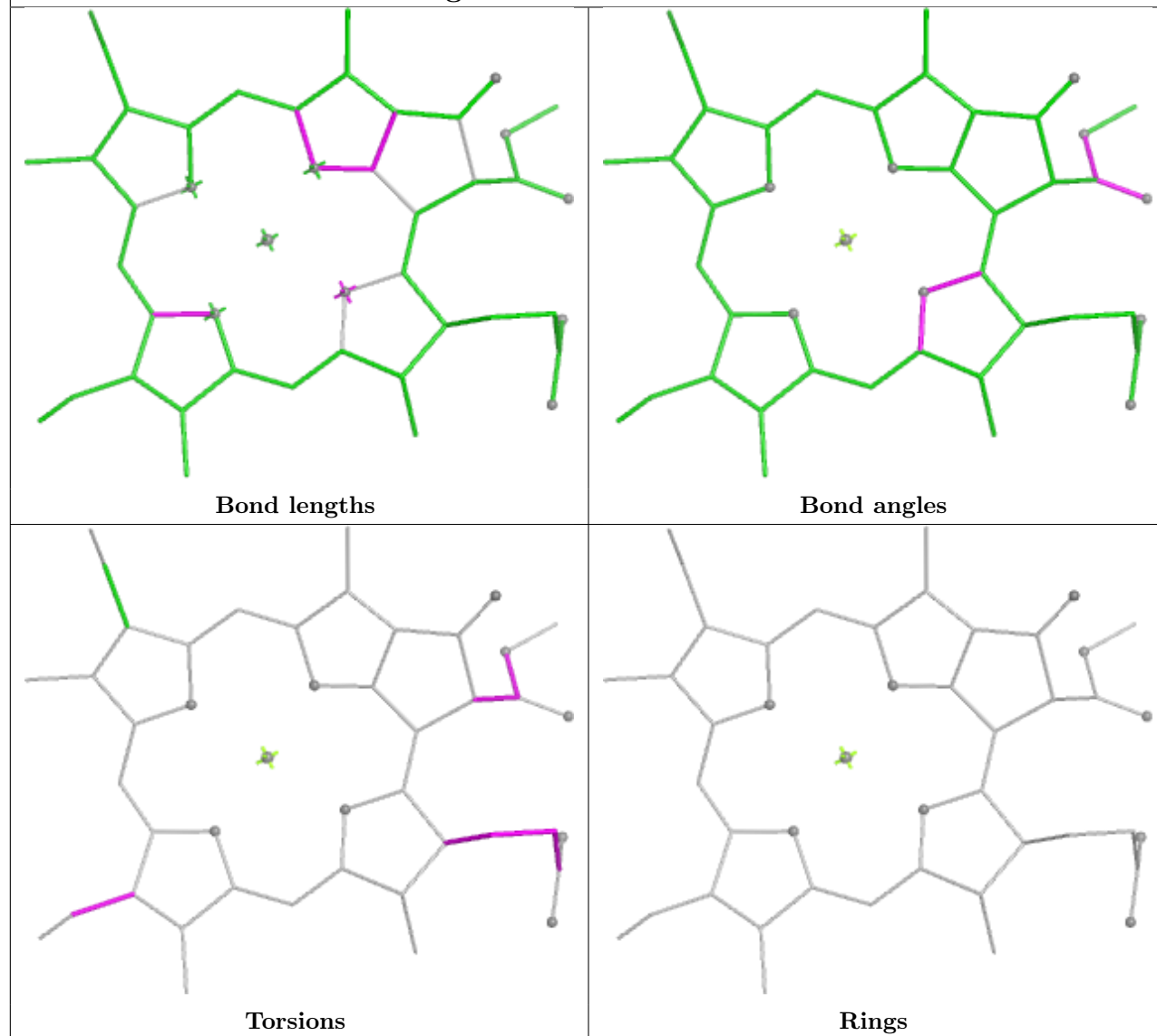


Torsions

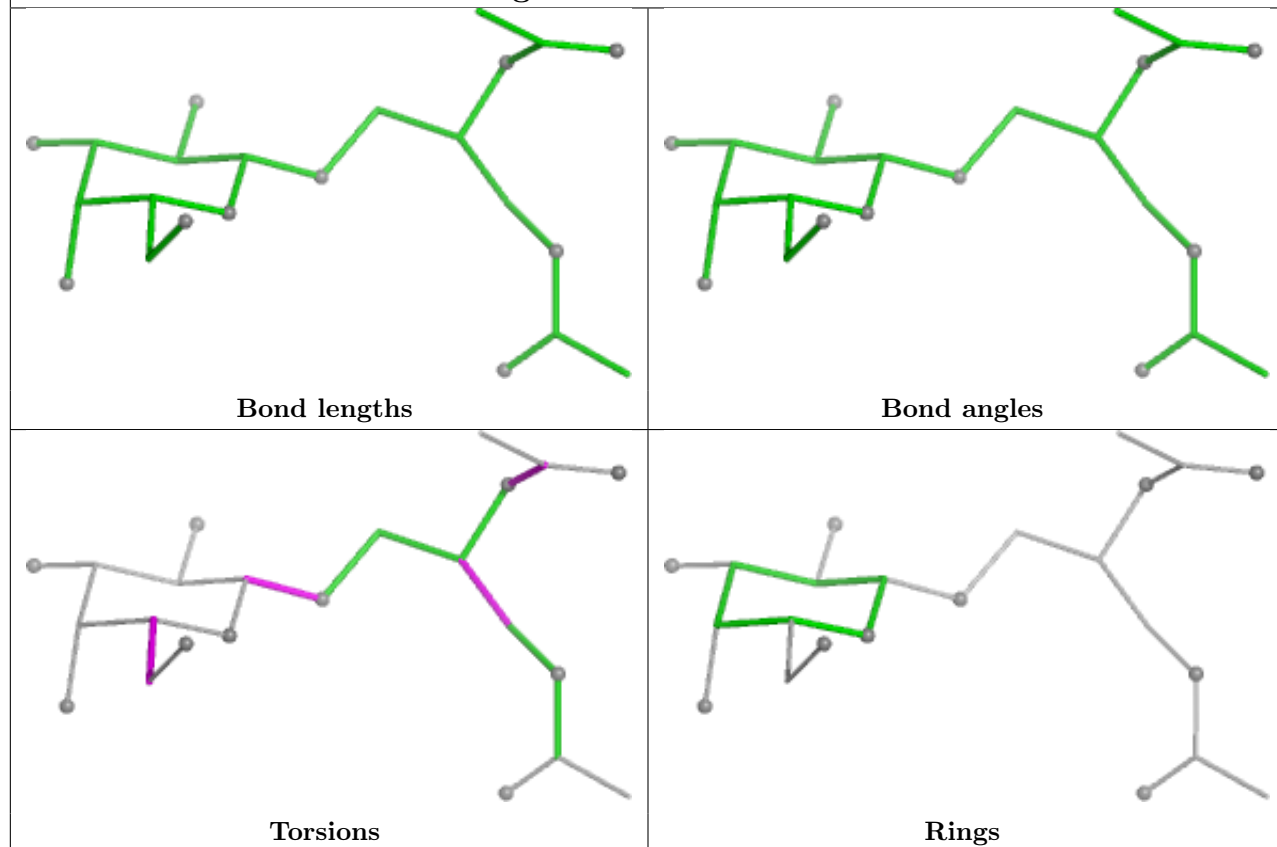


Rings

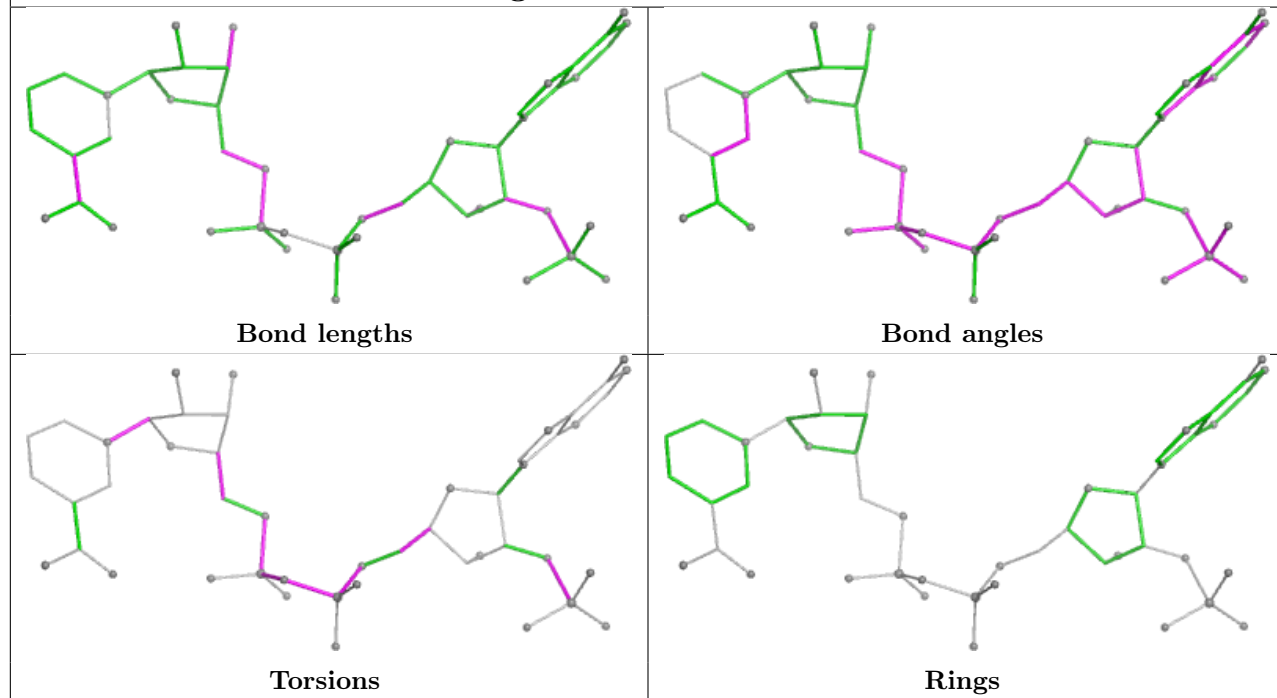
Ligand A1JWG YY 503



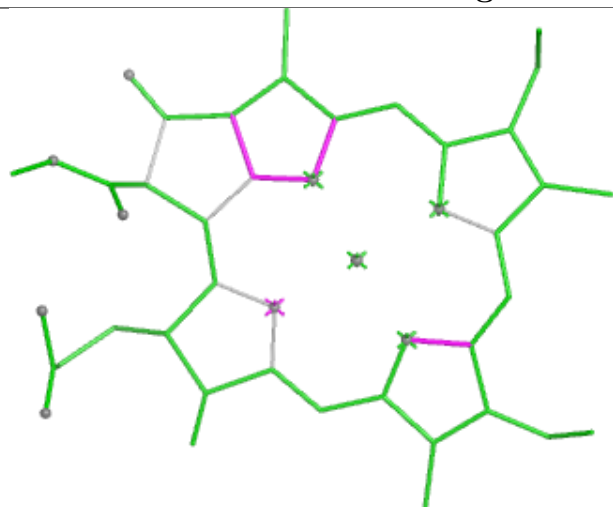
Ligand LMG YV 502



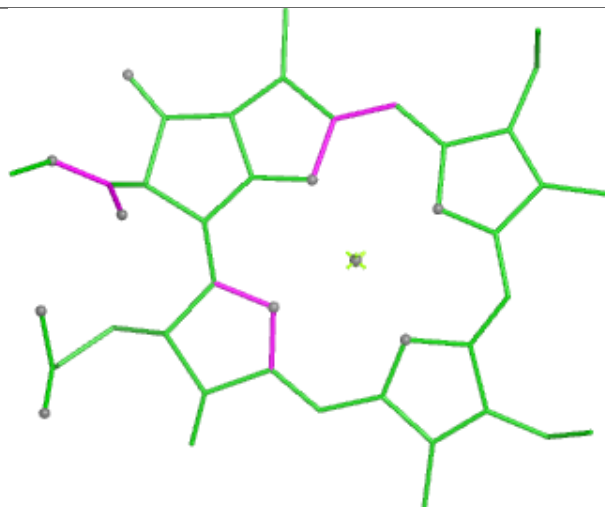
Ligand NDP YW 501



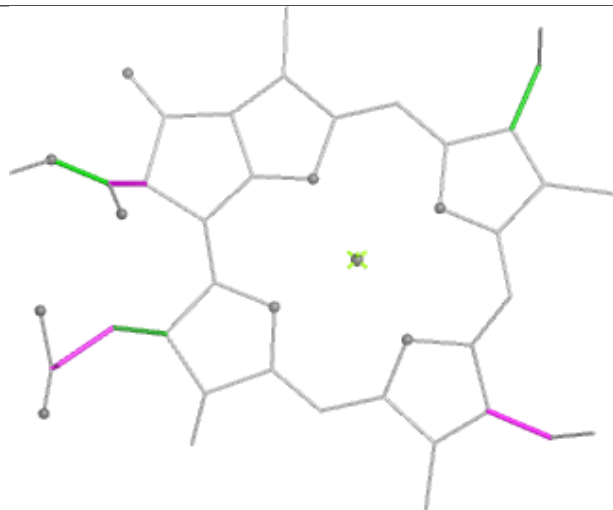
Ligand A1JWG ZV 503



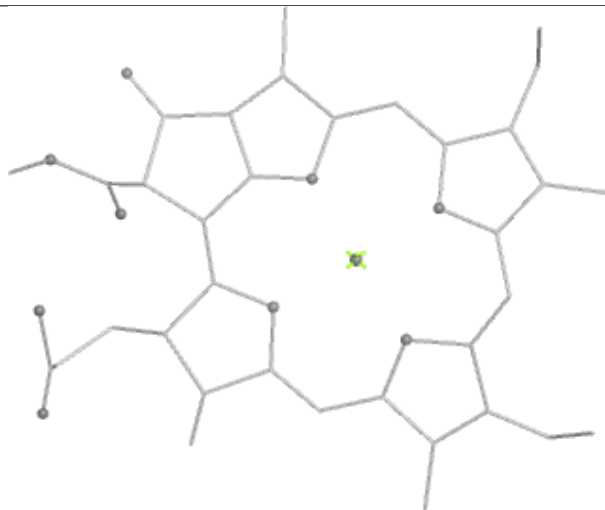
Bond lengths



Bond angles

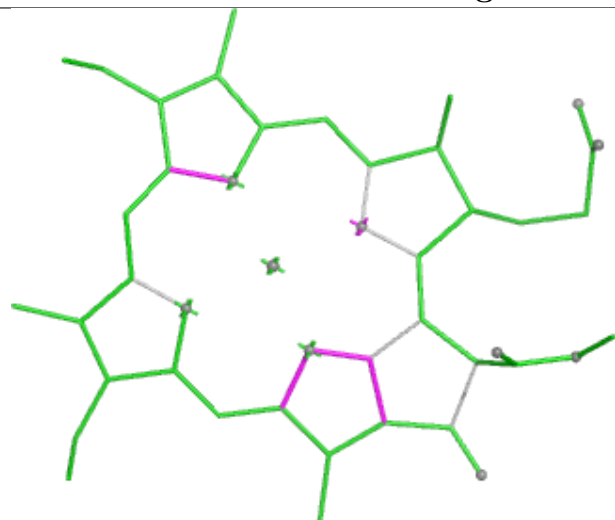


Torsions

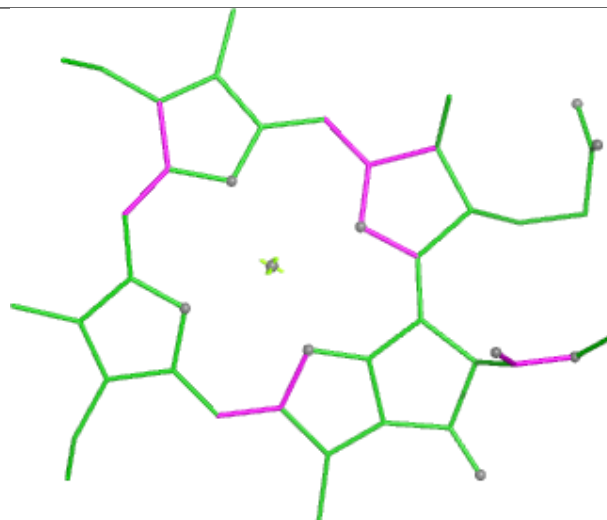


Rings

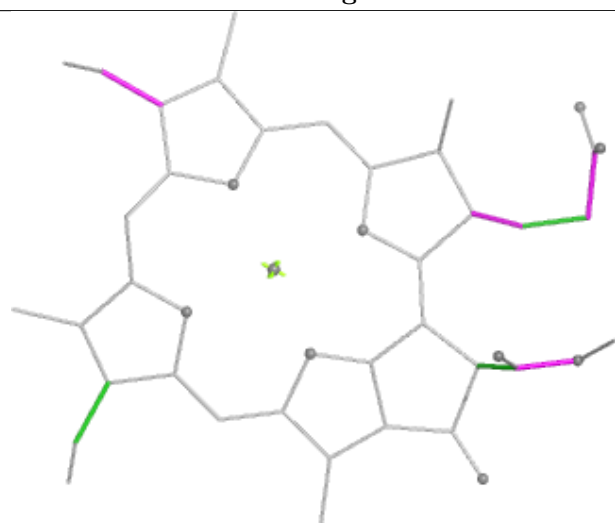
Ligand A1JWG YR 503



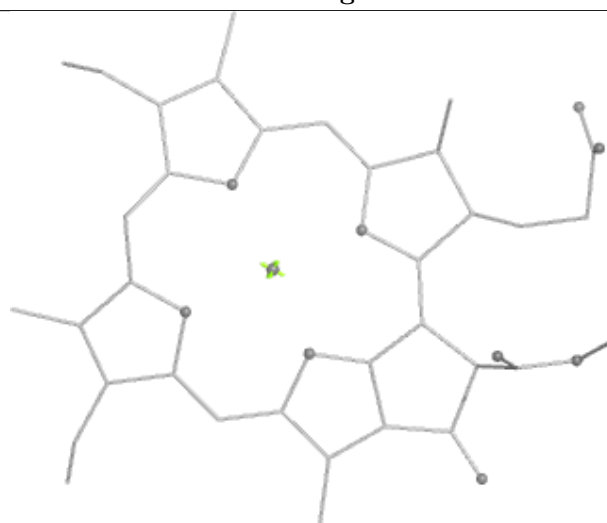
Bond lengths



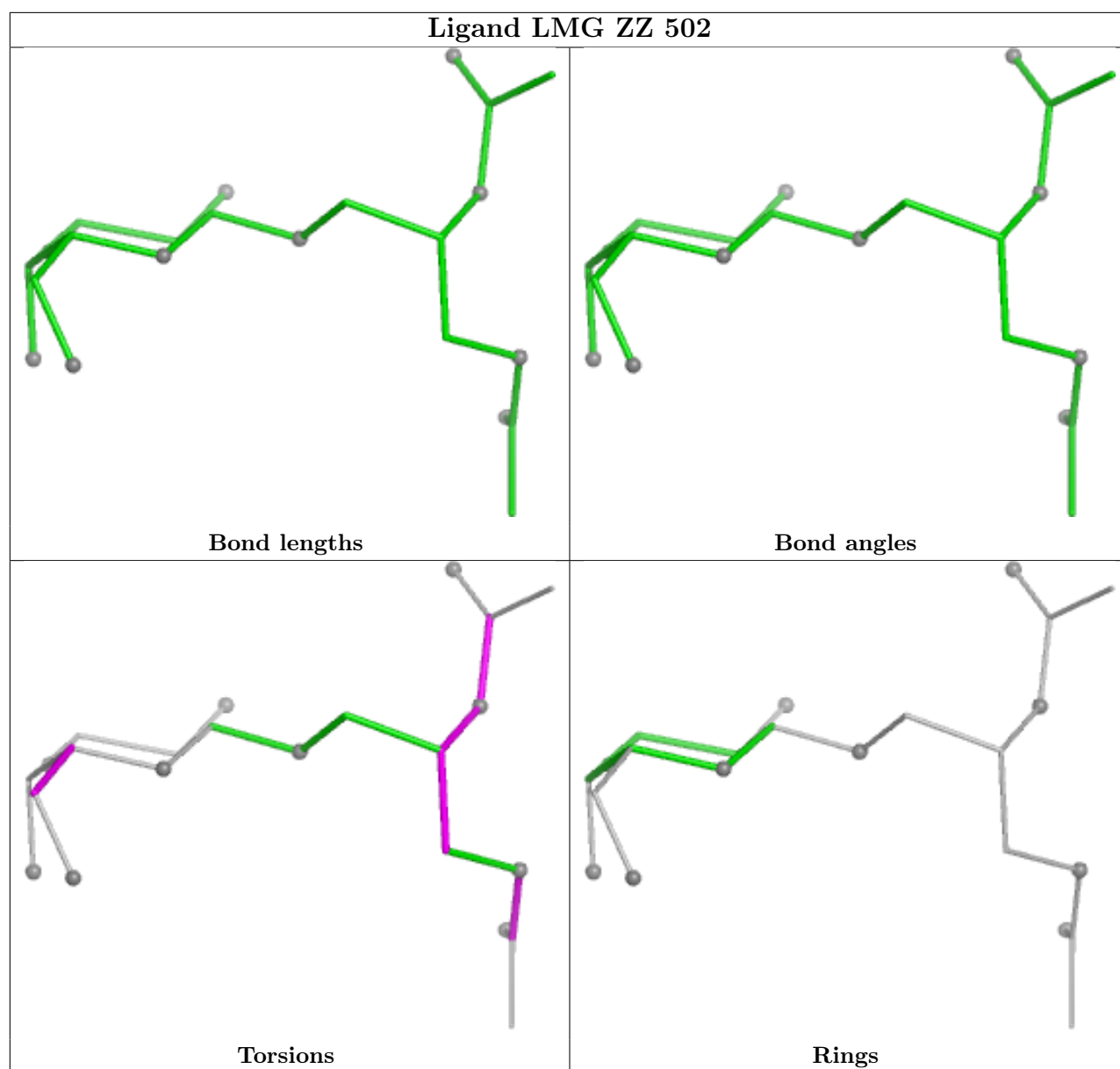
Bond angles

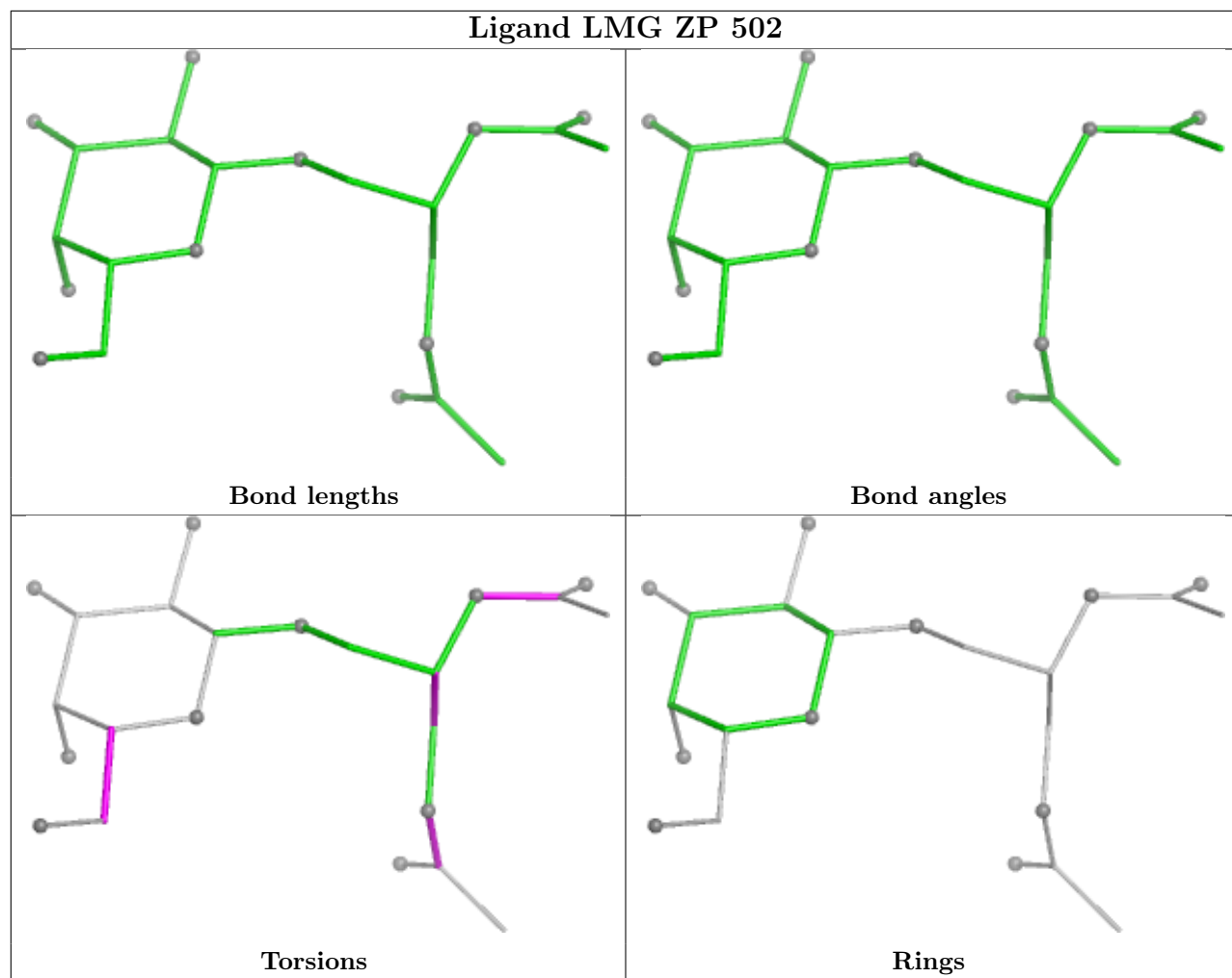


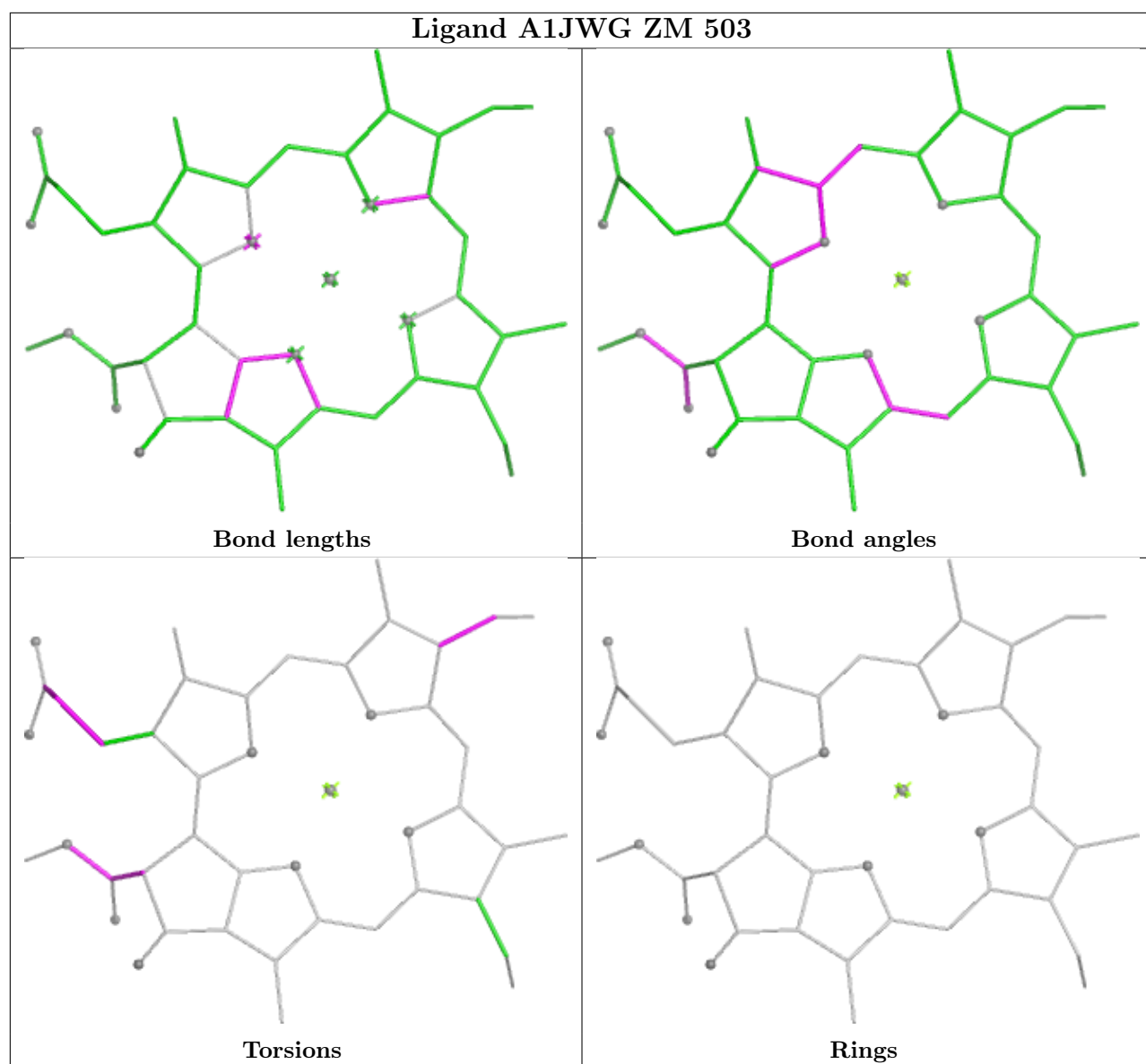
Torsions



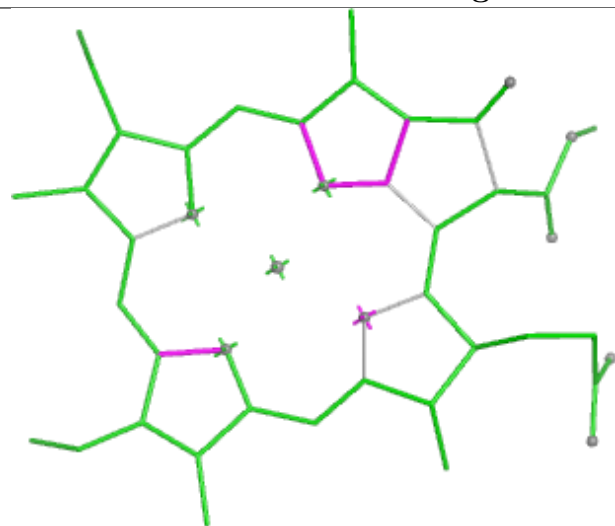
Rings



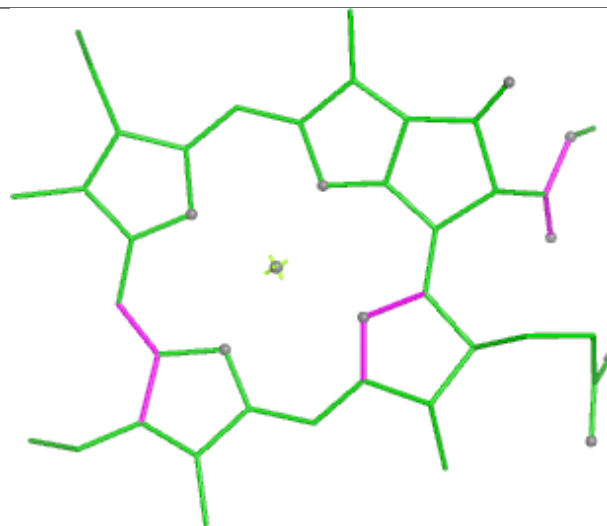




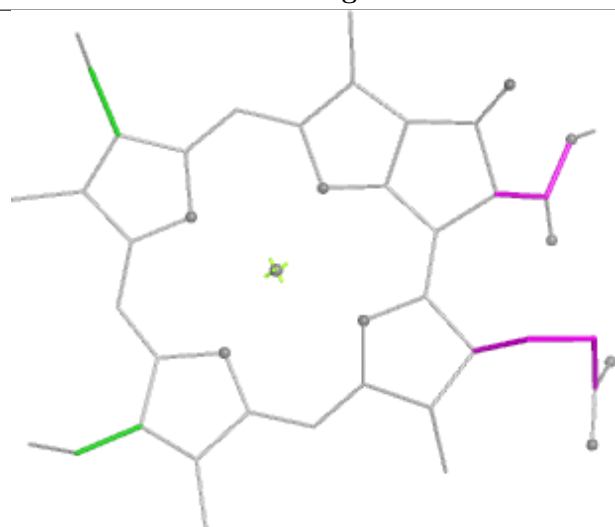
Ligand A1JWG YM 503



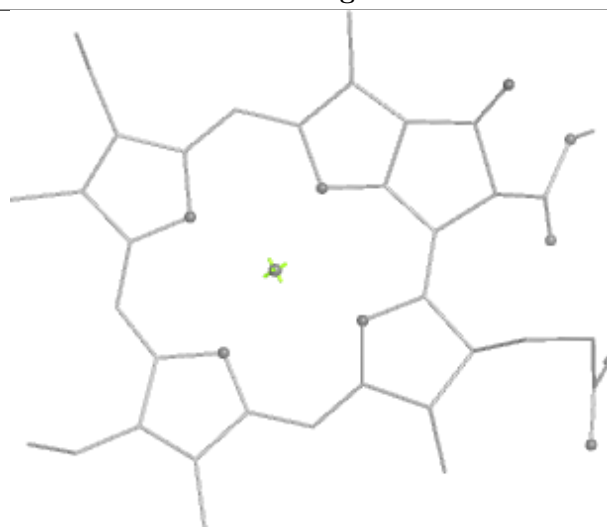
Bond lengths



Bond angles

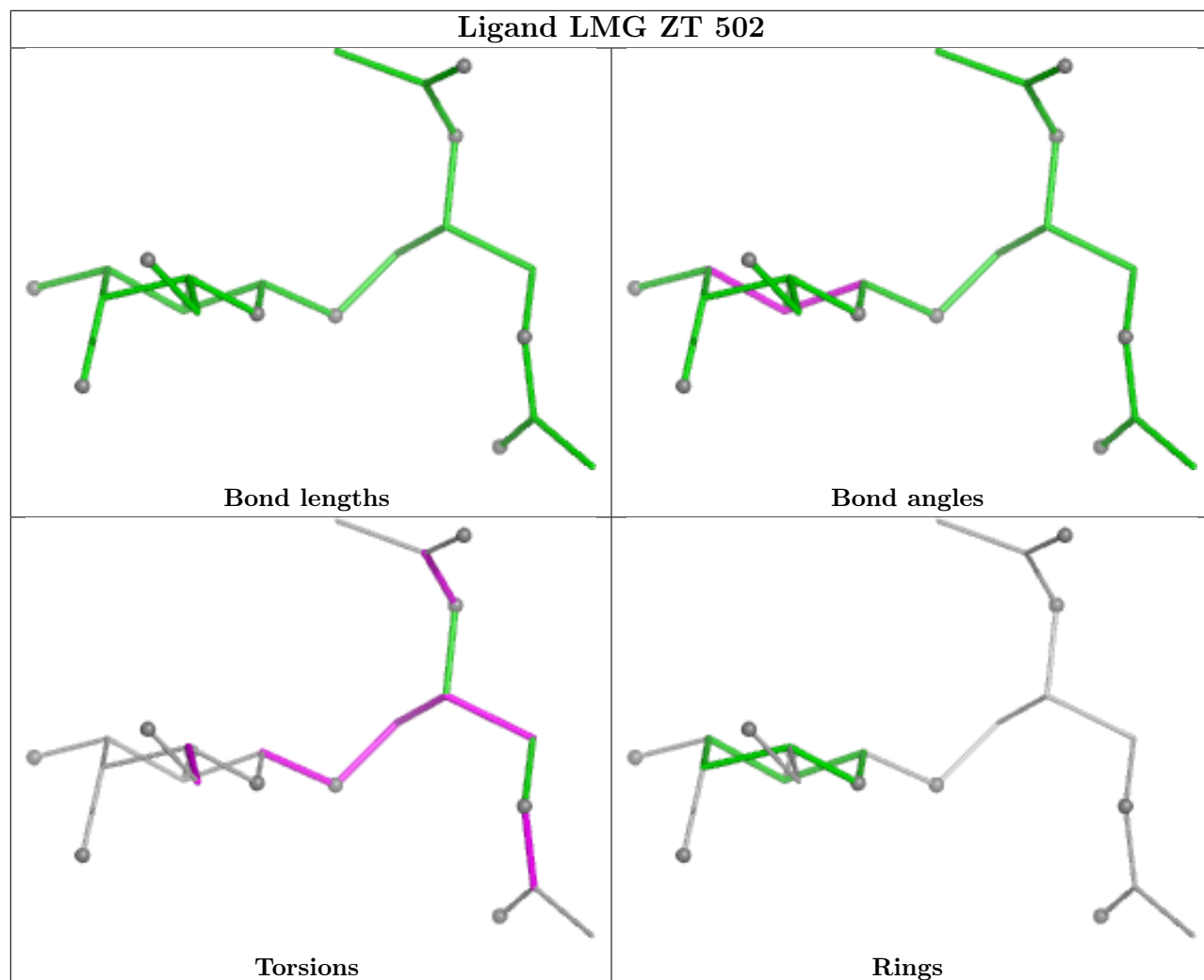


Torsions

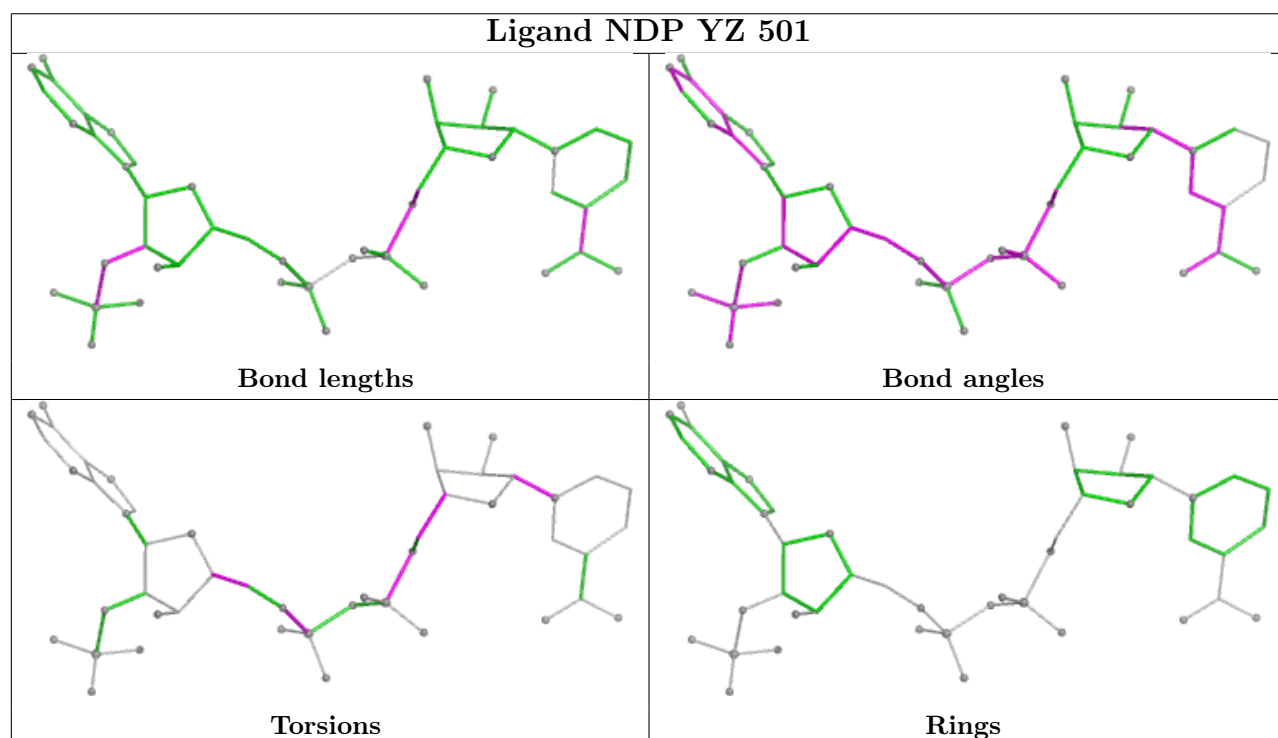


Rings

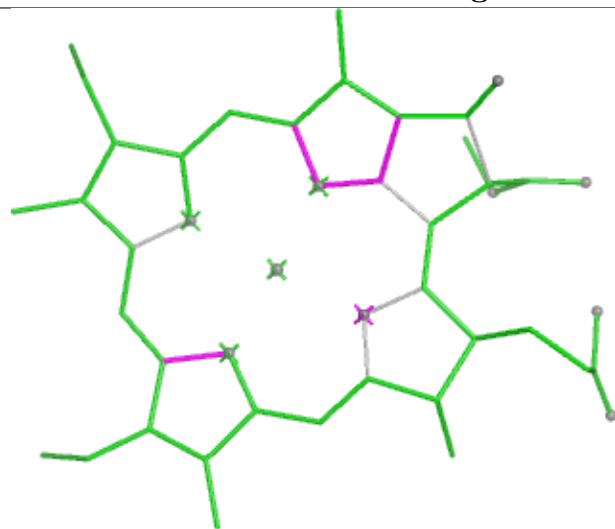
Ligand LMG ZT 502



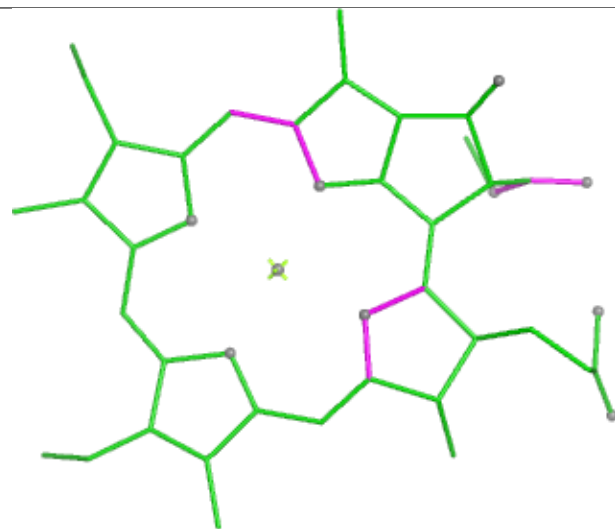
Ligand NDP YZ 501



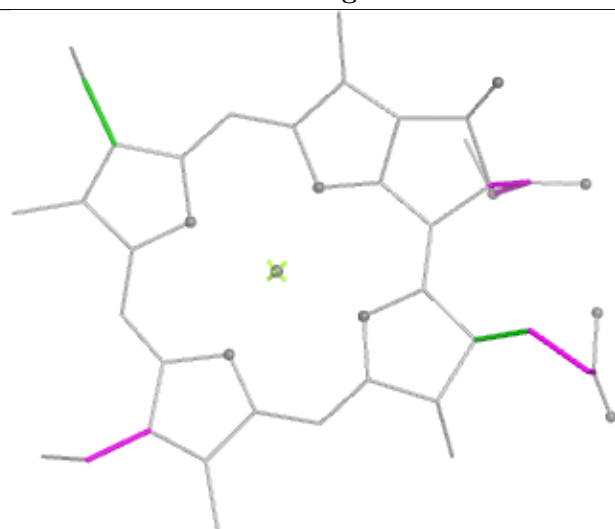
Ligand A1JWG YW 503



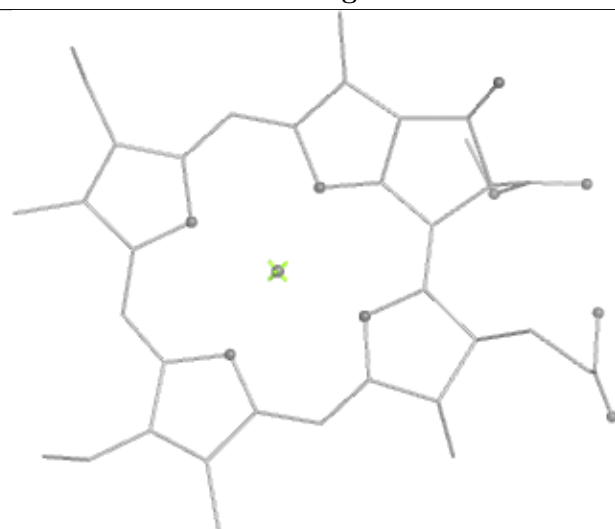
Bond lengths



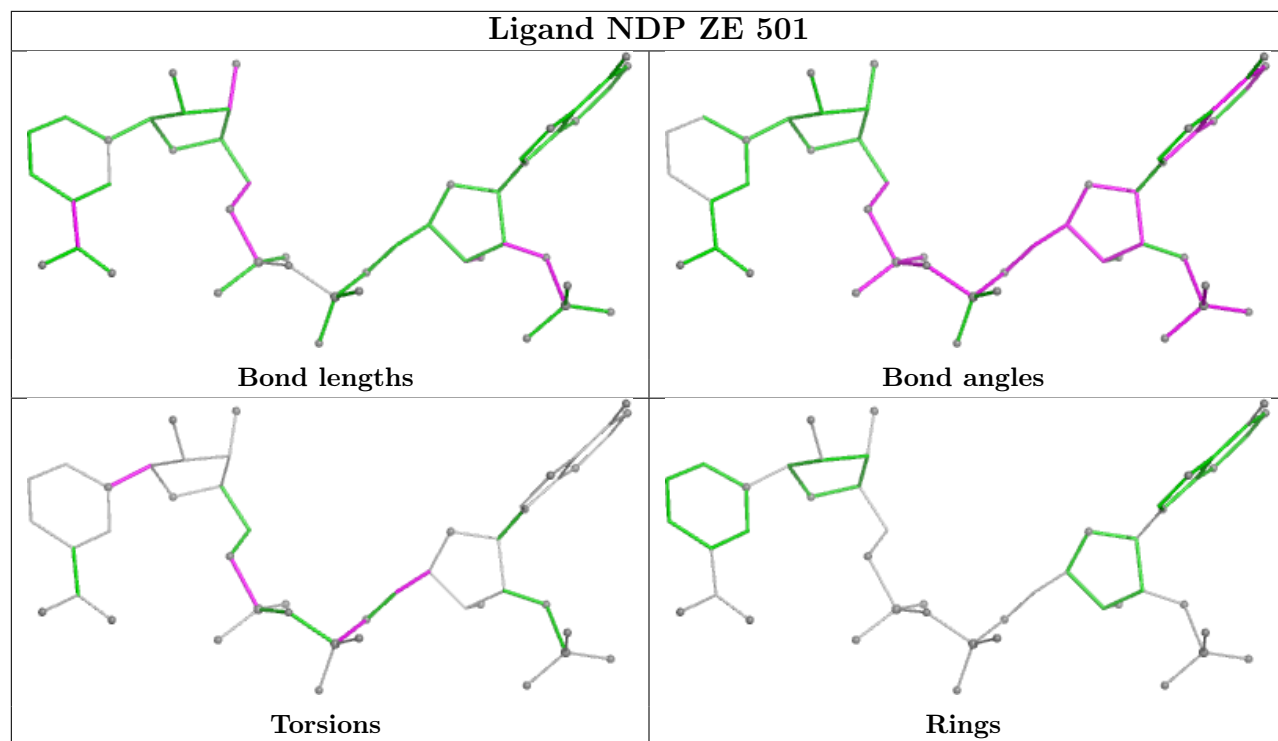
Bond angles

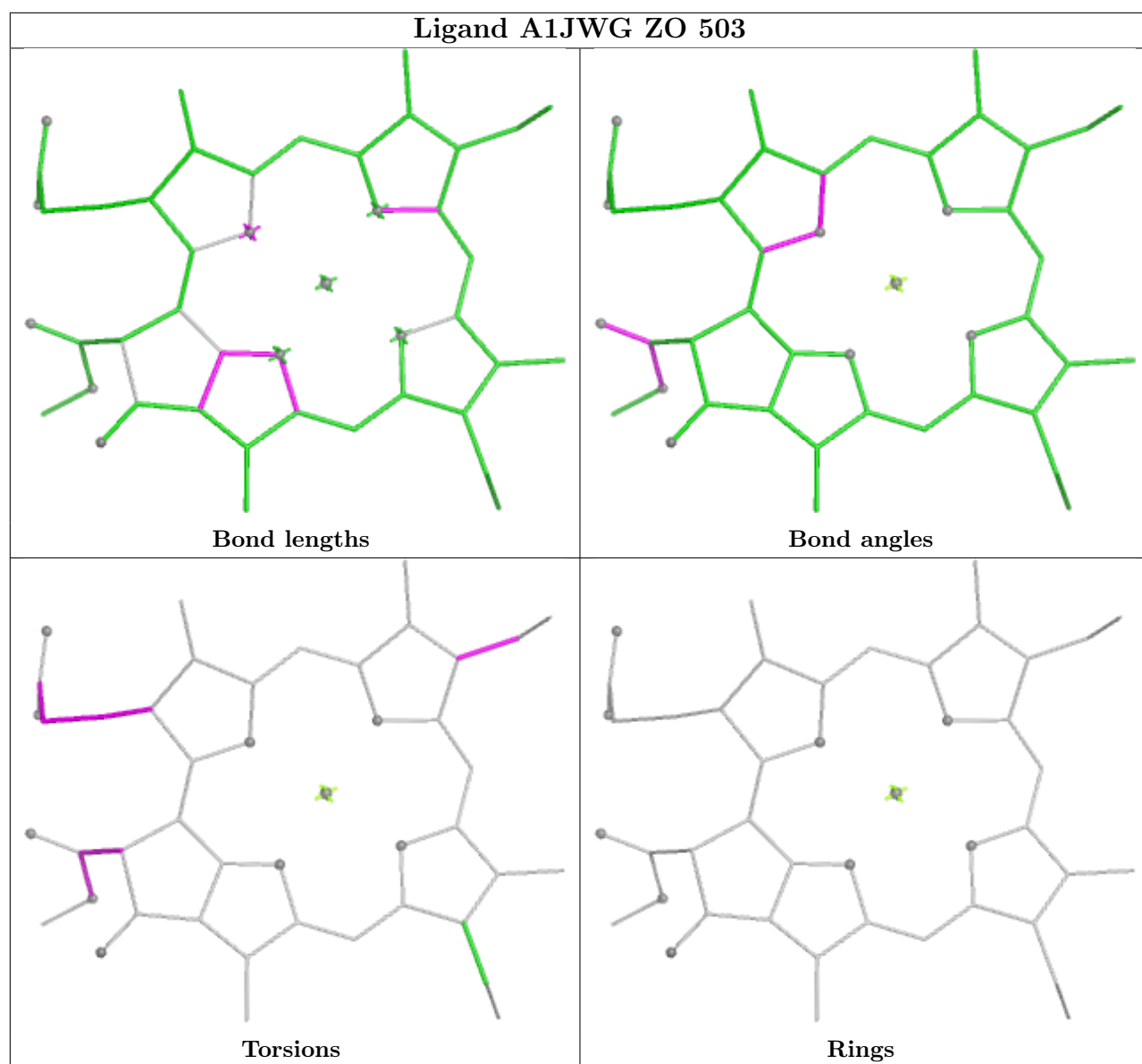


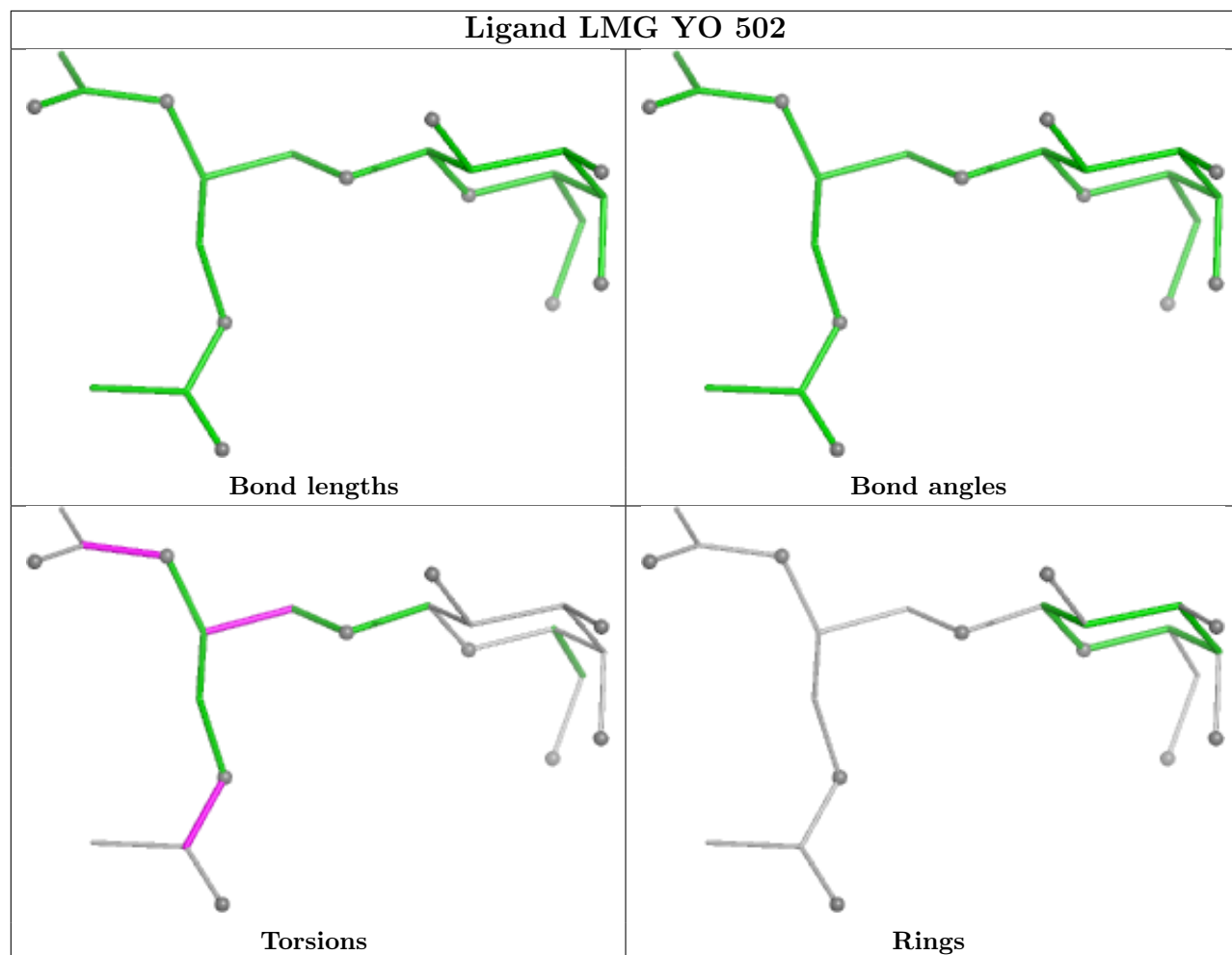
Torsions

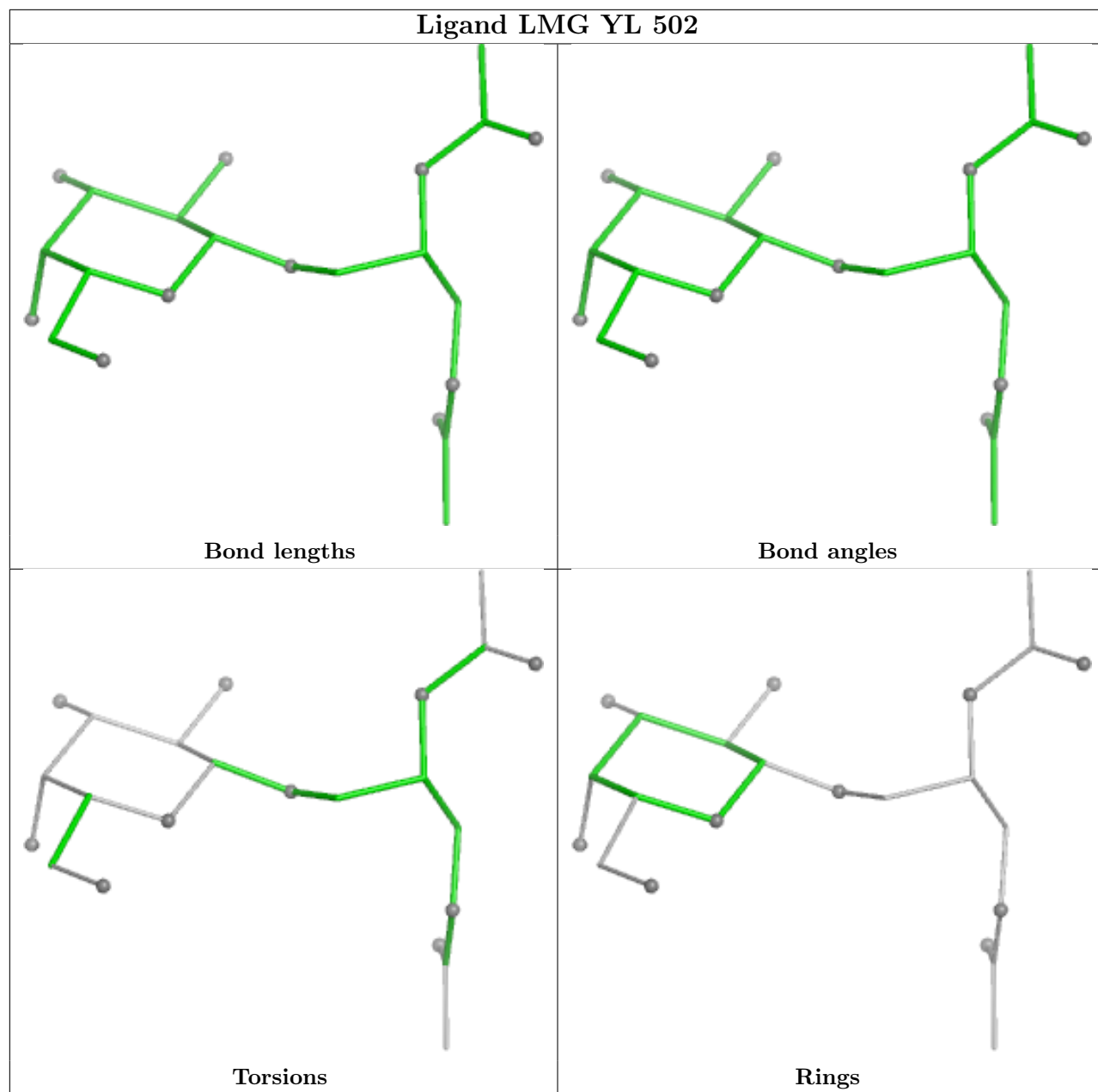


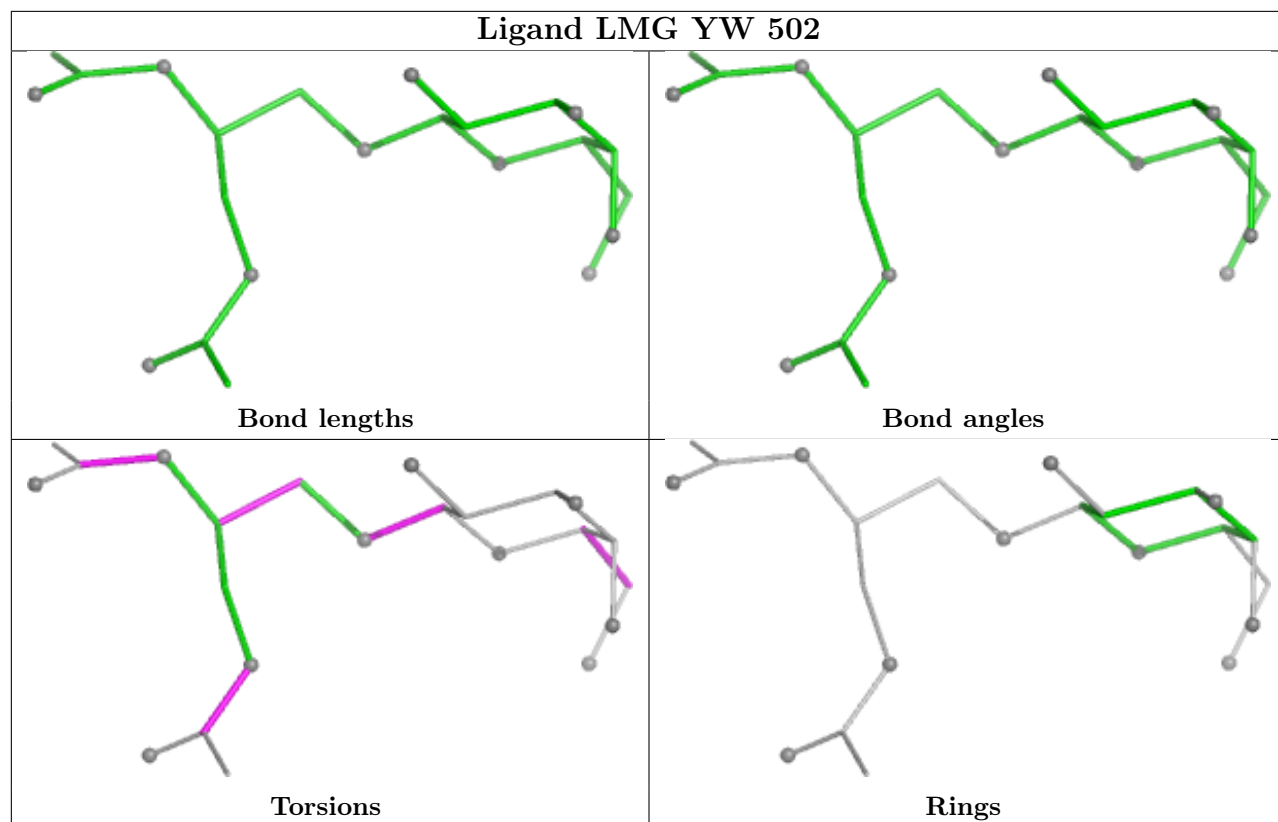
Rings



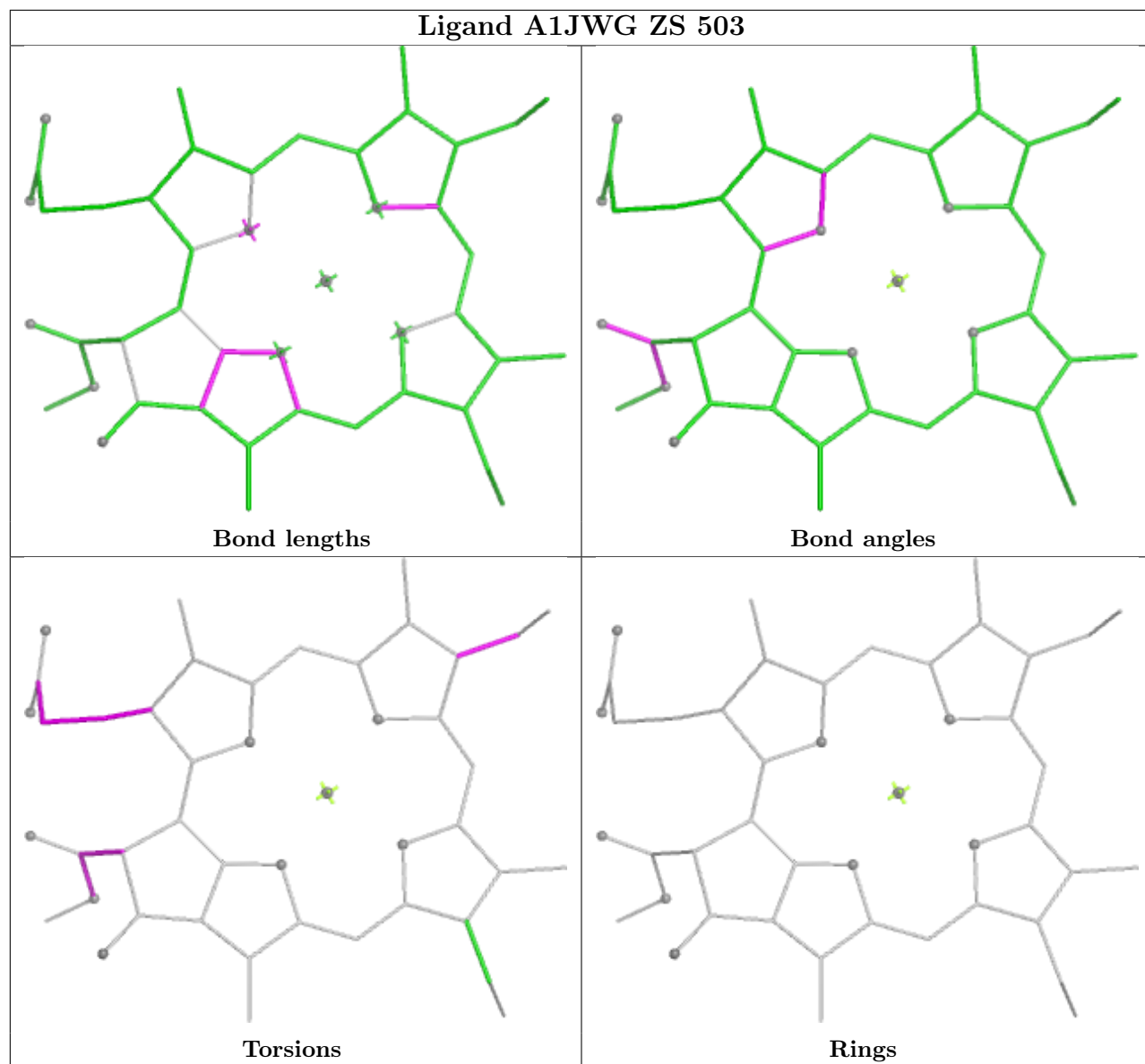




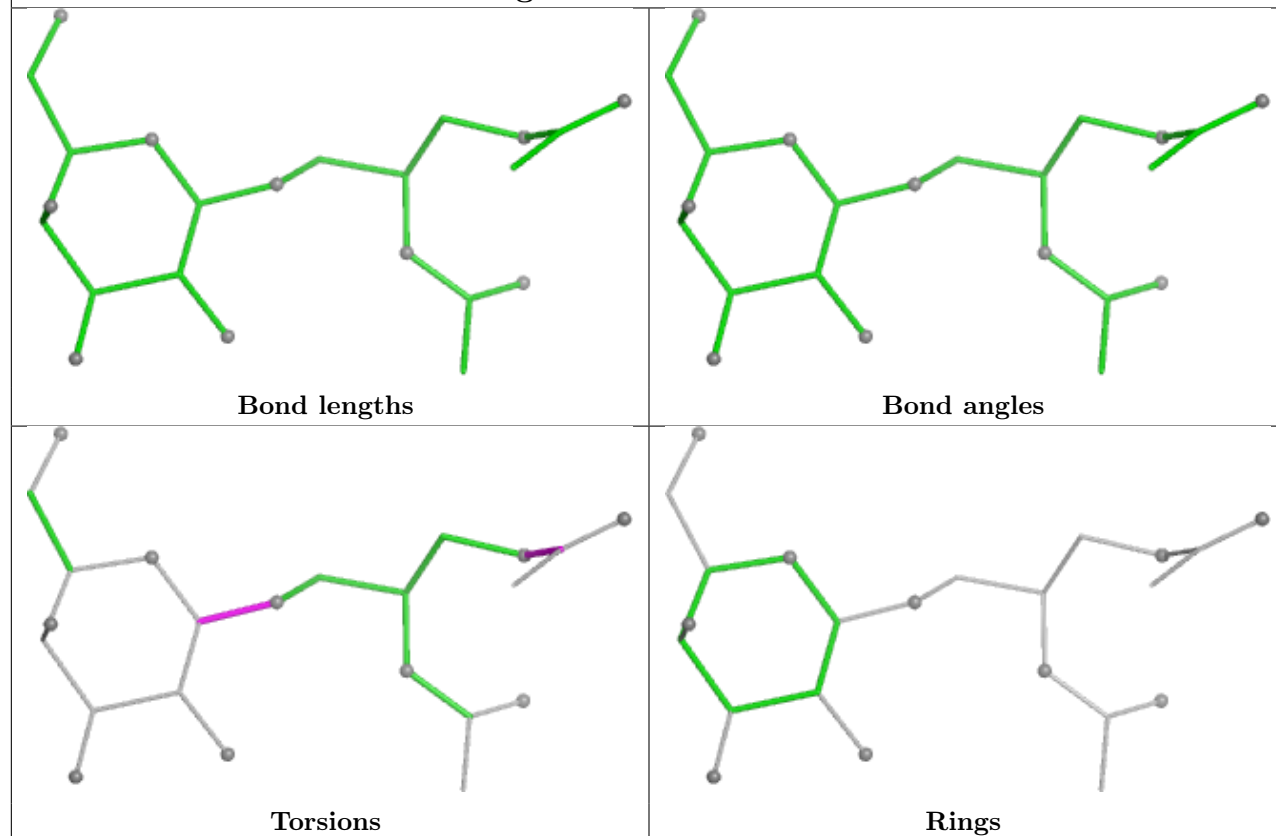




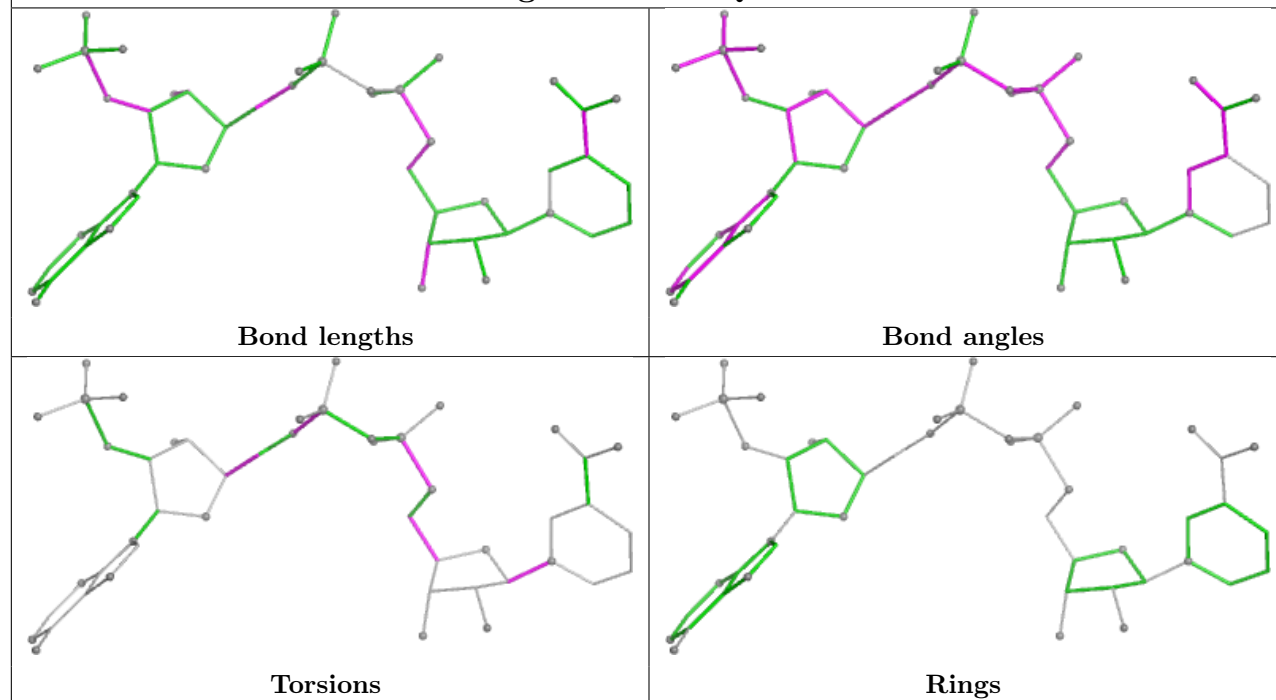
Ligand A1JWG ZS 503

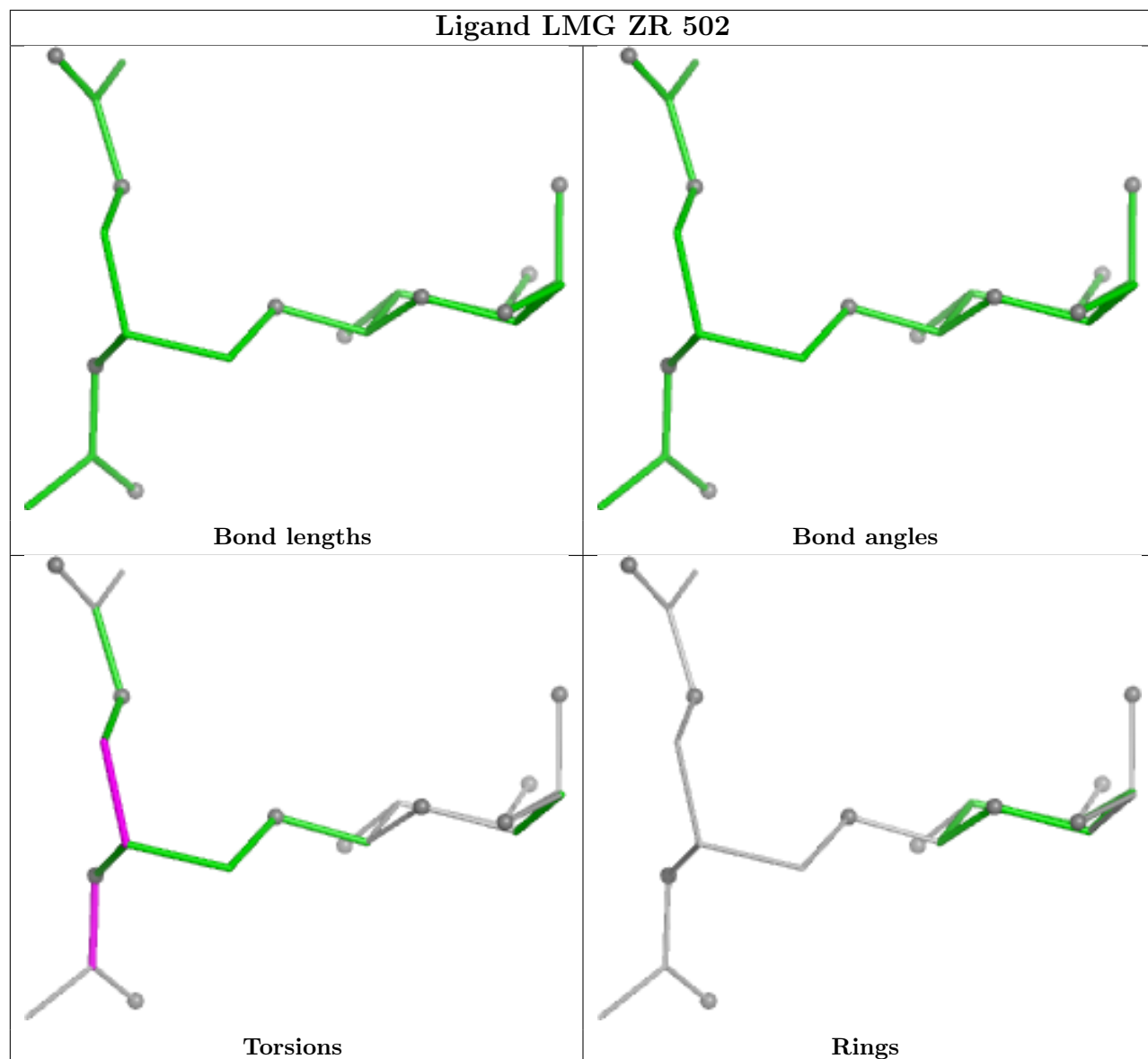


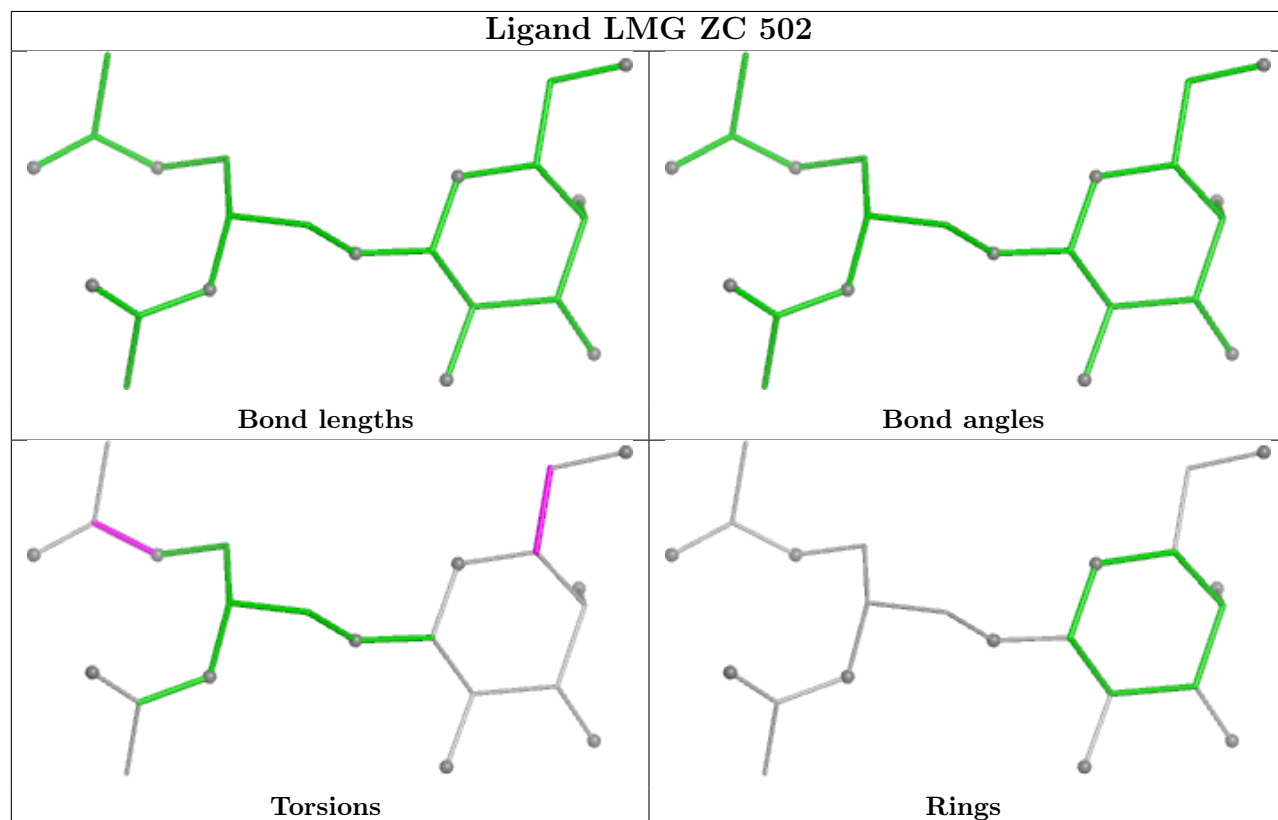
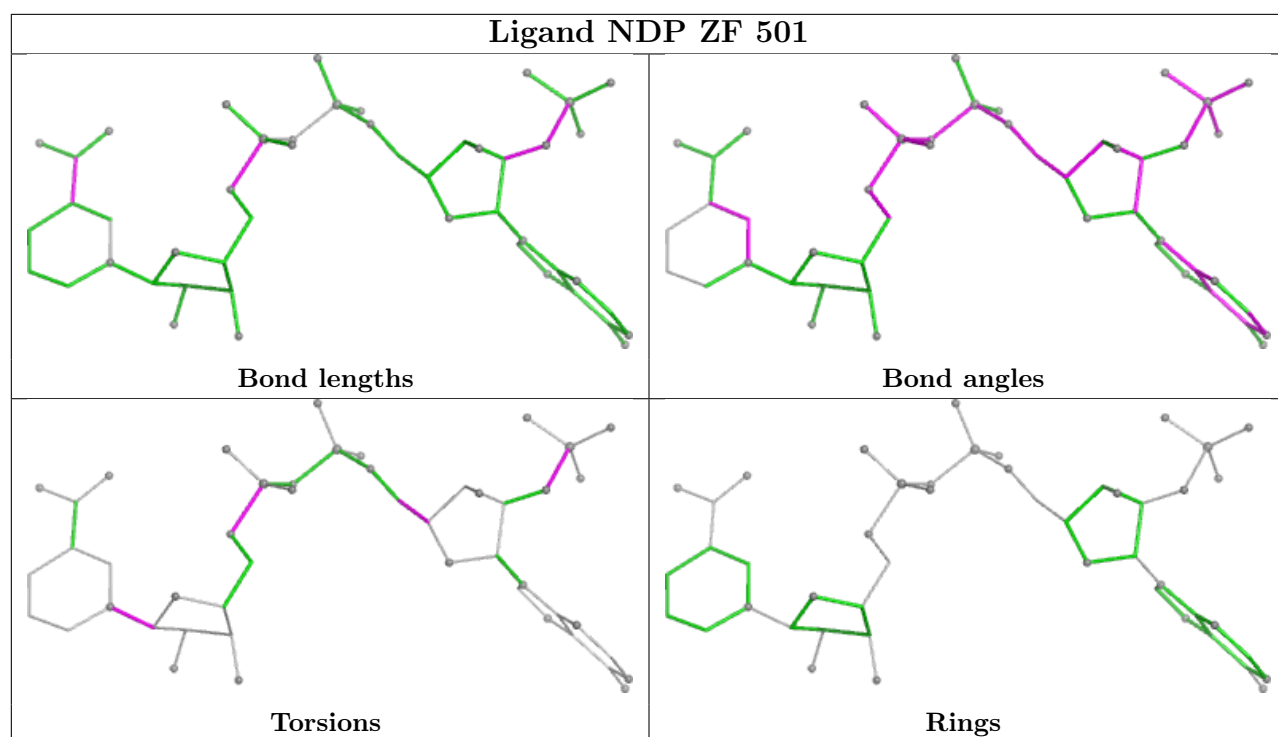
Ligand LMG YZ 502



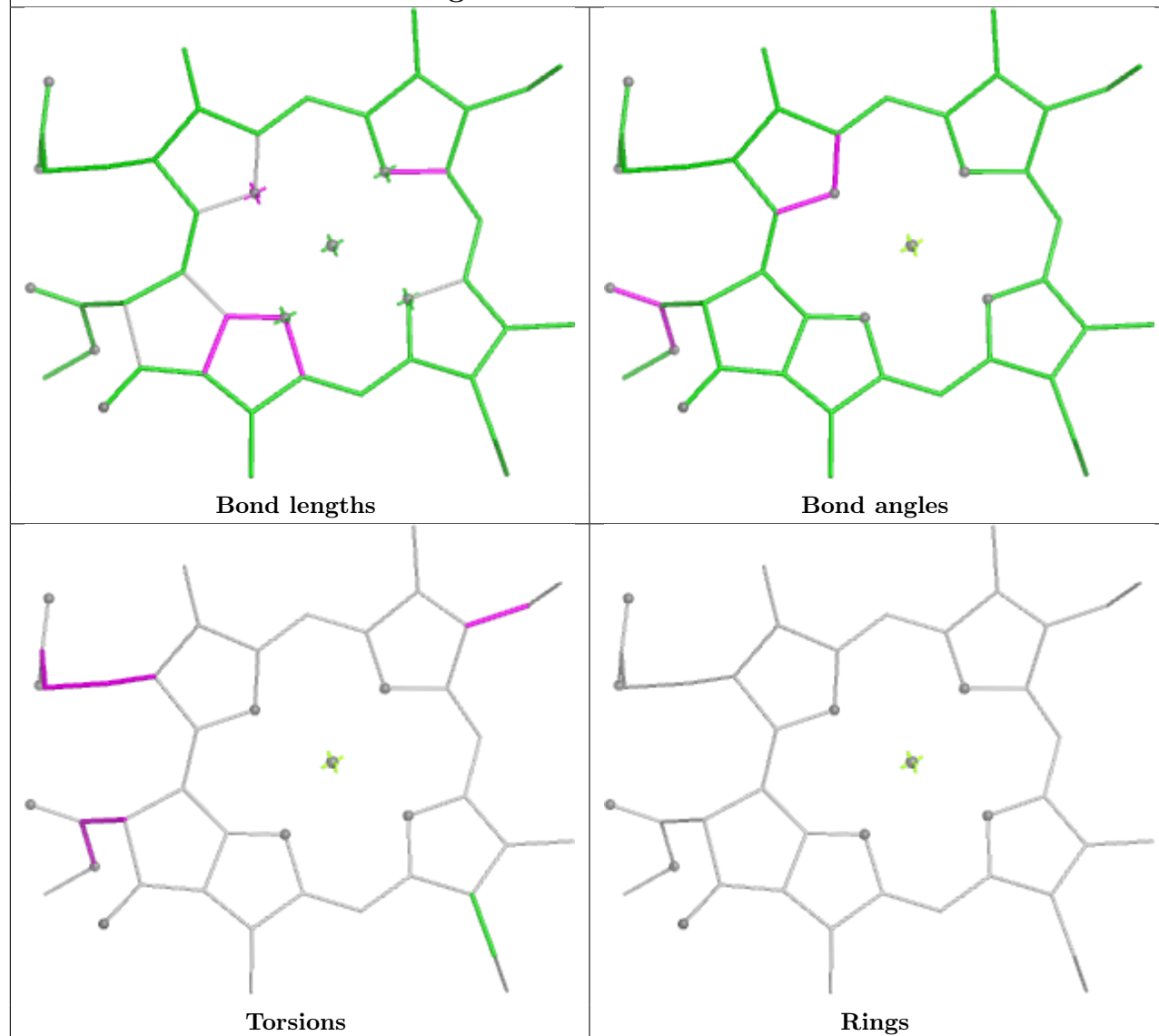
Ligand NDP ZQ 501

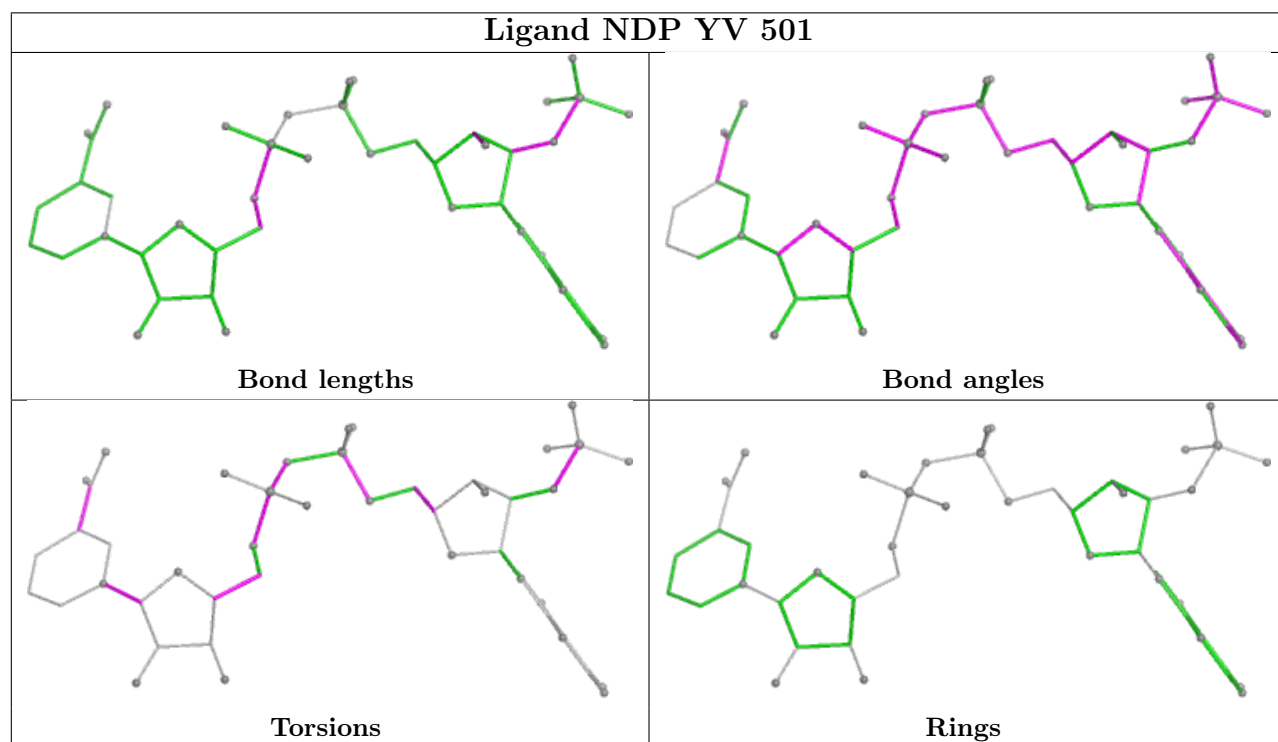
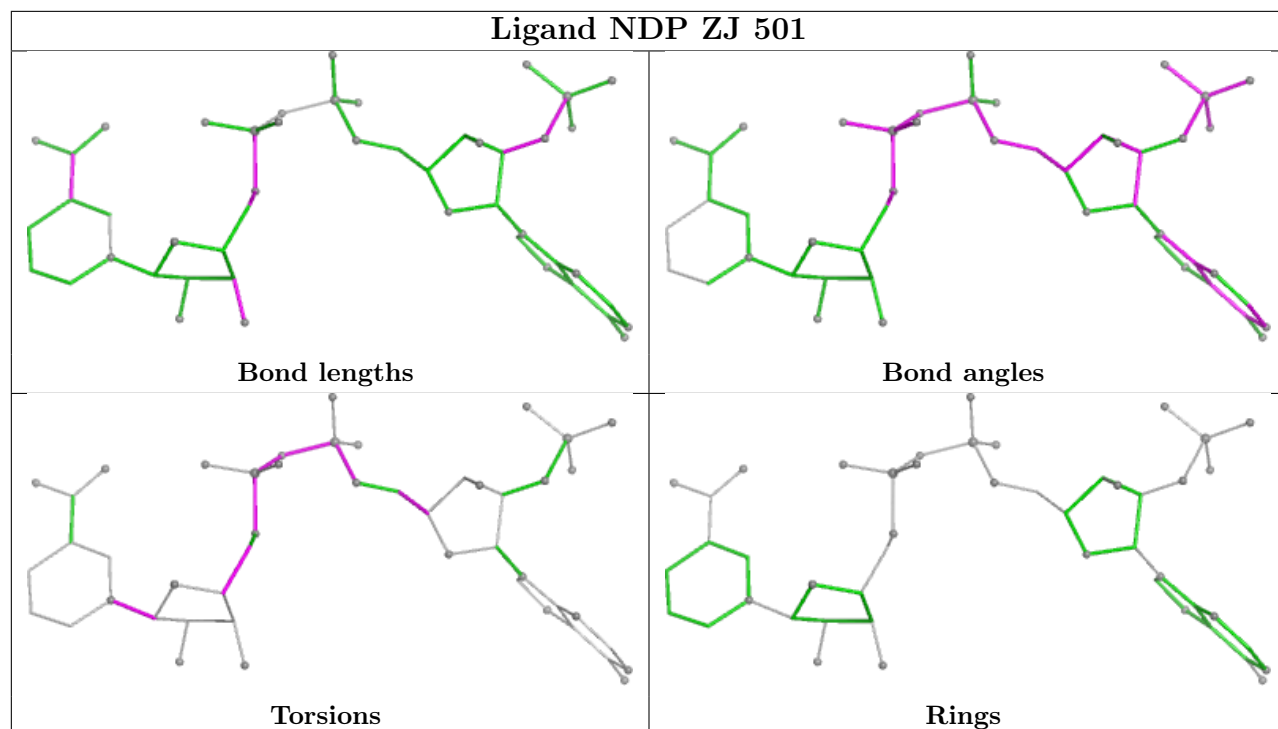




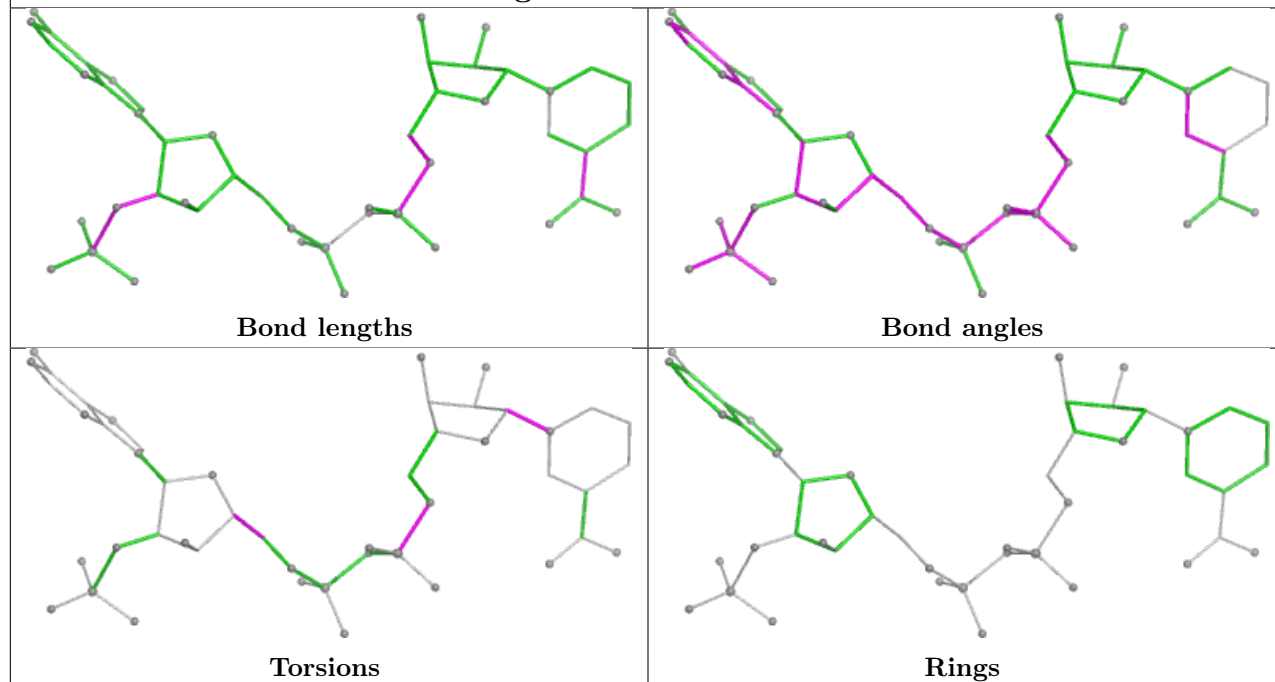


Ligand A1JWG ZK 503

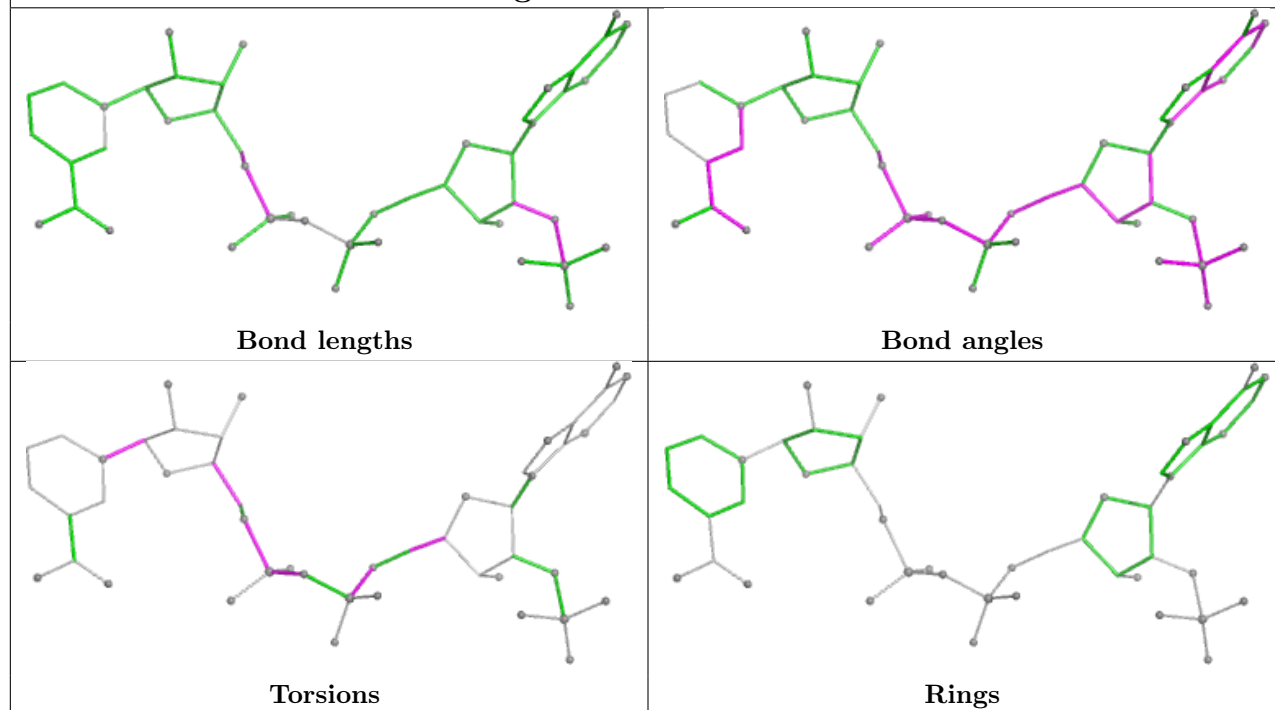




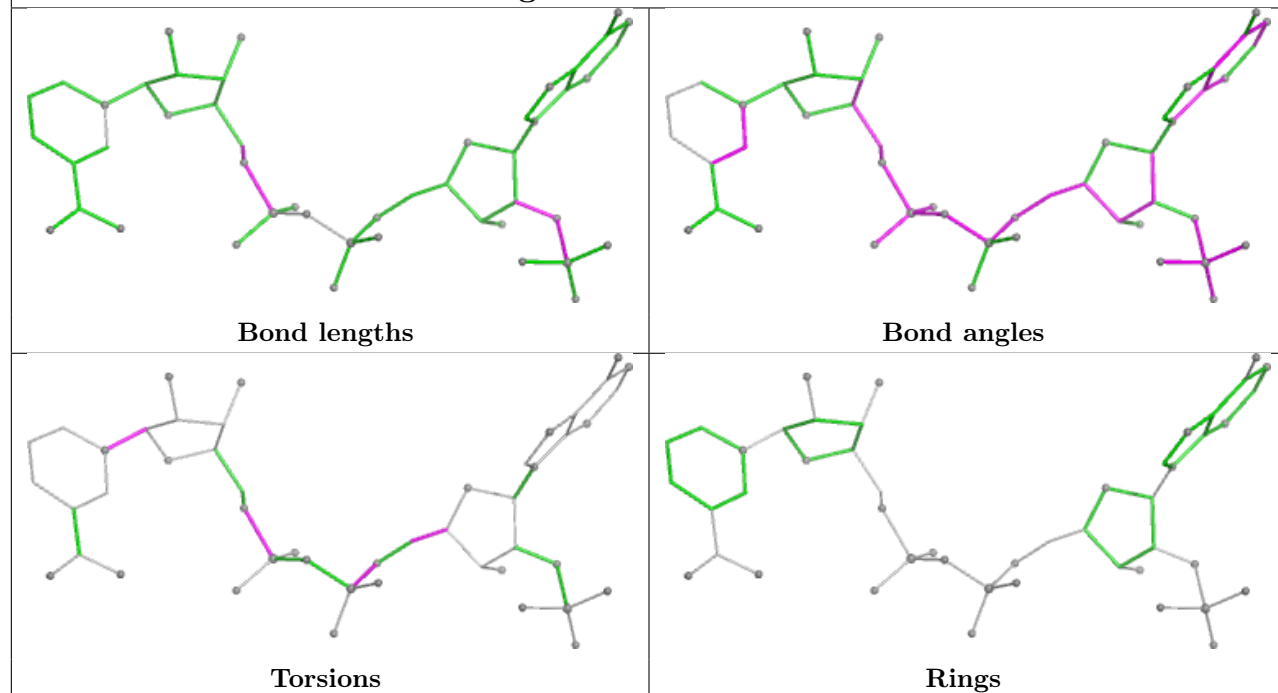
Ligand NDP YX 501



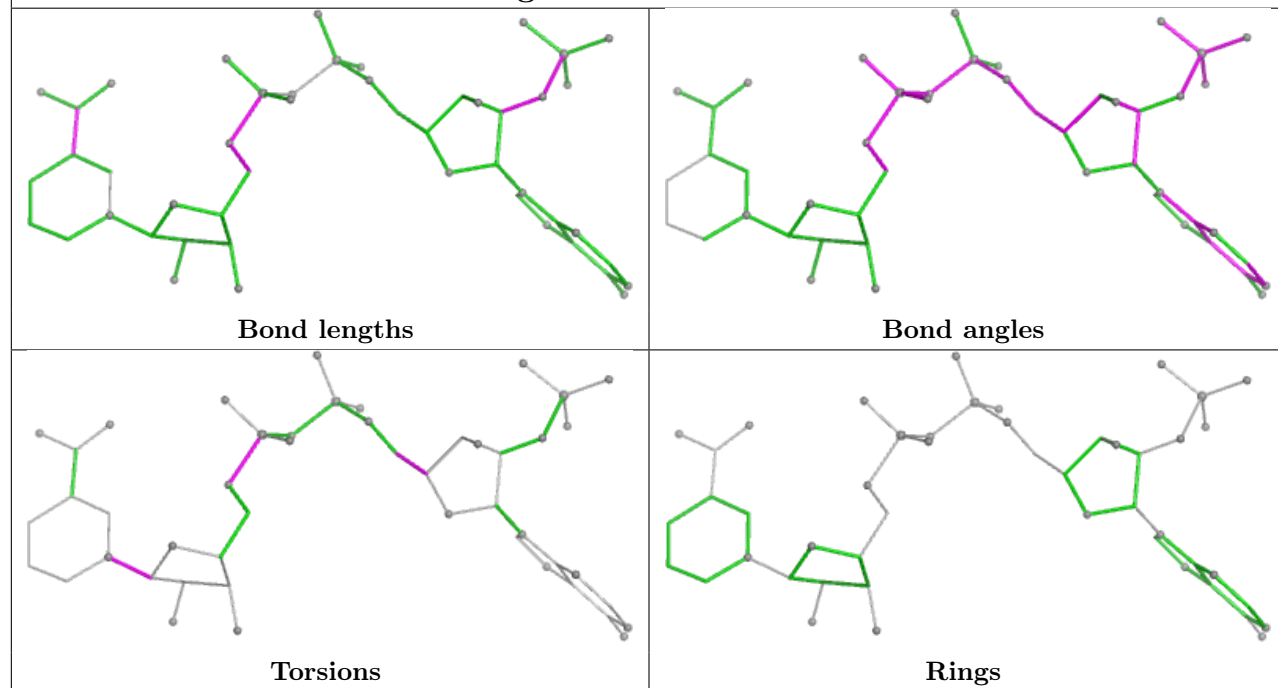
Ligand NDP YU 501

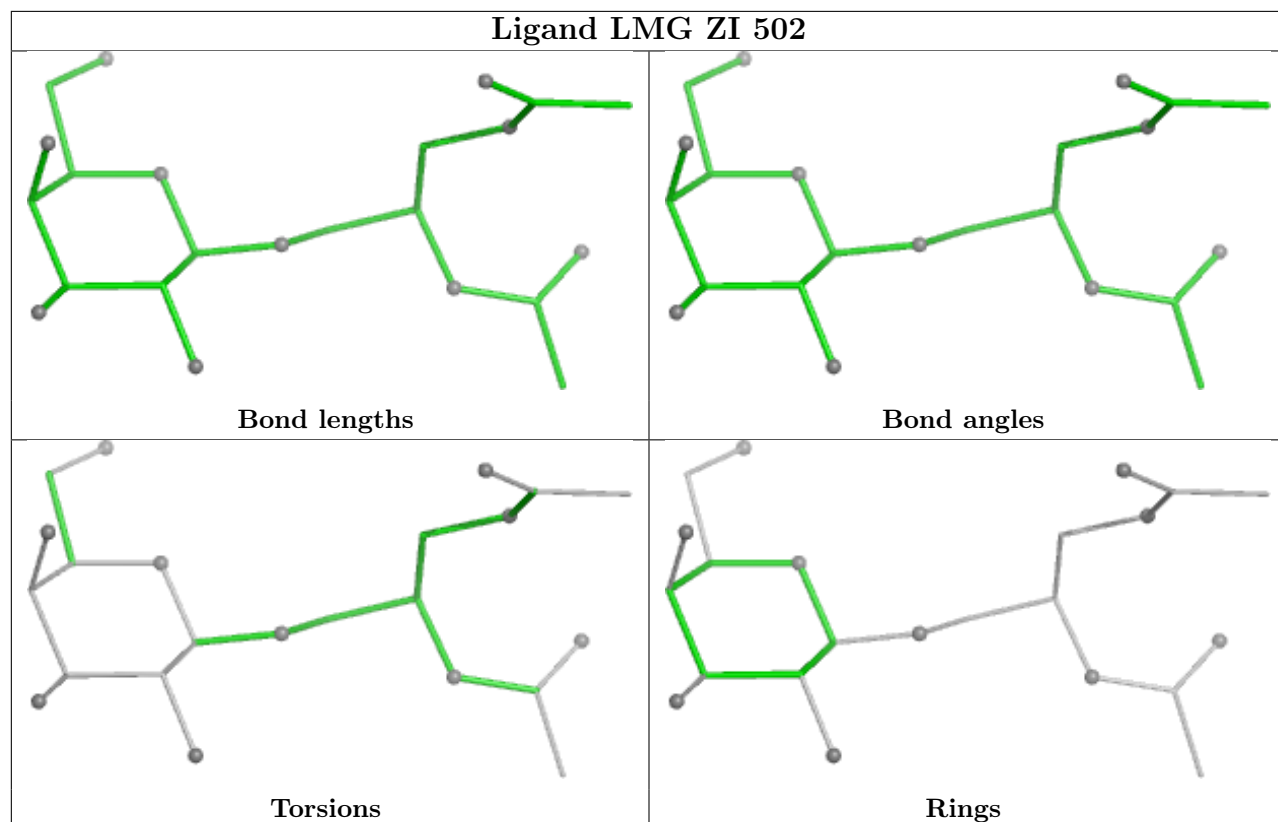
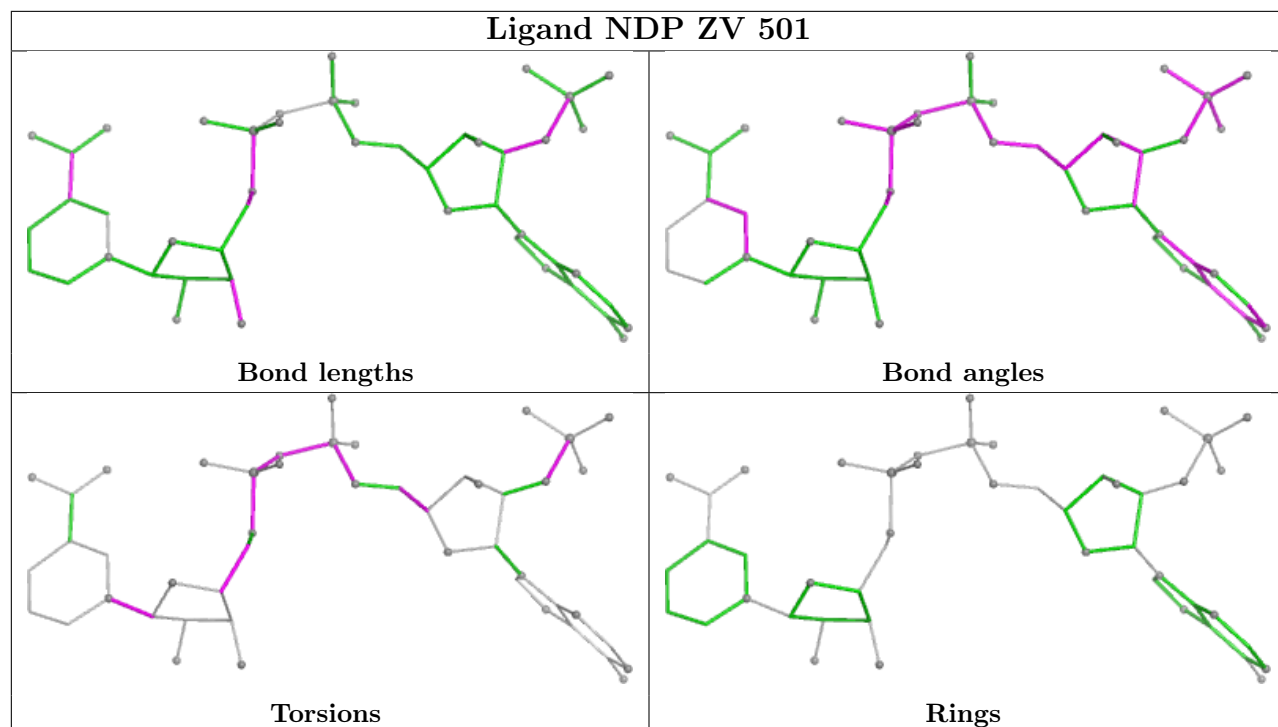


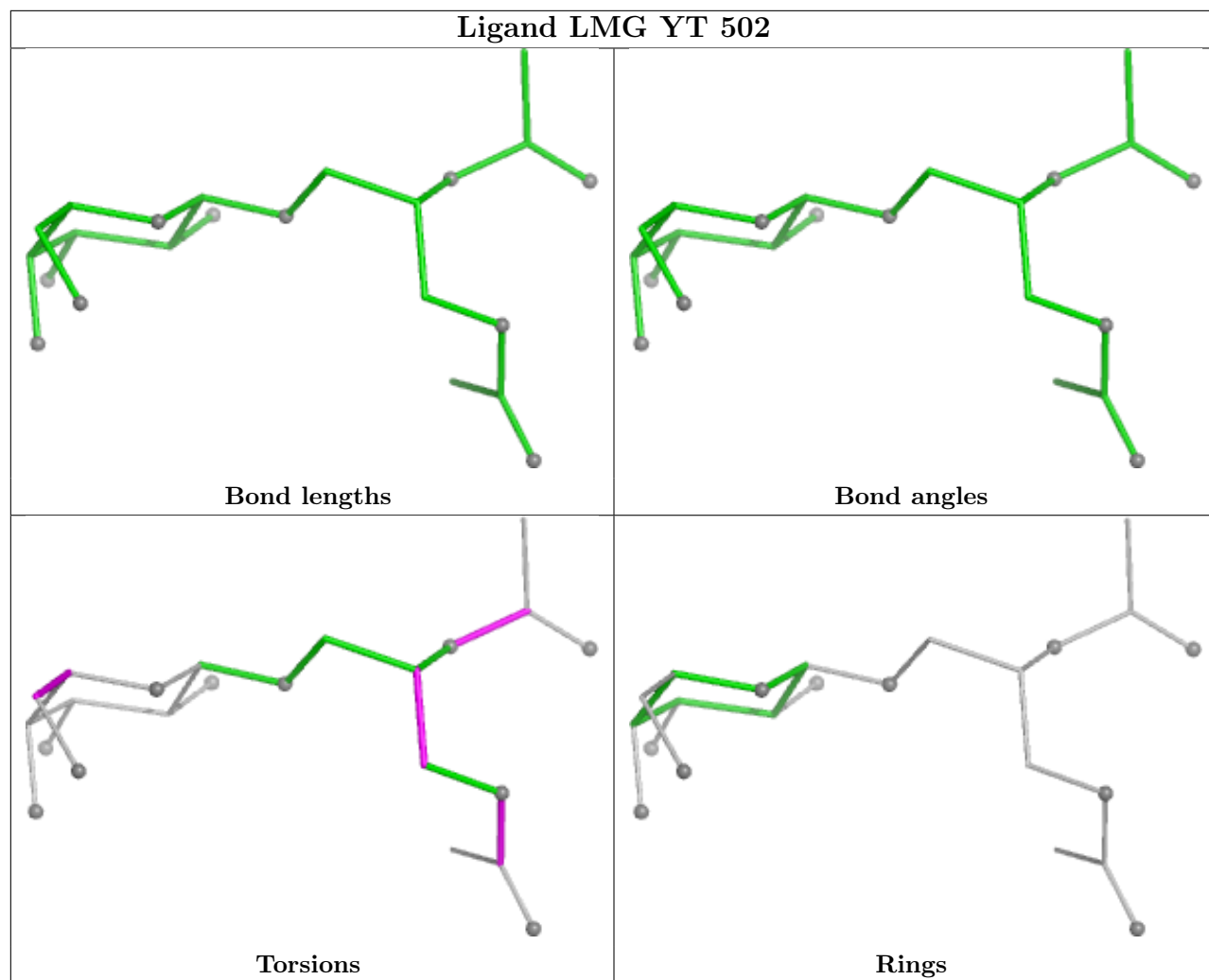
Ligand NDP ZC 501

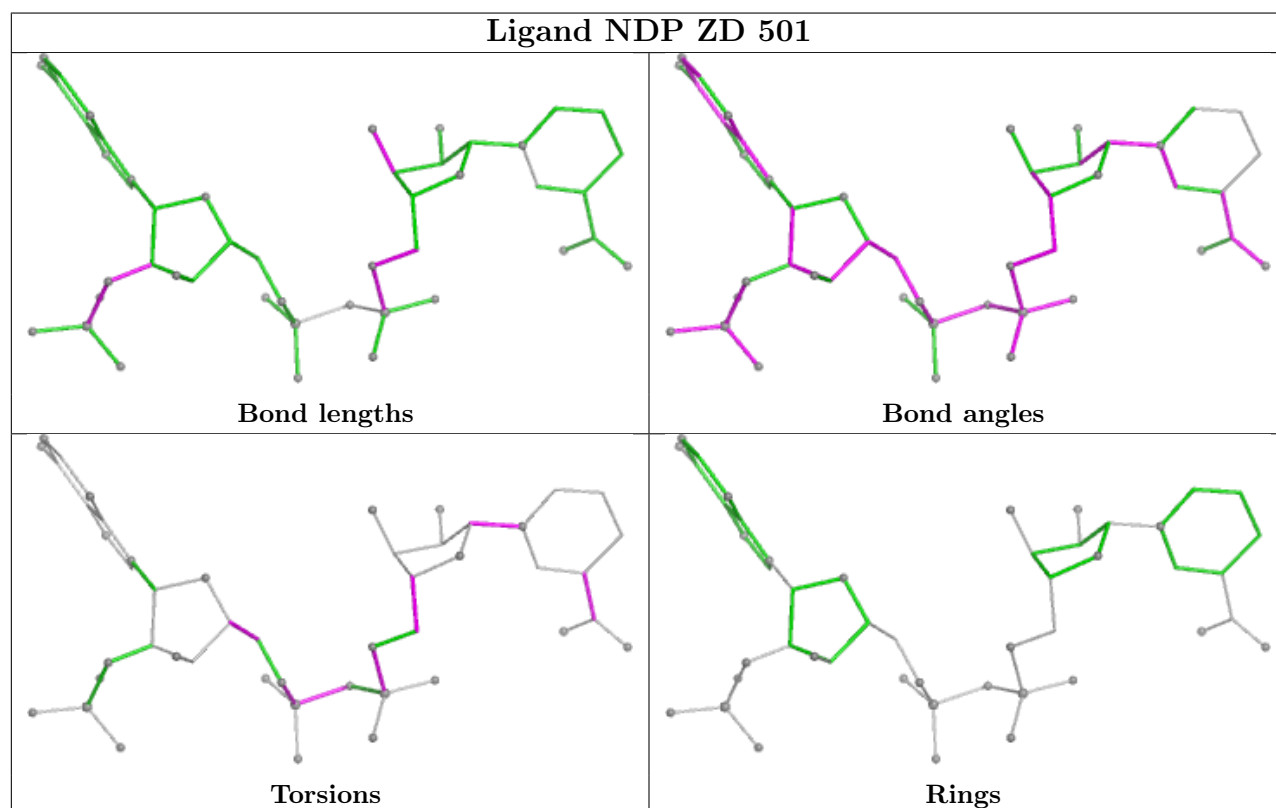
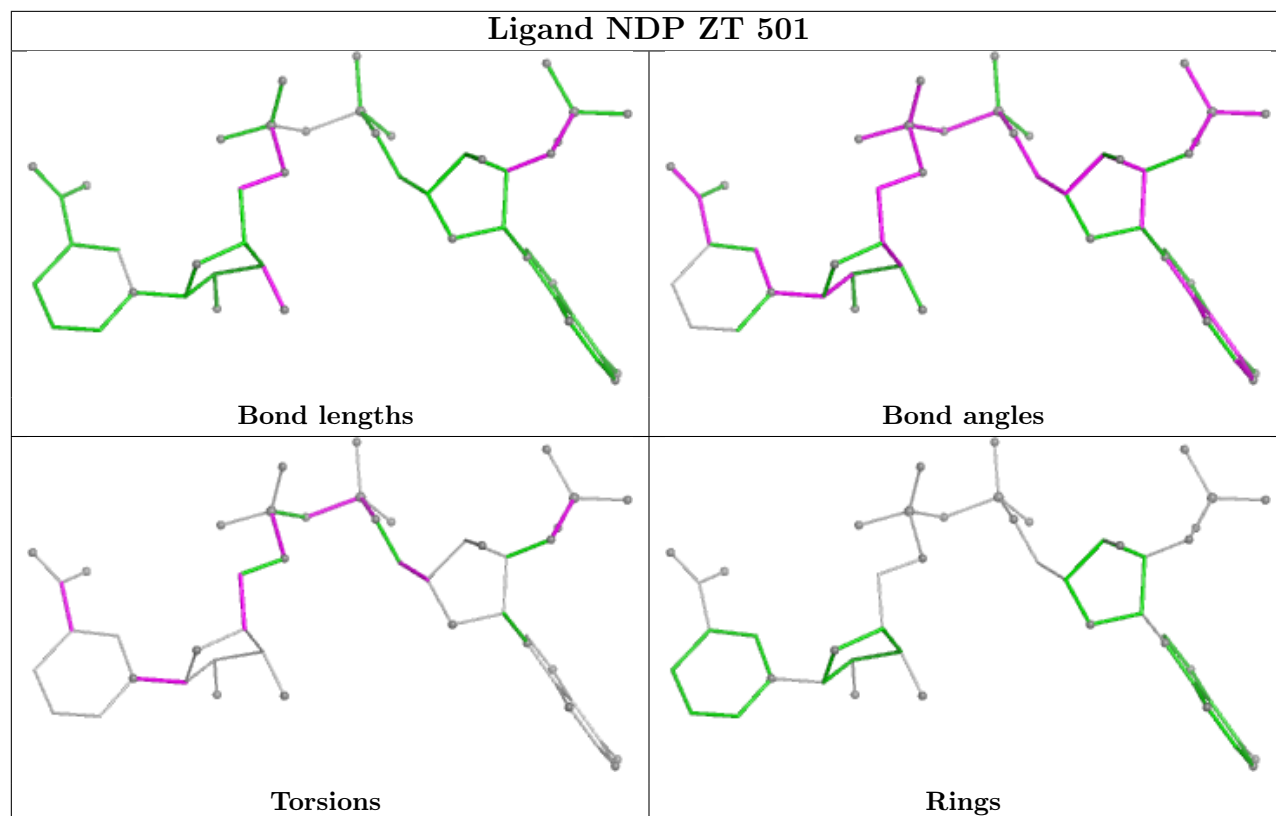


Ligand NDP ZN 501

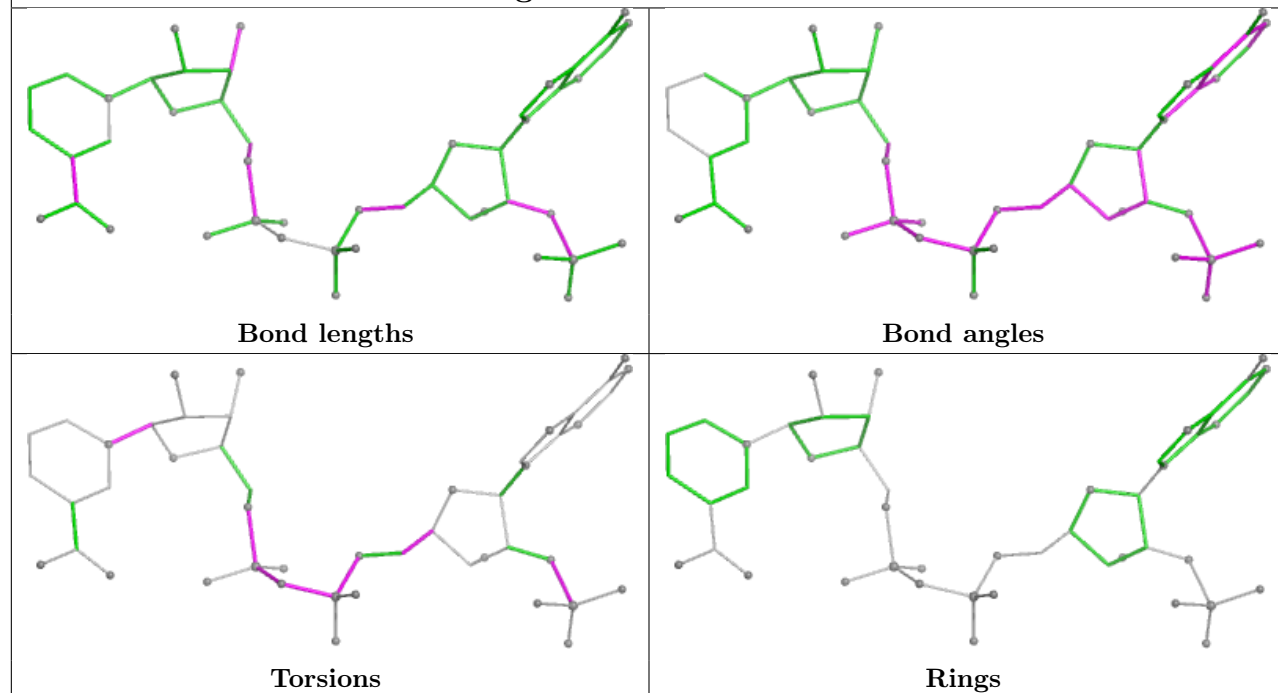




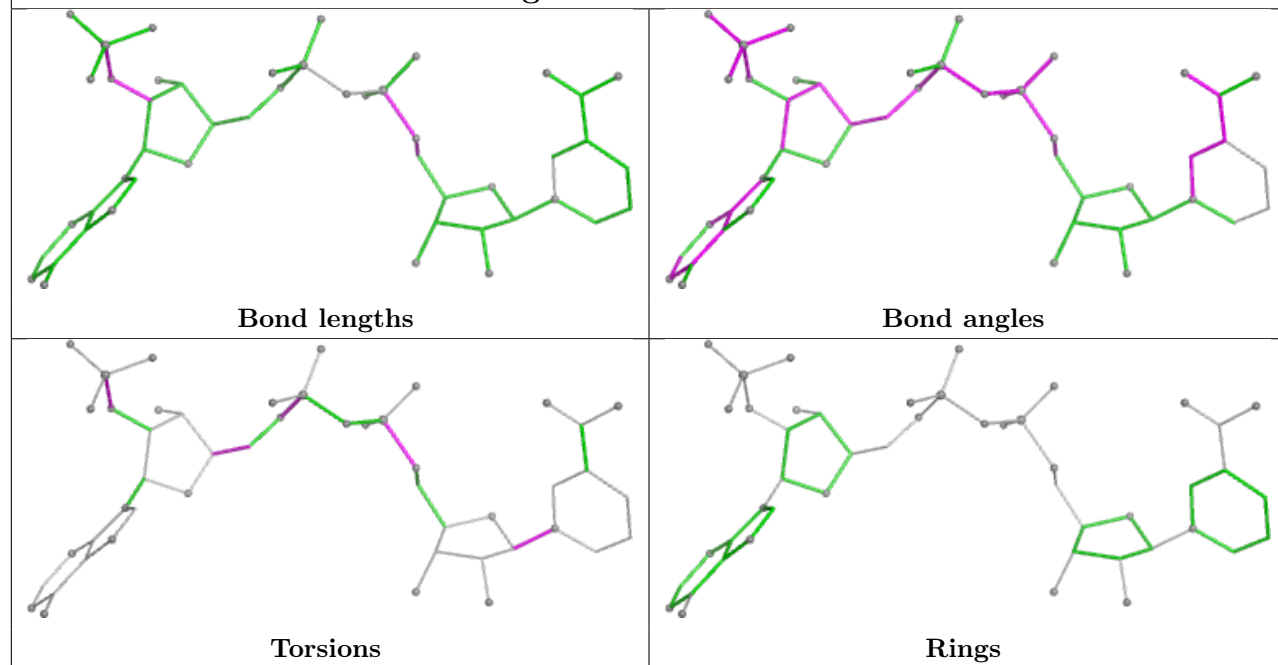


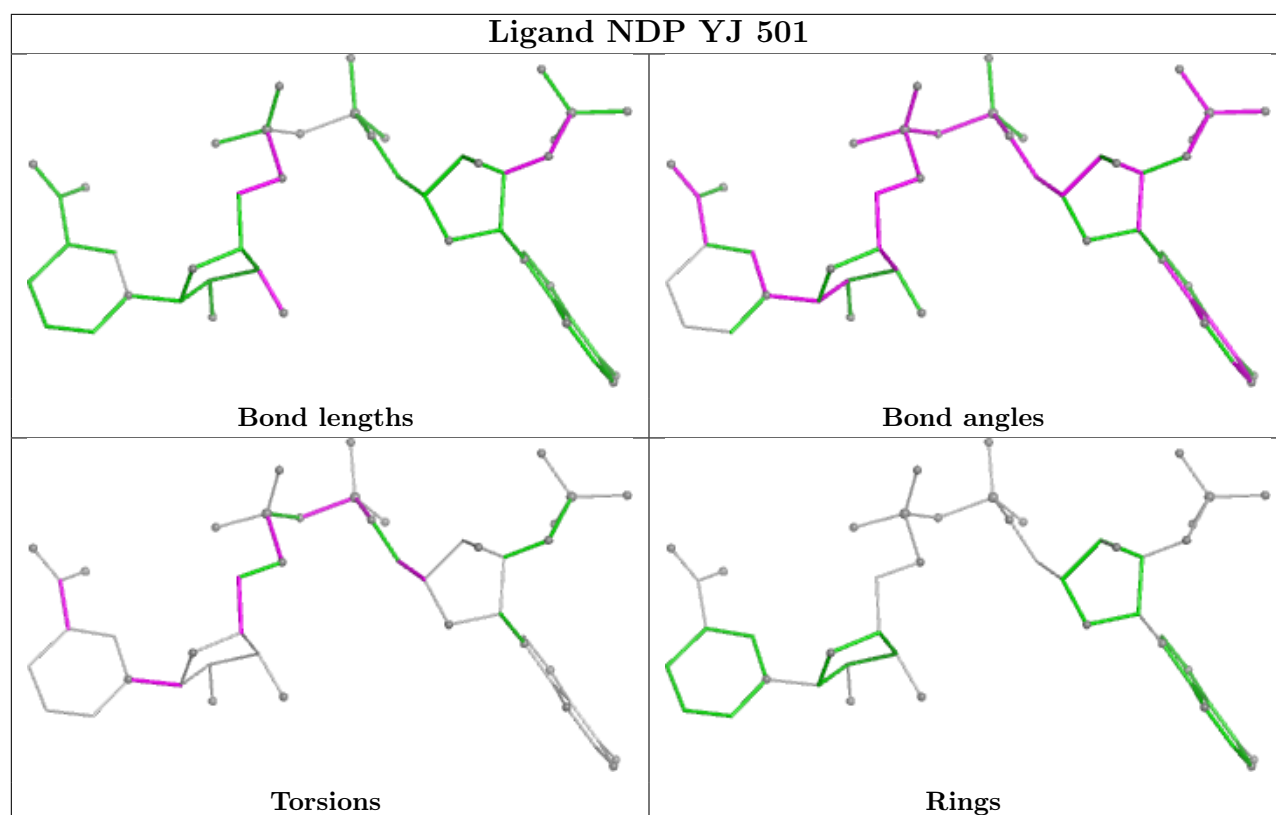


Ligand NDP YO 501



Ligand NDP YI 501





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

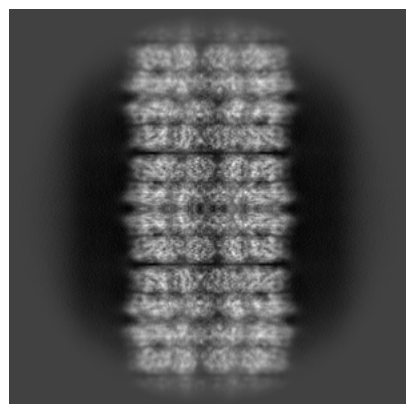
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-56048. These allow visual inspection of the internal detail of the map and identification of artifacts.

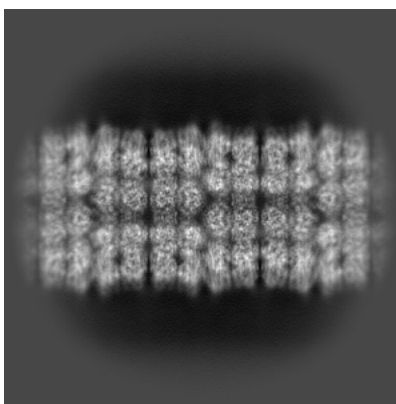
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

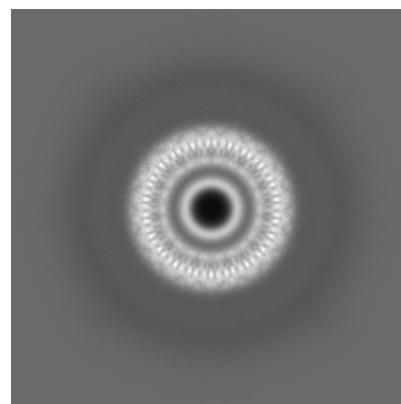
6.1.1 Primary map



X

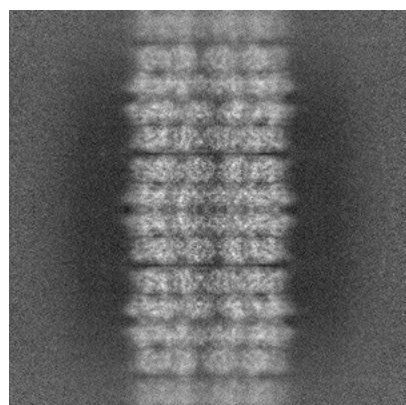


Y

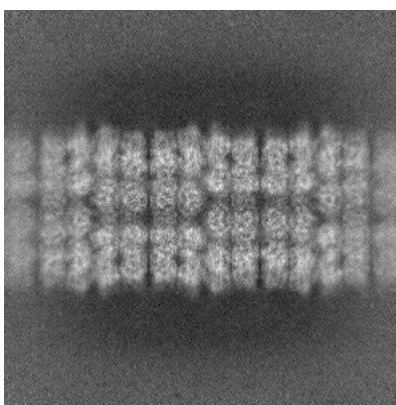


Z

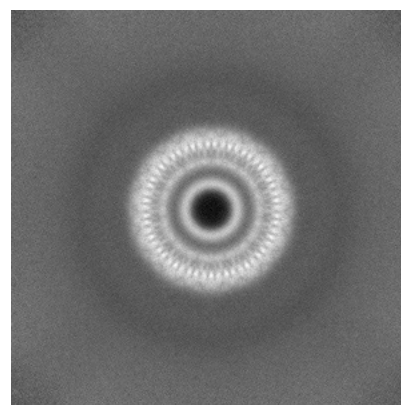
6.1.2 Raw map



X



Y

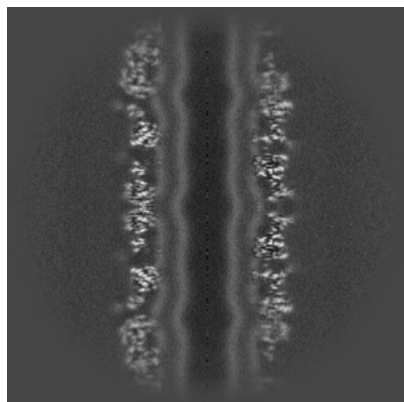


Z

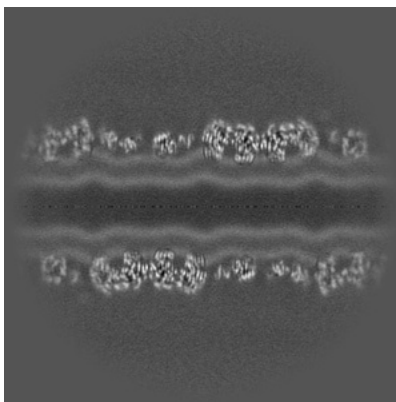
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

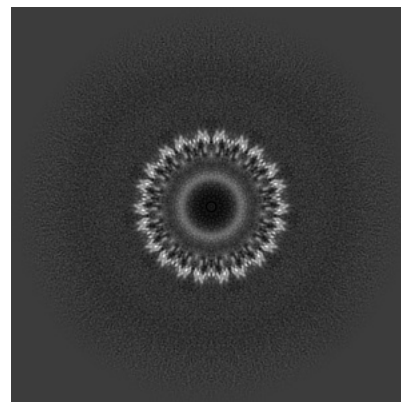
6.2.1 Primary map



X Index: 320

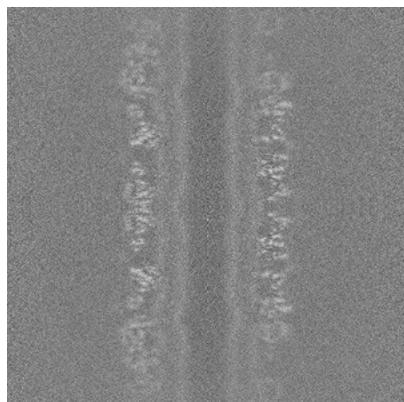


Y Index: 320

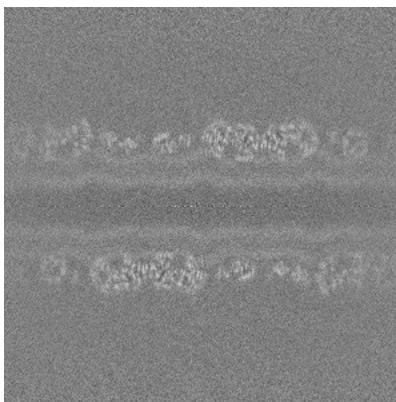


Z Index: 320

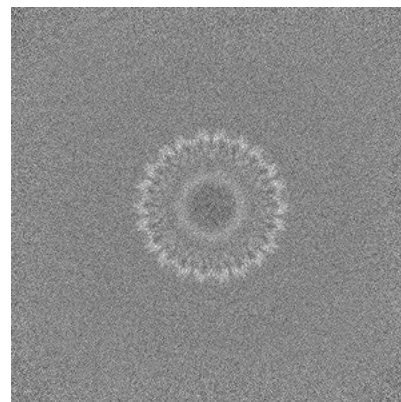
6.2.2 Raw map



X Index: 320



Y Index: 320

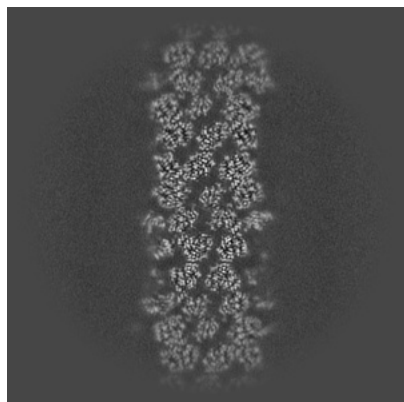


Z Index: 320

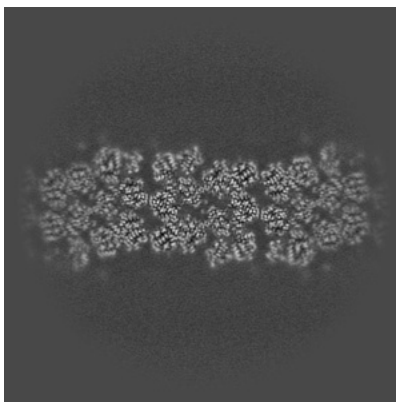
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

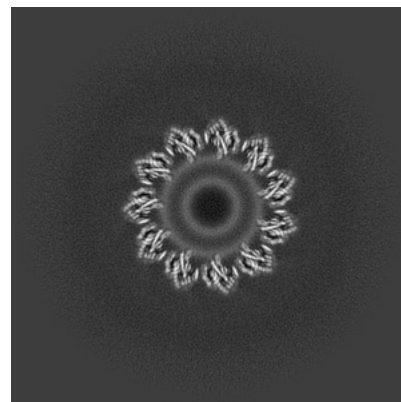
6.3.1 Primary map



X Index: 415

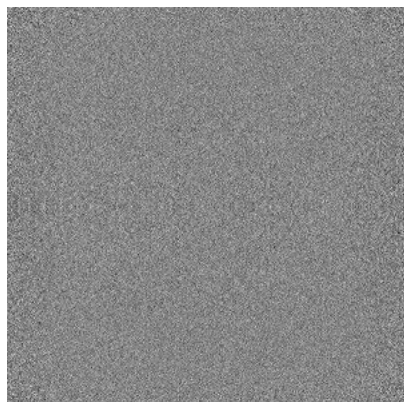


Y Index: 419

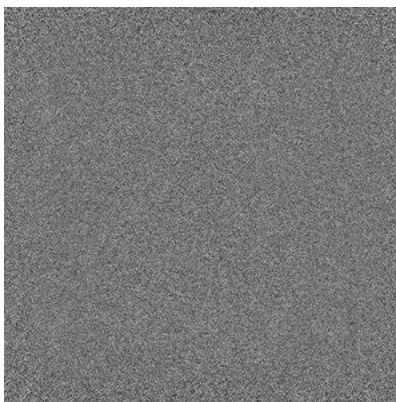


Z Index: 293

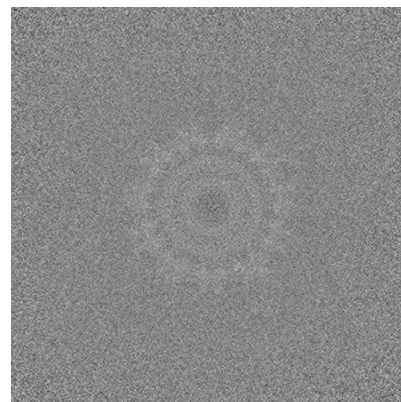
6.3.2 Raw map



X Index: 0



Y Index: 0

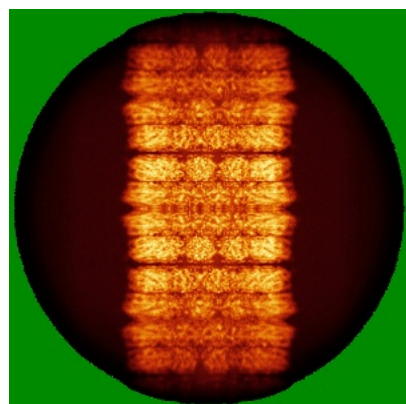


Z Index: 0

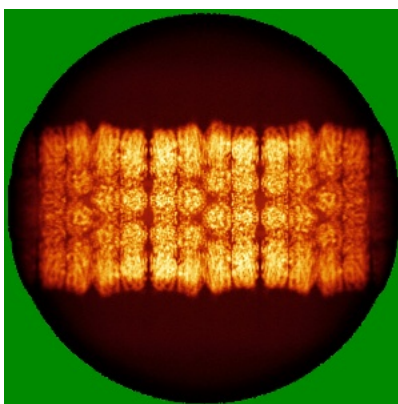
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

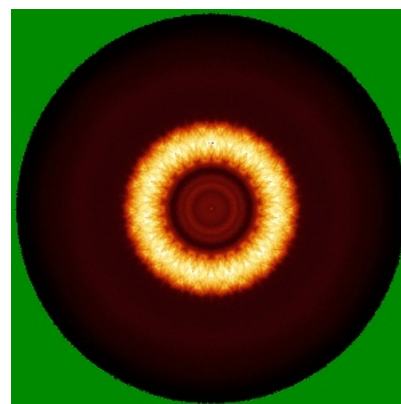
6.4.1 Primary map



X

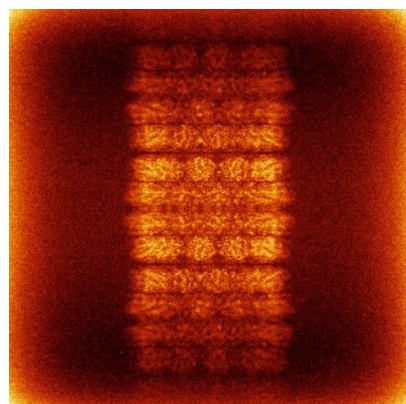


Y

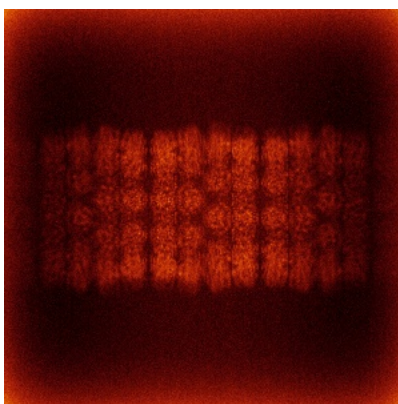


Z

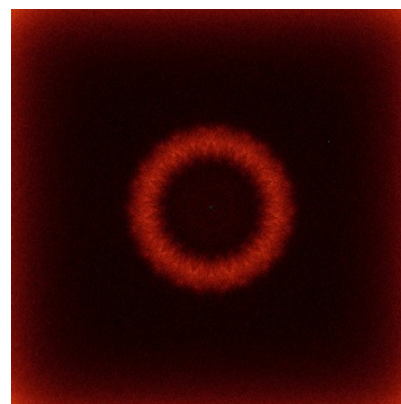
6.4.2 Raw map



X



Y

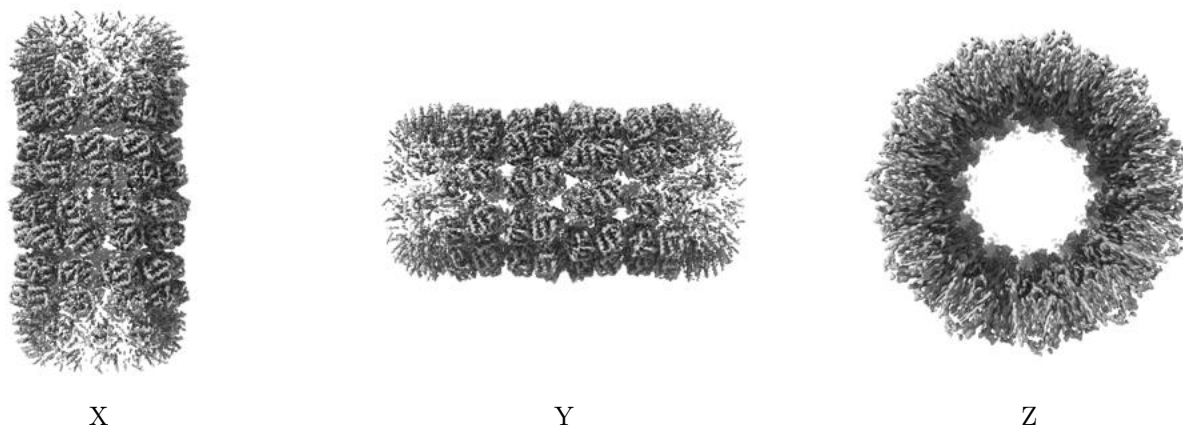


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

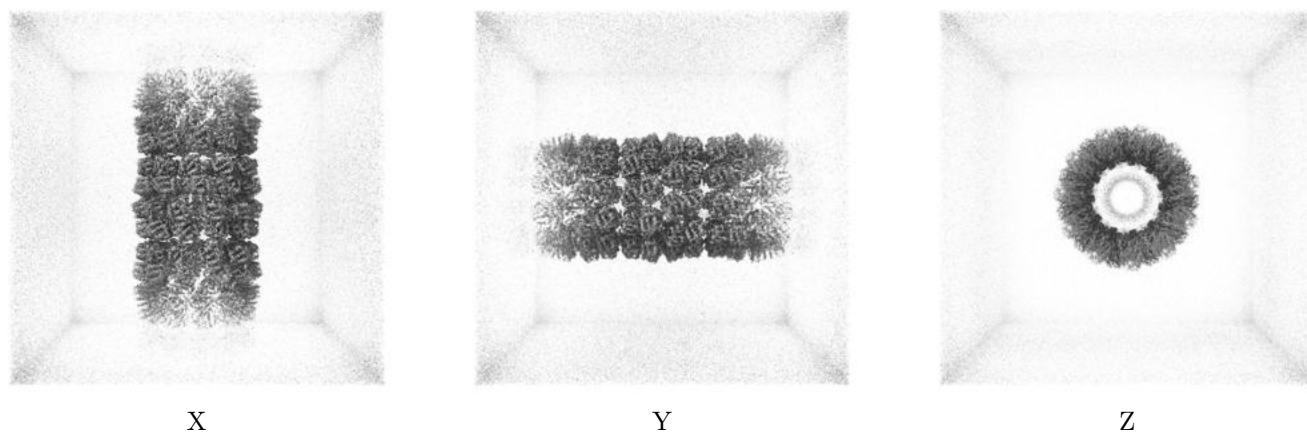
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.104. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

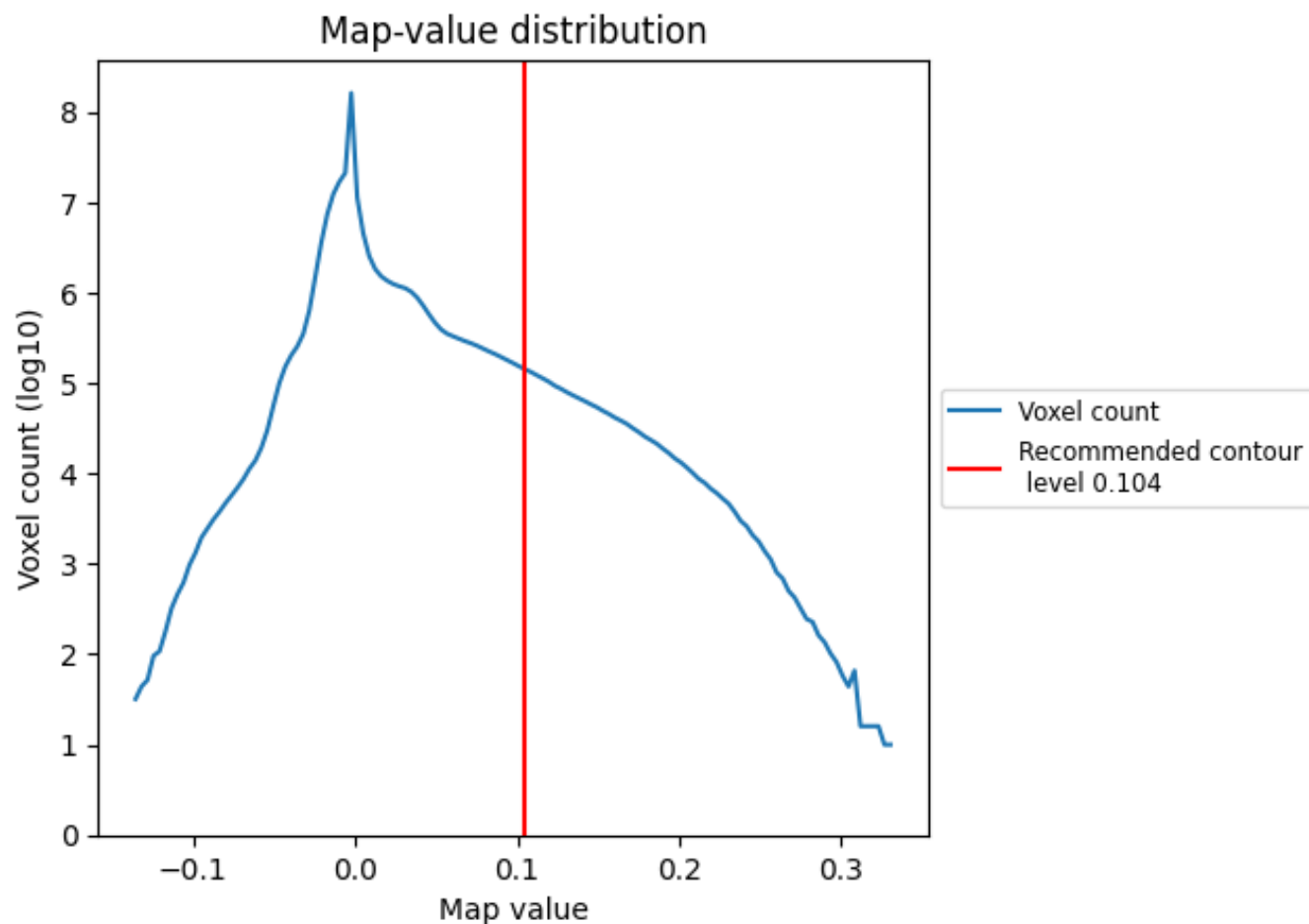
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

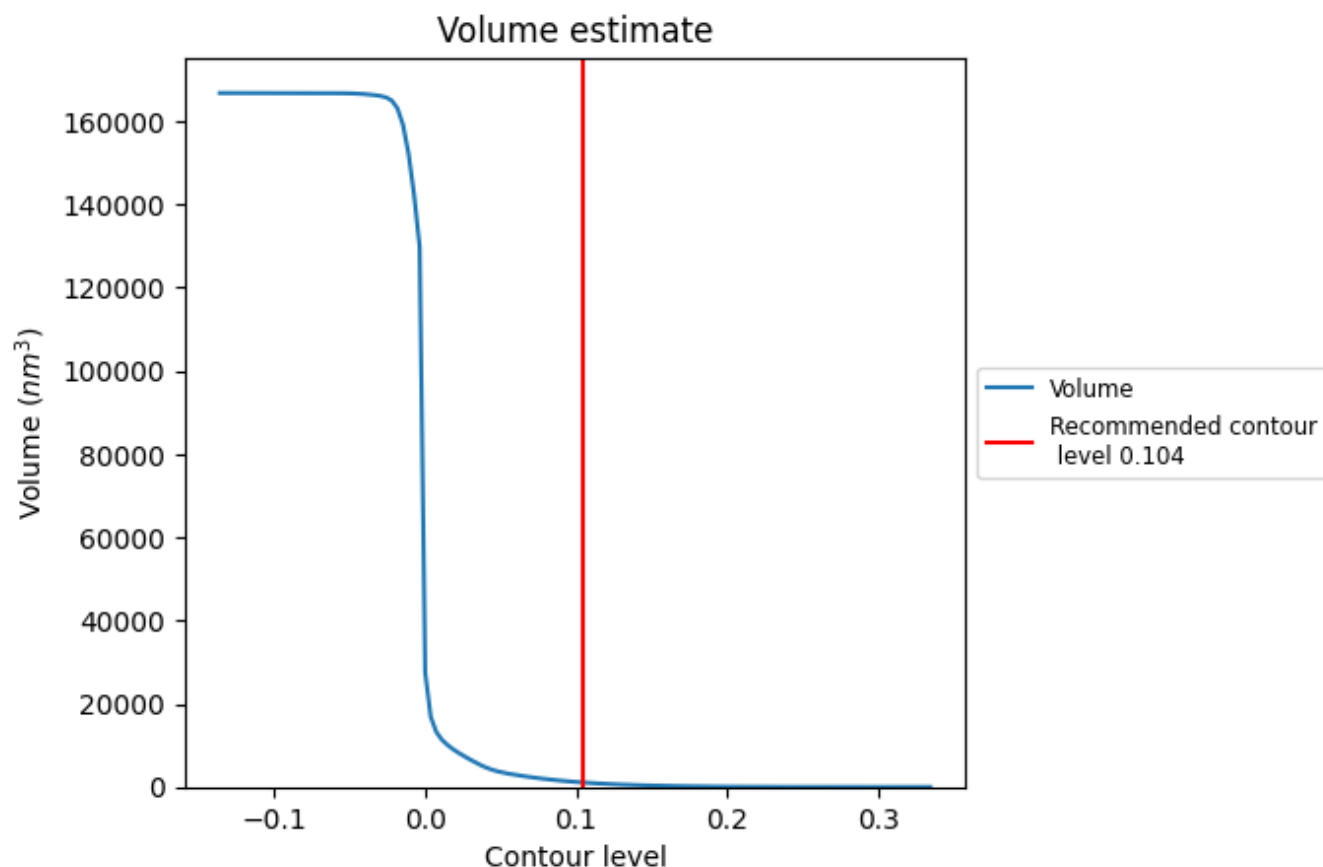
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

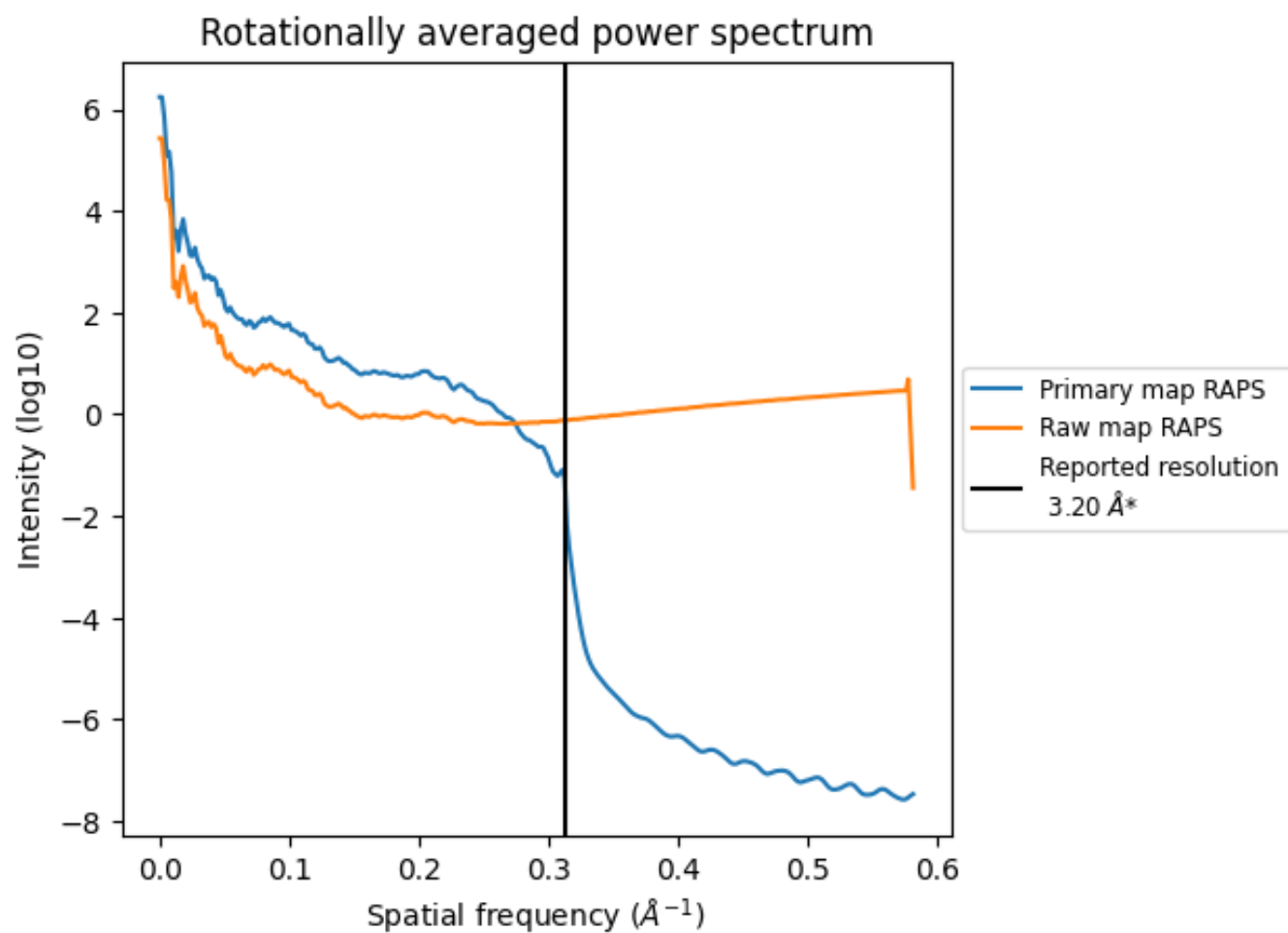
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1085 nm^3 ; this corresponds to an approximate mass of 980 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum [i](#)

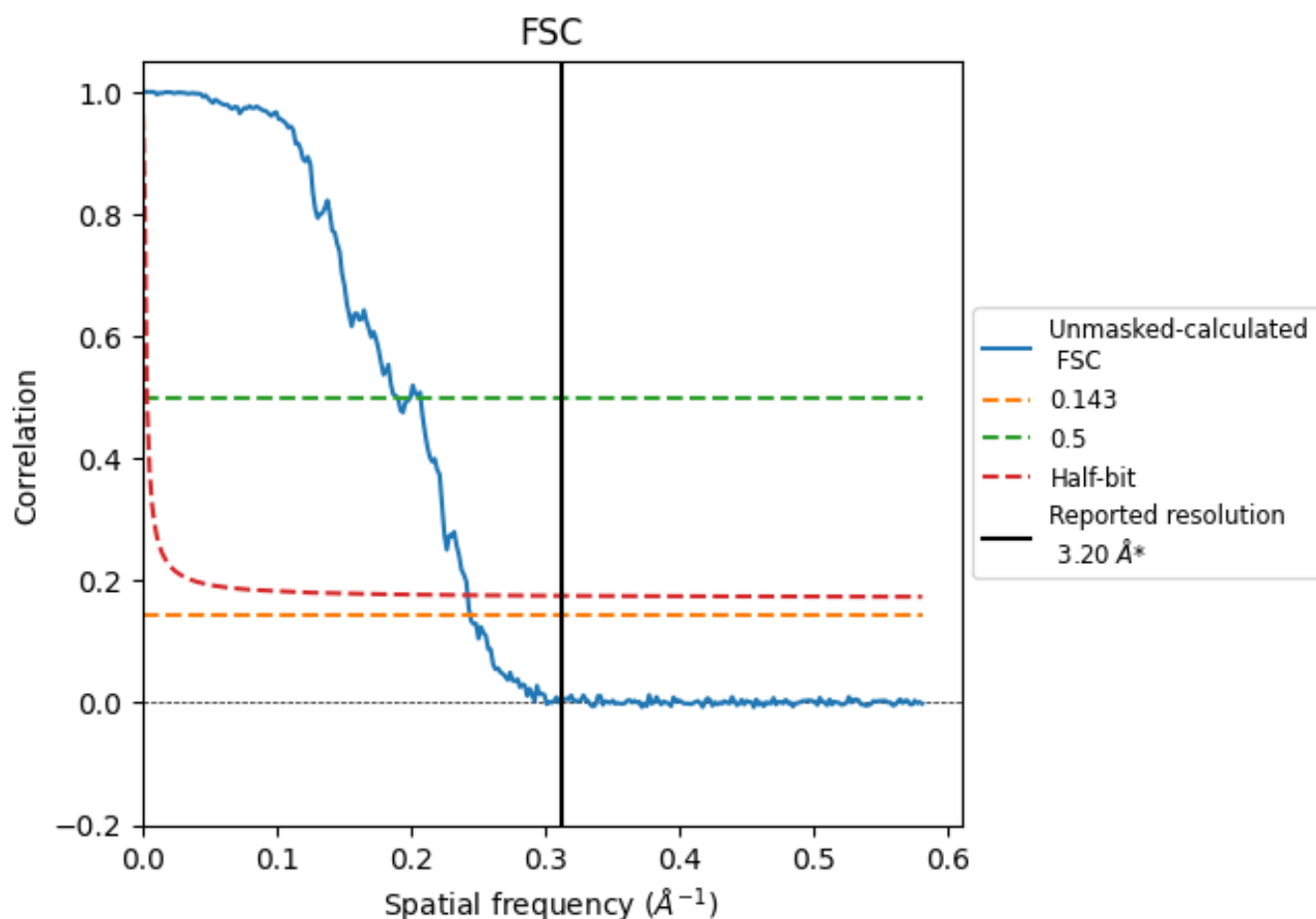


*Reported resolution corresponds to spatial frequency of $0.312 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.312 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

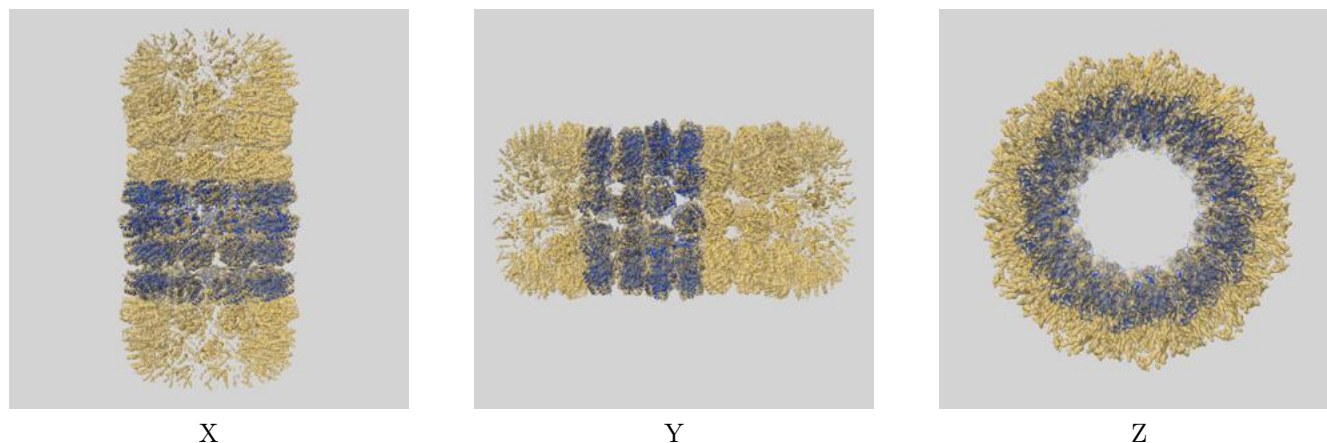
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.20	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.09	5.27	4.12

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.09 differs from the reported value 3.2 by more than 10 %

9 Map-model fit [i](#)

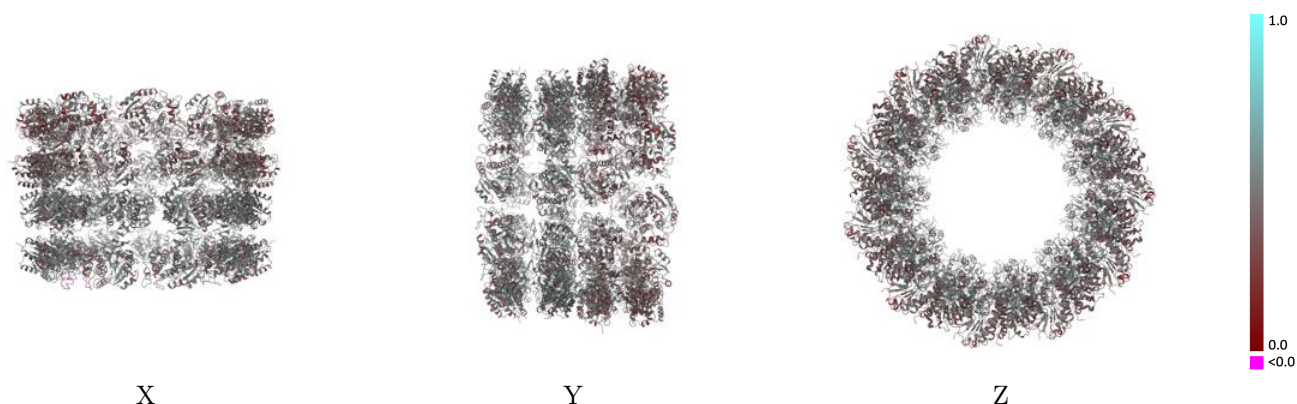
This section contains information regarding the fit between EMDB map EMD-56048 and PDB model 9TLB. Per-residue inclusion information can be found in section 3 on page 36.

9.1 Map-model overlay [i](#)



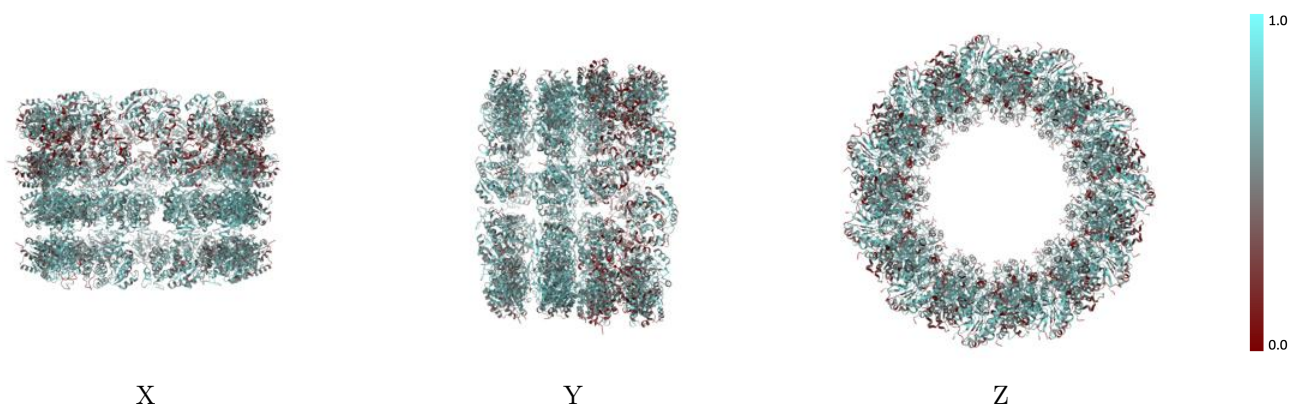
The images above show the 3D surface view of the map at the recommended contour level 0.104 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



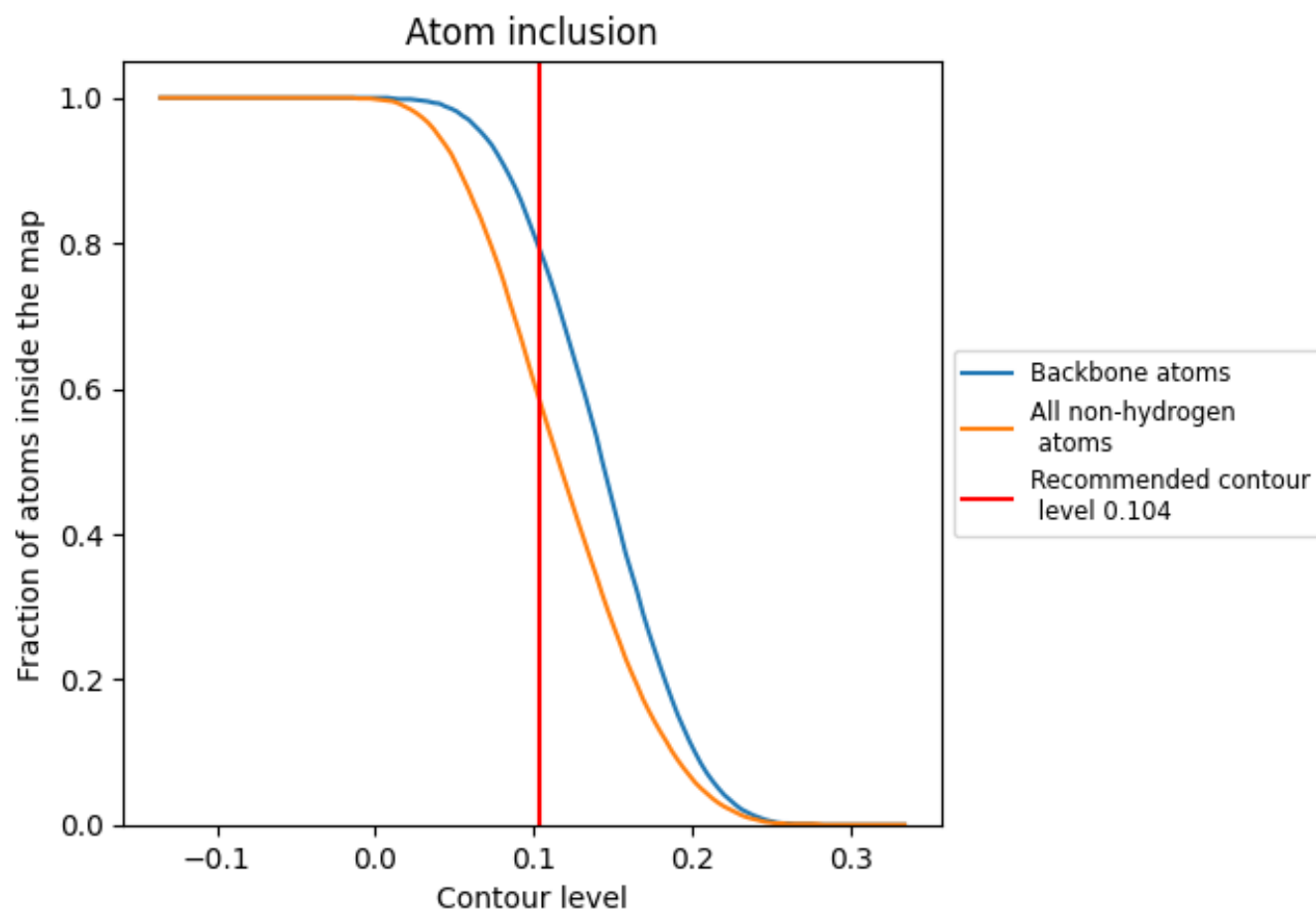
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.104).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 58% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.104) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5820	 0.4450
YI	 0.5180	 0.4120
YJ	 0.5320	 0.4310
YK	 0.6770	 0.4800
YL	 0.6080	 0.4650
YM	 0.5270	 0.4150
YN	 0.5290	 0.4300
YO	 0.6750	 0.4820
YP	 0.6040	 0.4660
YQ	 0.5180	 0.4090
YR	 0.5320	 0.4330
YS	 0.6720	 0.4820
YT	 0.6090	 0.4630
YU	 0.5200	 0.4130
YV	 0.5350	 0.4320
YW	 0.6730	 0.4830
YX	 0.6050	 0.4660
YY	 0.5230	 0.4100
YZ	 0.5170	 0.4110
ZA	 0.6730	 0.4830
ZB	 0.6070	 0.4620
ZC	 0.5210	 0.4130
ZD	 0.5310	 0.4320
ZE	 0.6770	 0.4820
ZF	 0.6070	 0.4660
ZG	 0.5200	 0.4130
ZH	 0.5320	 0.4310
ZI	 0.6720	 0.4830
ZJ	 0.6070	 0.4670
ZK	 0.5250	 0.4110
ZL	 0.5160	 0.4120
ZM	 0.6730	 0.4830
ZN	 0.5890	 0.4480
ZO	 0.5200	 0.4140
ZP	 0.5350	 0.4310



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Chain	Atom inclusion	Q-score
ZQ	 0.6760	 0.4810
ZR	 0.5830	 0.4410
ZS	 0.5280	 0.4160
ZT	 0.5290	 0.4310
ZU	 0.6750	 0.4820
ZV	 0.6050	 0.4660
ZW	 0.5240	 0.4110
ZX	 0.5150	 0.4090
ZY	 0.6740	 0.4820
ZZ	 0.6090	 0.4650