Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 09:23 PM EDT

PDB ID	:	2G87
Title	:	Crystallographic model of bathorhodopsin
Authors	:	Nakamichi, H.; Okada, T.
Deposited on	:	2006-03-02
Resolution	:	2.60 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

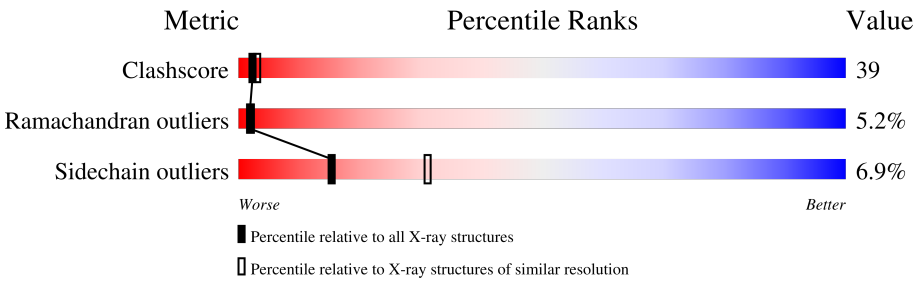
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	349	48% 45% 7%
1	B	349	46% 44% 10%
2	C	3	33% 33% 33%
3	D	2	50% 50%
3	F	2	100%
4	E	4	25% 75%

2 Entry composition (i)

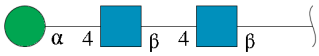
There are 11 unique types of molecules in this entry. The entry contains 5948 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Rhodopsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	349	2749	1818	424	481	26	0	0	0
1	B	349	2749	1818	424	481	26	0	0	0

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	3	39	22	2	15	0	0	0

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	D	2	28	16	2	10	0	0	0
3	F	2	28	16	2	10	0	0	0

- Molecule 4 is an oligosaccharide called beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	E	4	50	28	2	20	0	0	0

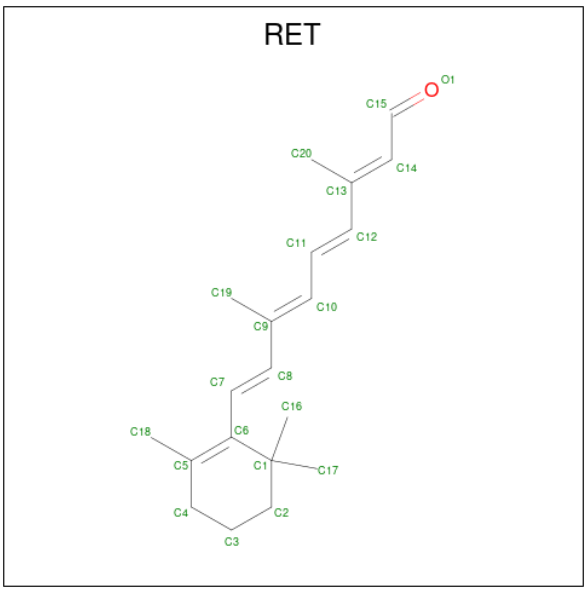
- Molecule 5 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Hg		
5	A	3	3	3	0	0
5	B	3	3	3	0	0

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

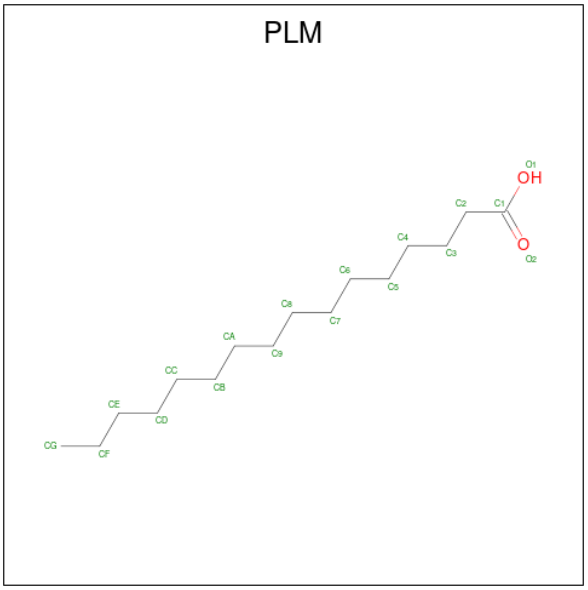
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
6	A	4	4	4	0	0
6	B	3	3	3	0	0

- Molecule 7 is RETINAL (three-letter code: RET) (formula: $C_{20}H_{28}O$).



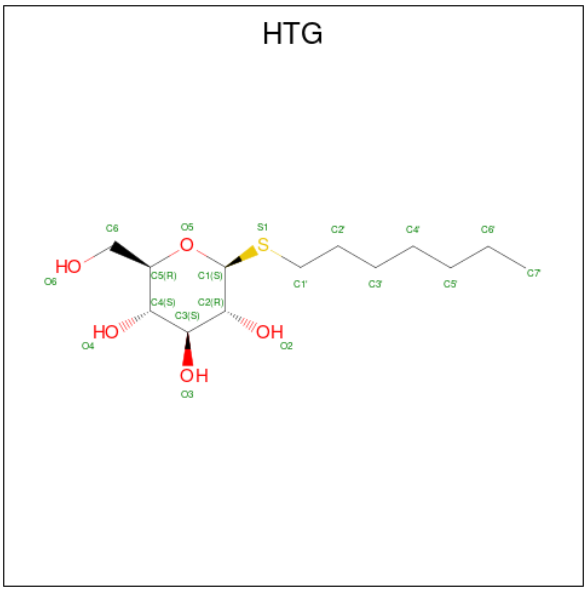
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C 20 20	0	0
7	B	1	Total C 20 20	0	0

- Molecule 8 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



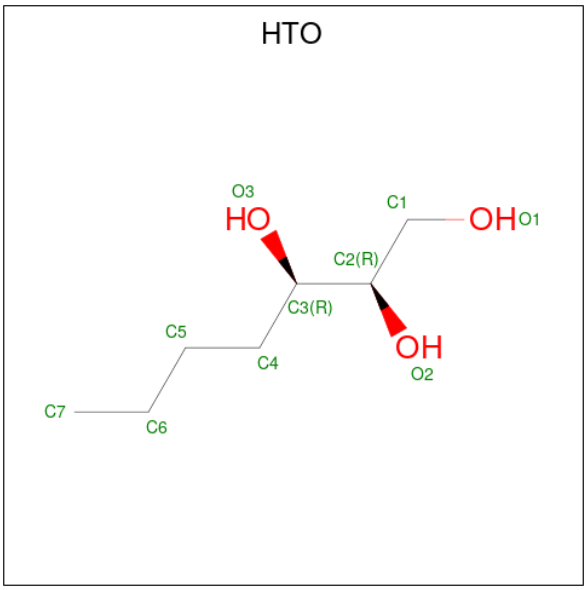
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 17 16 1	0	0
8	A	1	Total C O 17 16 1	0	0
8	A	1	Total C 16 16	0	0
8	B	1	Total C O 17 16 1	0	0
8	B	1	Total C O 17 16 1	0	0
8	B	1	Total C 16 16	0	0

- Molecule 9 is heptyl 1-thio-beta-D-glucopyranoside (three-letter code: HTG) (formula: $C_{13}H_{26}O_5S$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	O	S	0	0
			19	13	5	1		
9	B	1	Total	C	O	S	0	0
			19	13	5	1		
9	B	1	Total	C	O	S	0	0
			19	13	5	1		
9	B	1	Total	C	O	S	0	0
			19	13	5	1		

- Molecule 10 is HEPTANE-1,2,3-TRIOL (three-letter code: HTO) (formula: $C_7H_{16}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
10	B	1	10	7	3	0	0

- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
11	A	37	37	37	0	0
			Total	O		
11	B	29	29	29	0	0

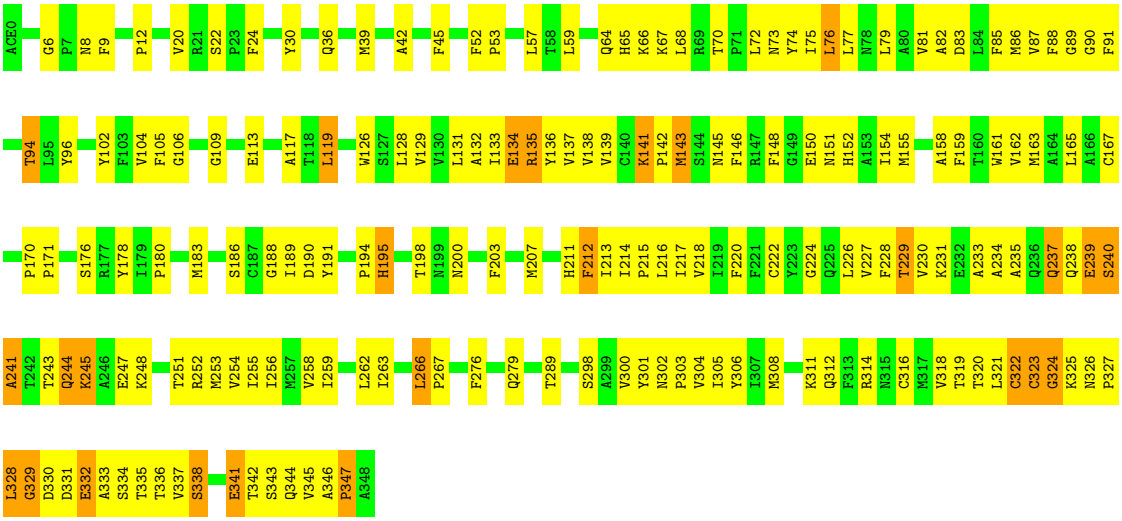
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

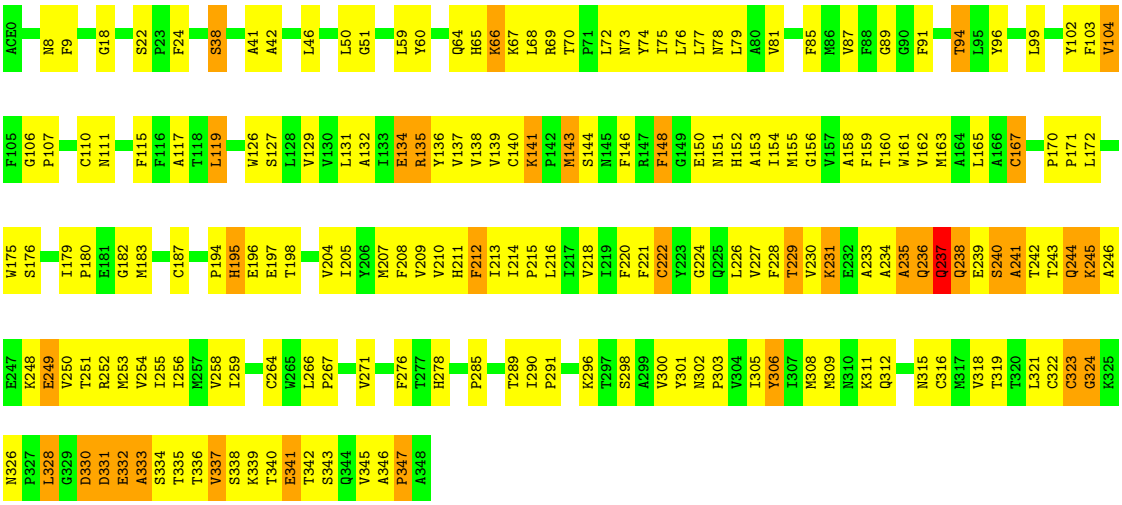
• Molecule 1: Rhodopsin

Chain A: 48% 45% 7%

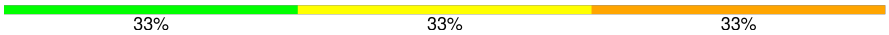


• Molecule 1: Rhodopsin

Chain B: 46% 44% 10%



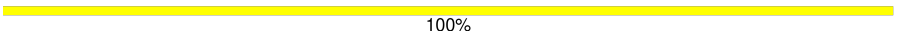
• Molecule 2: alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C: 

• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: 

• Molecule 4: beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α, β, γ	96.49Å 96.49Å 150.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (50.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.178 , 0.181	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5948	wwPDB-VP
Average B, all atoms (Å^2)	65.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: HG, BMA, ZN, RET, HTO, NAG, PLM, HTG, MAN, ACE

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	A	0.61	0/2831	0.67	0/3859
1	B	0.60	0/2831	0.68	0/3859
All	All	0.60	0/5662	0.67	0/7718

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	306	TYR	Sidechain

5.2 Too-close contacts (i)

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	A	2749	0	2706	217	0
1	B	2749	0	2707	239	0
2	C	39	0	34	1	0
3	D	28	0	25	1	0
3	F	28	0	25	0	0
4	E	50	0	43	2	0
5	A	3	0	0	0	0
5	B	3	0	0	0	0
6	A	4	0	0	0	0
6	B	3	0	0	0	0
7	A	20	0	27	3	0
7	B	20	0	27	4	0
8	A	50	0	89	4	0
8	B	50	0	89	5	0
9	A	19	0	26	0	0
9	B	57	0	78	5	0
10	B	10	0	16	1	0
11	A	37	0	0	3	0
11	B	29	0	0	4	0
All	All	5948	0	5892	454	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 39.

All (454) 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:B:312:GLN:HA	1:B:332:GLU:HG2	1.32	1.08
1:A:65:HIS:HB3	1:A:337:VAL:HG22	1.37	1.03
1:A:345:VAL:HG12	1:A:347:PRO:HD3	1.37	1.03
1:B:239:GLU:HB2	1:B:245:LYS:HD2	1.40	1.03
1:B:245:LYS:HE3	1:B:245:LYS:HA	1.39	1.02
1:B:245:LYS:NZ	1:B:248:LYS:HD2	1.76	1.01
1:A:316:CYS:SG	1:A:337:VAL:HG13	2.05	0.96
1:A:311:LYS:HD2	1:A:314:ARG:NH2	1.79	0.96
1:A:326:ASN:O	1:A:328:LEU:HD22	1.71	0.91
1:A:88:PHE:HB3	8:B:1323:PLM:HB1	1.54	0.90
1:B:346:ALA:N	1:B:347:PRO:HD3	1.86	0.90
1:A:94:THR:HB	1:A:113:GLU:OE2	1.73	0.89
1:B:322:CYS:HA	8:B:1322:PLM:O1	1.73	0.88
1:B:72:LEU:HD22	1:B:250:VAL:HG13	1.55	0.87
1:B:239:GLU:HB3	1:B:244:GLN:HE22	1.40	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:PHE:HA	1:A:94:THR:CG2	2.06	0.85
1:B:337:VAL:HB	1:B:343:SER:HA	1.59	0.85
1:A:326:ASN:HD21	1:A:328:LEU:HD11	1.41	0.85
1:B:238:GLN:HG3	1:B:241:ALA:H	1.41	0.84
1:B:245:LYS:HZ1	1:B:248:LYS:HD2	1.37	0.84
1:A:70:THR:H	1:A:73:ASN:HD22	1.26	0.83
1:B:170:PRO:HB2	1:B:171:PRO:HD3	1.60	0.82
1:A:341:GLU:OE1	1:A:342:THR:HG22	1.78	0.82
1:B:230:VAL:HG23	1:B:248:LYS:HD3	1.62	0.82
7:A:1296:RET:H8	7:A:1296:RET:H181	1.60	0.82
1:B:75:ILE:HG13	1:B:131:LEU:CD1	2.10	0.82
1:B:91:PHE:HA	1:B:94:THR:CG2	2.09	0.81
1:A:329:GLY:HA2	1:B:96:TYR:CE1	2.18	0.79
1:B:67:LYS:HD3	1:B:337:VAL:HG13	1.65	0.79
1:A:308:MET:HE3	1:B:99:LEU:HD21	1.62	0.79
1:A:96:TYR:HE2	1:A:104:VAL:HG21	1.49	0.78
1:A:345:VAL:HG12	1:A:347:PRO:CD	2.13	0.78
1:B:183:MET:HE3	1:B:289:THR:HG21	1.64	0.77
7:B:1296:RET:H181	7:B:1296:RET:H8	1.66	0.77
1:B:332:GLU:HB3	1:B:335:THR:O	1.85	0.77
1:A:67:LYS:HB2	1:A:337:VAL:HB	1.65	0.77
1:A:341:GLU:CG	1:A:342:THR:H	1.97	0.76
1:A:341:GLU:HG3	1:A:342:THR:N	2.01	0.76
1:B:139:VAL:HG11	1:B:230:VAL:HG12	1.66	0.75
1:A:308:MET:CE	1:B:99:LEU:HD21	2.16	0.75
1:A:59:LEU:HD12	1:A:77:LEU:HD11	1.68	0.75
1:A:304:VAL:O	1:A:308:MET:HG2	1.86	0.75
1:B:337:VAL:HG23	1:B:338:SER:N	2.02	0.75
1:B:214:ILE:HB	1:B:215:PRO:HD3	1.67	0.74
1:A:329:GLY:HA2	1:B:96:TYR:HE1	1.52	0.74
1:A:341:GLU:HG3	1:A:342:THR:H	1.53	0.73
1:B:67:LYS:NZ	1:B:312:GLN:HG3	2.03	0.73
1:B:75:ILE:HG13	1:B:131:LEU:HD11	1.72	0.72
1:A:239:GLU:HB3	1:A:244:GLN:CD	2.09	0.71
1:A:170:PRO:HB2	1:A:171:PRO:HD3	1.72	0.71
1:B:139:VAL:CG1	1:B:230:VAL:HG12	2.20	0.71
1:B:208:PHE:O	1:B:213:ILE:HG13	1.90	0.71
1:A:245:LYS:NZ	1:A:245:LYS:HA	2.06	0.71
1:B:70:THR:H	1:B:73:ASN:HD22	1.39	0.70
1:A:230:VAL:HG23	1:A:248:LYS:HE2	1.73	0.70
1:B:59:LEU:HD13	1:B:77:LEU:HD11	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:LEU:HB3	1:A:266:LEU:HD22	1.73	0.70
1:B:65:HIS:HB2	1:B:68:LEU:HD12	1.74	0.70
1:A:150:GLU:O	1:A:154:ILE:HG13	1.91	0.70
1:A:245:LYS:HA	1:A:245:LYS:HZ2	1.55	0.70
1:B:9:PHE:HA	1:B:179:ILE:HD11	1.72	0.69
1:B:337:VAL:HA	1:B:342:THR:HG23	1.73	0.69
1:B:234:ALA:HB1	1:B:245:LYS:HE2	1.74	0.69
1:B:67:LYS:HB3	1:B:337:VAL:CG2	2.23	0.69
1:A:332:GLU:C	1:A:334:SER:H	1.95	0.69
1:B:64:GLN:HG3	1:B:339:LYS:HE2	1.74	0.69
1:A:135:ARG:HD3	1:A:251:THR:HA	1.75	0.68
1:A:65:HIS:ND1	1:A:338:SER:HA	2.07	0.68
1:A:183:MET:HE3	1:A:289:THR:HG21	1.76	0.68
1:A:239:GLU:HG2	1:A:248:LYS:NZ	2.08	0.68
1:A:325:LYS:HE3	1:A:341:GLU:OE1	1.93	0.68
1:B:239:GLU:HB3	1:B:244:GLN:NE2	2.07	0.68
1:B:231:LYS:NZ	1:B:233:ALA:HB3	2.09	0.68
1:A:65:HIS:CG	1:A:338:SER:HA	2.27	0.68
1:B:96:TYR:HE2	1:B:104:VAL:HG21	1.59	0.68
1:B:69:ARG:N	1:B:73:ASN:HD22	1.92	0.68
1:B:132:ALA:O	1:B:222:CYS:SG	2.52	0.67
1:B:237:GLN:O	1:B:238:GLN:HB3	1.94	0.67
1:A:75:ILE:HG13	1:A:131:LEU:CD1	2.23	0.67
1:A:239:GLU:HB3	1:A:244:GLN:NE2	2.10	0.67
1:A:314:ARG:O	1:A:318:VAL:HG23	1.93	0.67
1:B:70:THR:H	1:B:73:ASN:ND2	1.93	0.67
1:A:76:LEU:HD22	1:A:306:TYR:CG	2.29	0.67
1:A:311:LYS:HD2	1:A:314:ARG:HH21	1.57	0.67
1:B:255:ILE:O	1:B:259:ILE:HG12	1.94	0.67
1:A:137:VAL:O	1:A:141:LYS:HA	1.94	0.67
1:B:267:PRO:HG2	11:B:2019:HOH:O	1.95	0.67
1:A:161:TRP:O	1:A:165:LEU:HD23	1.95	0.66
1:B:238:GLN:HG3	1:B:241:ALA:N	2.09	0.66
1:A:96:TYR:HE2	1:A:104:VAL:CG2	2.09	0.66
1:A:126:TRP:CH2	1:A:215:PRO:HG3	2.31	0.65
1:B:38:SER:HB3	9:B:1508:HTG:H4'1	1.78	0.65
1:B:183:MET:CE	1:B:289:THR:HG21	2.27	0.65
1:A:198:THR:HG23	1:A:200:ASN:OD1	1.96	0.65
1:A:235:ALA:HB3	1:A:239:GLU:OE1	1.97	0.65
1:A:59:LEU:CD1	1:A:77:LEU:HD11	2.26	0.64
1:B:136:TYR:HA	1:B:226:LEU:HD11	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:GLN:HG2	1:B:65:HIS:CD2	2.32	0.64
1:B:312:GLN:HA	1:B:332:GLU:CG	2.17	0.64
1:A:302:ASN:HB2	1:A:303:PRO:HD3	1.80	0.64
1:B:338:SER:HB2	1:B:341:GLU:CG	2.27	0.64
1:A:247:GLU:HA	1:A:247:GLU:OE1	1.96	0.64
1:B:67:LYS:HZ2	1:B:312:GLN:HG3	1.60	0.64
1:B:245:LYS:HE3	1:B:245:LYS:CA	2.21	0.64
1:B:341:GLU:HG3	1:B:342:THR:H	1.62	0.64
1:A:325:LYS:HE2	1:A:327:PRO:HD3	1.80	0.64
1:B:59:LEU:CD1	1:B:77:LEU:HD11	2.27	0.64
1:B:216:LEU:O	1:B:220:PHE:HD2	1.79	0.64
1:A:337:VAL:O	1:A:338:SER:OG	2.16	0.63
1:A:239:GLU:HG2	1:A:248:LYS:HZ2	1.62	0.63
1:B:271:VAL:HG21	1:B:291:PRO:HG3	1.81	0.63
1:A:64:GLN:HE22	1:A:320:THR:HG23	1.64	0.62
1:A:341:GLU:CG	1:A:342:THR:N	2.61	0.62
1:B:230:VAL:HG23	1:B:248:LYS:CD	2.28	0.62
1:B:337:VAL:CA	1:B:342:THR:HG23	2.30	0.62
1:B:239:GLU:CB	1:B:244:GLN:HE22	2.10	0.62
1:B:298:SER:HA	1:B:301:TYR:CD2	2.34	0.62
1:B:337:VAL:HA	1:B:342:THR:O	1.98	0.62
1:A:129:VAL:HG13	1:A:218:VAL:HG11	1.82	0.62
1:A:148:PHE:HA	1:A:152:HIS:ND1	2.15	0.62
1:A:239:GLU:HB3	1:A:244:GLN:OE1	2.00	0.61
1:B:253:MET:HA	1:B:309:MET:SD	2.40	0.61
1:A:189:ILE:HG22	1:A:190:ASP:N	2.15	0.61
1:A:251:THR:O	1:A:255:ILE:HG13	2.00	0.61
1:B:67:LYS:HB3	1:B:337:VAL:HG21	1.82	0.61
1:B:234:ALA:HA	1:B:245:LYS:NZ	2.16	0.61
1:A:298:SER:HA	1:A:301:TYR:CE2	2.36	0.61
1:B:91:PHE:HA	1:B:94:THR:HG22	1.82	0.61
1:B:248:LYS:O	1:B:252:ARG:HG3	2.00	0.61
1:A:312:GLN:HB2	1:A:332:GLU:HG3	1.83	0.60
1:B:298:SER:HA	1:B:301:TYR:CE2	2.34	0.60
1:A:195:HIS:HD1	1:A:198:THR:CG2	2.15	0.60
1:A:237:GLN:O	1:A:238:GLN:HB3	2.01	0.60
1:A:67:LYS:H	1:A:337:VAL:CG2	2.15	0.60
1:A:311:LYS:HZ1	1:A:330:ASP:HB2	1.67	0.60
1:B:67:LYS:HB3	1:B:337:VAL:CG1	2.31	0.60
1:A:65:HIS:HB3	1:A:337:VAL:CG2	2.22	0.59
1:A:212:PHE:O	1:A:216:LEU:HD23	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:LYS:HZ3	1:B:335:THR:C	2.06	0.59
1:A:326:ASN:OD1	1:A:328:LEU:HD13	2.03	0.59
1:B:65:HIS:CG	1:B:338:SER:HA	2.38	0.59
1:B:338:SER:HB2	1:B:341:GLU:HG3	1.84	0.59
1:A:237:GLN:CD	1:A:237:GLN:H	2.06	0.58
1:A:311:LYS:HD2	1:A:314:ARG:HH22	1.68	0.58
1:A:211:HIS:O	1:A:212:PHE:HB2	2.02	0.58
1:B:239:GLU:HB2	1:B:245:LYS:CD	2.23	0.58
1:A:227:VAL:HG13	1:A:228:PHE:N	2.18	0.58
1:B:238:GLN:HA	1:B:242:THR:HA	1.85	0.58
1:B:143:MET:HE2	1:B:143:MET:HA	1.86	0.57
1:B:64:GLN:HG2	1:B:65:HIS:HD2	1.68	0.57
1:B:245:LYS:HZ2	1:B:248:LYS:HD2	1.64	0.57
1:B:194:PRO:O	1:B:195:HIS:C	2.43	0.57
1:A:198:THR:CG2	1:A:200:ASN:OD1	2.53	0.57
1:B:209:VAL:O	1:B:214:ILE:HG13	2.05	0.57
1:A:267:PRO:HG2	11:A:964:HOH:O	2.03	0.57
1:B:9:PHE:CA	1:B:179:ILE:HD11	2.33	0.57
1:A:90:GLY:O	1:A:94:THR:HG22	2.03	0.57
1:B:235:ALA:HB3	1:B:239:GLU:OE1	2.05	0.57
1:A:311:LYS:NZ	1:A:330:ASP:HB2	2.20	0.57
2:C:1:NAG:H61	2:C:2:NAG:C7	2.35	0.57
1:A:75:ILE:HG13	1:A:131:LEU:HD11	1.85	0.56
1:A:342:THR:O	1:A:342:THR:HG23	2.03	0.56
1:B:119:LEU:HD21	1:B:165:LEU:HD22	1.85	0.56
1:B:96:TYR:HE2	1:B:104:VAL:CG2	2.18	0.56
1:A:139:VAL:HG11	1:A:226:LEU:HG	1.86	0.56
1:B:69:ARG:HG2	1:B:69:ARG:HH11	1.70	0.56
1:A:12:PRO:HD3	11:A:2064:HOH:O	2.04	0.56
8:B:1322:PLM:H52	8:B:1323:PLM:H61	1.86	0.56
1:A:254:VAL:O	1:A:258:VAL:HG23	2.06	0.56
1:B:22:SER:C	1:B:24:PHE:H	2.09	0.56
1:A:321:LEU:HD13	8:A:1322:PLM:H72	1.88	0.56
1:B:246:ALA:O	1:B:249:GLU:HG2	2.06	0.56
1:A:183:MET:CE	1:A:289:THR:HG21	2.35	0.56
1:B:126:TRP:CE2	1:B:163:MET:HB3	2.40	0.56
1:B:224:GLY:O	1:B:227:VAL:HG12	2.06	0.56
1:B:131:LEU:O	1:B:135:ARG:HB2	2.06	0.55
1:A:91:PHE:HA	1:A:94:THR:HG22	1.88	0.55
1:A:167:CYS:SG	1:A:207:MET:HG3	2.46	0.55
1:B:256:ILE:HG22	1:B:305:ILE:HD13	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:VAL:HA	1:B:342:THR:C	2.27	0.55
1:A:65:HIS:HB2	1:A:68:LEU:HD12	1.89	0.55
1:B:156:GLY:O	1:B:160:THR:HG23	2.07	0.55
1:B:345:VAL:C	1:B:347:PRO:HD3	2.26	0.55
1:A:322:CYS:O	1:A:324:GLY:N	2.39	0.55
1:A:244:GLN:OE1	1:A:244:GLN:C	2.45	0.55
1:B:135:ARG:HD3	1:B:251:THR:HA	1.89	0.54
1:B:332:GLU:O	1:B:333:ALA:HB3	2.07	0.54
1:A:135:ARG:HA	1:A:135:ARG:NE	2.22	0.54
8:A:1410:PLM:H81	1:B:46:LEU:HD11	1.88	0.54
1:B:278:HIS:NE2	11:B:2038:HOH:O	2.34	0.54
1:A:91:PHE:HA	1:A:94:THR:HG23	1.87	0.54
1:B:167:CYS:SG	1:B:211:HIS:CD2	3.01	0.54
1:B:330:ASP:O	1:B:331:ASP:O	2.26	0.54
1:A:321:LEU:C	1:A:323:CYS:H	2.10	0.54
1:B:143:MET:HG2	1:B:146:PHE:CB	2.38	0.53
1:B:226:LEU:N	1:B:226:LEU:HD12	2.23	0.53
1:B:311:LYS:HE3	9:B:1509:HTG:O3	2.09	0.53
1:B:161:TRP:O	1:B:165:LEU:HD23	2.09	0.53
1:B:216:LEU:O	1:B:220:PHE:CD2	2.60	0.53
1:A:65:HIS:ND1	1:A:337:VAL:O	2.36	0.53
1:A:82:ALA:O	1:A:86:MET:HG3	2.08	0.53
1:B:46:LEU:HD12	8:B:1407:PLM:HD2	1.89	0.53
1:B:167:CYS:O	1:B:170:PRO:HD2	2.09	0.53
1:B:51:GLY:HA2	1:B:300:VAL:HG12	1.90	0.53
1:B:68:LEU:HA	1:B:73:ASN:ND2	2.24	0.53
1:B:212:PHE:O	1:B:216:LEU:HD23	2.09	0.53
1:A:239:GLU:CG	1:A:244:GLN:HE22	2.21	0.53
1:B:332:GLU:CB	1:B:335:THR:O	2.54	0.53
1:A:67:LYS:HB2	1:A:337:VAL:CB	2.37	0.52
1:B:266:LEU:N	1:B:267:PRO:HD2	2.24	0.52
1:B:332:GLU:O	1:B:332:GLU:OE1	2.26	0.52
1:A:74:TYR:CE2	1:A:150:GLU:HG2	2.44	0.52
1:A:327:PRO:O	1:A:328:LEU:C	2.47	0.52
1:A:231:LYS:HZ1	1:A:233:ALA:HB3	1.75	0.52
1:B:150:GLU:O	1:B:154:ILE:CD1	2.58	0.52
1:A:326:ASN:HD21	1:A:328:LEU:CD1	2.17	0.52
1:B:143:MET:HA	1:B:143:MET:CE	2.40	0.51
1:B:159:PHE:O	1:B:163:MET:HG2	2.09	0.51
1:A:214:ILE:HB	1:A:215:PRO:HD3	1.91	0.51
1:B:139:VAL:HG21	1:B:226:LEU:HD23	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:GLY:O	1:A:113:GLU:HB2	2.10	0.51
1:A:314:ARG:HD2	1:B:99:LEU:HD13	1.91	0.51
1:B:67:LYS:HB3	1:B:337:VAL:HG11	1.91	0.51
1:B:254:VAL:O	1:B:258:VAL:HG23	2.09	0.51
1:A:135:ARG:HD3	1:A:251:THR:CA	2.41	0.51
1:A:224:GLY:O	1:A:227:VAL:HG12	2.11	0.51
1:B:9:PHE:C	1:B:179:ILE:HD11	2.30	0.51
1:A:87:VAL:O	1:A:91:PHE:HB2	2.10	0.51
1:A:136:TYR:HA	1:A:226:LEU:HD11	1.91	0.51
1:A:262:LEU:HB3	1:A:266:LEU:CD2	2.40	0.51
1:B:85:PHE:O	1:B:89:GLY:N	2.40	0.51
1:B:127:SER:O	1:B:131:LEU:HD13	2.11	0.51
1:B:67:LYS:N	1:B:337:VAL:HG21	2.26	0.51
1:B:139:VAL:CG1	1:B:226:LEU:HG	2.41	0.51
1:B:167:CYS:SG	1:B:207:MET:HG3	2.51	0.51
1:A:88:PHE:HB2	8:B:1323:PLM:HF2	1.93	0.50
1:A:346:ALA:N	1:A:347:PRO:HD3	2.26	0.50
1:B:67:LYS:NZ	1:B:332:GLU:HG3	2.26	0.50
1:B:129:VAL:HG13	1:B:218:VAL:HG11	1.94	0.50
1:A:94:THR:CB	1:A:113:GLU:OE2	2.54	0.50
1:A:176:SER:HA	1:A:200:ASN:OD1	2.12	0.50
1:B:67:LYS:HG2	1:B:337:VAL:HG22	1.92	0.50
1:B:298:SER:HB2	11:B:2036:HOH:O	2.12	0.50
1:A:24:PHE:CD1	1:A:106:GLY:HA2	2.46	0.50
1:A:325:LYS:HG2	1:A:327:PRO:HD3	1.93	0.50
1:B:321:LEU:C	1:B:323:CYS:H	2.15	0.50
1:B:302:ASN:HB2	11:B:2016:HOH:O	2.11	0.50
1:B:42:ALA:HB2	9:B:1508:HTG:H7'3	1.93	0.50
1:A:259:ILE:O	1:A:263:ILE:HG13	2.12	0.50
1:A:308:MET:CE	1:B:41:ALA:HB1	2.42	0.50
1:B:51:GLY:CA	1:B:300:VAL:HG12	2.42	0.50
1:A:83:ASP:O	1:A:87:VAL:HG23	2.12	0.49
1:A:240:SER:O	1:A:241:ALA:HB2	2.12	0.49
1:A:326:ASN:ND2	1:A:328:LEU:HD11	2.18	0.49
1:B:158:ALA:O	1:B:162:VAL:HG23	2.12	0.49
1:A:67:LYS:H	1:A:337:VAL:HG23	1.75	0.49
1:A:234:ALA:HB2	1:A:248:LYS:HG2	1.93	0.49
1:B:75:ILE:HG21	1:B:131:LEU:HD11	1.95	0.49
1:B:143:MET:HG2	1:B:146:PHE:HB2	1.94	0.49
1:A:139:VAL:HG21	1:A:230:VAL:HG11	1.94	0.49
1:B:195:HIS:HD1	1:B:198:THR:CG2	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:PHE:HA	1:B:179:ILE:CD1	2.41	0.49
1:A:231:LYS:NZ	1:A:233:ALA:HB3	2.28	0.49
1:A:42:ALA:O	1:A:45:PHE:HB3	2.13	0.49
1:A:134:GLU:HG2	1:A:148:PHE:CD2	2.48	0.49
1:B:346:ALA:N	1:B:347:PRO:CD	2.66	0.49
1:A:96:TYR:CE2	1:A:104:VAL:HG21	2.40	0.49
1:A:338:SER:HB2	1:A:341:GLU:HG2	1.94	0.49
1:A:6:GLY:HA3	1:A:9:PHE:CZ	2.48	0.48
1:A:189:ILE:CG2	1:A:190:ASP:N	2.76	0.48
1:B:107:PRO:O	1:B:111:ASN:ND2	2.46	0.48
1:A:70:THR:O	1:A:74:TYR:HD2	1.96	0.48
1:A:239:GLU:OE2	1:A:248:LYS:HD2	2.14	0.48
1:A:161:TRP:O	1:A:165:LEU:HB2	2.13	0.48
1:A:311:LYS:HB3	1:A:332:GLU:HG2	1.94	0.48
1:A:318:VAL:O	1:A:322:CYS:SG	2.71	0.48
1:A:195:HIS:HD1	1:A:198:THR:HG22	1.79	0.48
1:A:72:LEU:HD21	1:A:253:MET:CE	2.44	0.48
1:B:240:SER:O	1:B:241:ALA:HB2	2.14	0.48
1:A:227:VAL:CG1	1:A:228:PHE:N	2.76	0.48
1:B:134:GLU:O	1:B:138:VAL:HG13	2.14	0.48
1:B:322:CYS:O	1:B:324:GLY:N	2.46	0.48
1:B:337:VAL:CG2	1:B:338:SER:N	2.70	0.48
1:A:326:ASN:ND2	1:A:328:LEU:CD1	2.77	0.47
1:B:302:ASN:HB2	1:B:303:PRO:HD3	1.96	0.47
1:B:319:THR:HG21	1:B:341:GLU:OE1	2.14	0.47
1:A:152:HIS:O	1:A:155:MET:HB2	2.14	0.47
1:B:221:PHE:O	1:B:224:GLY:N	2.42	0.47
1:A:85:PHE:O	1:A:89:GLY:N	2.47	0.47
1:A:337:VAL:CG2	1:A:338:SER:N	2.76	0.47
1:B:238:GLN:HG3	1:B:241:ALA:CA	2.44	0.47
1:A:329:GLY:HA2	1:B:96:TYR:CZ	2.49	0.47
1:A:342:THR:OG1	1:A:344:GLN:CD	2.53	0.47
1:B:196:GLU:C	1:B:198:THR:H	2.18	0.47
1:A:66:LYS:HD2	1:A:66:LYS:N	2.29	0.47
1:A:132:ALA:O	1:A:222:CYS:SG	2.73	0.47
1:A:151:ASN:ND2	1:A:152:HIS:CD2	2.82	0.47
1:A:70:THR:H	1:A:73:ASN:ND2	2.04	0.47
1:A:298:SER:HA	1:A:301:TYR:CD2	2.50	0.47
1:B:170:PRO:CB	1:B:171:PRO:HD3	2.38	0.47
1:B:236:GLN:O	1:B:238:GLN:N	2.48	0.47
1:B:290:ILE:HB	1:B:291:PRO:CD	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:LYS:HD3	1:B:343:SER:OG	2.15	0.47
1:B:227:VAL:HG13	1:B:228:PHE:N	2.29	0.47
1:A:194:PRO:O	1:A:195:HIS:C	2.54	0.47
1:A:302:ASN:HB2	11:A:2017:HOH:O	2.15	0.47
1:A:322:CYS:HA	8:A:1322:PLM:O1	2.14	0.47
1:A:326:ASN:ND2	1:A:328:LEU:HD21	2.30	0.47
1:B:204:VAL:HG11	1:B:276:PHE:CD1	2.50	0.47
1:B:237:GLN:O	1:B:238:GLN:CB	2.62	0.47
1:B:115:PHE:CD1	1:B:172:LEU:HD11	2.50	0.47
1:A:178:TYR:HA	1:A:188:GLY:O	2.15	0.46
1:B:140:CYS:O	1:B:141:LYS:C	2.53	0.46
1:B:143:MET:CB	1:B:146:PHE:HB3	2.46	0.46
1:B:240:SER:O	1:B:241:ALA:CB	2.63	0.46
1:A:302:ASN:CB	1:A:303:PRO:HD3	2.45	0.46
1:B:238:GLN:CG	1:B:241:ALA:H	2.21	0.46
1:B:256:ILE:HD13	1:B:256:ILE:HA	1.72	0.46
1:B:76:LEU:HD22	1:B:306:TYR:CG	2.51	0.46
1:B:137:VAL:HG12	1:B:137:VAL:O	2.15	0.46
1:A:131:LEU:O	1:A:135:ARG:HB2	2.15	0.46
1:A:137:VAL:HA	1:A:142:PRO:CD	2.46	0.46
1:A:180:PRO:HA	1:A:186:SER:O	2.16	0.46
1:B:239:GLU:OE1	1:B:239:GLU:HA	2.16	0.46
1:A:143:MET:HB2	1:A:146:PHE:HB3	1.98	0.46
1:B:170:PRO:HB2	1:B:171:PRO:CD	2.39	0.46
1:A:300:VAL:O	1:A:303:PRO:HD2	2.16	0.45
1:A:75:ILE:HG21	1:A:131:LEU:HD11	1.97	0.45
1:A:137:VAL:HA	1:A:142:PRO:HD3	1.98	0.45
1:B:150:GLU:O	1:B:154:ILE:HD12	2.16	0.45
1:B:227:VAL:CG1	1:B:228:PHE:N	2.79	0.45
1:A:230:VAL:O	1:A:230:VAL:HG13	2.16	0.45
1:B:70:THR:O	1:B:74:TYR:HD2	2.00	0.45
1:B:75:ILE:O	1:B:79:LEU:HD13	2.17	0.45
1:B:315:ASN:OD1	1:B:331:ASP:O	2.34	0.45
1:B:340:THR:HG23	1:B:341:GLU:N	2.31	0.45
1:A:342:THR:HA	1:A:344:GLN:OE1	2.16	0.45
1:B:64:GLN:HE21	1:B:65:HIS:CD2	2.33	0.45
1:B:134:GLU:OE1	1:B:135:ARG:NH2	2.50	0.45
1:B:150:GLU:O	1:B:153:ALA:HB3	2.17	0.45
1:A:77:LEU:O	1:A:81:VAL:HG23	2.17	0.45
1:A:266:LEU:N	1:A:267:PRO:HD2	2.31	0.45
1:B:148:PHE:CE1	1:B:152:HIS:HB3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:PHE:O	1:B:213:ILE:CG1	2.62	0.45
1:A:67:LYS:H	1:A:337:VAL:HG21	1.82	0.45
1:B:234:ALA:HA	1:B:245:LYS:HZ3	1.82	0.45
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.78	0.45
1:A:325:LYS:HE2	1:A:327:PRO:CD	2.45	0.45
1:B:75:ILE:O	1:B:78:ASN:HB3	2.17	0.45
1:B:216:LEU:HD12	1:B:220:PHE:HE2	1.82	0.45
1:A:158:ALA:O	1:A:162:VAL:HG23	2.17	0.44
1:B:78:ASN:O	1:B:81:VAL:N	2.49	0.44
1:B:221:PHE:CD1	1:B:221:PHE:C	2.90	0.44
1:B:300:VAL:O	1:B:303:PRO:HD2	2.17	0.44
1:A:245:LYS:NZ	1:A:245:LYS:CA	2.77	0.44
1:B:148:PHE:CD1	1:B:152:HIS:HB2	2.53	0.44
1:B:205:ILE:O	1:B:209:VAL:HG23	2.17	0.44
1:A:143:MET:HG2	1:A:146:PHE:HB2	1.98	0.44
1:B:332:GLU:OE2	1:B:334:SER:HB2	2.17	0.44
1:A:67:LYS:HB2	1:A:337:VAL:CG2	2.48	0.44
1:A:213:ILE:O	1:A:217:ILE:HG13	2.17	0.44
1:B:60:TYR:C	1:B:60:TYR:CD1	2.91	0.44
1:A:191:TYR:OH	7:A:1296:RET:H192	2.18	0.44
1:B:22:SER:C	1:B:24:PHE:N	2.70	0.44
1:B:126:TRP:CH2	1:B:215:PRO:HG3	2.53	0.44
1:B:231:LYS:HB2	1:B:231:LYS:HE3	1.81	0.44
1:A:337:VAL:C	1:A:338:SER:OG	2.56	0.44
1:A:39:MET:CE	1:A:39:MET:HA	2.48	0.43
1:A:329:GLY:HA2	1:B:96:TYR:OH	2.18	0.43
1:A:74:TYR:OH	1:A:150:GLU:HG2	2.18	0.43
1:A:319:THR:OG1	1:A:327:PRO:HB3	2.18	0.43
1:B:102:TYR:CE2	1:B:104:VAL:HA	2.54	0.43
1:B:305:ILE:O	1:B:309:MET:N	2.48	0.43
1:B:139:VAL:HG11	1:B:226:LEU:HG	2.00	0.43
1:A:138:VAL:HG23	1:A:139:VAL:N	2.32	0.43
1:A:231:LYS:NZ	1:A:231:LYS:HB2	2.32	0.43
1:A:102:TYR:CE2	1:A:104:VAL:HA	2.53	0.43
1:B:244:GLN:O	1:B:248:LYS:HG3	2.18	0.43
1:B:301:TYR:CD1	1:B:301:TYR:C	2.91	0.43
1:B:315:ASN:O	1:B:318:VAL:HG12	2.18	0.43
1:B:18:GLY:HA2	4:E:1:NAG:O5	2.19	0.43
1:B:117:ALA:CB	7:B:1296:RET:H14	2.49	0.43
1:B:231:LYS:CE	1:B:233:ALA:HB3	2.48	0.43
1:B:234:ALA:HA	1:B:245:LYS:HZ1	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:216:LEU:HD12	1:B:220:PHE:CE2	2.54	0.43
1:B:311:LYS:HD2	9:B:1509:HTG:O2	2.19	0.43
1:A:117:ALA:CB	7:A:1296:RET:H14	2.49	0.42
1:A:52:PHE:HB3	1:A:53:PRO:CD	2.50	0.42
1:B:332:GLU:O	1:B:333:ALA:CB	2.67	0.42
1:A:133:ILE:O	1:A:136:TYR:N	2.52	0.42
1:A:332:GLU:C	1:A:334:SER:N	2.65	0.42
1:B:103:PHE:CZ	1:B:187:CYS:SG	3.12	0.42
1:B:226:LEU:N	1:B:226:LEU:CD1	2.83	0.42
1:B:312:GLN:O	1:B:316:CYS:SG	2.78	0.42
1:B:69:ARG:HG2	1:B:69:ARG:NH1	2.32	0.42
1:A:129:VAL:O	1:A:133:ILE:HG12	2.18	0.42
1:B:67:LYS:H	1:B:337:VAL:CG2	2.31	0.42
1:B:210:VAL:O	1:B:215:PRO:HD3	2.19	0.42
10:B:1401:HTO:H72	9:B:1509:HTG:H3'2	2.01	0.42
1:A:6:GLY:HA3	1:A:9:PHE:CE1	2.54	0.42
1:A:135:ARG:O	1:A:138:VAL:HG22	2.19	0.42
1:A:159:PHE:O	1:A:163:MET:HG2	2.20	0.42
1:B:196:GLU:C	1:B:198:THR:N	2.73	0.42
1:B:238:GLN:HG2	1:B:239:GLU:N	2.34	0.42
3:D:1:NAG:H61	3:D:2:NAG:H82	2.01	0.42
1:A:155:MET:HA	1:A:155:MET:CE	2.50	0.42
1:A:104:VAL:CG2	1:A:105:PHE:N	2.83	0.42
1:B:141:LYS:NZ	1:B:144:SER:O	2.38	0.42
1:B:326:ASN:C	1:B:328:LEU:HD22	2.39	0.42
1:A:22:SER:C	1:A:24:PHE:H	2.23	0.42
1:A:239:GLU:CB	1:A:244:GLN:OE1	2.68	0.42
1:A:321:LEU:HD13	8:A:1322:PLM:H51	2.02	0.41
1:B:18:GLY:HA2	4:E:1:NAG:C1	2.50	0.41
1:B:151:ASN:O	1:B:155:MET:HG2	2.20	0.41
1:B:224:GLY:O	1:B:227:VAL:CG1	2.67	0.41
1:B:150:GLU:O	1:B:154:ILE:HD13	2.20	0.41
1:B:243:THR:HG23	1:B:244:GLN:N	2.34	0.41
1:B:99:LEU:HD23	1:B:99:LEU:HA	1.92	0.41
1:B:330:ASP:C	1:B:331:ASP:O	2.58	0.41
1:A:20:VAL:HA	1:A:30:TYR:CZ	2.56	0.41
1:A:119:LEU:HD23	1:A:119:LEU:HA	1.89	0.41
1:A:216:LEU:HD12	1:A:220:PHE:CE2	2.56	0.41
1:A:256:ILE:HG22	1:A:305:ILE:HD13	2.01	0.41
1:B:9:PHE:HA	1:B:179:ILE:CG1	2.50	0.41
1:B:67:LYS:HE2	1:B:336:THR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:326:ASN:HB3	1:B:328:LEU:CD2	2.51	0.41
1:A:326:ASN:C	1:A:328:LEU:H	2.23	0.41
1:B:243:THR:CG2	1:B:244:GLN:N	2.83	0.41
1:A:36:GLN:O	1:A:39:MET:HB2	2.20	0.41
1:A:241:ALA:C	1:A:243:THR:N	2.74	0.41
1:A:332:GLU:O	1:A:333:ALA:HB3	2.20	0.41
1:B:182:GLY:HA2	1:B:285:PRO:O	2.21	0.41
1:A:170:PRO:CB	1:A:171:PRO:HD3	2.45	0.41
1:B:87:VAL:O	1:B:91:PHE:HB2	2.20	0.41
1:A:67:LYS:HG2	1:A:343:SER:OG	2.21	0.41
1:A:134:GLU:HG2	1:A:148:PHE:CE2	2.55	0.41
1:A:170:PRO:HB2	1:A:203:PHE:CE1	2.56	0.41
1:B:117:ALA:HB1	7:B:1296:RET:H14	2.03	0.41
1:A:64:GLN:NE2	1:A:320:THR:HG23	2.34	0.41
1:A:72:LEU:HD21	1:A:253:MET:HE1	2.03	0.41
1:B:59:LEU:HA	1:B:77:LEU:CD1	2.51	0.40
1:B:110:CYS:SG	1:B:180:PRO:HD3	2.61	0.40
1:A:231:LYS:NZ	1:A:231:LYS:CB	2.85	0.40
1:A:276:PHE:O	1:A:279:GLN:HG3	2.21	0.40
1:A:338:SER:CB	1:A:341:GLU:HG2	2.51	0.40
1:B:67:LYS:CE	1:B:312:GLN:HG3	2.51	0.40
1:B:207:MET:CE	7:B:1296:RET:H161	2.51	0.40
1:A:85:PHE:O	1:A:89:GLY:HA3	2.21	0.40
1:A:126:TRP:CE2	1:A:163:MET:HB3	2.56	0.40
1:A:128:LEU:HD23	1:A:128:LEU:HA	1.77	0.40
1:A:229:THR:HG22	1:A:230:VAL:H	1.87	0.40
1:B:91:PHE:CZ	1:B:296:LYS:HB3	2.56	0.40
1:B:175:TRP:O	1:B:176:SER:HB3	2.22	0.40
1:B:238:GLN:HG3	1:B:241:ALA:C	2.42	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	A	347/349 (99%)	292 (84%)	39 (11%)	16 (5%)	2 3
1	B	347/349 (99%)	287 (83%)	40 (12%)	20 (6%)	1 1
All	All	694/698 (99%)	579 (83%)	79 (11%)	36 (5%)	2 2

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	HIS
1	A	212	PHE
1	A	240	SER
1	A	241	ALA
1	A	323	CYS
1	A	341	GLU
1	B	237	GLN
1	B	240	SER
1	B	241	ALA
1	B	331	ASP
1	B	337	VAL
1	B	347	PRO
1	A	324	GLY
1	A	328	LEU
1	A	329	GLY
1	A	347	PRO
1	B	195	HIS
1	B	229	THR
1	B	323	CYS
1	B	328	LEU
1	B	341	GLU
1	A	335	THR
1	A	338	SER
1	B	238	GLN
1	A	145	ASN
1	A	322	CYS
1	B	235	ALA
1	B	324	GLY
1	A	229	THR
1	B	66	LYS
1	B	141	LYS
1	B	212	PHE
1	B	333	ALA
1	A	141	LYS
1	B	106	GLY

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Mol	Chain	Res	Type
1	B	197	GLU

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	A	296/296 (100%)	278 (94%)	18 (6%)	18	38
1	B	296/296 (100%)	273 (92%)	23 (8%)	12	25
All	All	592/592 (100%)	551 (93%)	41 (7%)	15	31

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	57	LEU
1	A	76	LEU
1	A	79	LEU
1	A	94	THR
1	A	119	LEU
1	A	134	GLU
1	A	135	ARG
1	A	143	MET
1	A	237	GLN
1	A	239	GLU
1	A	244	GLN
1	A	245	LYS
1	A	252	ARG
1	A	266	LEU
1	A	331	ASP
1	A	332	GLU
1	A	336	THR
1	B	8	ASN
1	B	38	SER
1	B	50	LEU
1	B	94	THR

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Mol	Chain	Res	Type
1	B	104	VAL
1	B	119	LEU
1	B	134	GLU
1	B	135	ARG
1	B	143	MET
1	B	148	PHE
1	B	167	CYS
1	B	222	CYS
1	B	229	THR
1	B	231	LYS
1	B	236	GLN
1	B	237	GLN
1	B	244	GLN
1	B	245	LYS
1	B	249	GLU
1	B	264	CYS
1	B	308	MET
1	B	330	ASP
1	B	332	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	64	GLN
1	A	73	ASN
1	A	100	HIS
1	A	111	ASN
1	A	225	GLN
1	A	237	GLN
1	A	326	ASN
1	B	8	ASN
1	B	64	GLN
1	B	73	ASN
1	B	111	ASN
1	B	151	ASN
1	B	237	GLN
1	B	244	GLN

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)

11 monosaccharides 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	$\#\|Z\| > 2$
2	NAG	C	1	1,2	14,14,15	0.61	0	17,19,21	0.77	0
2	NAG	C	2	2	14,14,15	0.83	0	17,19,21	0.91	1 (5%)
2	MAN	C	3	2	11,11,12	0.94	0	15,15,17	0.72	0
3	NAG	D	1	1,3	14,14,15	0.80	0	17,19,21	0.76	0
3	NAG	D	2	3	14,14,15	0.60	0	17,19,21	0.85	1 (5%)
4	NAG	E	1	1,4	14,14,15	0.55	0	17,19,21	0.69	0
4	NAG	E	2	4	14,14,15	0.43	0	17,19,21	1.14	1 (5%)
4	BMA	E	3	4	11,11,12	0.83	0	15,15,17	0.64	0
4	BMA	E	4	4	11,11,12	0.90	0	15,15,17	0.84	1 (6%)
3	NAG	F	1	1,3	14,14,15	0.61	0	17,19,21	0.83	1 (5%)
3	NAG	F	2	3	14,14,15	0.75	0	17,19,21	0.74	1 (5%)

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
2	NAG	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	C	3	2	-	2/2/19/22	0/1/1/1
3	NAG	D	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
4	NAG	E	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	4/6/23/26	0/1/1/1
4	BMA	E	3	4	-	2/2/19/22	0/1/1/1
4	BMA	E	4	4	-	0/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	E	2	NAG	C2-N2-C7	-3.90	117.67	122.90
3	D	2	NAG	C2-N2-C7	-2.71	119.27	122.90
2	C	2	NAG	C8-C7-N2	2.35	120.02	116.12
3	F	2	NAG	C2-N2-C7	-2.15	120.02	122.90
4	E	4	BMA	C1-O5-C5	2.12	115.03	112.19
3	F	1	NAG	C2-N2-C7	-2.01	120.21	122.90

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
3	D	1	NAG	C8-C7-N2-C2
3	D	1	NAG	O7-C7-N2-C2
3	F	1	NAG	C8-C7-N2-C2
3	F	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
2	C	3	MAN	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
4	E	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
4	E	3	BMA	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
4	E	1	NAG	C8-C7-N2-C2
4	E	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C4-C5-C6-O6
2	C	3	MAN	C4-C5-C6-O6
4	E	2	NAG	C4-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
4	E	2	NAG	C8-C7-N2-C2
4	E	2	NAG	O7-C7-N2-C2

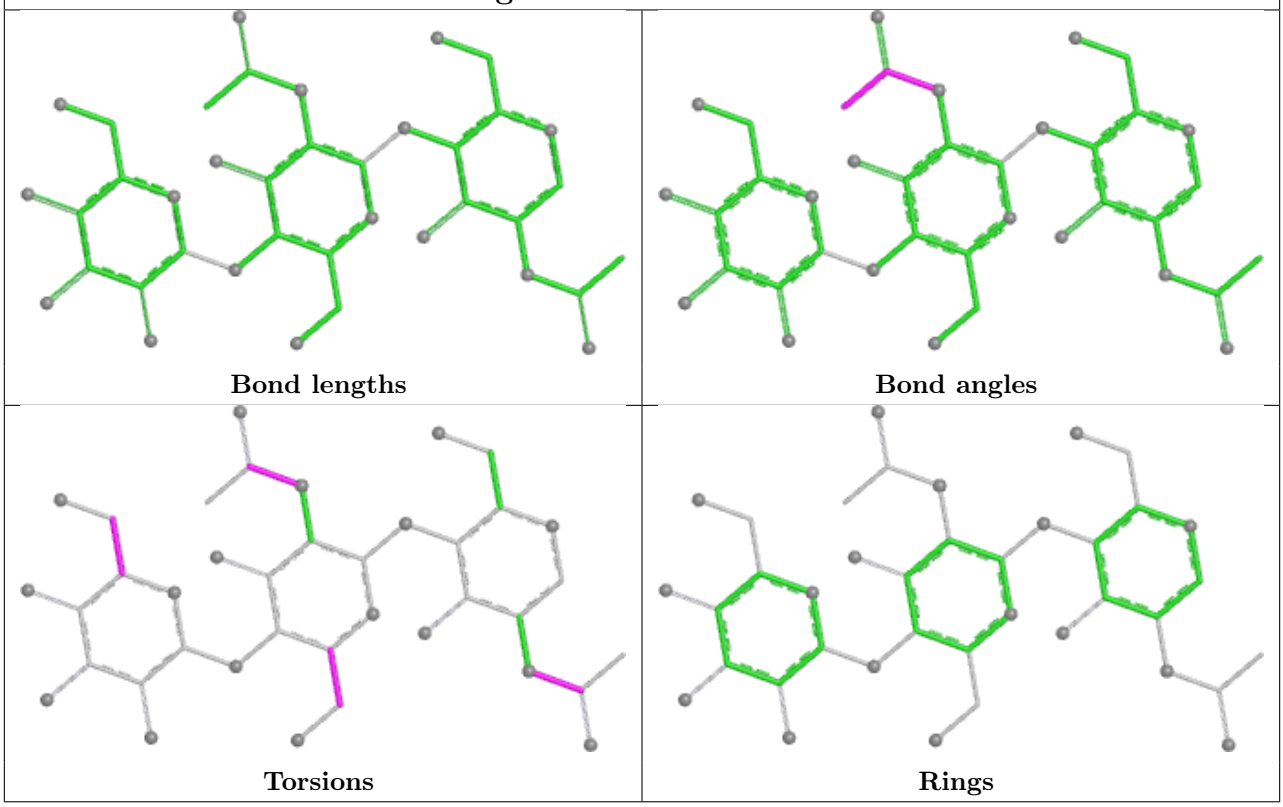
There are no ring outliers.

5 monomers are involved in 4 short contacts:

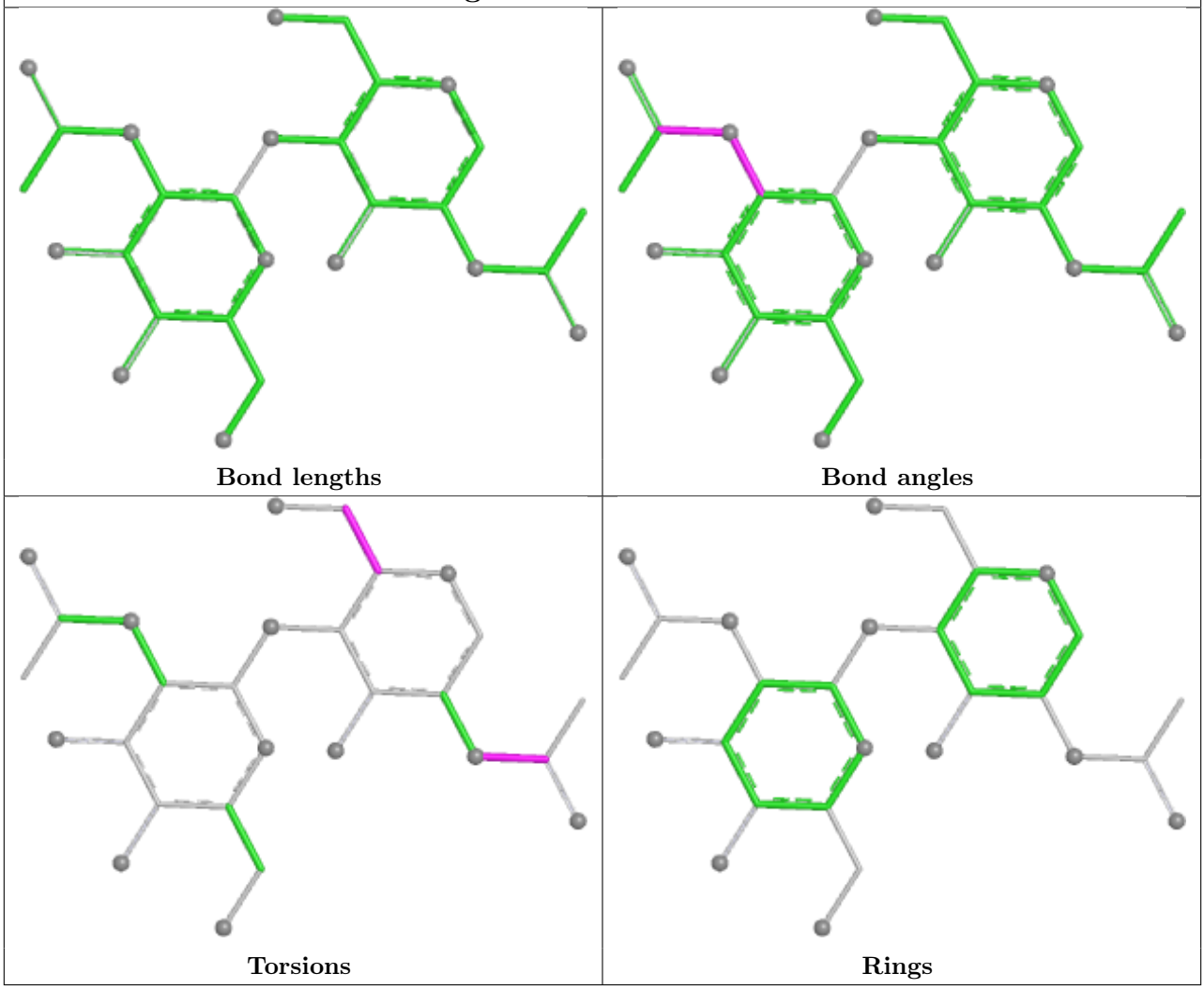
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1	NAG	1	0
2	C	2	NAG	1	0
4	E	1	NAG	2	0
3	D	2	NAG	1	0
2	C	1	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

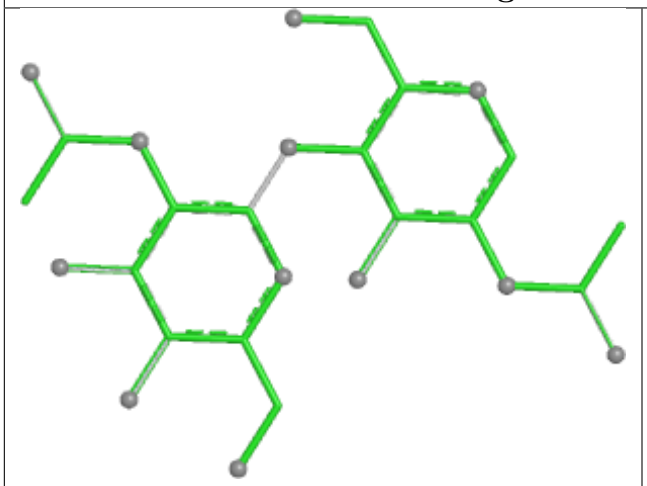
Oligosaccharide Chain C



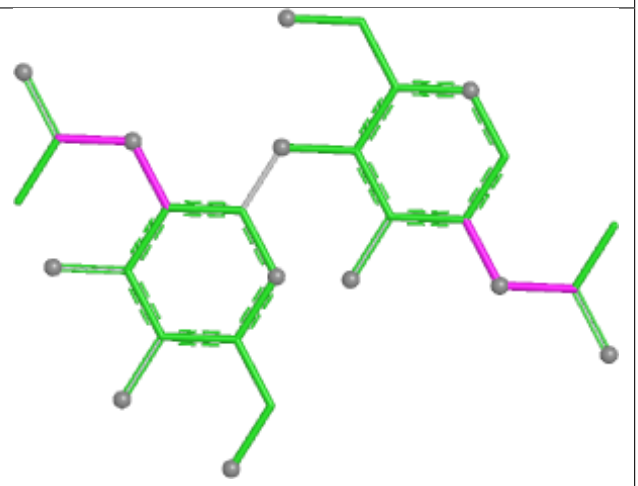
Oligosaccharide Chain D



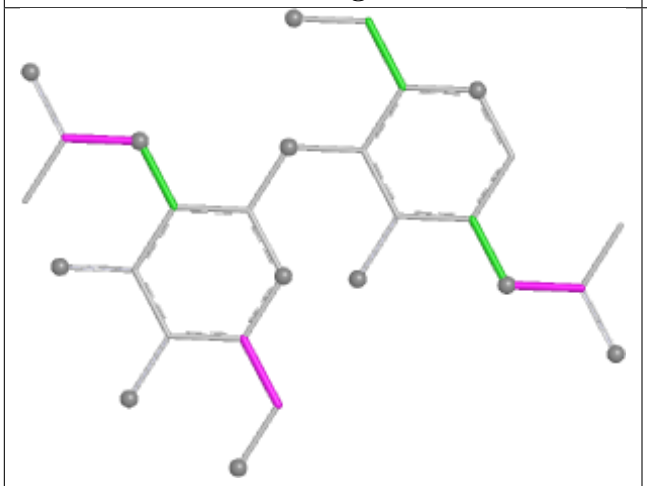
Oligosaccharide Chain F



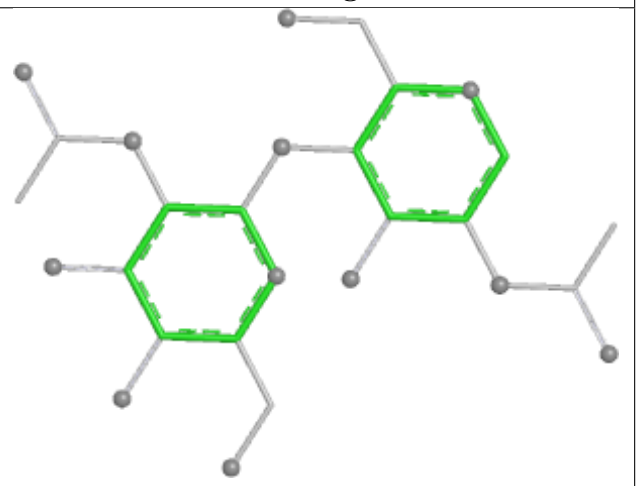
Bond lengths



Bond angles

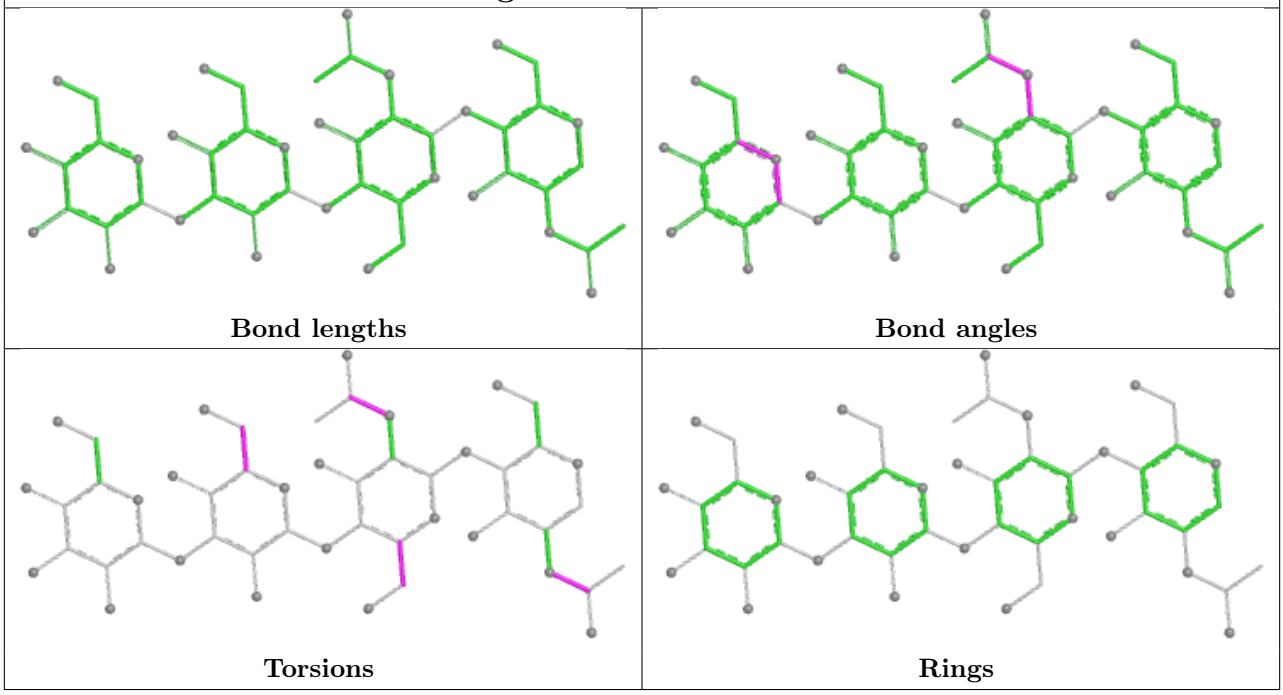


Torsions



Rings

Oligosaccharide Chain E



5.6 Ligand geometry (i)

Of 26 ligands modelled in this entry, 13 are monoatomic - leaving 13 for Mogul analysis.

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	$\#\|Z\| > 2$
9	HTG	B	1506	-	19,19,19	2.85	8 (42%)	23,24,24	1.74	1 (4%)
8	PLM	B	1322	1	16,16,17	1.07	1 (6%)	15,15,17	0.42	0
8	PLM	A	1322	1	16,16,17	0.98	1 (6%)	15,15,17	0.61	0
9	HTG	A	1507	-	19,19,19	2.95	8 (42%)	23,24,24	1.91	1 (4%)
9	HTG	B	1508	-	19,19,19	2.36	8 (42%)	23,24,24	2.34	2 (8%)
8	PLM	A	1323	1	16,16,17	0.99	1 (6%)	15,15,17	0.54	0
8	PLM	B	1323	1	16,16,17	0.98	1 (6%)	15,15,17	0.48	0
8	PLM	B	1407	-	15,15,17	2.00	3 (20%)	14,14,17	3.44	8 (57%)
9	HTG	B	1509	-	19,19,19	3.07	8 (42%)	23,24,24	1.99	2 (8%)
10	HTO	B	1401	-	9,9,9	2.19	3 (33%)	10,10,10	1.15	1 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
7	RET	A	1296	1	20,20,21	1.94	2 (10%)	27,27,28	1.95	8 (29%)
7	RET	B	1296	1	20,20,21	2.50	4 (20%)	27,27,28	1.95	6 (22%)
8	PLM	A	1410	-	15,15,17	2.23	3 (20%)	14,14,17	3.39	7 (50%)

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
9	HTG	B	1506	-	-	2/10/30/30	0/1/1/1
8	PLM	B	1322	1	-	10/14/14/15	-
8	PLM	A	1322	1	-	11/14/14/15	-
9	HTG	A	1507	-	-	4/10/30/30	0/1/1/1
9	HTG	B	1508	-	-	4/10/30/30	0/1/1/1
8	PLM	A	1323	1	-	8/14/14/15	-
8	PLM	B	1323	1	-	9/14/14/15	-
8	PLM	B	1407	-	-	5/13/13/15	-
9	HTG	B	1509	-	-	5/10/30/30	0/1/1/1
10	HTO	B	1401	-	-	2/10/10/10	-
7	RET	A	1296	1	-	10/13/30/31	0/1/1/1
7	RET	B	1296	1	-	10/13/30/31	0/1/1/1
8	PLM	A	1410	-	-	6/13/13/15	-

All (51) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1507	HTG	O5-C1	8.75	1.56	1.42
9	B	1509	HTG	O5-C1	8.51	1.55	1.42
9	B	1506	HTG	O5-C1	8.25	1.55	1.42
7	A	1296	RET	C5-C6	6.50	1.45	1.34
9	B	1508	HTG	O5-C1	6.32	1.52	1.42
7	B	1296	RET	C5-C6	6.29	1.45	1.34
7	B	1296	RET	C14-C13	5.88	1.37	1.33
8	A	1410	PLM	CC-CB	-5.86	1.22	1.51
7	B	1296	RET	C1-C6	5.50	1.60	1.53
10	B	1401	HTO	C3-C2	5.26	1.65	1.53
8	B	1407	PLM	CC-CB	-5.18	1.26	1.51
9	B	1509	HTG	C4-C5	5.14	1.64	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	1509	HTG	C1-C2	4.66	1.61	1.53
9	B	1506	HTG	C4-C5	4.56	1.62	1.53
7	A	1296	RET	C1-C6	4.54	1.59	1.53
8	A	1410	PLM	CB-CA	-4.24	1.30	1.51
9	A	1507	HTG	C1-S1	4.16	1.87	1.80
9	B	1506	HTG	C1-C2	3.92	1.60	1.53
9	A	1507	HTG	C1-C2	3.90	1.60	1.53
9	B	1508	HTG	C4-C5	3.86	1.61	1.53
8	B	1322	PLM	O1-C1	-3.81	1.22	1.42
8	A	1410	PLM	CD-CC	-3.81	1.32	1.51
8	A	1322	PLM	O1-C1	-3.75	1.23	1.42
9	B	1509	HTG	O5-C5	3.70	1.53	1.44
8	A	1323	PLM	O1-C1	-3.68	1.23	1.42
9	B	1509	HTG	C1-S1	3.64	1.86	1.80
8	B	1323	PLM	O1-C1	-3.53	1.24	1.42
8	B	1407	PLM	CB-CA	-3.53	1.34	1.51
9	B	1508	HTG	O5-C5	3.45	1.52	1.44
9	A	1507	HTG	C4-C3	3.43	1.61	1.52
9	A	1507	HTG	O5-C5	3.37	1.52	1.44
9	B	1506	HTG	C1-S1	3.35	1.86	1.80
9	A	1507	HTG	C4-C5	3.34	1.60	1.53
9	B	1509	HTG	C4-C3	3.29	1.60	1.52
9	B	1508	HTG	C1-C2	3.28	1.59	1.53
9	A	1507	HTG	C3-C2	3.24	1.60	1.52
8	B	1407	PLM	CD-CC	-3.18	1.36	1.51
9	B	1509	HTG	C3-C2	3.16	1.60	1.52
9	B	1506	HTG	C4-C3	3.10	1.60	1.52
9	B	1506	HTG	C3-C2	3.10	1.60	1.52
9	B	1506	HTG	O5-C5	2.99	1.51	1.44
9	B	1508	HTG	C3-C2	2.90	1.59	1.52
9	B	1509	HTG	C6-C5	2.75	1.61	1.51
9	B	1506	HTG	C1'-S1	2.74	1.85	1.81
9	A	1507	HTG	C1'-S1	2.66	1.85	1.81
10	B	1401	HTO	C4-C3	2.61	1.57	1.52
7	B	1296	RET	C17-C1	2.59	1.58	1.53
9	B	1508	HTG	C4-C3	2.36	1.58	1.52
9	B	1508	HTG	C1-S1	2.29	1.84	1.80
10	B	1401	HTO	C5-C4	2.14	1.61	1.52
9	B	1508	HTG	C1'-S1	2.00	1.84	1.81

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	B	1508	HTG	C1'-S1-C1	10.29	122.68	100.45
9	B	1509	HTG	C1'-S1-C1	8.46	118.72	100.45
8	B	1407	PLM	CD-CC-CB	8.30	156.34	114.37
8	A	1410	PLM	CD-CC-CB	8.20	155.84	114.37
9	A	1507	HTG	C1'-S1-C1	8.01	117.75	100.45
9	B	1506	HTG	C1'-S1-C1	7.49	116.62	100.45
8	A	1410	PLM	CA-C9-C8	7.43	151.91	114.37
8	B	1407	PLM	CA-C9-C8	7.41	151.83	114.37
7	B	1296	RET	C8-C9-C10	-5.44	110.45	119.01
7	A	1296	RET	C18-C5-C6	5.01	129.95	124.48
7	B	1296	RET	C18-C5-C6	4.53	129.43	124.48
7	A	1296	RET	C8-C9-C10	-4.32	112.22	119.01
7	B	1296	RET	C19-C9-C10	3.68	128.78	122.82
7	A	1296	RET	C19-C9-C10	3.49	128.47	122.82
8	B	1407	PLM	CC-CB-CA	3.41	131.61	114.37
8	A	1410	PLM	CC-CB-CA	3.16	130.36	114.37
7	A	1296	RET	C7-C8-C9	3.02	130.70	126.23
7	A	1296	RET	C2-C1-C6	2.72	114.39	110.44
8	A	1410	PLM	C5-C4-C3	2.61	127.57	114.37
8	B	1407	PLM	CE-CD-CC	2.56	127.31	114.37
7	A	1296	RET	C18-C5-C4	-2.52	108.23	113.60
7	B	1296	RET	C10-C11-C12	-2.47	116.05	123.20
8	B	1407	PLM	C5-C4-C3	2.42	126.60	114.37
7	B	1296	RET	C2-C1-C6	2.40	113.92	110.44
7	A	1296	RET	C10-C11-C12	-2.39	116.28	123.20
7	B	1296	RET	C18-C5-C4	-2.37	108.55	113.60
8	A	1410	PLM	CF-CE-CD	-2.34	94.70	115.25
8	A	1410	PLM	CE-CD-CC	2.29	125.96	114.37
8	B	1407	PLM	C9-C8-C7	2.28	125.87	114.37
10	B	1401	HTO	O3-C3-C4	-2.26	104.01	109.14
7	A	1296	RET	C1-C6-C5	-2.25	119.56	122.64
8	B	1407	PLM	CG-CF-CE	2.21	128.32	113.36
8	A	1410	PLM	CG-CF-CE	2.14	127.81	113.36
8	B	1407	PLM	CF-CE-CD	-2.12	96.64	115.25
9	B	1508	HTG	O5-C5-C6	2.06	111.55	106.44
9	B	1509	HTG	O5-C5-C6	2.04	111.50	106.44

There are no chirality outliers.

All (86) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1296	RET	C5-C6-C7-C8
7	A	1296	RET	C7-C8-C9-C10

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Mol	Chain	Res	Type	Atoms
7	A	1296	RET	C11-C12-C13-C14
7	A	1296	RET	C11-C12-C13-C20
7	B	1296	RET	C7-C8-C9-C10
7	B	1296	RET	C11-C12-C13-C14
7	B	1296	RET	C11-C12-C13-C20
8	B	1407	PLM	CA-CB-CC-CD
9	B	1509	HTG	O5-C5-C6-O6
9	A	1507	HTG	O5-C5-C6-O6
9	B	1509	HTG	C4-C5-C6-O6
7	B	1296	RET	C7-C8-C9-C19
7	A	1296	RET	C6-C7-C8-C9
9	B	1508	HTG	C1'-C2'-C3'-C4'
7	A	1296	RET	C11-C10-C9-C8
7	B	1296	RET	C11-C10-C9-C8
8	A	1323	PLM	C1-C2-C3-C4
9	B	1506	HTG	C1'-C2'-C3'-C4'
8	A	1323	PLM	CB-CC-CD-CE
8	B	1322	PLM	C5-C6-C7-C8
9	A	1507	HTG	C2'-C3'-C4'-C5'
8	A	1410	PLM	CA-CB-CC-CD
8	B	1323	PLM	CA-CB-CC-CD
8	A	1322	PLM	CC-CD-CE-CF
8	B	1323	PLM	C9-CA-CB-CC
8	B	1323	PLM	C8-C9-CA-CB
8	A	1410	PLM	C5-C6-C7-C8
8	A	1323	PLM	C4-C5-C6-C7
8	B	1323	PLM	C3-C4-C5-C6
9	B	1508	HTG	C2'-C3'-C4'-C5'
9	A	1507	HTG	C4-C5-C6-O6
8	B	1322	PLM	CC-CD-CE-CF
7	A	1296	RET	C1-C6-C7-C8
7	B	1296	RET	C1-C6-C7-C8
7	B	1296	RET	C5-C6-C7-C8
8	B	1323	PLM	C6-C7-C8-C9
8	A	1322	PLM	C8-C9-CA-CB
8	A	1323	PLM	C7-C8-C9-CA
7	A	1296	RET	C10-C11-C12-C13
7	B	1296	RET	C10-C11-C12-C13
7	B	1296	RET	C6-C7-C8-C9
8	B	1407	PLM	C5-C6-C7-C8
9	B	1509	HTG	C2'-C3'-C4'-C5'
8	A	1323	PLM	C3-C4-C5-C6

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Mol	Chain	Res	Type	Atoms
8	A	1410	PLM	C7-C8-C9-CA
8	B	1323	PLM	CB-CC-CD-CE
8	A	1410	PLM	C9-CA-CB-CC
8	B	1322	PLM	C9-CA-CB-CC
8	B	1322	PLM	CB-CC-CD-CE
8	A	1322	PLM	C7-C8-C9-CA
9	B	1506	HTG	C2'-C3'-C4'-C5'
8	B	1407	PLM	C3-C4-C5-C6
8	B	1322	PLM	C8-C9-CA-CB
8	A	1322	PLM	C4-C5-C6-C7
8	A	1322	PLM	C3-C4-C5-C6
9	B	1509	HTG	C3'-C4'-C5'-C6'
8	B	1407	PLM	C7-C8-C9-CA
10	B	1401	HTO	C4-C5-C6-C7
9	B	1508	HTG	C4'-C5'-C6'-C7'
8	A	1323	PLM	CA-CB-CC-CD
8	A	1322	PLM	C5-C6-C7-C8
8	B	1323	PLM	C4-C5-C6-C7
8	B	1407	PLM	C2-C3-C4-C5
8	A	1322	PLM	C9-CA-CB-CC
8	A	1322	PLM	CB-CC-CD-CE
8	B	1323	PLM	C5-C6-C7-C8
7	A	1296	RET	C11-C10-C9-C19
7	B	1296	RET	C11-C10-C9-C19
9	A	1507	HTG	C1'-C2'-C3'-C4'
8	B	1322	PLM	C4-C5-C6-C7
8	B	1322	PLM	C1-C2-C3-C4
8	B	1322	PLM	CA-CB-CC-CD
8	A	1322	PLM	C2-C3-C4-C5
8	A	1322	PLM	C1-C2-C3-C4
8	A	1410	PLM	CC-CD-CE-CF
9	B	1509	HTG	C4'-C5'-C6'-C7'
7	A	1296	RET	C7-C8-C9-C19
8	A	1323	PLM	C2-C3-C4-C5
8	B	1322	PLM	O1-C1-C2-C3
9	B	1508	HTG	C3'-C4'-C5'-C6'
8	A	1410	PLM	C2-C3-C4-C5
8	B	1322	PLM	C3-C4-C5-C6
8	A	1323	PLM	C8-C9-CA-CB
8	B	1323	PLM	O1-C1-C2-C3
10	B	1401	HTO	C1-C2-C3-O3
8	A	1322	PLM	O1-C1-C2-C3

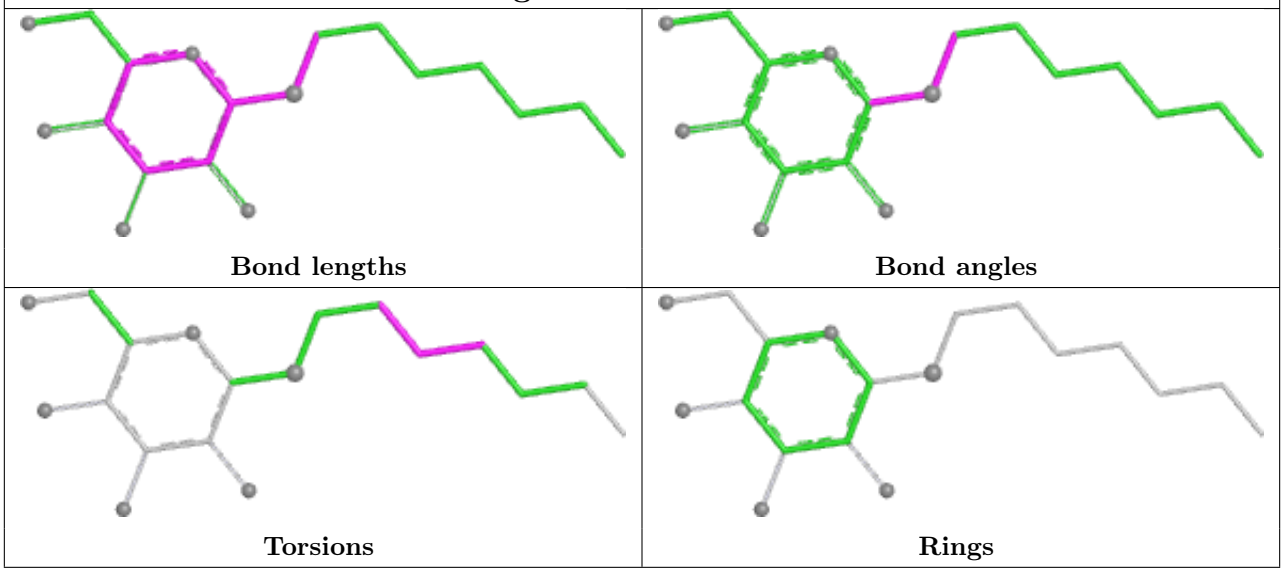
There are no ring outliers.

10 monomers are involved in 21 short contacts:

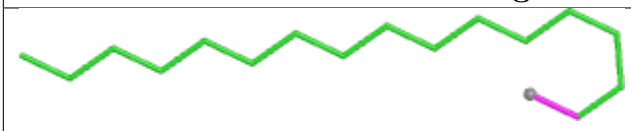
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	1322	PLM	2	0
8	A	1322	PLM	3	0
9	B	1508	HTG	2	0
8	B	1323	PLM	3	0
8	B	1407	PLM	1	0
9	B	1509	HTG	3	0
10	B	1401	HTO	1	0
7	A	1296	RET	3	0
7	B	1296	RET	4	0
8	A	1410	PLM	1	0

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 then 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.

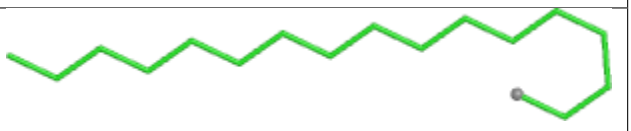
Ligand HTG B 1506



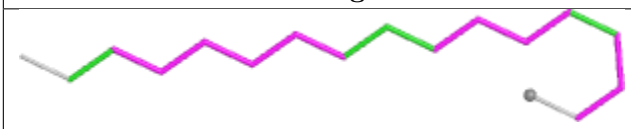
Ligand PLM B 1322



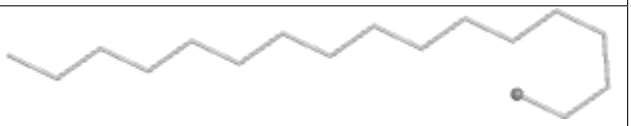
Bond lengths



Bond angles



Torsions

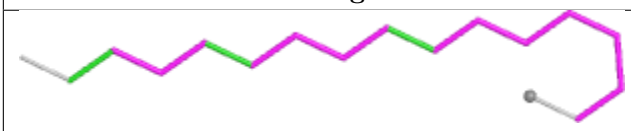


Rings

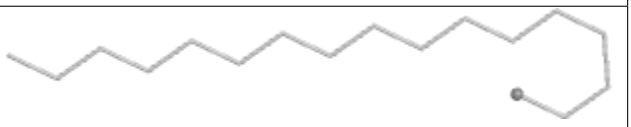
Ligand PLM A 1322

Bond lengths

Bond angles

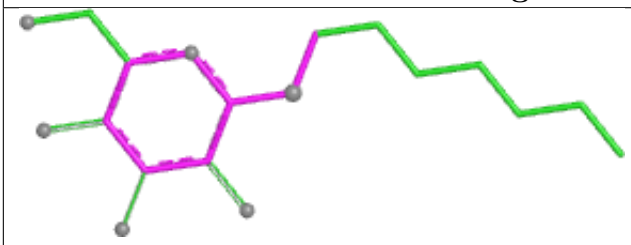


Torsions

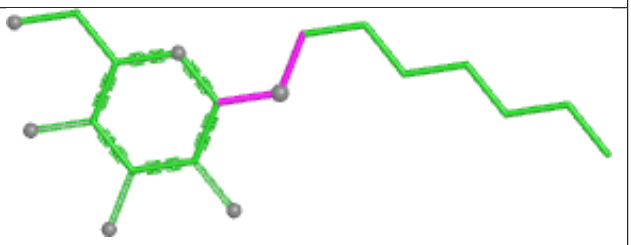


Rings

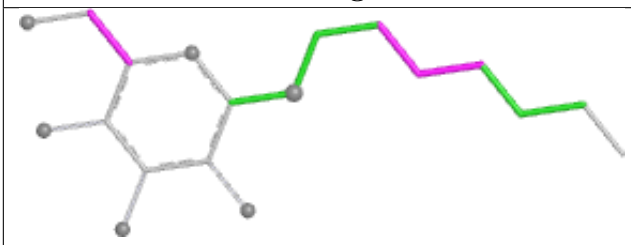
Ligand HTG A 1507



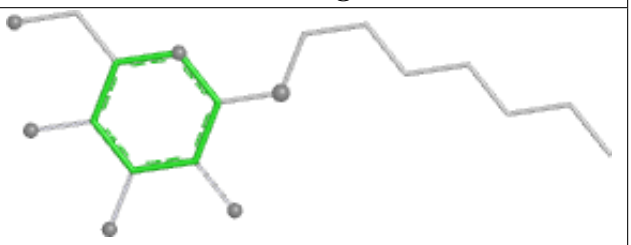
Bond lengths



Bond angles

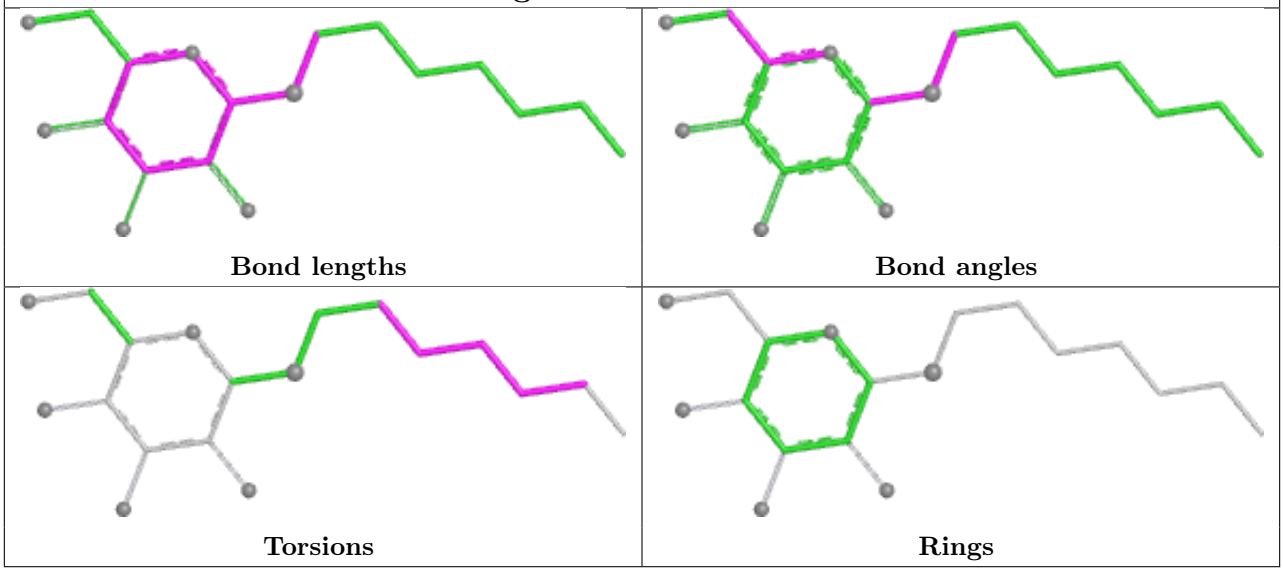


Torsions

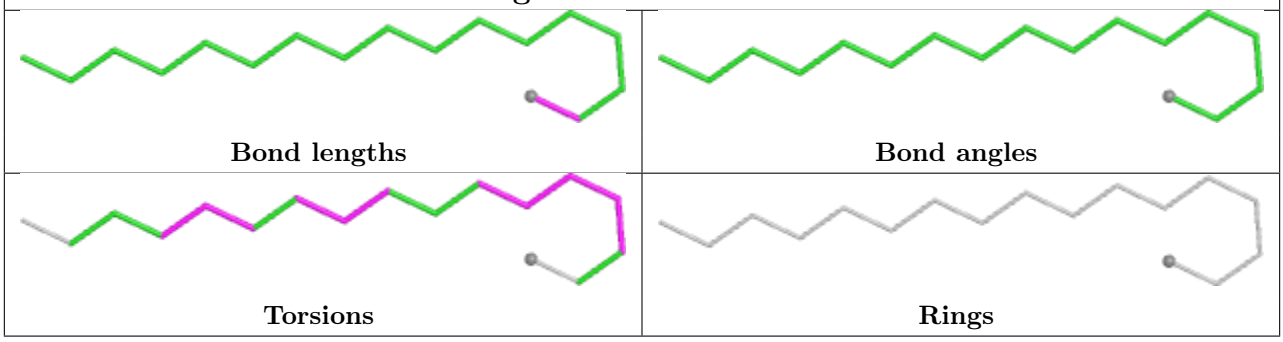


Rings

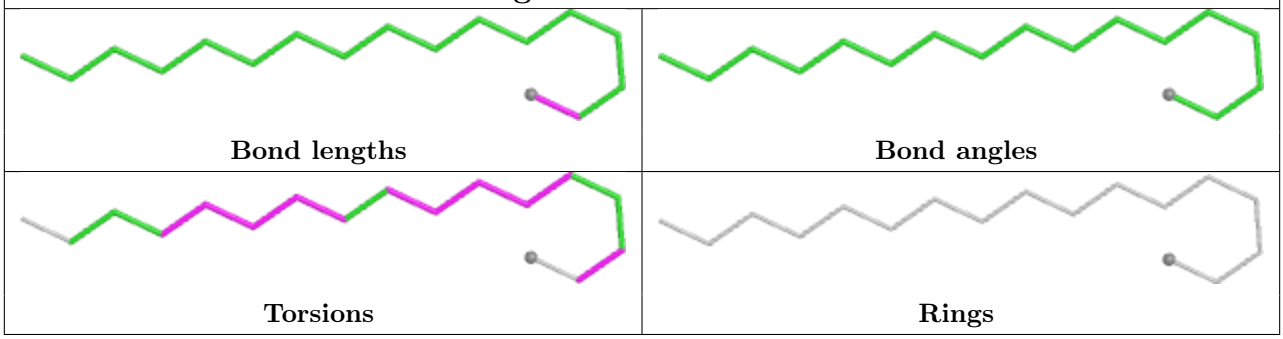
Ligand HTG B 1508



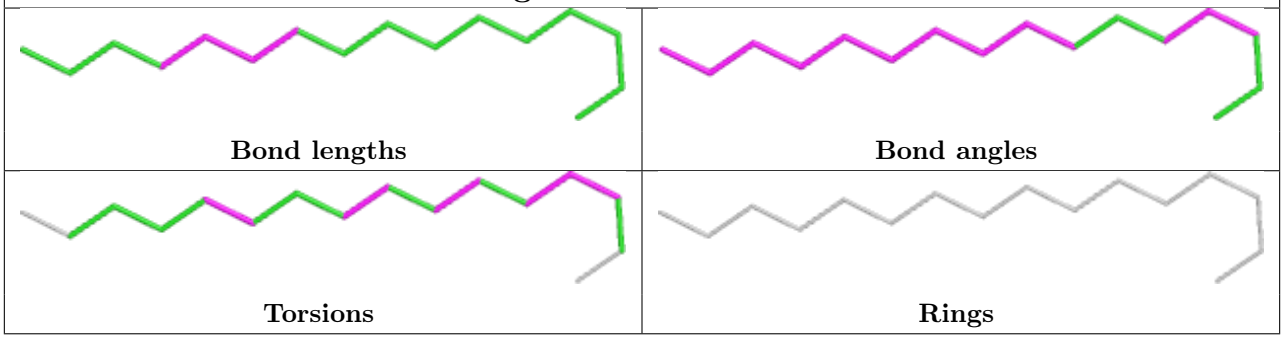
Ligand PLM A 1323



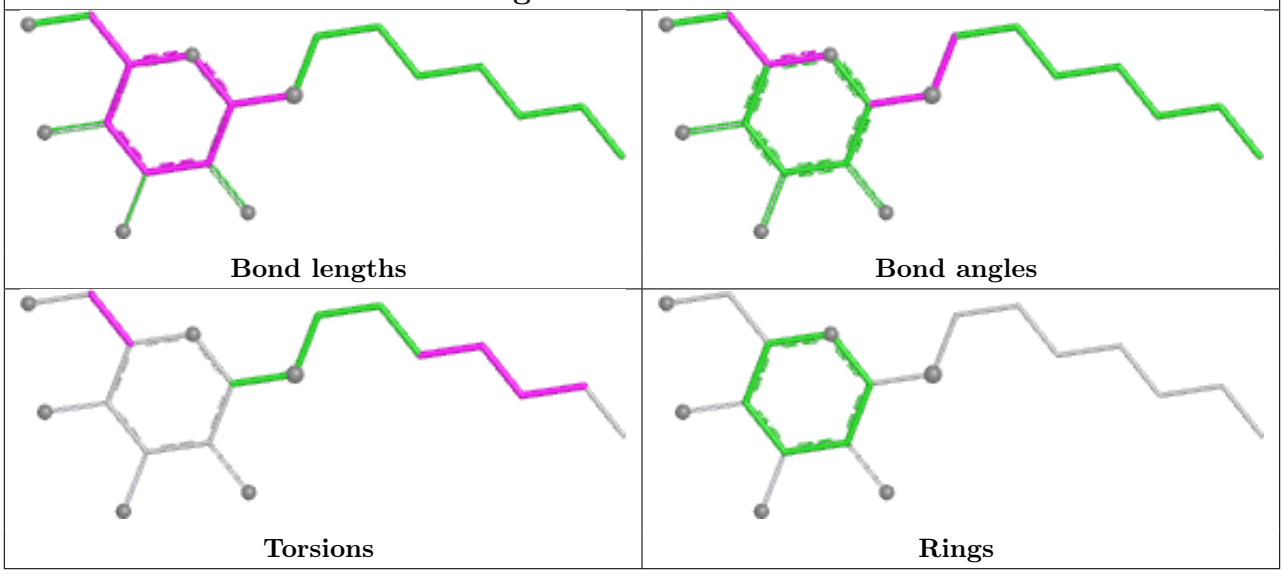
Ligand PLM B 1323



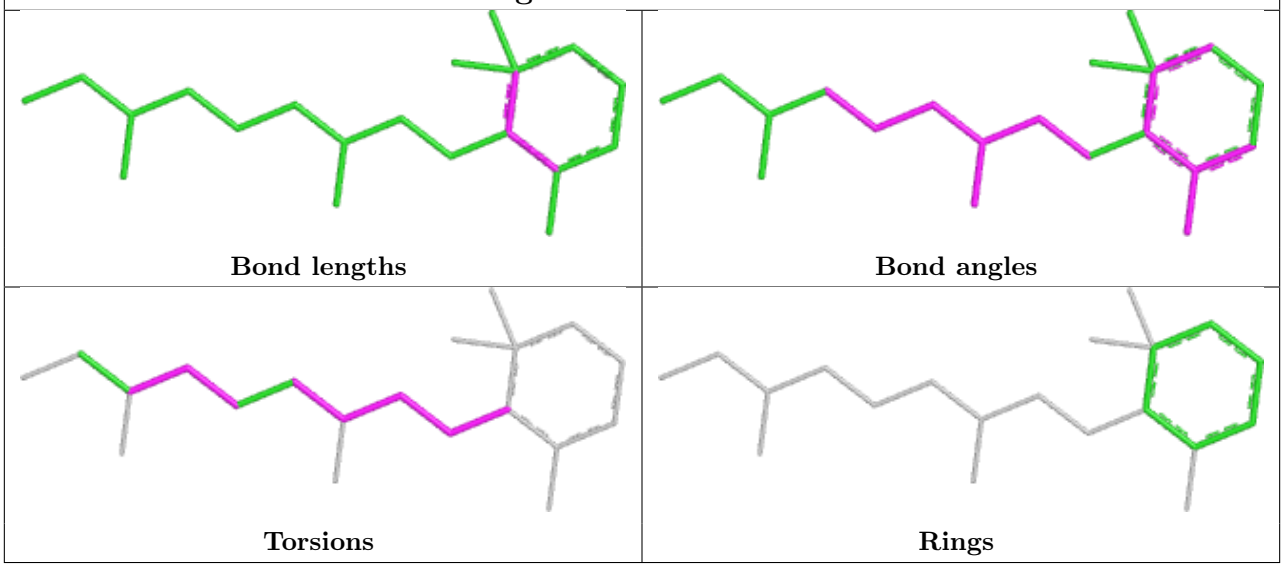
Ligand PLM B 1407



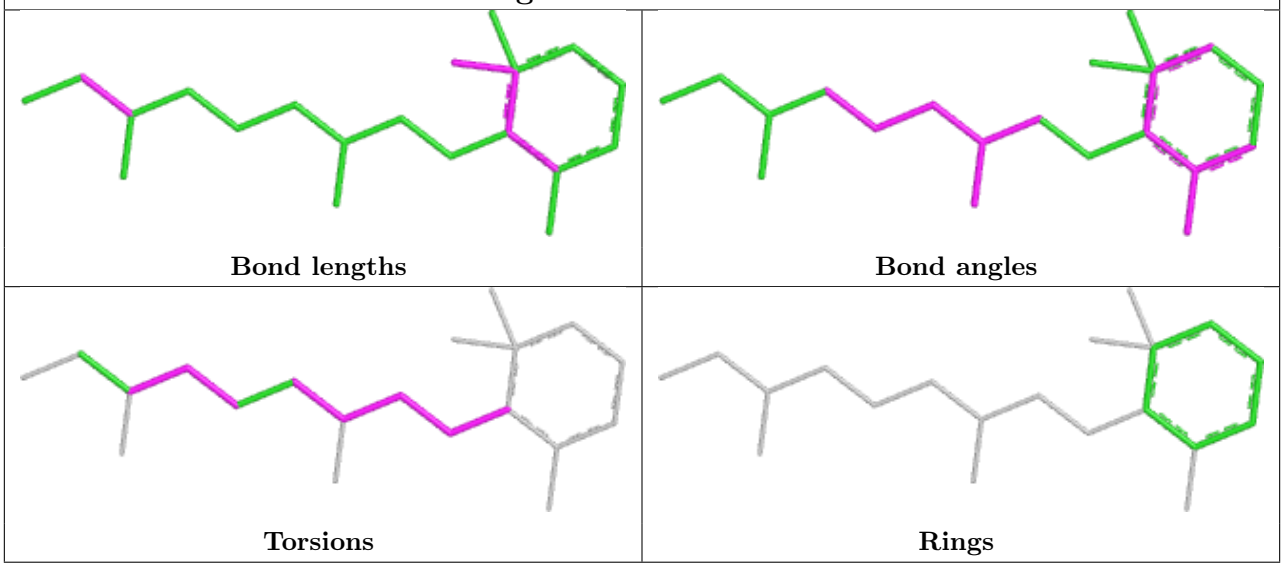
Ligand HTG B 1509



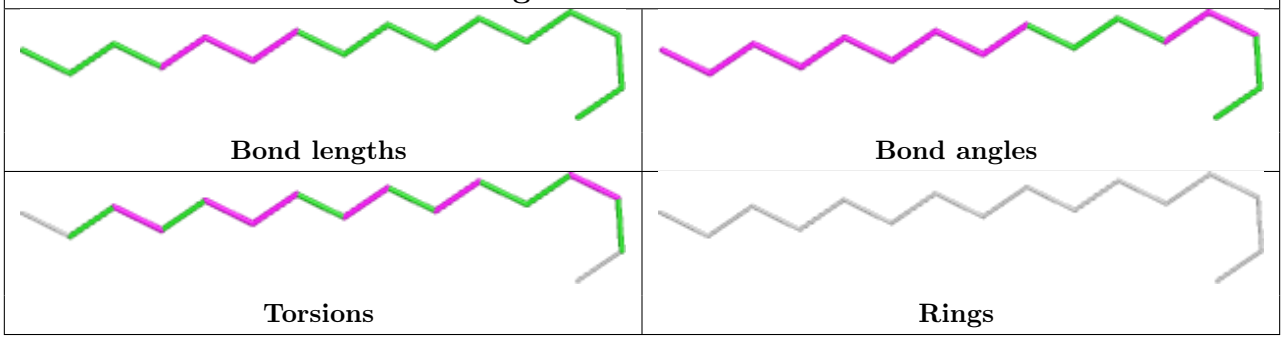
Ligand RET A 1296



Ligand RET B 1296



Ligand PLM A 1410



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 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.