Full wwPDB EM Validation Report (i)

Mar 4, 2024 – 07:11 AM EST

PDB ID	:	6CJU
EMDB ID	:	EMD-7484
Title	:	Structure of the SthK cyclic nucleotide-gated potassium channel in complex with cAMP
Authors	:	Nimigean, C.M.; Rheinberger, J.
Deposited on	:	2018-02-26
Resolution	:	3.35 Å(reported)

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 (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:

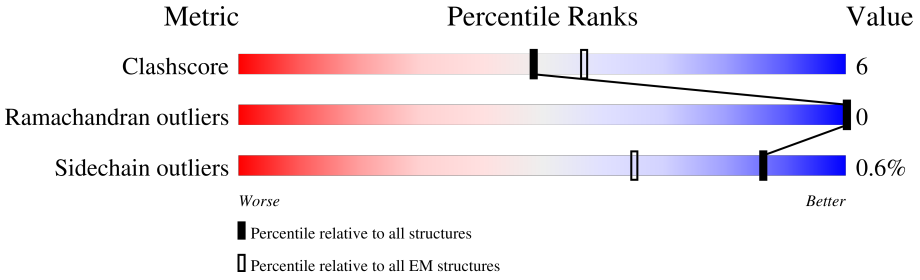
EMDB validation analysis	:	0.0.1.dev70
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.35 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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 <=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	A	456	5% 73% 13% 14%
1	B	456	5% 73% 12% 14%
1	C	456	5% 73% 12% 14%
1	D	456	5% 73% 12% 14%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 13076 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 SthK cyclic nucleotide-gated potassium channel.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	390	3102	2036	515	544	7	0	0
1	B	390	3102	2036	515	544	7	0	0
1	C	390	3102	2036	515	544	7	0	0
1	D	390	3102	2036	515	544	7	0	0

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP G0GA88
A	-17	ALA	-	expression tag	UNP G0GA88
A	-16	LYS	-	expression tag	UNP G0GA88
A	-15	ASP	-	expression tag	UNP G0GA88
A	-14	ILE	-	expression tag	UNP G0GA88
A	-13	GLY	-	expression tag	UNP G0GA88
A	-12	ILE	-	expression tag	UNP G0GA88
A	-11	ASN	-	expression tag	UNP G0GA88
A	-10	SER	-	expression tag	UNP G0GA88
A	-9	ASP	-	expression tag	UNP G0GA88
A	-8	PRO	-	expression tag	UNP G0GA88
A	-7	ASN	-	expression tag	UNP G0GA88
A	-6	SER	-	expression tag	UNP G0GA88
A	-5	SER	-	expression tag	UNP G0GA88
A	-4	SER	-	expression tag	UNP G0GA88
A	-3	VAL	-	expression tag	UNP G0GA88
A	-2	ASP	-	expression tag	UNP G0GA88
A	-1	LYS	-	expression tag	UNP G0GA88
A	0	LEU	-	expression tag	UNP G0GA88
A	421	LEU	-	expression tag	UNP G0GA88
A	422	GLU	-	expression tag	UNP G0GA88
A	423	SER	-	expression tag	UNP G0GA88

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Chain	Residue	Modelled	Actual	Comment	Reference
A	424	SER	-	expression tag	UNP G0GA88
A	425	GLY	-	expression tag	UNP G0GA88
A	426	LEU	-	expression tag	UNP G0GA88
A	427	VAL	-	expression tag	UNP G0GA88
A	428	PRO	-	expression tag	UNP G0GA88
A	429	ARG	-	expression tag	UNP G0GA88
A	430	GLY	-	expression tag	UNP G0GA88
A	431	SER	-	expression tag	UNP G0GA88
A	432	VAL	-	expression tag	UNP G0GA88
A	433	LYS	-	expression tag	UNP G0GA88
A	434	HIS	-	expression tag	UNP G0GA88
A	435	HIS	-	expression tag	UNP G0GA88
A	436	HIS	-	expression tag	UNP G0GA88
A	437	HIS	-	expression tag	UNP G0GA88
B	-18	MET	-	initiating methionine	UNP G0GA88
B	-17	ALA	-	expression tag	UNP G0GA88
B	-16	LYS	-	expression tag	UNP G0GA88
B	-15	ASP	-	expression tag	UNP G0GA88
B	-14	ILE	-	expression tag	UNP G0GA88
B	-13	GLY	-	expression tag	UNP G0GA88
B	-12	ILE	-	expression tag	UNP G0GA88
B	-11	ASN	-	expression tag	UNP G0GA88
B	-10	SER	-	expression tag	UNP G0GA88
B	-9	ASP	-	expression tag	UNP G0GA88
B	-8	PRO	-	expression tag	UNP G0GA88
B	-7	ASN	-	expression tag	UNP G0GA88
B	-6	SER	-	expression tag	UNP G0GA88
B	-5	SER	-	expression tag	UNP G0GA88
B	-4	SER	-	expression tag	UNP G0GA88
B	-3	VAL	-	expression tag	UNP G0GA88
B	-2	ASP	-	expression tag	UNP G0GA88
B	-1	LYS	-	expression tag	UNP G0GA88
B	0	LEU	-	expression tag	UNP G0GA88
B	421	LEU	-	expression tag	UNP G0GA88
B	422	GLU	-	expression tag	UNP G0GA88
B	423	SER	-	expression tag	UNP G0GA88
B	424	SER	-	expression tag	UNP G0GA88
B	425	GLY	-	expression tag	UNP G0GA88
B	426	LEU	-	expression tag	UNP G0GA88
B	427	VAL	-	expression tag	UNP G0GA88
B	428	PRO	-	expression tag	UNP G0GA88
B	429	ARG	-	expression tag	UNP G0GA88

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Chain	Residue	Modelled	Actual	Comment	Reference
B	430	GLY	-	expression tag	UNP G0GA88
B	431	SER	-	expression tag	UNP G0GA88
B	432	VAL	-	expression tag	UNP G0GA88
B	433	LYS	-	expression tag	UNP G0GA88
B	434	HIS	-	expression tag	UNP G0GA88
B	435	HIS	-	expression tag	UNP G0GA88
B	436	HIS	-	expression tag	UNP G0GA88
B	437	HIS	-	expression tag	UNP G0GA88
C	-18	MET	-	initiating methionine	UNP G0GA88
C	-17	ALA	-	expression tag	UNP G0GA88
C	-16	LYS	-	expression tag	UNP G0GA88
C	-15	ASP	-	expression tag	UNP G0GA88
C	-14	ILE	-	expression tag	UNP G0GA88
C	-13	GLY	-	expression tag	UNP G0GA88
C	-12	ILE	-	expression tag	UNP G0GA88
C	-11	ASN	-	expression tag	UNP G0GA88
C	-10	SER	-	expression tag	UNP G0GA88
C	-9	ASP	-	expression tag	UNP G0GA88
C	-8	PRO	-	expression tag	UNP G0GA88
C	-7	ASN	-	expression tag	UNP G0GA88
C	-6	SER	-	expression tag	UNP G0GA88
C	-5	SER	-	expression tag	UNP G0GA88
C	-4	SER	-	expression tag	UNP G0GA88
C	-3	VAL	-	expression tag	UNP G0GA88
C	-2	ASP	-	expression tag	UNP G0GA88
C	-1	LYS	-	expression tag	UNP G0GA88
C	0	LEU	-	expression tag	UNP G0GA88
C	421	LEU	-	expression tag	UNP G0GA88
C	422	GLU	-	expression tag	UNP G0GA88
C	423	SER	-	expression tag	UNP G0GA88
C	424	SER	-	expression tag	UNP G0GA88
C	425	GLY	-	expression tag	UNP G0GA88
C	426	LEU	-	expression tag	UNP G0GA88
C	427	VAL	-	expression tag	UNP G0GA88
C	428	PRO	-	expression tag	UNP G0GA88
C	429	ARG	-	expression tag	UNP G0GA88
C	430	GLY	-	expression tag	UNP G0GA88
C	431	SER	-	expression tag	UNP G0GA88
C	432	VAL	-	expression tag	UNP G0GA88
C	433	LYS	-	expression tag	UNP G0GA88
C	434	HIS	-	expression tag	UNP G0GA88
C	435	HIS	-	expression tag	UNP G0GA88

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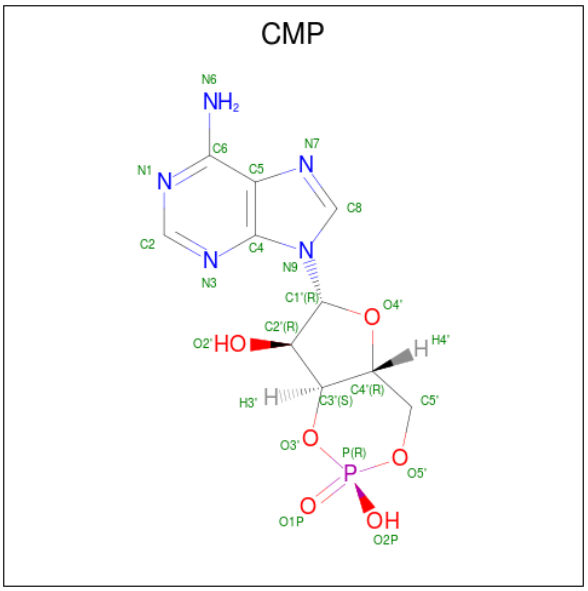
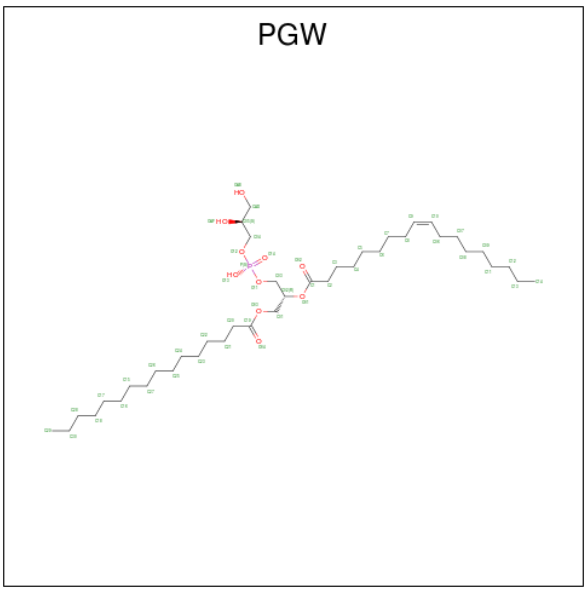
Chain	Residue	Modelled	Actual	Comment	Reference
C	436	HIS	-	expression tag	UNP G0GA88
C	437	HIS	-	expression tag	UNP G0GA88
D	-18	MET	-	initiating methionine	UNP G0GA88
D	-17	ALA	-	expression tag	UNP G0GA88
D	-16	LYS	-	expression tag	UNP G0GA88
D	-15	ASP	-	expression tag	UNP G0GA88
D	-14	ILE	-	expression tag	UNP G0GA88
D	-13	GLY	-	expression tag	UNP G0GA88
D	-12	ILE	-	expression tag	UNP G0GA88
D	-11	ASN	-	expression tag	UNP G0GA88
D	-10	SER	-	expression tag	UNP G0GA88
D	-9	ASP	-	expression tag	UNP G0GA88
D	-8	PRO	-	expression tag	UNP G0GA88
D	-7	ASN	-	expression tag	UNP G0GA88
D	-6	SER	-	expression tag	UNP G0GA88
D	-5	SER	-	expression tag	UNP G0GA88
D	-4	SER	-	expression tag	UNP G0GA88
D	-3	VAL	-	expression tag	UNP G0GA88
D	-2	ASP	-	expression tag	UNP G0GA88
D	-1	LYS	-	expression tag	UNP G0GA88
D	0	LEU	-	expression tag	UNP G0GA88
D	421	LEU	-	expression tag	UNP G0GA88
D	422	GLU	-	expression tag	UNP G0GA88
D	423	SER	-	expression tag	UNP G0GA88
D	424	SER	-	expression tag	UNP G0GA88
D	425	GLY	-	expression tag	UNP G0GA88
D	426	LEU	-	expression tag	UNP G0GA88
D	427	VAL	-	expression tag	UNP G0GA88
D	428	PRO	-	expression tag	UNP G0GA88
D	429	ARG	-	expression tag	UNP G0GA88
D	430	GLY	-	expression tag	UNP G0GA88
D	431	SER	-	expression tag	UNP G0GA88
D	432	VAL	-	expression tag	UNP G0GA88
D	433	LYS	-	expression tag	UNP G0GA88
D	434	HIS	-	expression tag	UNP G0GA88
D	435	HIS	-	expression tag	UNP G0GA88
D	436	HIS	-	expression tag	UNP G0GA88
D	437	HIS	-	expression tag	UNP G0GA88

- Molecule 2 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (three-letter code: CMP) (formula: $C_{10}H_{12}N_5O_6P$).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
2	A	1	22	10	5	6	1	0
2	B	1	22	10	5	6	1	0
2	C	1	22	10	5	6	1	0
2	D	1	22	10	5	6	1	0

- Molecule 3 is (1R)-2-{[(S)-{[(2S)-2,3-dihydroxypropyl]oxy}(hydroxy)phosphoryl]oxy}-1-[(hexadecanoyloxy)methyl]ethyl (9Z)-octadec-9-enoate (three-letter code: PGW) (formula: $C_{40}H_{77}O_{10}P$).



Mol	Chain	Residues	Atoms	AltConf
3	A	1	Total C O P 43 32 10 1	0
3	A	1	Total C 10 10	0
3	A	1	Total C 16 16	0
3	A	1	Total C 8 8	0
3	A	1	Total C 9 9	0
3	A	1	Total C O 29 25 4	0
3	A	1	Total C 14 14	0
3	A	1	Total C 16 16	0
3	B	1	Total C O P 43 32 10 1	0
3	B	1	Total C 10 10	0
3	B	1	Total C 16 16	0
3	B	1	Total C 8 8	0
3	B	1	Total C 9 9	0
3	B	1	Total C O 29 25 4	0
3	B	1	Total C 14 14	0
3	B	1	Total C 16 16	0
3	C	1	Total C O P 43 32 10 1	0
3	C	1	Total C 10 10	0
3	C	1	Total C 16 16	0
3	C	1	Total C 8 8	0
3	C	1	Total C 9 9	0
3	C	1	Total C O 29 25 4	0

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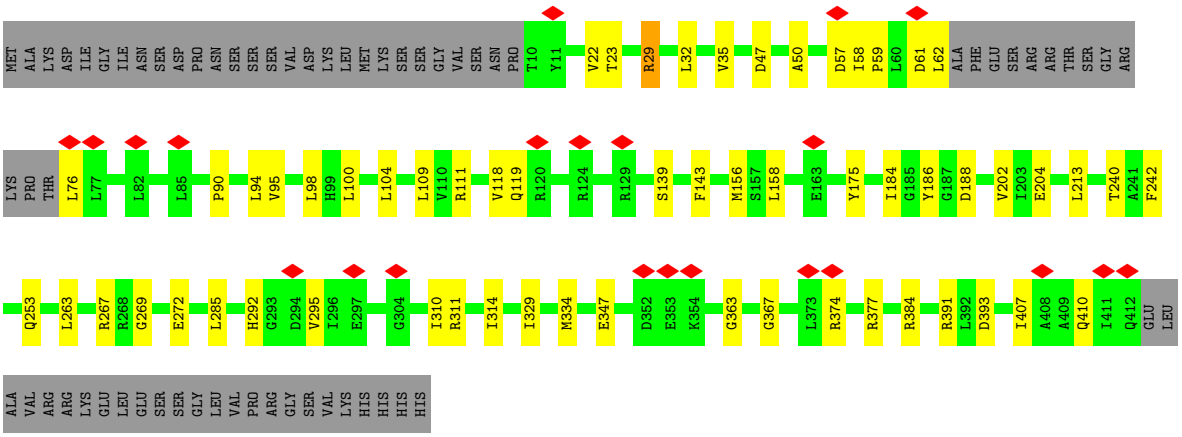
Mol	Chain	Residues	Atoms	AltConf
3	C	1	Total C 14 14	0
3	C	1	Total C 16 16	0
3	D	1	Total C O P 43 32 10 1	0
3	D	1	Total C 10 10	0
3	D	1	Total C 16 16	0
3	D	1	Total C 8 8	0
3	D	1	Total C 9 9	0
3	D	1	Total C O 29 25 4	0
3	D	1	Total C 14 14	0
3	D	1	Total C 16 16	0

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 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.

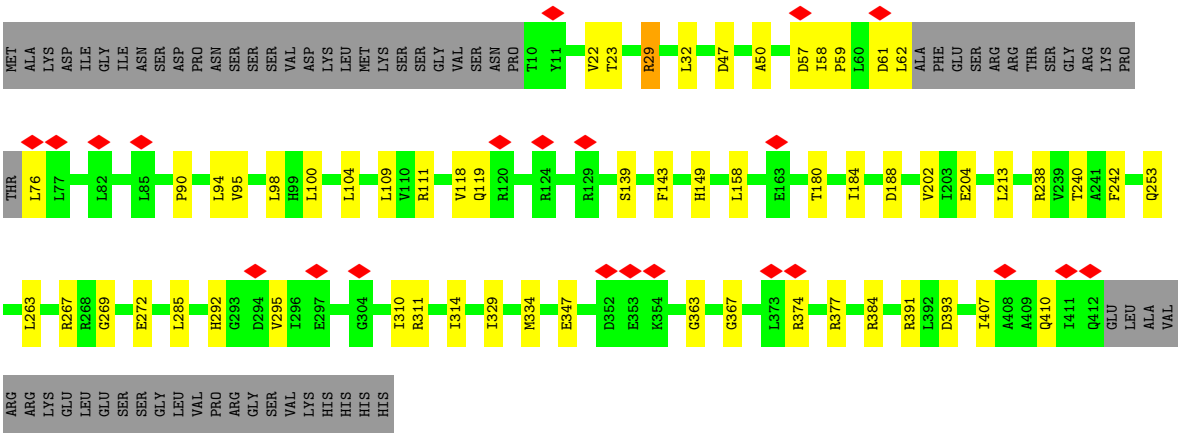
• Molecule 1: SthK cyclic nucleotide-gated potassium channel

Chain A:



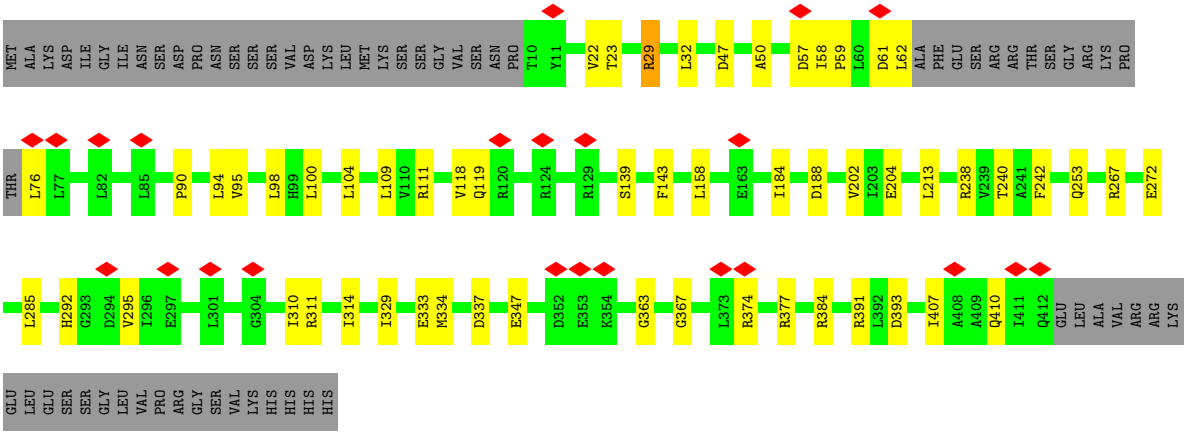
• Molecule 1: SthK cyclic nucleotide-gated potassium channel

Chain B:



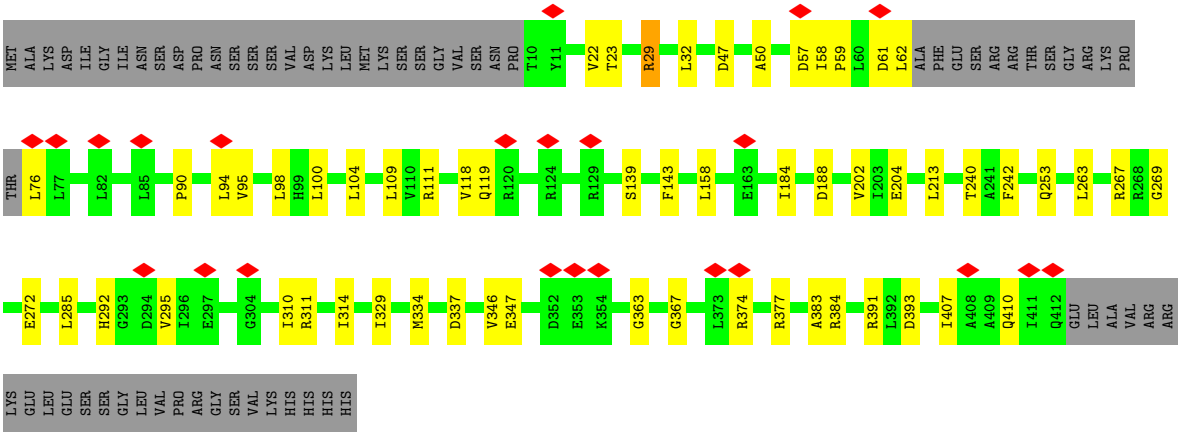
• Molecule 1: SthK cyclic nucleotide-gated potassium channel

Chain C:



• Molecule 1: SthK cyclic nucleotide-gated potassium channel

Chain D:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	81501	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	72	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.393	Depositor
Minimum map value	-0.225	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.010	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	274.752, 274.752, 274.752	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.07325, 1.07325, 1.07325	Depositor

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: CMP, PGW

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	0.47	0/3173	0.68	0/4325
1	B	0.47	0/3173	0.68	0/4325
1	C	0.47	0/3173	0.68	0/4325
1	D	0.47	0/3173	0.68	0/4325
All	All	0.47	0/12692	0.68	0/17300

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3102	0	3239	44	0
1	B	3102	0	3239	44	0
1	C	3102	0	3239	43	0
1	D	3102	0	3239	44	0
2	A	22	0	11	2	0
2	B	22	0	11	2	0
2	C	22	0	11	2	0
2	D	22	0	11	2	0
3	A	145	0	224	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	145	0	224	11	0
3	C	145	0	224	10	0
3	D	145	0	224	11	0
All	All	13076	0	13896	168	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 6.

All (168) 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)
2:C:500:CMP:H2	2:C:500:CMP:C2	0.97	1.50
2:D:500:CMP:H2	2:D:500:CMP:C2	0.97	1.49
2:A:500:CMP:H2	2:A:500:CMP:C2	0.97	1.48
2:B:500:CMP:H2	2:B:500:CMP:C2	0.97	1.48
1:A:143:PHE:CE1	3:A:506:PGW:C29	2.48	0.97
1:C:143:PHE:CE1	3:C:506:PGW:C29	2.48	0.96
1:D:143:PHE:CE1	3:D:506:PGW:C29	2.48	0.96
1:B:143:PHE:CE1	3:B:506:PGW:C29	2.48	0.96
1:D:143:PHE:HE1	3:D:506:PGW:C29	2.01	0.73
1:A:143:PHE:HE1	3:A:506:PGW:C29	2.01	0.73
1:B:143:PHE:HE1	3:B:506:PGW:C29	2.01	0.72
1:C:143:PHE:CE1	3:C:506:PGW:H29A	2.24	0.72
1:D:143:PHE:CE1	3:D:506:PGW:H29A	2.24	0.71
1:B:143:PHE:CE1	3:B:506:PGW:H29A	2.24	0.71
1:D:143:PHE:HE1	3:D:506:PGW:H29A	1.56	0.70
1:A:143:PHE:HE1	3:A:506:PGW:H29A	1.56	0.70
1:A:143:PHE:CE1	3:A:506:PGW:H29A	2.24	0.70
1:C:143:PHE:HE1	3:C:506:PGW:C29	2.01	0.70
1:B:143:PHE:HE1	3:B:506:PGW:H29A	1.56	0.69
1:D:143:PHE:CE1	3:D:506:PGW:H29B	2.28	0.68
1:A:143:PHE:CE1	3:A:506:PGW:H29B	2.28	0.66
1:C:143:PHE:HE1	3:C:506:PGW:H29A	1.56	0.66
1:C:143:PHE:CE1	3:C:506:PGW:H29B	2.28	0.66
1:B:143:PHE:CE1	3:B:506:PGW:H29B	2.28	0.64
1:D:62:LEU:HD11	1:D:76:LEU:HD13	1.81	0.62
1:A:62:LEU:HD11	1:A:76:LEU:HD13	1.81	0.62
1:B:62:LEU:HD11	1:B:76:LEU:HD13	1.81	0.61
1:C:62:LEU:HD11	1:C:76:LEU:HD13	1.81	0.61
1:D:95:VAL:HA	1:D:100:LEU:HD13	1.83	0.61
1:C:95:VAL:HA	1:C:100:LEU:HD13	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:VAL:HA	1:A:100:LEU:HD13	1.83	0.60
1:B:95:VAL:HA	1:B:100:LEU:HD13	1.83	0.59
1:C:90:PRO:HA	1:C:111:ARG:HH11	1.68	0.58
1:B:90:PRO:HA	1:B:111:ARG:HH11	1.68	0.58
1:A:90:PRO:HA	1:A:111:ARG:HH11	1.68	0.57
1:C:202:VAL:HG11	3:C:507:PGW:H15	1.86	0.57
1:D:90:PRO:HA	1:D:111:ARG:HH11	1.68	0.57
1:D:202:VAL:HG11	3:D:507:PGW:H15	1.86	0.57
1:C:347:GLU:HB3	1:C:384:ARG:HD2	1.87	0.57
1:B:202:VAL:HG11	3:B:507:PGW:H15	1.86	0.57
1:A:347:GLU:HB3	1:A:384:ARG:HD2	1.87	0.56
1:B:347:GLU:HB3	1:B:384:ARG:HD2	1.87	0.56
1:A:202:VAL:HG11	3:A:507:PGW:H15	1.86	0.56
1:B:119:GLN:NE2	1:B:139:SER:OG	2.36	0.56
1:A:119:GLN:NE2	1:A:139:SER:OG	2.36	0.56
1:D:347:GLU:HB3	1:D:384:ARG:HD2	1.87	0.56
1:C:242:PHE:HZ	1:D:272:GLU:HG2	1.72	0.55
1:A:285:LEU:HD21	1:A:311:ARG:HD2	1.90	0.54
1:B:242:PHE:HZ	1:C:272:GLU:HG2	1.72	0.54
1:B:285:LEU:HD21	1:B:311:ARG:HD2	1.90	0.54
1:A:204:GLU:HG2	1:D:184:ILE:CD1	2.38	0.54
1:A:242:PHE:HZ	1:B:272:GLU:HG2	1.72	0.54
1:C:119:GLN:NE2	1:C:139:SER:OG	2.36	0.54
1:A:272:GLU:HG2	1:D:242:PHE:HZ	1.72	0.54
1:C:184:ILE:CD1	1:D:204:GLU:HG2	2.38	0.54
1:D:119:GLN:NE2	1:D:139:SER:OG	2.36	0.53
1:A:213:LEU:HD21	3:D:508:PGW:H23A	1.91	0.53
3:B:508:PGW:H23A	1:C:213:LEU:HD21	1.91	0.53
3:A:508:PGW:H23A	1:B:213:LEU:HD21	1.91	0.53
1:B:184:ILE:CD1	1:C:204:GLU:HG2	2.38	0.53
1:A:184:ILE:CD1	1:B:204:GLU:HG2	2.38	0.53
3:C:508:PGW:H23A	1:D:213:LEU:HD21	1.91	0.53
1:D:285:LEU:HD21	1:D:311:ARG:HD2	1.90	0.53
1:C:285:LEU:HD21	1:C:311:ARG:HD2	1.90	0.53
1:A:240:THR:HG23	1:A:253:GLN:HE21	1.74	0.53
1:D:240:THR:HG23	1:D:253:GLN:HE21	1.74	0.53
1:C:240:THR:HG23	1:C:253:GLN:HE21	1.74	0.52
1:A:367:GLY:HA2	2:A:500:CMP:O1P	2.10	0.51
1:B:367:GLY:HA2	2:B:500:CMP:O1P	2.10	0.51
1:B:240:THR:HG23	1:B:253:GLN:HE21	1.74	0.51
1:C:367:GLY:HA2	2:C:500:CMP:O1P	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:367:GLY:HA2	2:D:500:CMP:O1P	2.10	0.51
1:B:23:THR:HG21	1:B:118:VAL:HG21	1.94	0.50
1:D:292:HIS:HB2	1:D:314:ILE:HD11	1.94	0.50
1:C:292:HIS:HB2	1:C:314:ILE:HD11	1.94	0.50
1:B:292:HIS:HB2	1:B:314:ILE:HD11	1.94	0.50
1:A:292:HIS:HB2	1:A:314:ILE:HD11	1.94	0.49
1:C:23:THR:HG21	1:C:118:VAL:HG21	1.94	0.49
1:C:109:LEU:HD21	3:C:506:PGW:H27	1.95	0.49
1:D:109:LEU:HD21	3:D:506:PGW:H27	1.95	0.49
1:A:58:ILE:HG23	1:A:59:PRO:HD3	1.95	0.49
1:D:58:ILE:HG23	1:D:59:PRO:HD3	1.95	0.49
1:A:23:THR:HG21	1:A:118:VAL:HG21	1.94	0.49
1:D:23:THR:HG21	1:D:118:VAL:HG21	1.94	0.49
1:A:186:TYR:HH	1:B:180:THR:HG1	1.52	0.49
1:B:58:ILE:HG23	1:B:59:PRO:HD3	1.95	0.49
1:C:58:ILE:HG23	1:C:59:PRO:HD3	1.95	0.49
1:B:109:LEU:HD21	3:B:506:PGW:H27	1.95	0.48
1:C:158:LEU:HD23	3:C:506:PGW:H2A	1.95	0.48
1:D:158:LEU:HD23	3:D:506:PGW:H2A	1.95	0.48
1:C:391:ARG:NH1	1:C:393:ASP:OD1	2.46	0.48
1:A:109:LEU:HD21	3:A:506:PGW:H27	1.95	0.48
1:B:391:ARG:NH1	1:B:393:ASP:OD1	2.46	0.48
1:A:391:ARG:NH1	1:A:393:ASP:OD1	2.46	0.47
1:A:158:LEU:HD23	3:A:506:PGW:H2A	1.95	0.47
1:B:158:LEU:HD23	3:B:506:PGW:H2A	1.95	0.47
1:D:391:ARG:NH1	1:D:393:ASP:OD1	2.46	0.47
1:A:295:VAL:HG22	1:A:363:GLY:HA2	1.97	0.47
1:A:94:LEU:O	1:A:98:LEU:HB2	2.16	0.46
1:B:94:LEU:O	1:B:98:LEU:HB2	2.16	0.46
1:B:295:VAL:HG22	1:B:363:GLY:HA2	1.97	0.46
1:C:22:VAL:HG13	1:C:50:ALA:HB1	1.98	0.46
1:C:94:LEU:O	1:C:98:LEU:HB2	2.16	0.46
1:C:337:ASP:OD1	1:C:337:ASP:N	2.49	0.46
1:A:22:VAL:HG13	1:A:50:ALA:HB1	1.98	0.46
1:C:295:VAL:HG22	1:C:363:GLY:HA2	1.97	0.45
1:D:295:VAL:HG22	1:D:363:GLY:HA2	1.97	0.45
1:D:94:LEU:O	1:D:98:LEU:HB2	2.16	0.45
1:B:22:VAL:HG13	1:B:50:ALA:HB1	1.98	0.44
1:D:22:VAL:HG13	1:D:50:ALA:HB1	1.98	0.44
1:D:337:ASP:OD1	1:D:337:ASP:N	2.49	0.44
1:B:100:LEU:HD23	1:B:104:LEU:HG	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:100:LEU:HD23	1:C:104:LEU:HG	1.99	0.44
1:A:95:VAL:HG22	1:A:100:LEU:HD22	2.00	0.44
1:B:95:VAL:HG22	1:B:100:LEU:HD22	2.00	0.44
1:A:188:ASP:OD1	1:A:188:ASP:N	2.50	0.43
1:D:100:LEU:HD23	1:D:104:LEU:HG	1.99	0.43
1:A:100:LEU:HD23	1:A:104:LEU:HG	1.99	0.43
1:A:242:PHE:CZ	1:B:272:GLU:HG2	2.53	0.43
1:C:95:VAL:HG22	1:C:100:LEU:HD22	2.00	0.43
1:B:242:PHE:CZ	1:C:272:GLU:HG2	2.53	0.43
1:D:188:ASP:OD1	1:D:188:ASP:N	2.50	0.43
1:A:32:LEU:HD11	3:A:503:PGW:H15	2.01	0.43
1:B:188:ASP:OD1	1:B:188:ASP:N	2.50	0.43
1:B:32:LEU:HD11	3:B:503:PGW:H15	2.01	0.43
1:C:32:LEU:HD11	3:C:503:PGW:H15	2.01	0.42
1:D:32:LEU:HD11	3:D:503:PGW:H15	2.01	0.42
1:D:95:VAL:HG22	1:D:100:LEU:HD22	2.00	0.42
3:A:501:PGW:O12	3:A:501:PGW:OAE	2.33	0.42
1:C:242:PHE:CZ	1:D:272:GLU:HG2	2.53	0.42
1:D:57:ASP:O	1:D:61:ASP:N	2.53	0.42
1:C:188:ASP:OD1	1:C:188:ASP:N	2.50	0.42
1:A:57:ASP:O	1:A:61:ASP:N	2.53	0.42
1:C:334:MET:HG3	1:C:374:ARG:HH12	1.84	0.42
1:A:272:GLU:HG2	1:D:242:PHE:CZ	2.53	0.42
1:B:334:MET:HG3	1:B:374:ARG:HH12	1.84	0.42
3:D:501:PGW:O12	3:D:501:PGW:OAE	2.33	0.42
1:A:334:MET:HG3	1:A:374:ARG:HH12	1.84	0.41
3:B:501:PGW:O12	3:B:501:PGW:OAE	2.33	0.41
1:B:57:ASP:O	1:B:61:ASP:N	2.53	0.41
1:C:407:ILE:HA	1:C:410:GLN:HB3	2.03	0.41
1:A:407:ILE:HA	1:A:410:GLN:HB3	2.03	0.41
1:B:29:ARG:NH1	1:B:47:ASP:OD1	2.54	0.41
1:B:407:ILE:HA	1:B:410:GLN:HB3	2.03	0.41
1:D:407:ILE:HA	1:D:410:GLN:HB3	2.03	0.41
1:A:29:ARG:NH1	1:A:47:ASP:OD1	2.54	0.41
1:D:329:ILE:HG21	1:D:377:ARG:HD2	2.02	0.41
1:C:57:ASP:O	1:C:61:ASP:N	2.53	0.41
1:D:334:MET:HG3	1:D:374:ARG:HH12	1.84	0.41
1:A:175:TYR:OH	1:A:188:ASP:OD2	2.28	0.41
1:C:238:ARG:HH11	1:C:238:ARG:HD3	1.76	0.41
1:A:35:VAL:HG11	1:A:156:MET:HE2	2.03	0.41
1:A:263:LEU:HD23	1:A:269:GLY:HA2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:329:ILE:HG21	1:B:377:ARG:HD2	2.02	0.41
1:C:29:ARG:NH1	1:C:47:ASP:OD1	2.54	0.41
1:D:29:ARG:NH1	1:D:47:ASP:OD1	2.54	0.41
1:D:346:VAL:HG23	1:D:383:ALA:HA	2.03	0.41
1:B:263:LEU:HD23	1:B:269:GLY:HA2	2.03	0.41
1:A:310:ILE:O	1:A:314:ILE:HB	2.22	0.40
1:B:310:ILE:O	1:B:314:ILE:HB	2.22	0.40
1:C:310:ILE:O	1:C:314:ILE:HB	2.22	0.40
1:C:329:ILE:HG21	1:C:377:ARG:HD2	2.02	0.40
1:D:310:ILE:O	1:D:314:ILE:HB	2.22	0.40
1:A:329:ILE:HG21	1:A:377:ARG:HD2	2.02	0.40
1:B:238:ARG:HH11	1:B:238:ARG:HD3	1.76	0.40
1:D:263:LEU:HD23	1:D:269:GLY:HA2	2.03	0.40
1:B:32:LEU:HD12	1:B:149:HIS:HE1	1.87	0.40
1:C:333:GLU:OE2	1:D:311:ARG:NH2	2.55	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 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	A	386/456 (85%)	358 (93%)	28 (7%)	0	100	100
1	B	386/456 (85%)	358 (93%)	28 (7%)	0	100	100
1	C	386/456 (85%)	358 (93%)	28 (7%)	0	100	100
1	D	386/456 (85%)	358 (93%)	28 (7%)	0	100	100
All	All	1544/1824 (85%)	1432 (93%)	112 (7%)	0	100	100

There are no Ramachandran outliers to report.

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 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	A	335/393 (85%)	333 (99%)	2 (1%)	86	93
1	B	335/393 (85%)	333 (99%)	2 (1%)	86	93
1	C	335/393 (85%)	333 (99%)	2 (1%)	86	93
1	D	335/393 (85%)	333 (99%)	2 (1%)	86	93
All	All	1340/1572 (85%)	1332 (99%)	8 (1%)	86	93

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

Mol	Chain	Res	Type
1	A	29	ARG
1	A	267	ARG
1	B	29	ARG
1	B	267	ARG
1	C	29	ARG
1	C	267	ARG
1	D	29	ARG
1	D	267	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
1	A	253	GLN
1	B	119	GLN
1	B	253	GLN
1	C	119	GLN
1	C	253	GLN
1	D	119	GLN
1	D	253	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)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

36 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	$\#\|Z\| > 2$
3	PGW	A	504	-	7,7,50	0.19	0	6,6,56	0.55	0
3	PGW	D	501	-	42,42,50	0.90	3 (7%)	44,48,56	1.03	2 (4%)
3	PGW	B	506	-	28,28,50	0.66	1 (3%)	30,30,56	0.99	1 (3%)
3	PGW	C	506	-	28,28,50	0.66	1 (3%)	30,30,56	0.99	1 (3%)
3	PGW	D	504	-	7,7,50	0.19	0	6,6,56	0.55	0
3	PGW	D	508	-	15,15,50	0.24	0	14,14,56	1.35	0
3	PGW	D	505	-	8,8,50	0.16	0	7,7,56	1.22	0
3	PGW	D	506	-	28,28,50	0.66	1 (3%)	30,30,56	0.99	1 (3%)
3	PGW	C	508	-	15,15,50	0.24	0	14,14,56	1.35	0
3	PGW	C	502	-	9,9,50	0.13	0	8,8,56	1.09	0
2	CMP	C	500	-	22,25,25	1.38	4 (18%)	24,39,39	2.14	4 (16%)
2	CMP	B	500	-	22,25,25	1.38	4 (18%)	24,39,39	2.14	4 (16%)
3	PGW	A	506	-	28,28,50	0.66	1 (3%)	30,30,56	0.99	1 (3%)
3	PGW	D	503	-	15,15,50	0.21	0	14,14,56	1.31	0
3	PGW	B	508	-	15,15,50	0.24	0	14,14,56	1.35	0
3	PGW	D	502	-	9,9,50	0.13	0	8,8,56	1.09	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	PGW	A	502	-	9,9,50	0.13	0	8,8,56	1.09	0
3	PGW	A	507	-	13,13,50	0.19	0	12,12,56	1.24	0
2	CMP	A	500	-	22,25,25	1.38	4 (18%)	24,39,39	2.14	4 (16%)
2	CMP	D	500	-	22,25,25	1.38	4 (18%)	24,39,39	2.14	4 (16%)
3	PGW	A	501	-	42,42,50	0.90	3 (7%)	44,48,56	1.03	2 (4%)
3	PGW	C	503	-	15,15,50	0.21	0	14,14,56	1.31	0
3	PGW	D	507	-	13,13,50	0.19	0	12,12,56	1.24	0
3	PGW	A	508	-	15,15,50	0.24	0	14,14,56	1.35	0
3	PGW	B	503	-	15,15,50	0.21	0	14,14,56	1.31	0
3	PGW	B	504	-	7,7,50	0.19	0	6,6,56	0.55	0
3	PGW	B	502	-	9,9,50	0.13	0	8,8,56	1.09	0
3	PGW	A	505	-	8,8,50	0.16	0	7,7,56	1.22	0
3	PGW	B	501	-	42,42,50	0.90	3 (7%)	44,48,56	1.03	2 (4%)
3	PGW	A	503	-	15,15,50	0.21	0	14,14,56	1.31	0
3	PGW	C	501	-	42,42,50	0.90	3 (7%)	44,48,56	1.03	2 (4%)
3	PGW	B	505	-	8,8,50	0.16	0	7,7,56	1.22	0
3	PGW	C	505	-	8,8,50	0.16	0	7,7,56	1.22	0
3	PGW	C	504	-	7,7,50	0.19	0	6,6,56	0.55	0
3	PGW	C	507	-	13,13,50	0.19	0	12,12,56	1.24	0
3	PGW	B	507	-	13,13,50	0.19	0	12,12,56	1.24	0

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
3	PGW	A	504	-	-	2/5/5/55	-
3	PGW	D	501	-	-	26/47/47/55	-
3	PGW	B	506	-	-	9/29/29/55	-
3	PGW	C	506	-	-	9/29/29/55	-
3	PGW	D	504	-	-	2/5/5/55	-
3	PGW	D	508	-	-	2/13/13/55	-
3	PGW	D	505	-	-	2/6/6/55	-
3	PGW	D	506	-	-	9/29/29/55	-
3	PGW	C	508	-	-	2/13/13/55	-
3	PGW	C	502	-	-	0/7/7/55	-
2	CMP	C	500	-	-	0/0/31/31	0/4/4/4
2	CMP	B	500	-	-	0/0/31/31	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PGW	A	506	-	-	9/29/29/55	-
3	PGW	D	503	-	-	5/13/13/55	-
3	PGW	B	508	-	-	2/13/13/55	-
3	PGW	D	502	-	-	0/7/7/55	-
3	PGW	A	502	-	-	0/7/7/55	-
3	PGW	A	507	-	-	6/11/11/55	-
2	CMP	A	500	-	-	0/0/31/31	0/4/4/4
2	CMP	D	500	-	-	0/0/31/31	0/4/4/4
3	PGW	A	501	-	-	26/47/47/55	-
3	PGW	C	503	-	-	5/13/13/55	-
3	PGW	D	507	-	-	6/11/11/55	-
3	PGW	A	508	-	-	2/13/13/55	-
3	PGW	B	503	-	-	5/13/13/55	-
3	PGW	B	504	-	-	2/5/5/55	-
3	PGW	B	502	-	-	0/7/7/55	-
3	PGW	A	505	-	-	2/6/6/55	-
3	PGW	B	501	-	-	26/47/47/55	-
3	PGW	A	503	-	-	5/13/13/55	-
3	PGW	C	501	-	-	26/47/47/55	-
3	PGW	B	505	-	-	2/6/6/55	-
3	PGW	C	505	-	-	2/6/6/55	-
3	PGW	C	504	-	-	2/5/5/55	-
3	PGW	C	507	-	-	6/11/11/55	-
3	PGW	B	507	-	-	6/11/11/55	-

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	CMP	P-O5'	2.91	1.61	1.57
2	B	500	CMP	P-O5'	2.91	1.61	1.57
2	C	500	CMP	P-O5'	2.91	1.61	1.57
2	D	500	CMP	P-O5'	2.91	1.61	1.57
3	A	506	PGW	O01-C02	-2.57	1.42	1.47
3	B	506	PGW	O01-C02	-2.57	1.42	1.47
3	C	506	PGW	O01-C02	-2.57	1.42	1.47
3	D	506	PGW	O01-C02	-2.57	1.42	1.47
3	A	501	PGW	P-O12	2.57	1.69	1.59
3	B	501	PGW	P-O12	2.57	1.69	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	501	PGW	P-O12	2.57	1.69	1.59
3	D	501	PGW	P-O12	2.57	1.69	1.59
2	A	500	CMP	C5-C4	2.33	1.47	1.40
2	B	500	CMP	C5-C4	2.33	1.47	1.40
2	C	500	CMP	C5-C4	2.33	1.47	1.40
2	D	500	CMP	C5-C4	2.33	1.47	1.40
2	A	500	CMP	O3'-C3'	-2.20	1.41	1.44
2	B	500	CMP	O3'-C3'	-2.20	1.41	1.44
2	C	500	CMP	O3'-C3'	-2.20	1.41	1.44
2	D	500	CMP	O3'-C3'	-2.20	1.41	1.44
3	A	501	PGW	O11-C03	-2.10	1.36	1.44
3	B	501	PGW	O11-C03	-2.10	1.36	1.44
3	C	501	PGW	O11-C03	-2.10	1.36	1.44
3	D	501	PGW	O11-C03	-2.10	1.36	1.44
2	A	500	CMP	O5'-C5'	-2.08	1.43	1.46
2	B	500	CMP	O5'-C5'	-2.08	1.43	1.46
2	C	500	CMP	O5'-C5'	-2.08	1.43	1.46
2	D	500	CMP	O5'-C5'	-2.08	1.43	1.46
3	A	501	PGW	P-O11	2.00	1.67	1.59
3	B	501	PGW	P-O11	2.00	1.67	1.59
3	C	501	PGW	P-O11	2.00	1.67	1.59
3	D	501	PGW	P-O11	2.00	1.67	1.59

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	CMP	O3'-C3'-C4'	-8.07	104.61	110.71
2	B	500	CMP	O3'-C3'-C4'	-8.07	104.61	110.71
2	C	500	CMP	O3'-C3'-C4'	-8.07	104.61	110.71
2	D	500	CMP	O3'-C3'-C4'	-8.07	104.61	110.71
2	A	500	CMP	N3-C2-N1	-3.22	123.64	128.68
2	B	500	CMP	N3-C2-N1	-3.22	123.64	128.68
2	C	500	CMP	N3-C2-N1	-3.22	123.64	128.68
2	D	500	CMP	N3-C2-N1	-3.22	123.64	128.68
2	A	500	CMP	O2P-P-O1P	3.16	118.63	108.73
2	B	500	CMP	O2P-P-O1P	3.16	118.63	108.73
2	C	500	CMP	O2P-P-O1P	3.16	118.63	108.73
2	D	500	CMP	O2P-P-O1P	3.16	118.63	108.73
3	A	501	PGW	O13-P-O14	2.91	126.62	112.24
3	B	501	PGW	O13-P-O14	2.91	126.62	112.24
3	C	501	PGW	O13-P-O14	2.91	126.62	112.24
3	D	501	PGW	O13-P-O14	2.91	126.62	112.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	500	CMP	C4-C5-N7	-2.63	106.66	109.40
2	B	500	CMP	C4-C5-N7	-2.63	106.66	109.40
2	C	500	CMP	C4-C5-N7	-2.63	106.66	109.40
2	D	500	CMP	C4-C5-N7	-2.63	106.66	109.40
3	A	506	PGW	O01-C1-C2	2.27	116.40	111.50
3	B	506	PGW	O01-C1-C2	2.27	116.40	111.50
3	C	506	PGW	O01-C1-C2	2.27	116.40	111.50
3	D	506	PGW	O01-C1-C2	2.27	116.40	111.50
3	A	501	PGW	O01-C1-C2	2.02	115.86	111.50
3	B	501	PGW	O01-C1-C2	2.02	115.86	111.50
3	C	501	PGW	O01-C1-C2	2.02	115.86	111.50
3	D	501	PGW	O01-C1-C2	2.02	115.86	111.50

There are no chirality outliers.

All (208) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	PGW	C03-O11-P-O14
3	A	501	PGW	C04-O12-P-O14
3	A	501	PGW	C2-C1-O01-C02
3	B	501	PGW	C03-O11-P-O14
3	B	501	PGW	C04-O12-P-O14
3	B	501	PGW	C2-C1-O01-C02
3	C	501	PGW	C03-O11-P-O14
3	C	501	PGW	C04-O12-P-O14
3	C	501	PGW	C2-C1-O01-C02
3	D	501	PGW	C03-O11-P-O14
3	D	501	PGW	C04-O12-P-O14
3	D	501	PGW	C2-C1-O01-C02
3	A	501	PGW	O04-C19-O03-C01
3	B	501	PGW	O04-C19-O03-C01
3	C	501	PGW	O04-C19-O03-C01
3	D	501	PGW	O04-C19-O03-C01
3	A	501	PGW	C20-C19-O03-C01
3	B	501	PGW	C20-C19-O03-C01
3	C	501	PGW	C20-C19-O03-C01
3	D	501	PGW	C20-C19-O03-C01
3	A	501	PGW	O02-C1-O01-C02
3	B	501	PGW	O02-C1-O01-C02
3	C	501	PGW	O02-C1-O01-C02
3	D	501	PGW	O02-C1-O01-C02
3	A	501	PGW	C1-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
3	A	506	PGW	C1-C2-C3-C4
3	B	501	PGW	C1-C2-C3-C4
3	B	506	PGW	C1-C2-C3-C4
3	C	501	PGW	C1-C2-C3-C4
3	C	506	PGW	C1-C2-C3-C4
3	D	501	PGW	C1-C2-C3-C4
3	D	506	PGW	C1-C2-C3-C4
3	A	501	PGW	C03-O11-P-O12
3	A	501	PGW	C04-O12-P-O11
3	B	501	PGW	C03-O11-P-O12
3	B	501	PGW	C04-O12-P-O11
3	C	501	PGW	C03-O11-P-O12
3	C	501	PGW	C04-O12-P-O11
3	D	501	PGW	C03-O11-P-O12
3	D	501	PGW	C04-O12-P-O11
3	A	501	PGW	C23-C24-C25-C26
3	B	501	PGW	C23-C24-C25-C26
3	C	501	PGW	C23-C24-C25-C26
3	D	501	PGW	C23-C24-C25-C26
3	A	501	PGW	C20-C21-C22-C23
3	A	503	PGW	C27-C15-C16-C17
3	B	501	PGW	C20-C21-C22-C23
3	B	503	PGW	C27-C15-C16-C17
3	C	501	PGW	C20-C21-C22-C23
3	C	503	PGW	C27-C15-C16-C17
3	D	501	PGW	C20-C21-C22-C23
3	D	503	PGW	C27-C15-C16-C17
3	A	503	PGW	C23-C24-C25-C26
3	B	503	PGW	C23-C24-C25-C26
3	C	503	PGW	C23-C24-C25-C26
3	D	503	PGW	C23-C24-C25-C26
3	A	507	PGW	C25-C26-C27-C15
3	B	507	PGW	C25-C26-C27-C15
3	C	507	PGW	C25-C26-C27-C15
3	D	507	PGW	C25-C26-C27-C15
3	A	501	PGW	C21-C22-C23-C24
3	B	501	PGW	C21-C22-C23-C24
3	C	501	PGW	C21-C22-C23-C24
3	D	501	PGW	C21-C22-C23-C24
3	A	501	PGW	C24-C25-C26-C27
3	B	501	PGW	C24-C25-C26-C27
3	C	501	PGW	C24-C25-C26-C27

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Mol	Chain	Res	Type	Atoms
3	D	501	PGW	C24-C25-C26-C27
3	A	508	PGW	C17-C18-C28-C30
3	B	508	PGW	C17-C18-C28-C30
3	C	508	PGW	C17-C18-C28-C30
3	D	508	PGW	C17-C18-C28-C30
3	A	501	PGW	C2-C3-C4-C5
3	B	501	PGW	C2-C3-C4-C5
3	C	501	PGW	C2-C3-C4-C5
3	D	501	PGW	C2-C3-C4-C5
3	A	507	PGW	C16-C15-C27-C26
3	B	507	PGW	C16-C15-C27-C26
3	C	507	PGW	C16-C15-C27-C26
3	D	507	PGW	C16-C15-C27-C26
3	A	501	PGW	C07-C06-C10-C9
3	B	501	PGW	C07-C06-C10-C9
3	C	501	PGW	C07-C06-C10-C9
3	D	501	PGW	C07-C06-C10-C9
3	A	506	PGW	C20-C19-O03-C01
3	B	506	PGW	C20-C19-O03-C01
3	C	506	PGW	C20-C19-O03-C01
3	D	506	PGW	C20-C19-O03-C01
3	A	506	PGW	C15-C16-C17-C18
3	B	506	PGW	C15-C16-C17-C18
3	C	506	PGW	C15-C16-C17-C18
3	D	506	PGW	C15-C16-C17-C18
3	A	506	PGW	C21-C22-C23-C24
3	B	506	PGW	C21-C22-C23-C24
3	C	506	PGW	C21-C22-C23-C24
3	D	506	PGW	C21-C22-C23-C24
3	A	506	PGW	C16-C17-C18-C28
3	B	506	PGW	C16-C17-C18-C28
3	C	506	PGW	C16-C17-C18-C28
3	D	506	PGW	C16-C17-C18-C28
3	A	507	PGW	C21-C22-C23-C24
3	B	507	PGW	C21-C22-C23-C24
3	C	507	PGW	C21-C22-C23-C24
3	D	507	PGW	C21-C22-C23-C24
3	A	507	PGW	C22-C23-C24-C25
3	B	507	PGW	C22-C23-C24-C25
3	C	507	PGW	C22-C23-C24-C25
3	D	507	PGW	C22-C23-C24-C25
3	A	501	PGW	C15-C16-C17-C18

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Mol	Chain	Res	Type	Atoms
3	B	501	PGW	C15-C16-C17-C18
3	C	501	PGW	C15-C16-C17-C18
3	D	501	PGW	C15-C16-C17-C18
3	A	506	PGW	O04-C19-O03-C01
3	B	506	PGW	O04-C19-O03-C01
3	C	506	PGW	O04-C19-O03-C01
3	D	506	PGW	O04-C19-O03-C01
3	A	503	PGW	C18-C28-C30-C29
3	B	503	PGW	C18-C28-C30-C29
3	C	503	PGW	C18-C28-C30-C29
3	D	503	PGW	C18-C28-C30-C29
3	A	505	PGW	C4-C5-C6-C7
3	B	505	PGW	C4-C5-C6-C7
3	C	505	PGW	C4-C5-C6-C7
3	D	505	PGW	C4-C5-C6-C7
3	A	501	PGW	C3-C4-C5-C6
3	B	501	PGW	C3-C4-C5-C6
3	C	501	PGW	C3-C4-C5-C6
3	D	501	PGW	C3-C4-C5-C6
3	A	501	PGW	O03-C01-C02-C03
3	B	501	PGW	O03-C01-C02-C03
3	C	501	PGW	O03-C01-C02-C03
3	D	501	PGW	O03-C01-C02-C03
3	A	508	PGW	C20-C21-C22-C23
3	B	508	PGW	C20-C21-C22-C23
3	C	508	PGW	C20-C21-C22-C23
3	D	508	PGW	C20-C21-C22-C23
3	A	501	PGW	O01-C02-C03-O11
3	B	501	PGW	O01-C02-C03-O11
3	C	501	PGW	O01-C02-C03-O11
3	D	501	PGW	O01-C02-C03-O11
3	A	501	PGW	O03-C01-C02-O01
3	B	501	PGW	O03-C01-C02-O01
3	C	501	PGW	O03-C01-C02-O01
3	D	501	PGW	O03-C01-C02-O01
3	A	503	PGW	C15-C16-C17-C18
3	B	503	PGW	C15-C16-C17-C18
3	C	503	PGW	C15-C16-C17-C18
3	D	503	PGW	C15-C16-C17-C18
3	A	506	PGW	C27-C15-C16-C17
3	A	507	PGW	C27-C15-C16-C17
3	B	506	PGW	C27-C15-C16-C17

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Mol	Chain	Res	Type	Atoms
3	B	507	PGW	C27-C15-C16-C17
3	C	506	PGW	C27-C15-C16-C17
3	C	507	PGW	C27-C15-C16-C17
3	D	506	PGW	C27-C15-C16-C17
3	D	507	PGW	C27-C15-C16-C17
3	A	501	PGW	C25-C26-C27-C15
3	A	505	PGW	C2-C3-C4-C5
3	B	501	PGW	C25-C26-C27-C15
3	B	505	PGW	C2-C3-C4-C5
3	C	501	PGW	C25-C26-C27-C15
3	C	505	PGW	C2-C3-C4-C5
3	D	501	PGW	C25-C26-C27-C15
3	D	505	PGW	C2-C3-C4-C5
3	A	501	PGW	C03-O11-P-O13
3	A	501	PGW	C04-O12-P-O13
3	B	501	PGW	C03-O11-P-O13
3	B	501	PGW	C04-O12-P-O13
3	C	501	PGW	C03-O11-P-O13
3	C	501	PGW	C04-O12-P-O13
3	D	501	PGW	C03-O11-P-O13
3	D	501	PGW	C04-O12-P-O13
3	A	501	PGW	C01-C02-C03-O11
3	B	501	PGW	C01-C02-C03-O11
3	C	501	PGW	C01-C02-C03-O11
3	D	501	PGW	C01-C02-C03-O11
3	A	504	PGW	C2-C3-C4-C5
3	B	504	PGW	C2-C3-C4-C5
3	C	504	PGW	C2-C3-C4-C5
3	D	504	PGW	C2-C3-C4-C5
3	A	501	PGW	C03-C02-O01-C1
3	B	501	PGW	C03-C02-O01-C1
3	C	501	PGW	C03-C02-O01-C1
3	D	501	PGW	C03-C02-O01-C1
3	A	503	PGW	C17-C18-C28-C30
3	B	503	PGW	C17-C18-C28-C30
3	C	503	PGW	C17-C18-C28-C30
3	D	503	PGW	C17-C18-C28-C30
3	A	506	PGW	O02-C1-O01-C02
3	B	506	PGW	O02-C1-O01-C02
3	C	506	PGW	O02-C1-O01-C02
3	D	506	PGW	O02-C1-O01-C02
3	A	504	PGW	C3-C4-C5-C6

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Mol	Chain	Res	Type	Atoms
3	B	504	PGW	C3-C4-C5-C6
3	C	504	PGW	C3-C4-C5-C6
3	D	504	PGW	C3-C4-C5-C6
3	A	507	PGW	C20-C21-C22-C23
3	B	507	PGW	C20-C21-C22-C23
3	C	507	PGW	C20-C21-C22-C23
3	D	507	PGW	C20-C21-C22-C23
3	A	501	PGW	C7-C8-C9-C10
3	B	501	PGW	C7-C8-C9-C10
3	C	501	PGW	C7-C8-C9-C10
3	D	501	PGW	C7-C8-C9-C10
3	A	506	PGW	C2-C1-O01-C02
3	B	506	PGW	C2-C1-O01-C02
3	C	506	PGW	C2-C1-O01-C02
3	D	506	PGW	C2-C1-O01-C02

There are no ring outliers.

23 monomers are involved in 51 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	501	PGW	1	0
3	B	506	PGW	7	0
3	C	506	PGW	7	0
3	D	508	PGW	1	0
3	D	506	PGW	7	0
3	C	508	PGW	1	0
2	C	500	CMP	2	0
2	B	500	CMP	2	0
3	A	506	PGW	7	0
3	D	503	PGW	1	0
3	B	508	PGW	1	0
3	A	507	PGW	1	0
2	A	500	CMP	2	0
2	D	500	CMP	2	0
3	A	501	PGW	1	0
3	C	503	PGW	1	0
3	D	507	PGW	1	0
3	A	508	PGW	1	0
3	B	503	PGW	1	0
3	B	501	PGW	1	0
3	A	503	PGW	1	0
3	C	507	PGW	1	0

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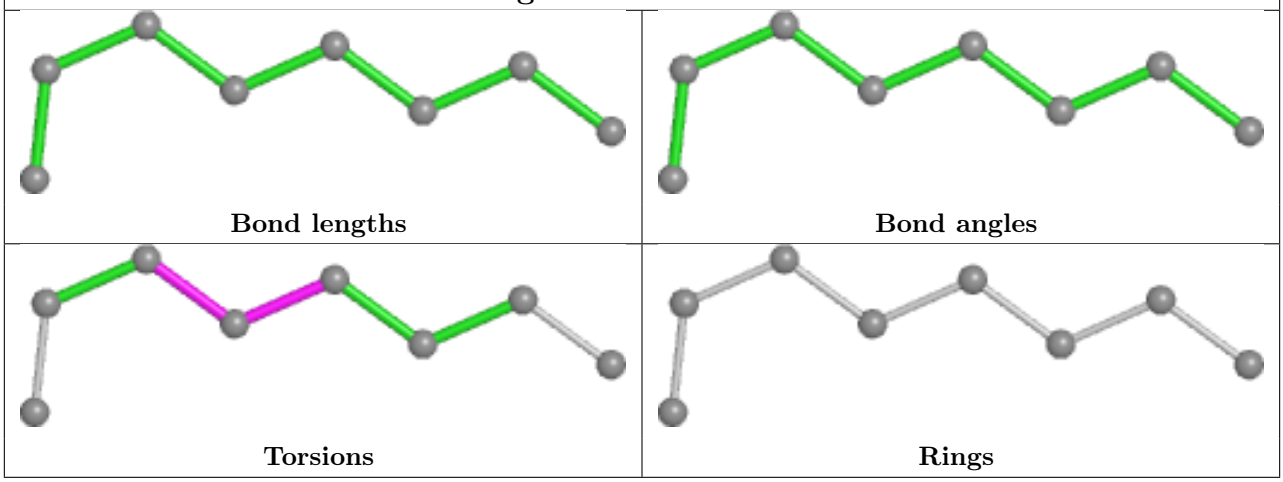
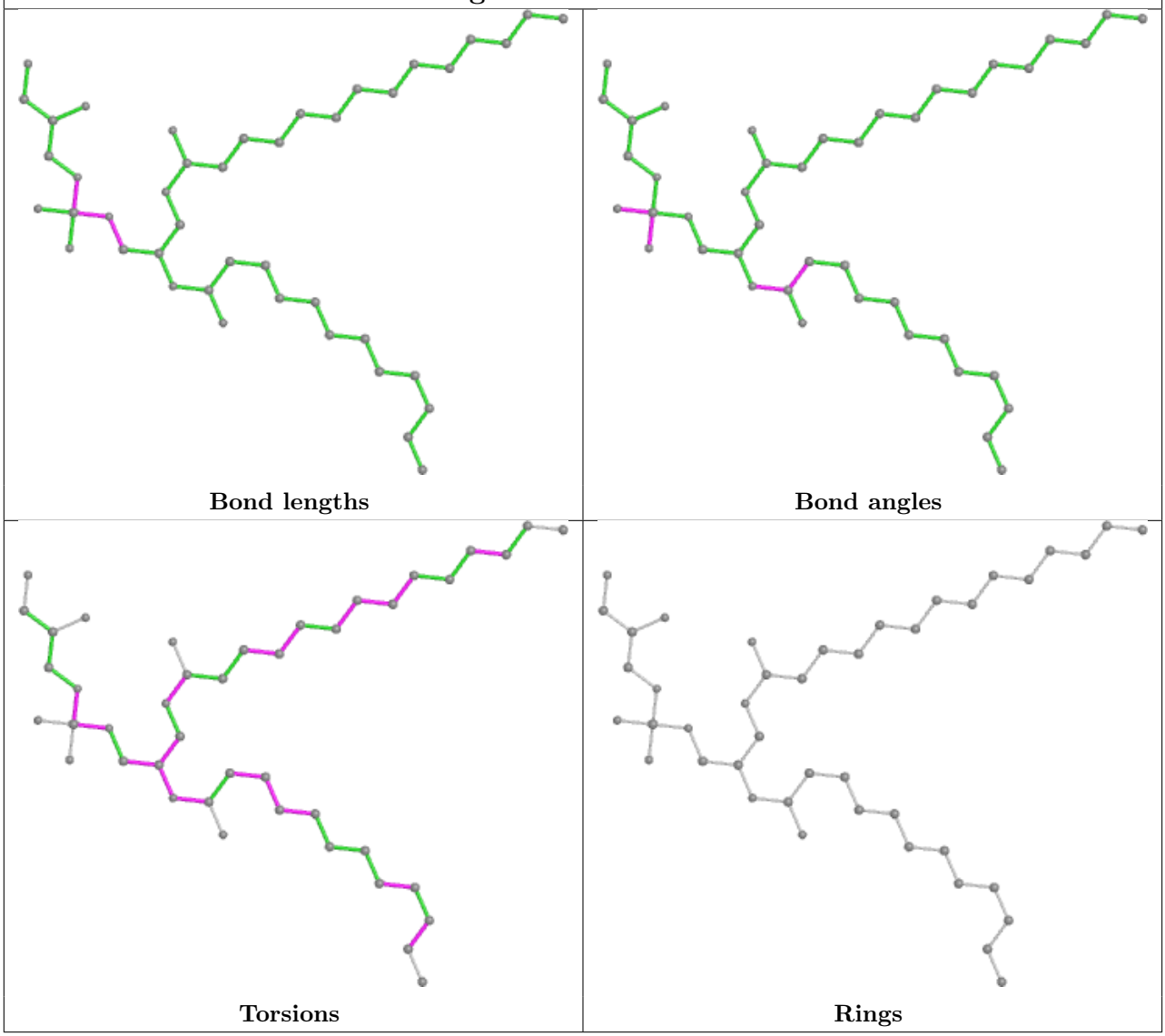
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	507	PGW	1	0

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.

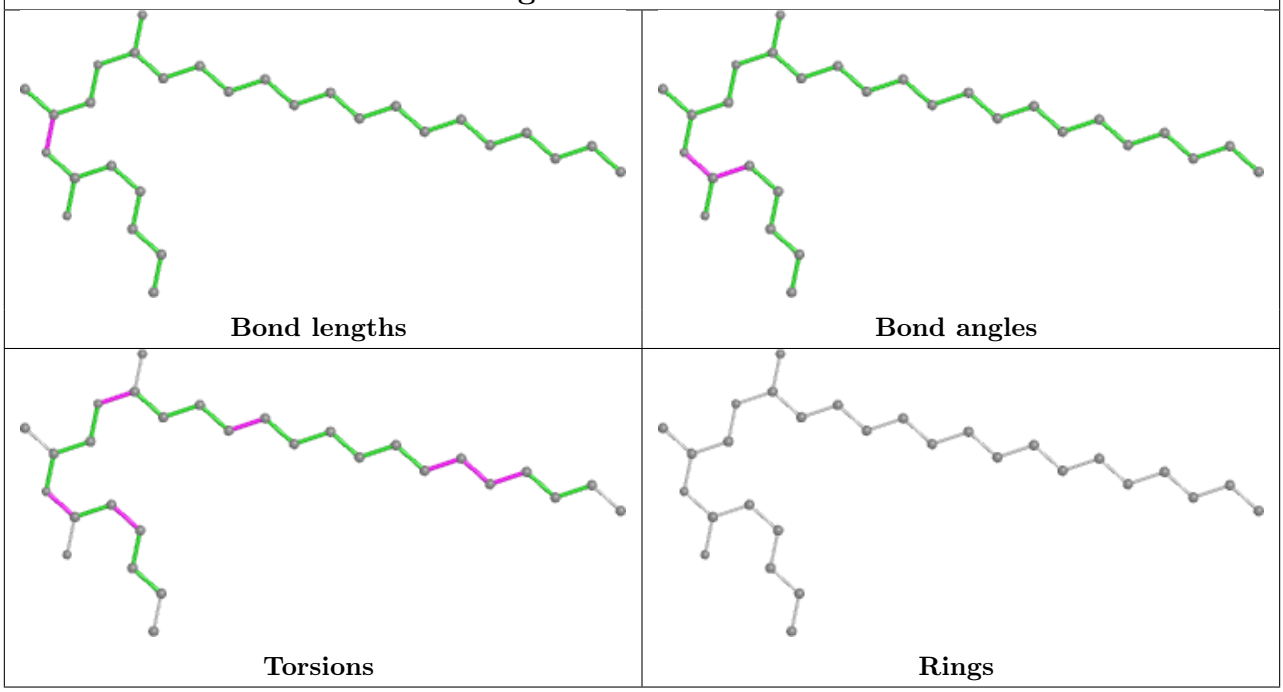
Ligand PGW A 504



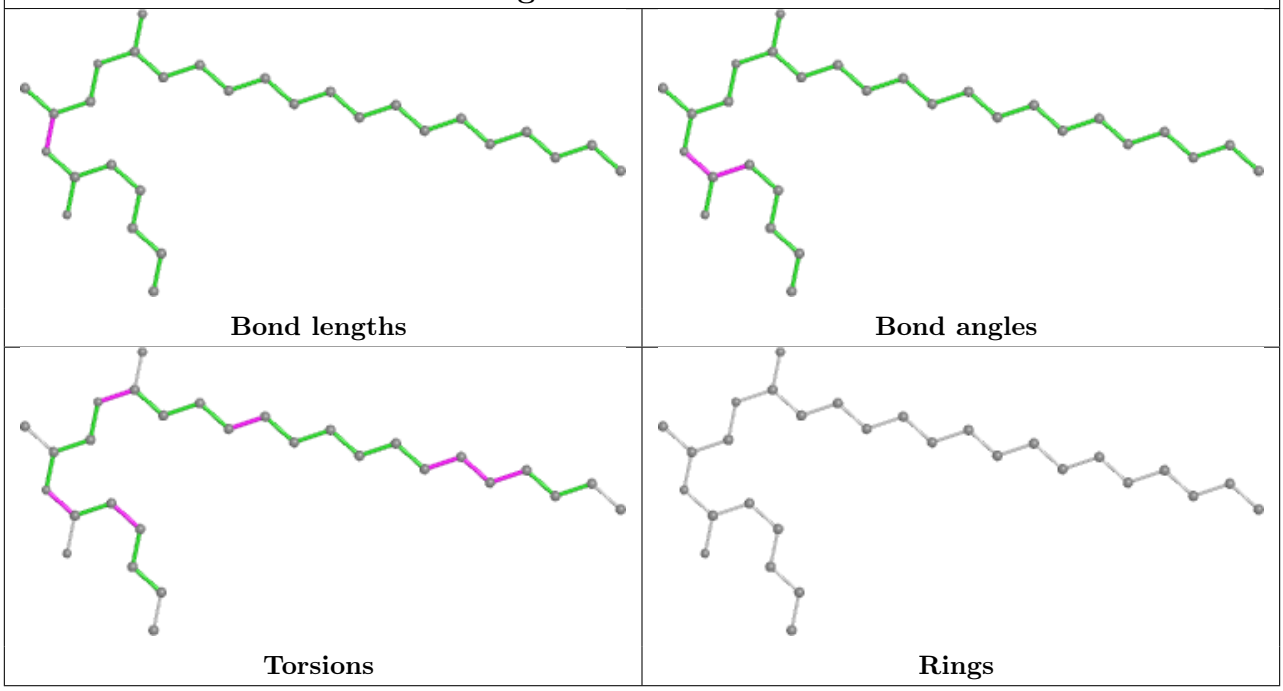
Ligand PGW D 501



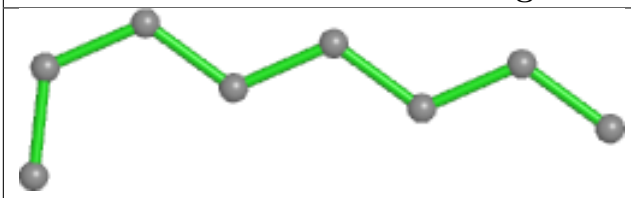
Ligand PGW B 506



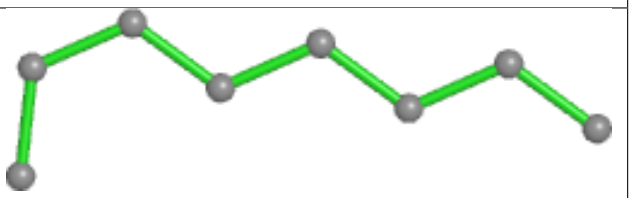
Ligand PGW C 506



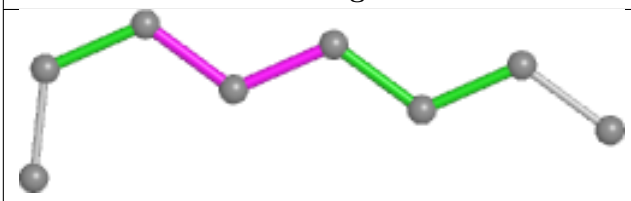
Ligand PGW D 504



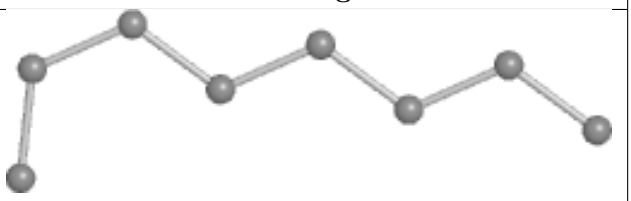
Bond lengths



Bond angles



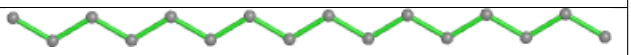
Torsions



Rings

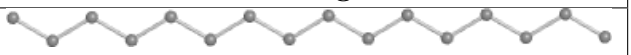
Ligand PGW D 508

Bond lengths



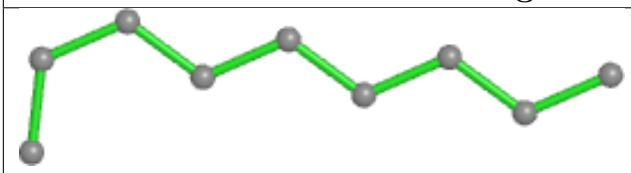
Bond angles

Torsions



Rings

Ligand PGW D 505



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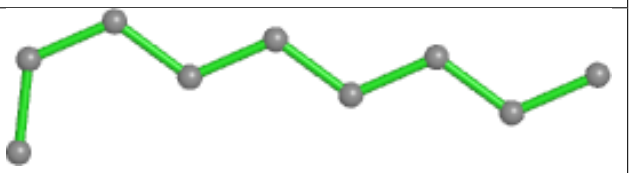
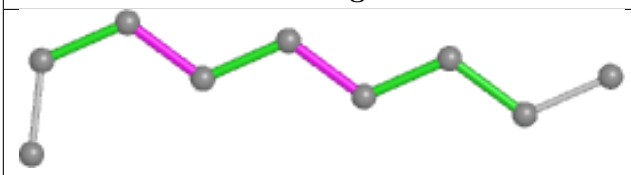


Bond angles



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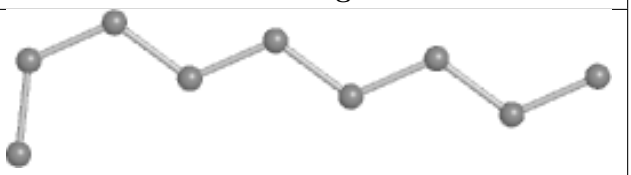

Rings

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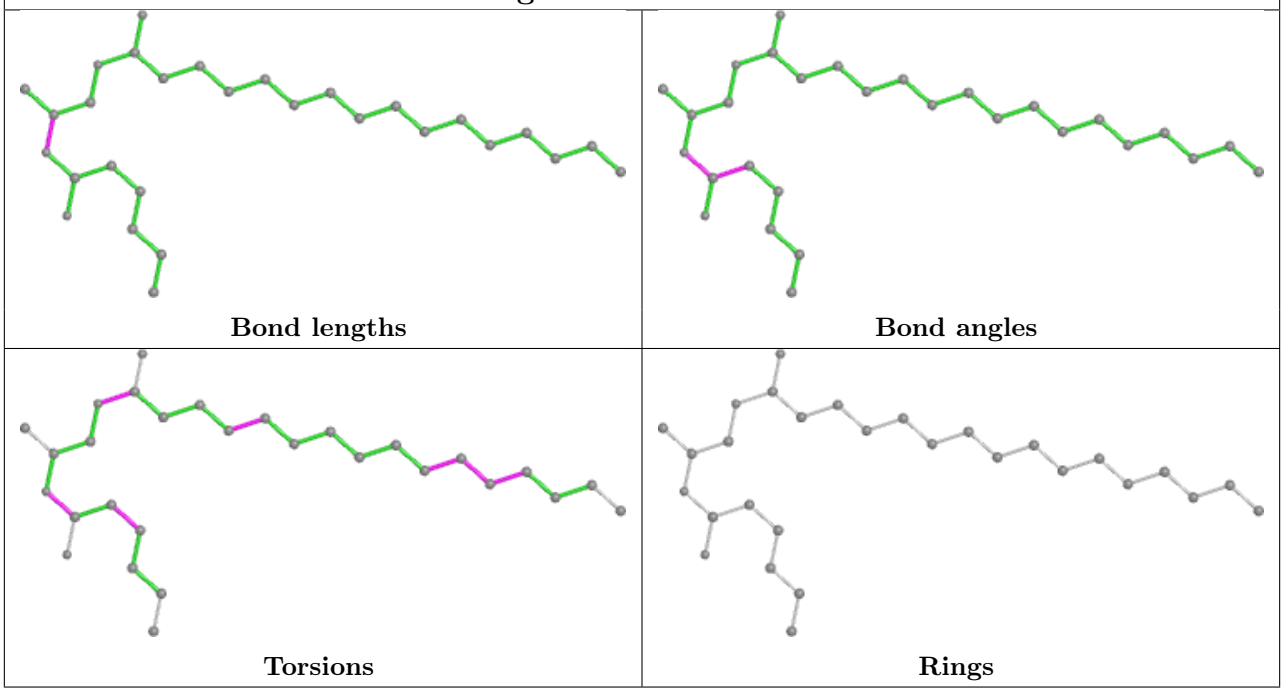
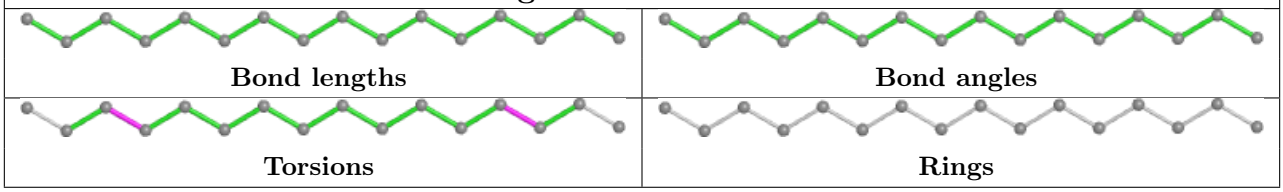
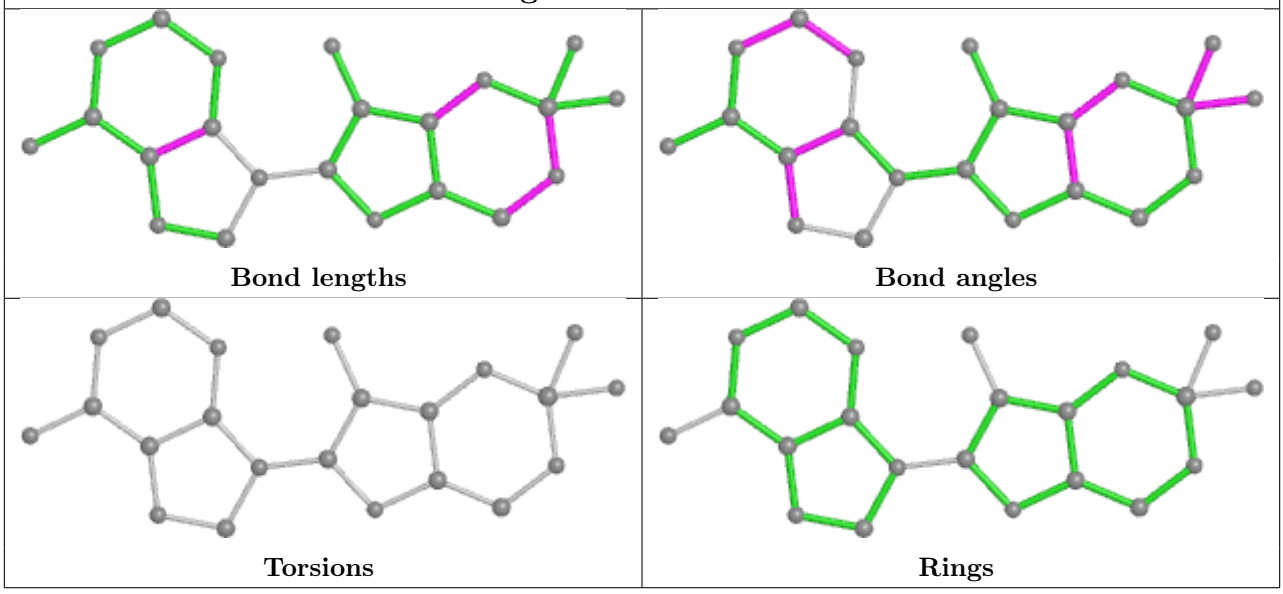


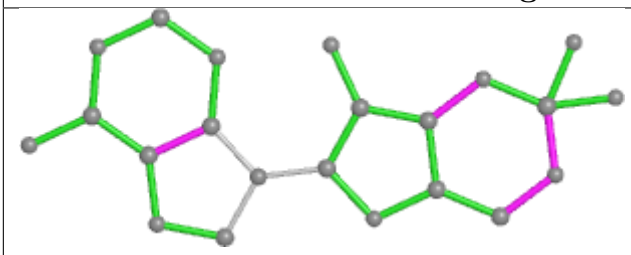
Ligand PGW C 508



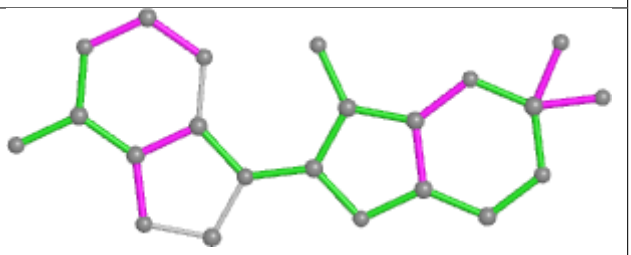
Ligand CMP C 500



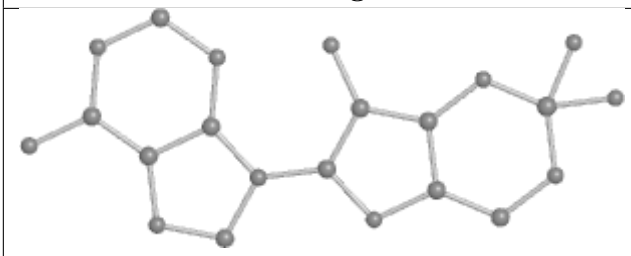
Ligand CMP B 500



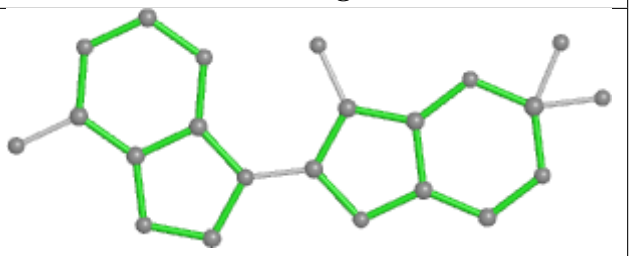
Bond lengths



Bond angles

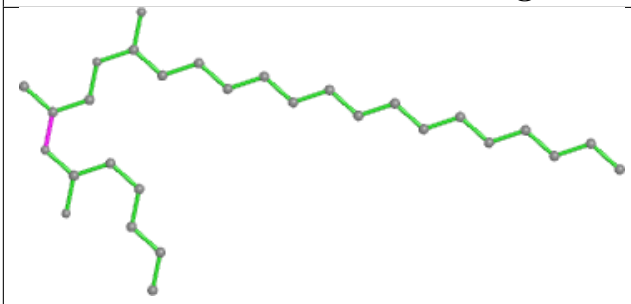


Torsions

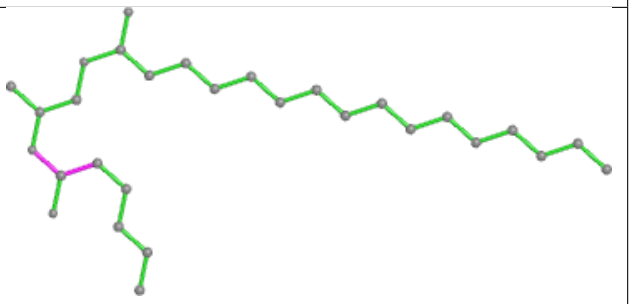


Rings

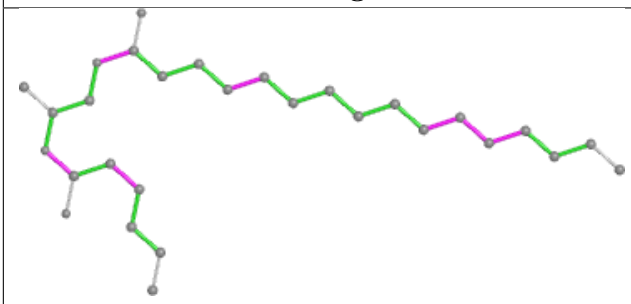
Ligand PGW A 506



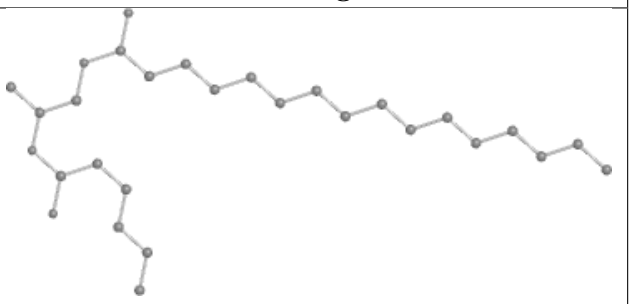
Bond lengths



Bond angles



Torsions



Rings

Ligand PGW D 503

Bond lengths

Bond angles

Torsions

Rings

Ligand PGW B 508

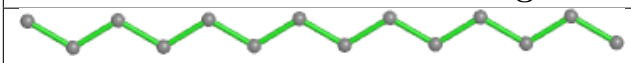
Bond lengths

Bond angles

Torsions

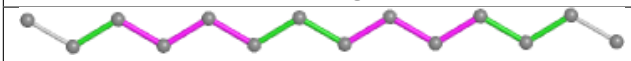
Rings

Ligand PGW A 507



Bond lengths

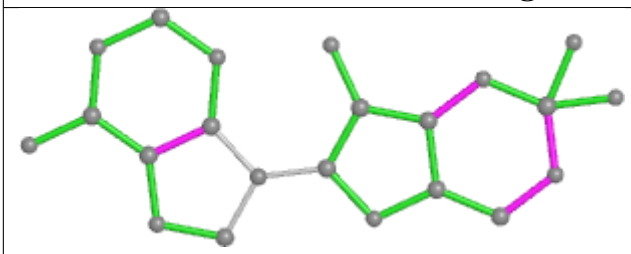
Bond angles



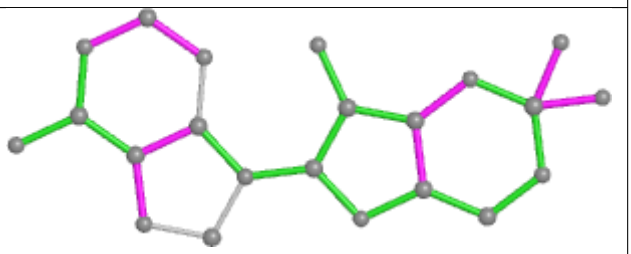
Torsions

Rings

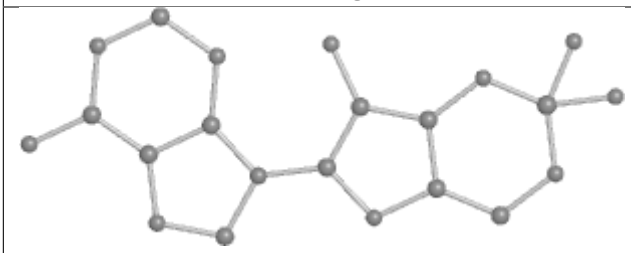
Ligand CMP A 500



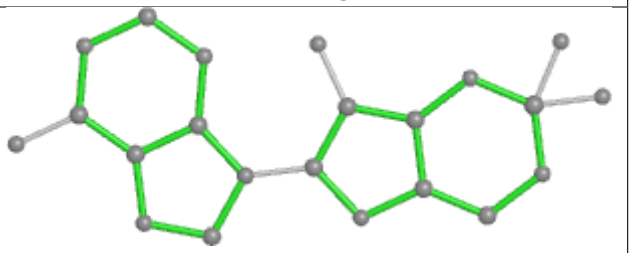
Bond lengths



Bond angles

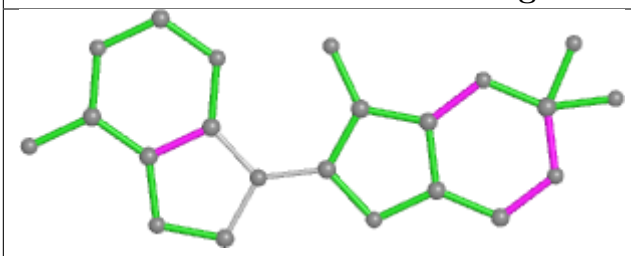


Torsions

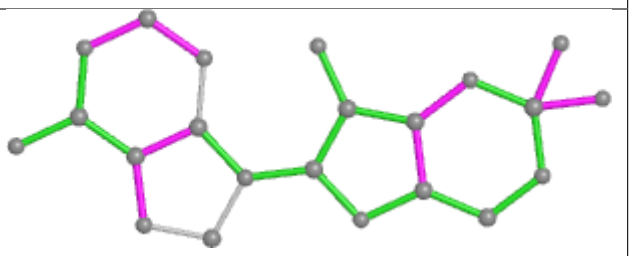


Rings

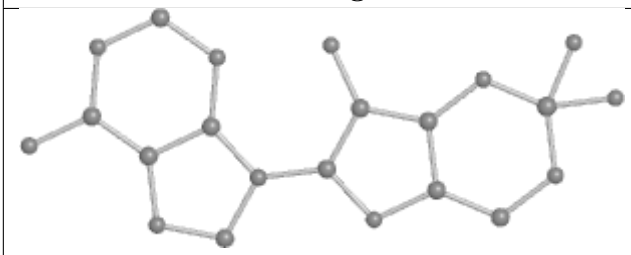
Ligand CMP D 500



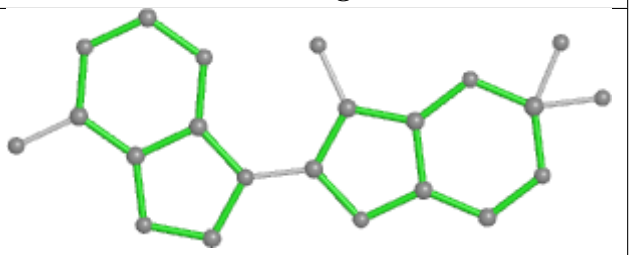
Bond lengths



Bond angles

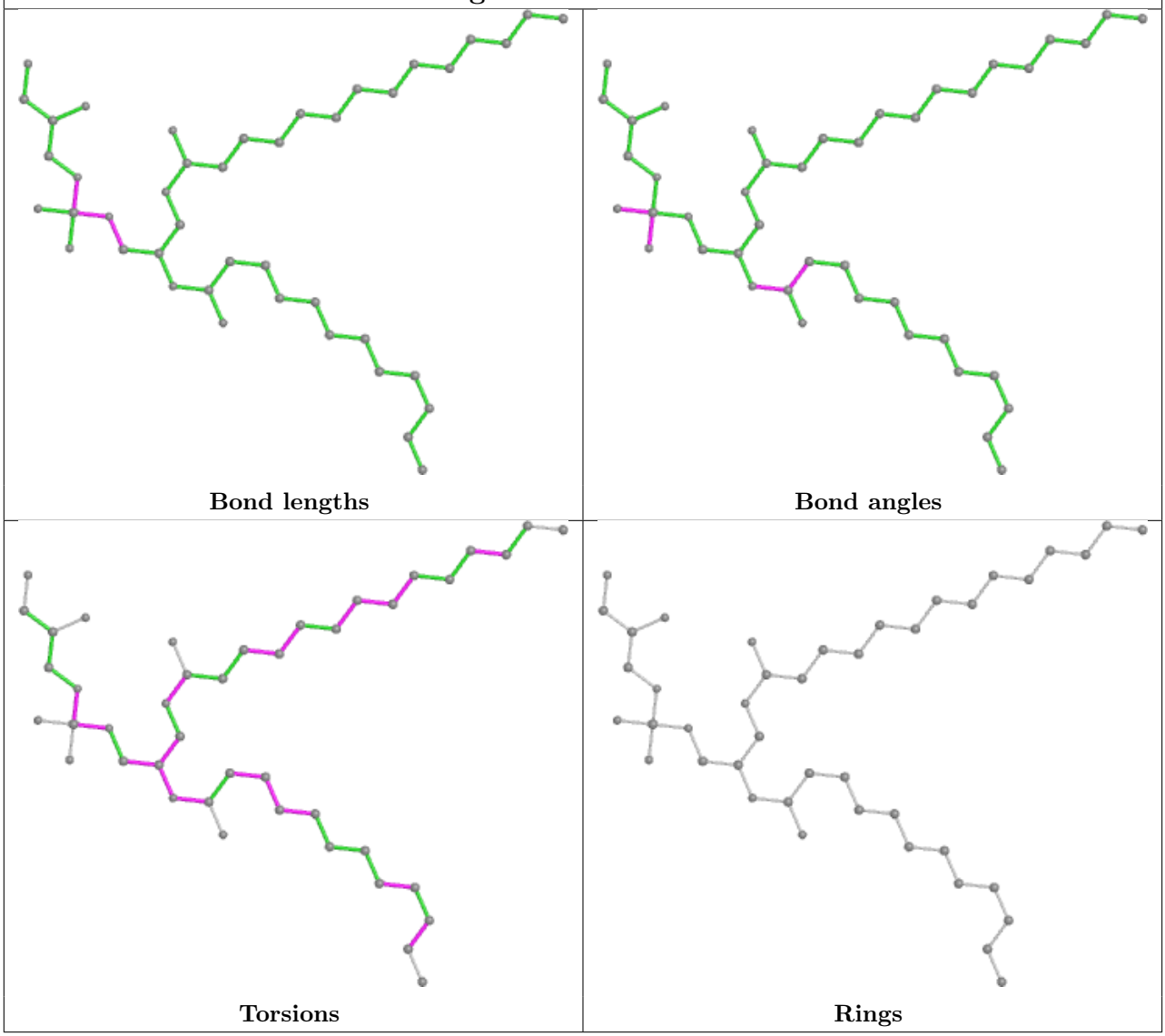


Torsions

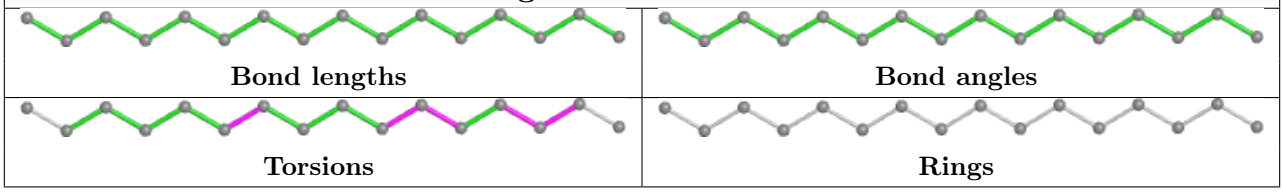


Rings

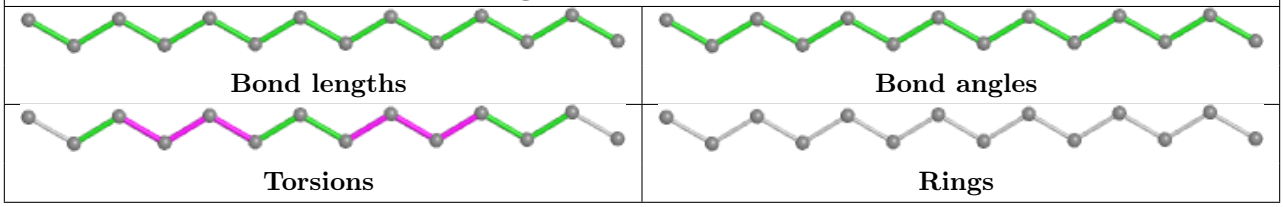
Ligand PGW A 501



Ligand PGW C 503



Ligand PGW D 507



Ligand PGW A 508

Bond lengths

Bond angles

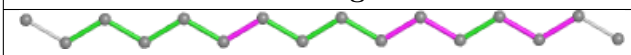
Torsions

Rings

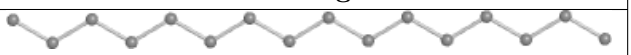
Ligand PGW B 503

Bond lengths

Bond angles

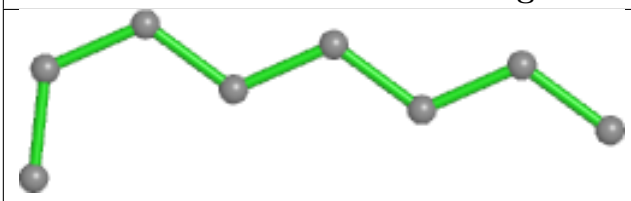


Torsions

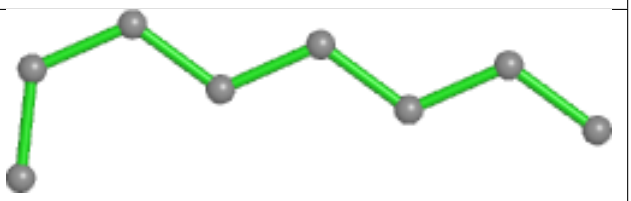


Rings

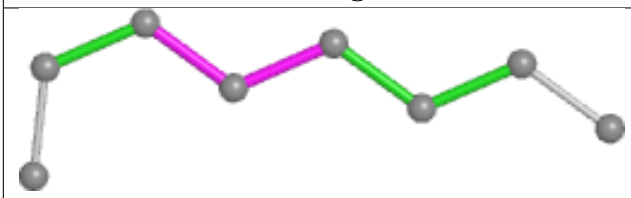
Ligand PGW B 504



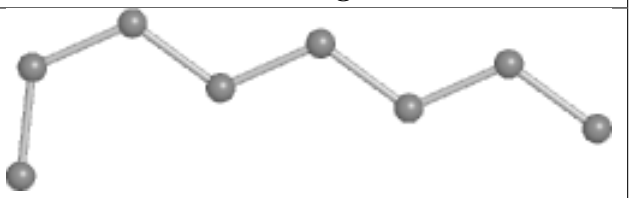
Bond lengths



Bond angles

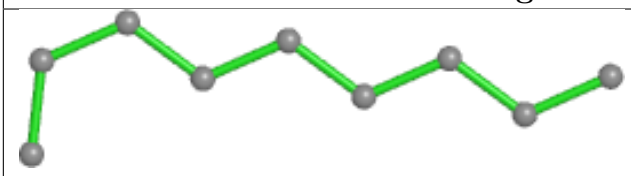


Torsions

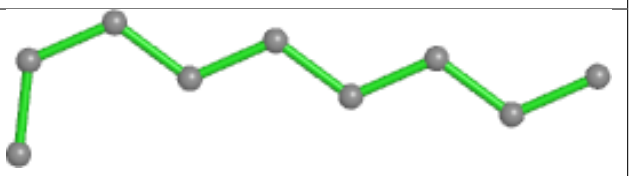


Rings

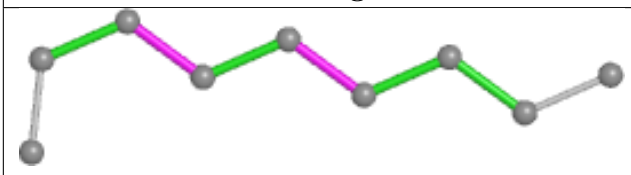
Ligand PGW A 505



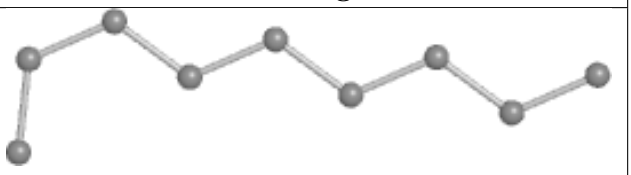
Bond lengths



Bond angles

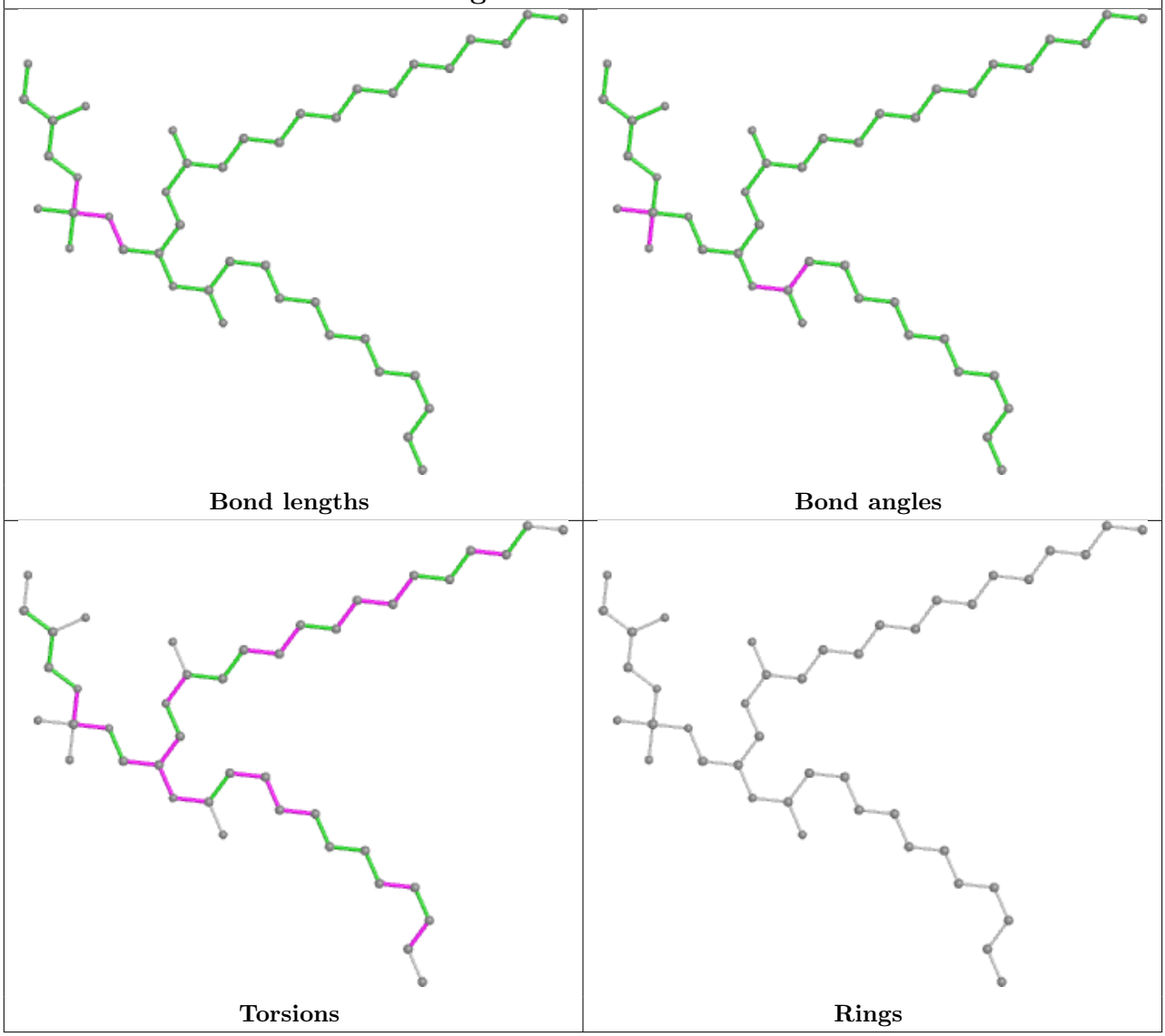


Torsions

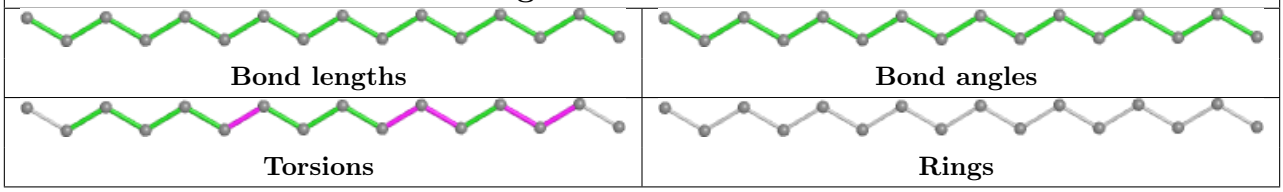


Rings

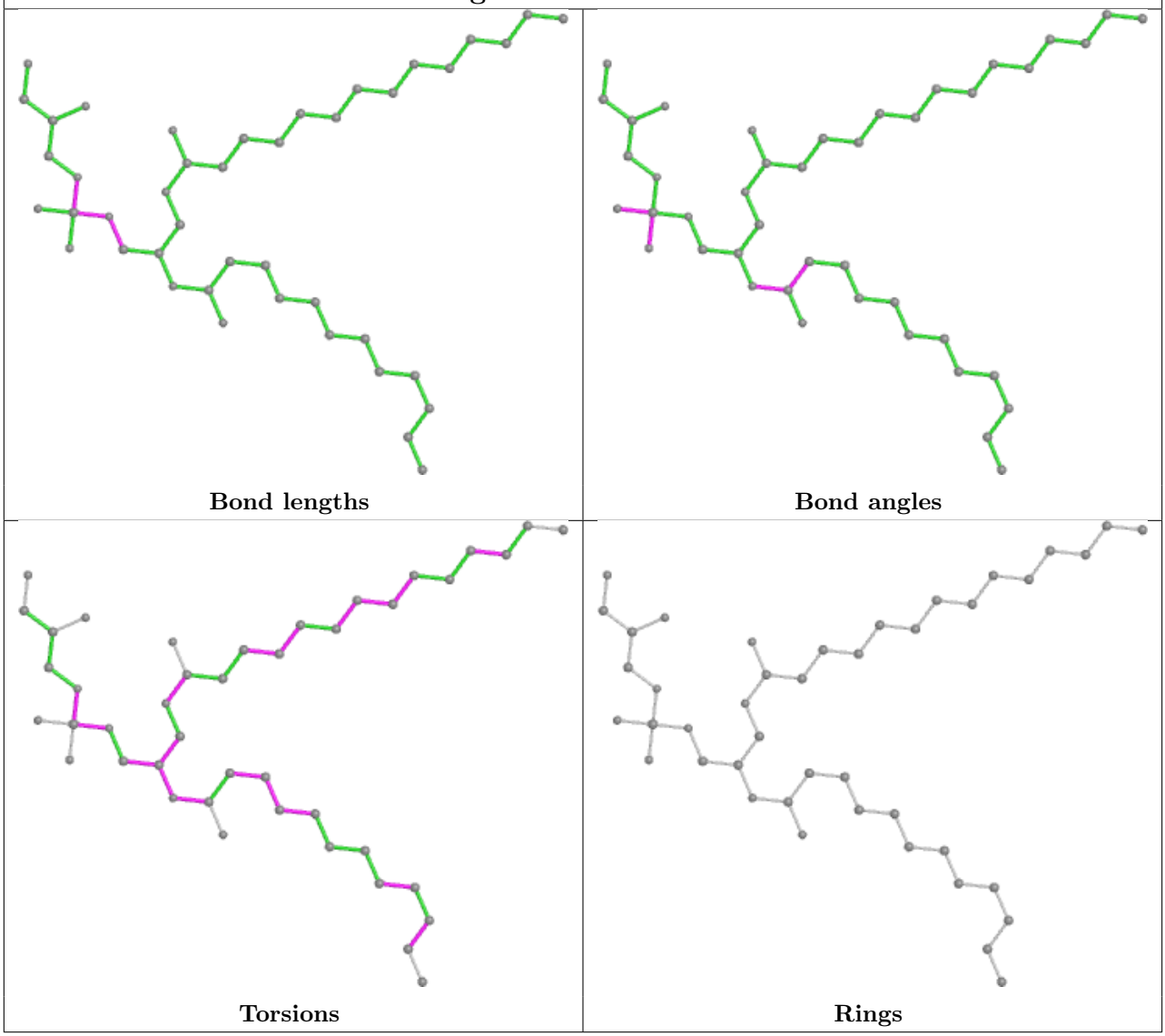
Ligand PGW B 501



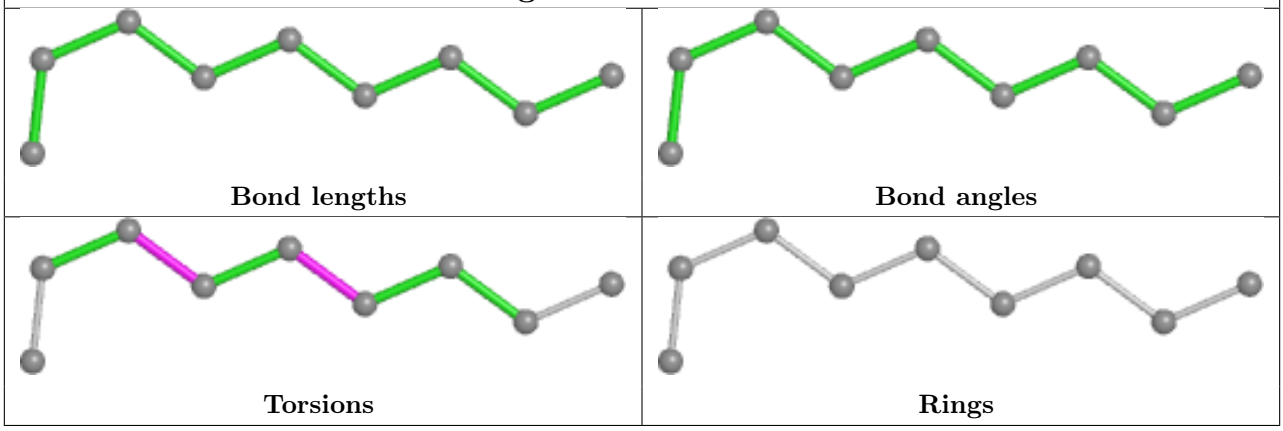
Ligand PGW A 503

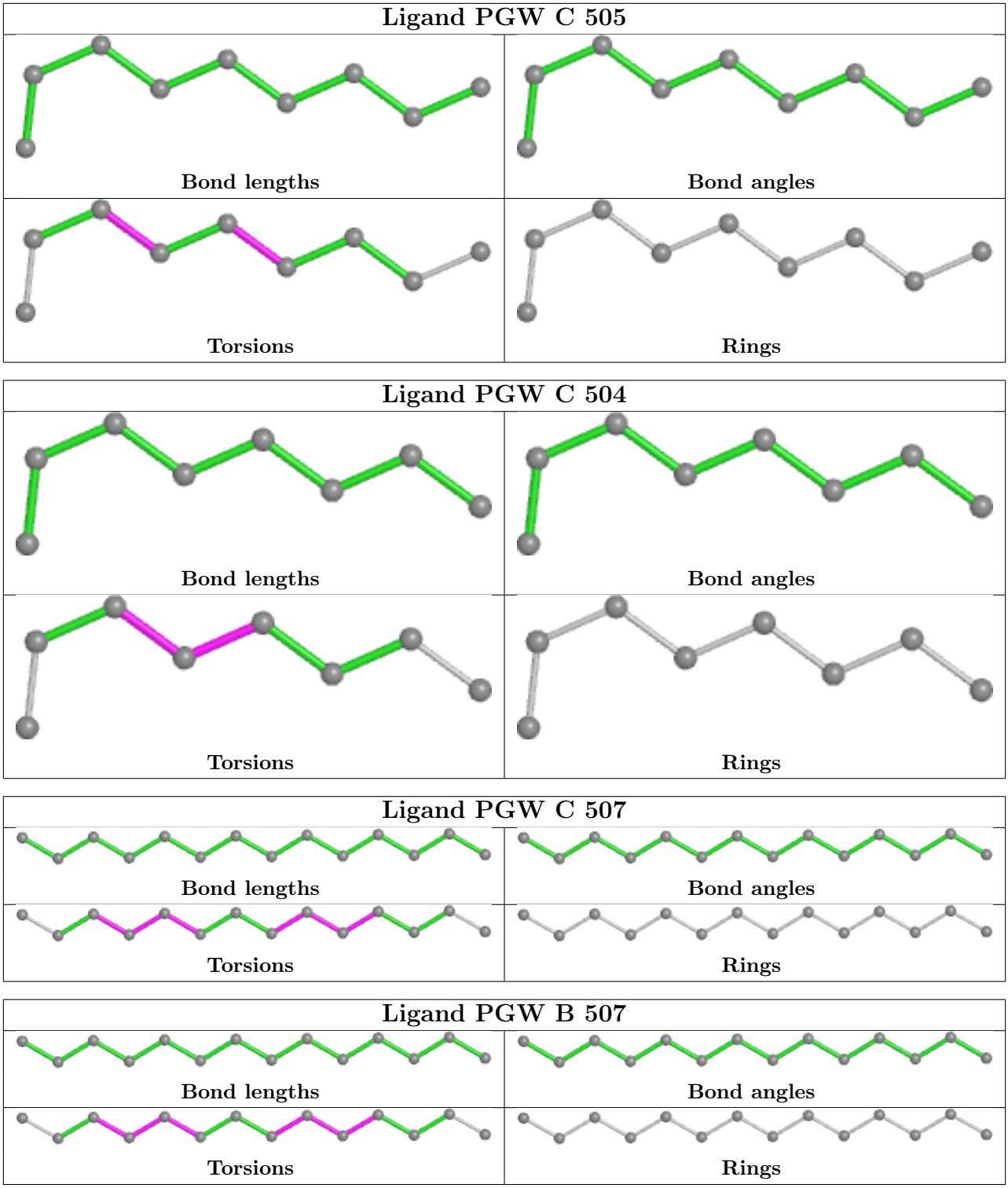


Ligand PGW C 501



Ligand PGW B 505





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.

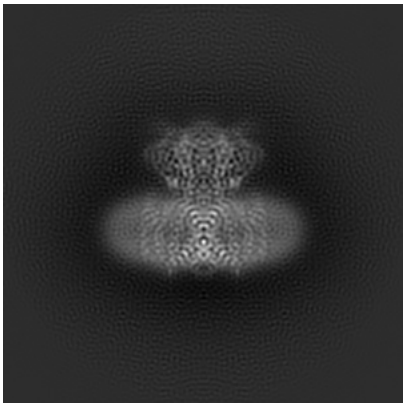
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-7484. These allow visual inspection of the internal detail of the map and identification of artifacts.

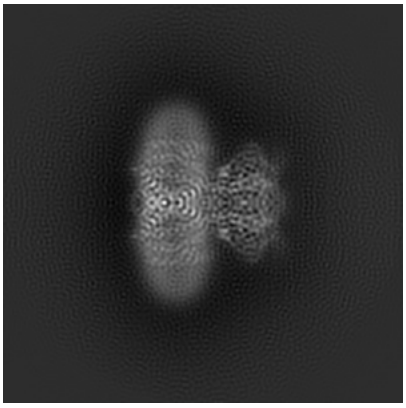
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

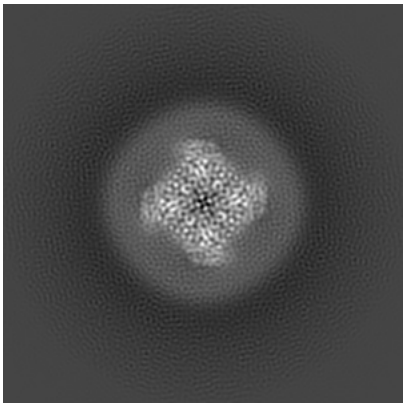
6.1.1 Primary map



X



Y

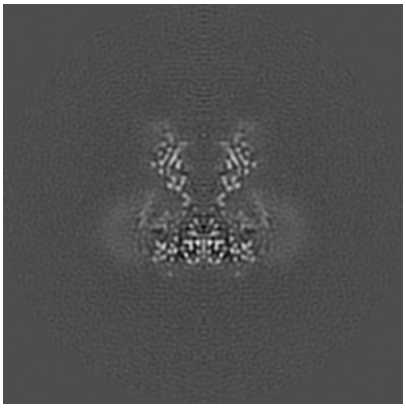


Z

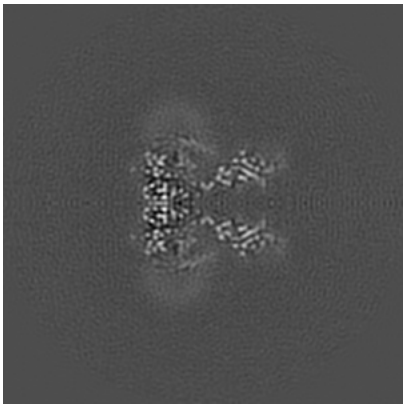
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

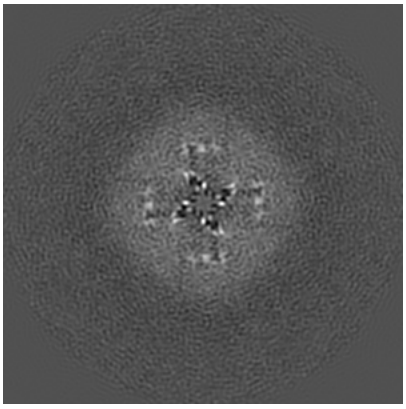
6.2.1 Primary map



X Index: 128



Y Index: 128

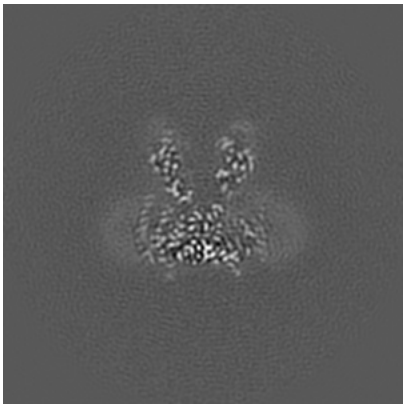


Z Index: 128

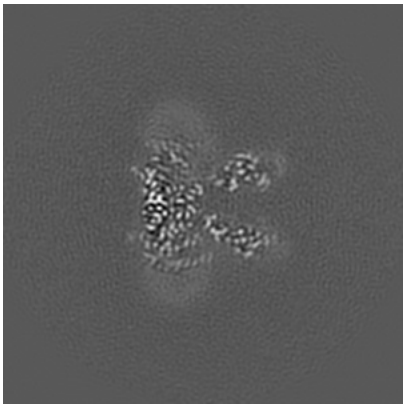
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

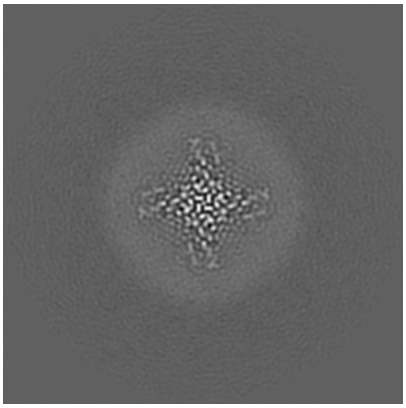
6.3.1 Primary map



X Index: 131



Y Index: 125

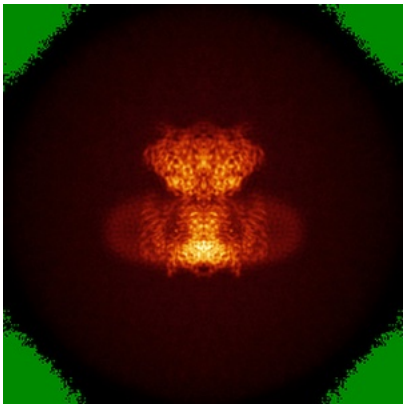


Z Index: 100

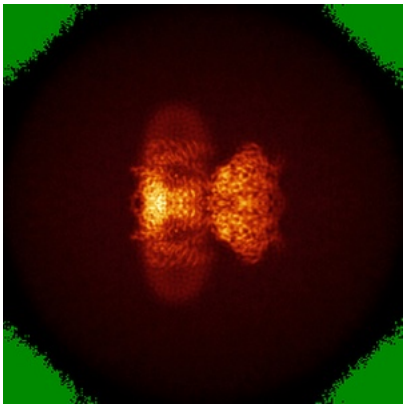
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

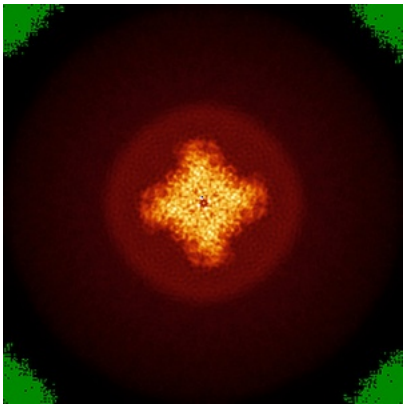
6.4.1 Primary map



X



Y



Z

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

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.06. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

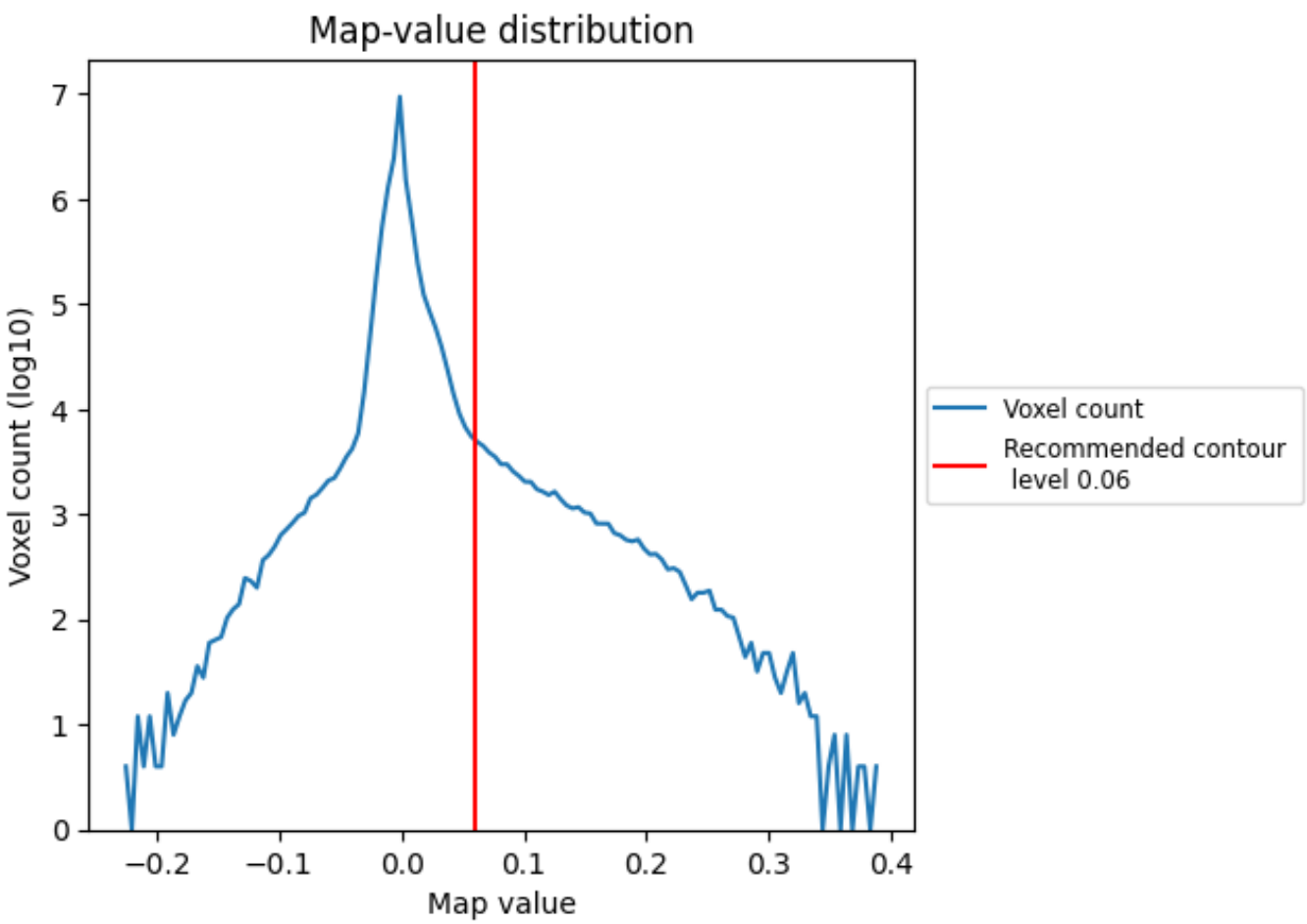
6.6 Mask visualisation (i)

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

7 Map analysis (i)

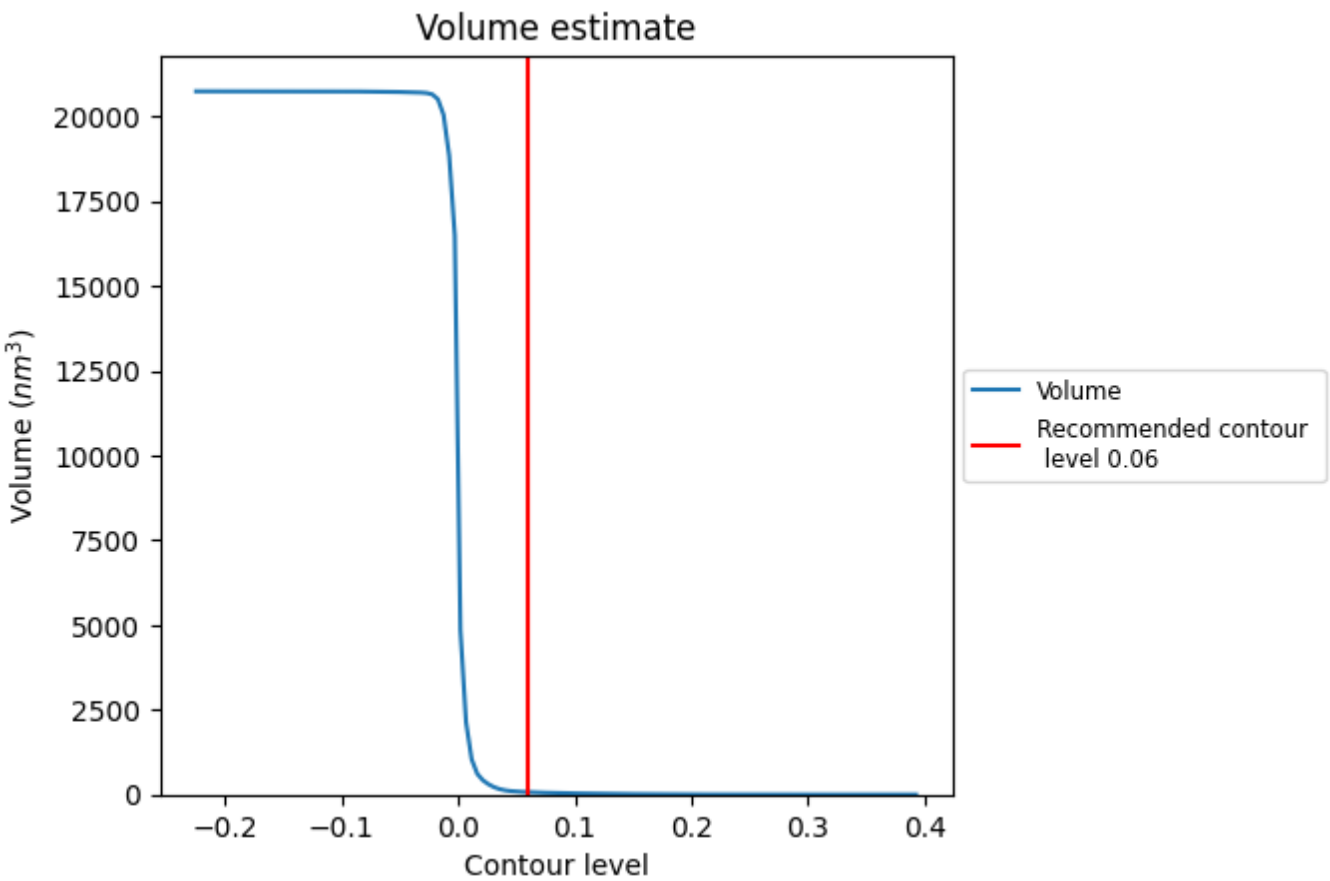
This section contains the results of statistical analysis of the map.

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.

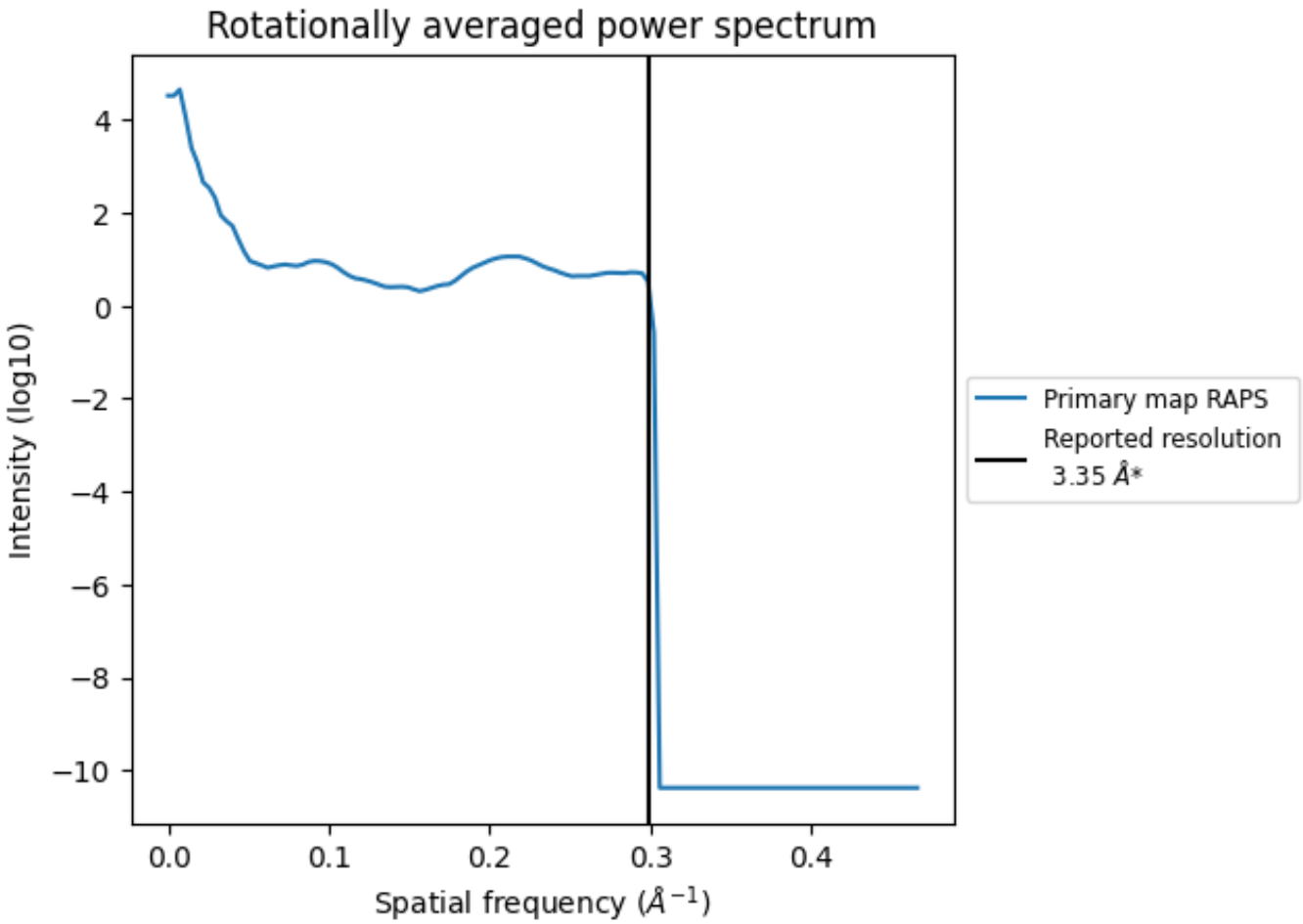
7.2 Volume estimate (i)



The volume at the recommended contour level is 71 nm^3; this corresponds to an approximate mass of 64 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)

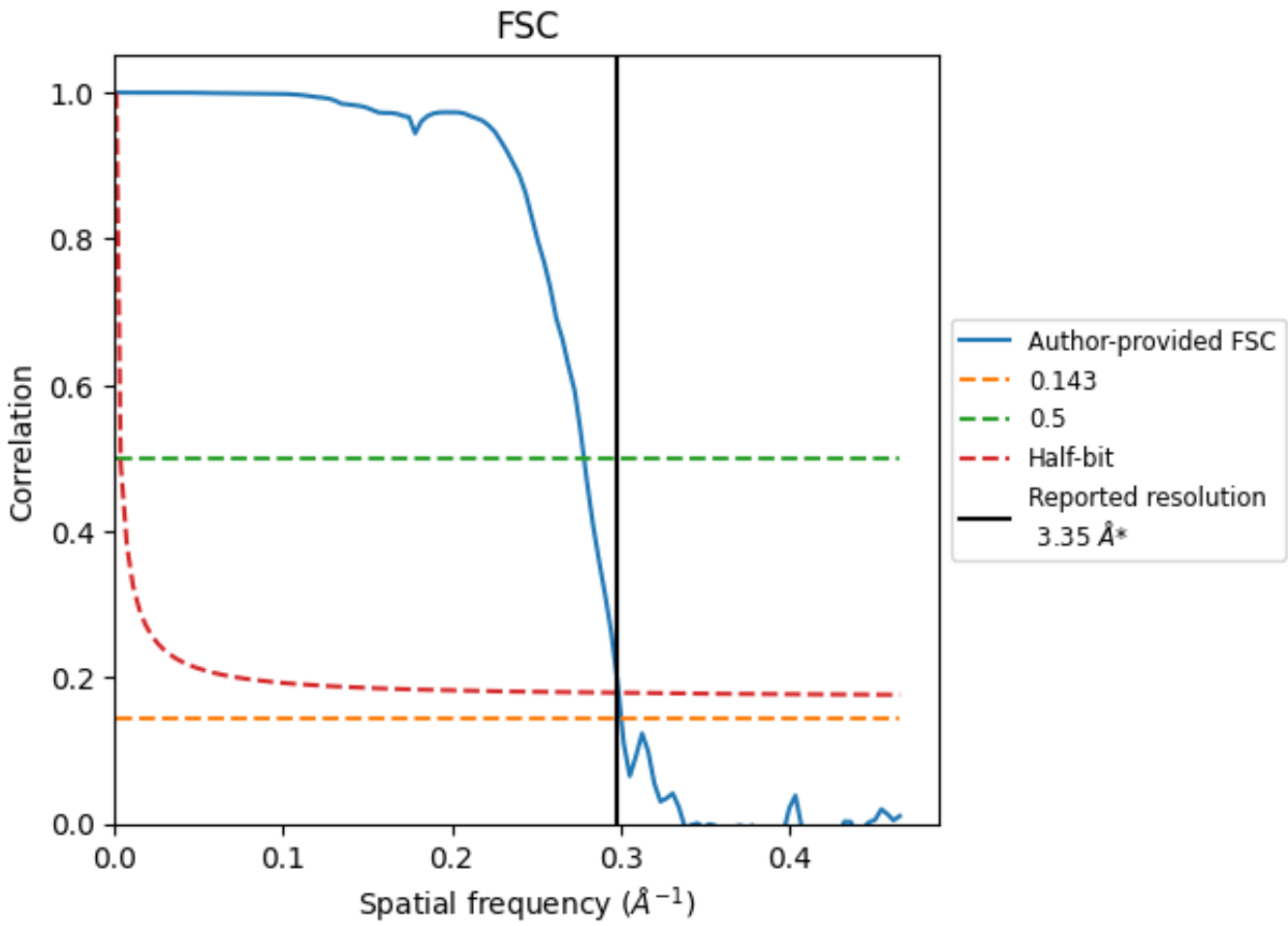


*Reported resolution corresponds to spatial frequency of 0.299 Å^{-1}

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 0.299 Å^{-1}

8.2 Resolution estimates (i)

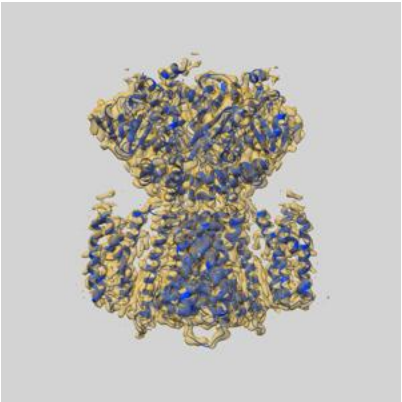
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.35	-	-
Author-provided FSC curve	3.33	3.59	3.34
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

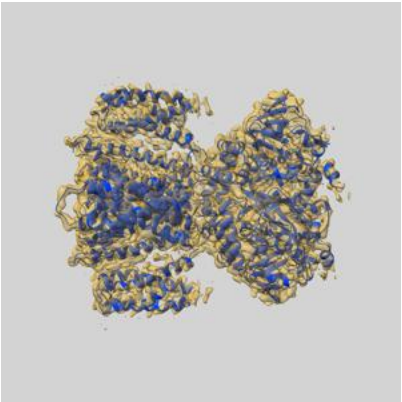
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-7484 and PDB model 6CJU. Per-residue inclusion information can be found in section 3 on page 10.

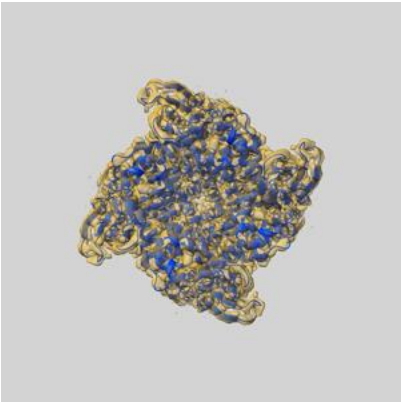
9.1 Map-model overlay (i)



X



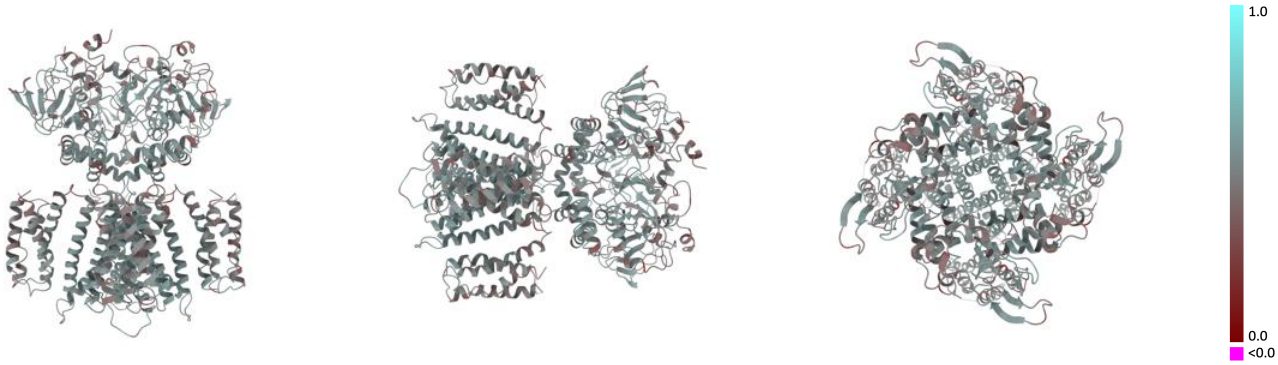
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.06 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.

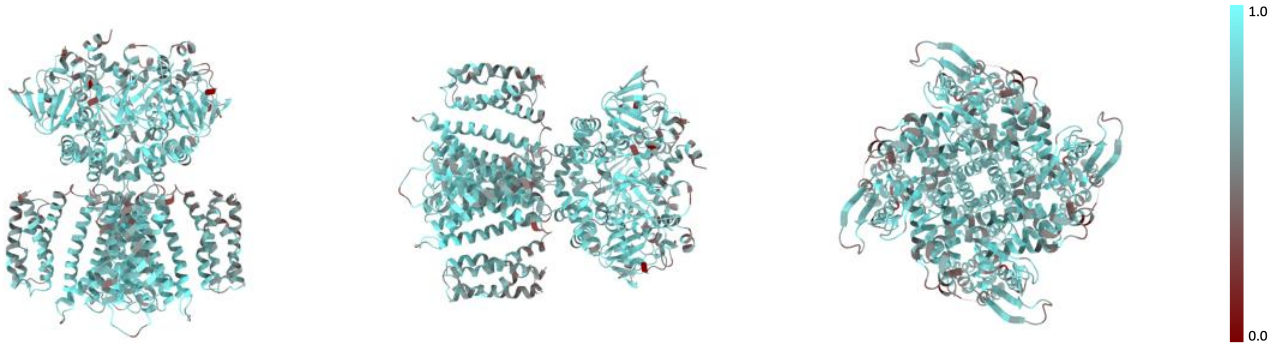
9.2 Q-score mapped to coordinate model (i)



X Y Z

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.

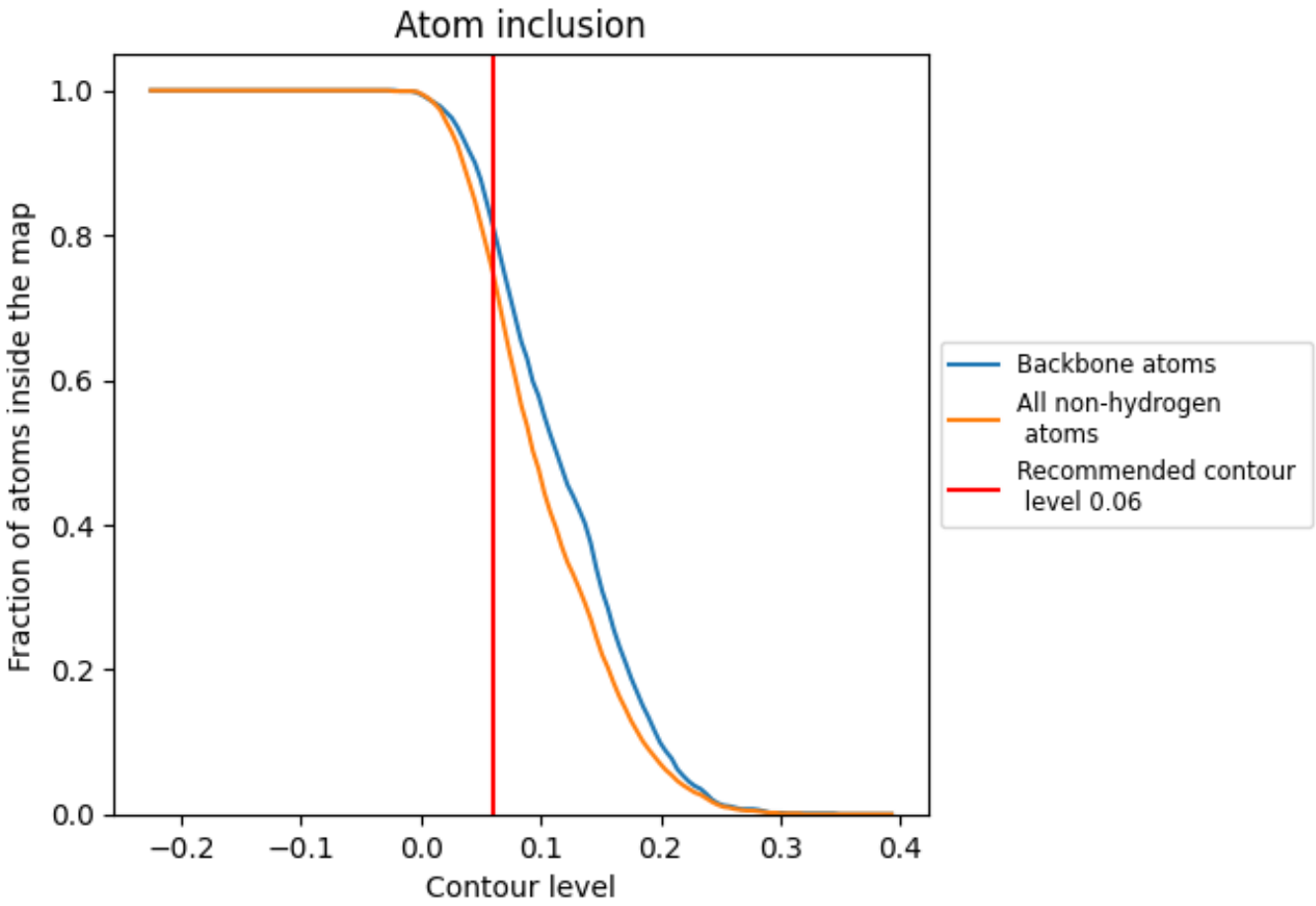
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.06).

9.4 Atom inclusion (i)



At the recommended contour level, 81% of all backbone atoms, 75% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7500	0.5050
A	0.7510	0.5040
B	0.7490	0.5050
C	0.7490	0.5070
D	0.7500	0.5050