



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 10:49 am BST

PDB ID : 5WPK
Title : Structure of the class II 3-hydroxy-3-methylglutaryl-CoA reductase from *Streptococcus pneumoniae* bound to HMG-CoA and in a partially closed conformation
Authors : Miller, B.R.; Kung, Y.
Deposited on : 2017-08-04
Resolution : 2.30 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

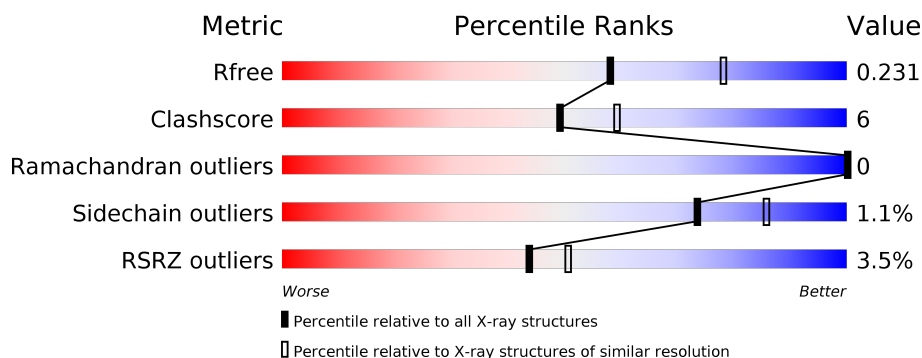
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 on 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	426	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>14%</div> <div>•</div> <div>11%</div> </div> </div>
1	B	426	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>12%</div> <div>•</div> </div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6617 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.

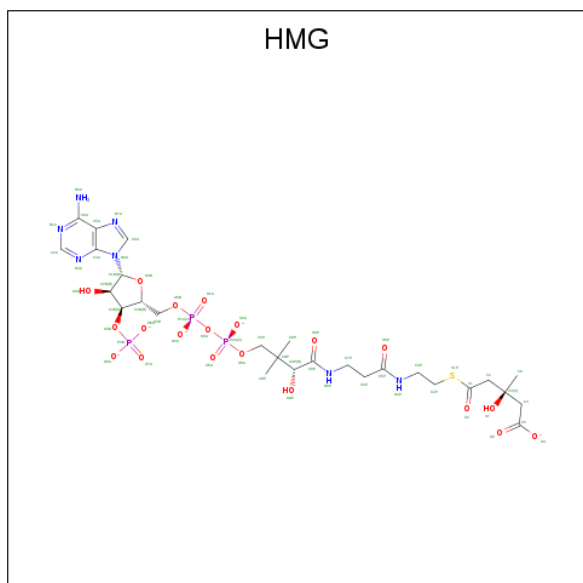
- Molecule 1 is a protein called 3-hydroxy-3-methylglutaryl coenzyme A reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	1	0
			2872	1810	498	551	13			
1	B	422	Total	C	N	O	S	0	3	0
			3162	1988	553	608	13			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A0D6J7E8
A	0	HIS	-	expression tag	UNP A0A0D6J7E8
A	355	GLU	VAL	conflict	UNP A0A0D6J7E8
B	-1	GLY	-	expression tag	UNP A0A0D6J7E8
B	0	HIS	-	expression tag	UNP A0A0D6J7E8
B	355	GLU	VAL	conflict	UNP A0A0D6J7E8

- Molecule 2 is 3-HYDROXY-3-METHYLGLUTARYL-COENZYME A (three-letter code: HMG) (formula: C₂₇H₃₉N₇O₂₀P₃S).



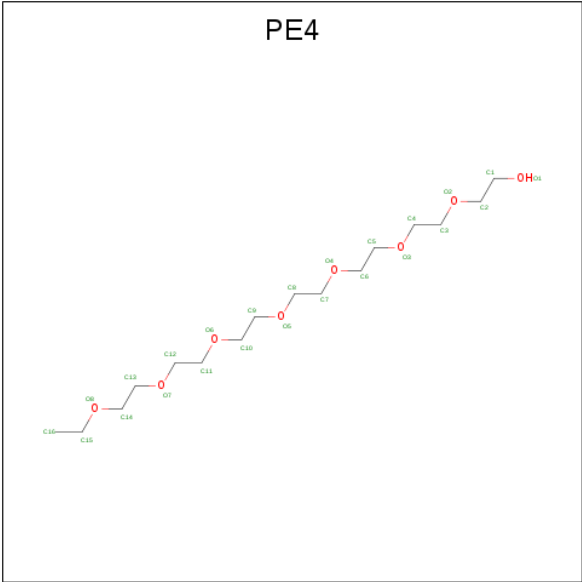
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			58	27	7	20	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			58	27	7	20	3	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is 2-{2-[2-(2-{2-[2-(2-ETHOXY-ETHOXY)-ETHOXY]-ETHOXY}-ETHOXY)-ETHOXY]-ETHOXY}-ETHANOL (three-letter code: PE4) (formula: C₁₆H₃₄O₈).

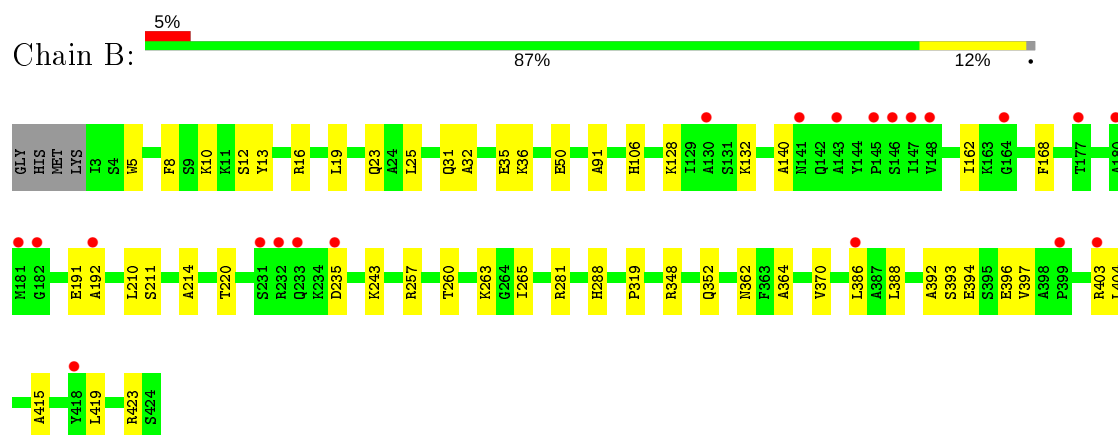
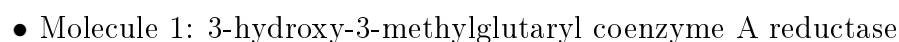


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			16	10	6		
4	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	230	Total	O	0	0
			230	230		
5	B	187	Total	O	0	0
			187	187		

- Molecule 1: 3-hydroxy-3-methylglutaryl coenzyme A reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.87Å 131.25Å 57.96Å 90.00° 102.54° 90.00°	Depositor
Resolution (Å)	19.24 – 2.30 19.24 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.3 (19.24-2.30) 97.9 (19.24-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.30Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.177 , 0.233 0.176 , 0.231	Depositor DCC
R_{free} test set	1778 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 30.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.216 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6617	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PE4, GOL, HMG

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	0.28	0/2922	0.47	0/3961
1	B	0.26	0/3221	0.45	0/4371
All	All	0.27	0/6143	0.46	0/8332

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

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	2872	0	2864	43	0
1	B	3162	0	3120	38	0
2	A	58	0	39	1	0
2	B	58	0	39	2	0
3	A	6	0	8	1	0
3	B	12	0	16	4	0
4	B	32	0	41	2	0
5	A	230	0	0	7	1
5	B	187	0	0	1	1
All	All	6617	0	6127	73	1

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 (73) 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:220:THR:HG21	3:B:503:GOL:H11	1.80	0.63
1:A:106:HIS:NE2	1:A:220:THR:HG22	2.14	0.62
1:A:151:GLY:O	1:A:155:ARG:NH2	2.30	0.61
1:A:297:TYR:O	5:A:601:HOH:O	2.17	0.60
1:B:10:LYS:HD2	1:B:394:GLU:HG3	1.84	0.59
1:A:139:LEU:HD23	1:A:193:LEU:HD23	1.84	0.59
1:A:199:GLU:O	1:A:202:GLN:NE2	2.34	0.59
1:B:162:ILE:CG2	1:B:168:PHE:HB2	2.33	0.58
1:A:91:ALA:HB2	1:A:364:ALA:HB2	1.84	0.58
1:A:249:GLN:NE2	5:A:613:HOH:O	2.37	0.58
1:B:25:LEU:HD21	4:B:505:PE4:H81	1.86	0.57
1:A:265:ILE:HG12	1:A:362:ASN:HB2	1.86	0.57
1:A:11:LYS:HZ2	1:A:19:LEU:HD12	1.70	0.57
1:A:128:LYS:O	1:A:132:LYS:HG3	2.04	0.57
1:A:51:ASN:OD1	2:B:502:HMG:H2A1	2.05	0.56
1:A:254:ASP:HB2	1:B:162:ILE:CD1	2.35	0.56
1:A:181:MET:HB3	3:A:501:GOL:H2	1.89	0.55
1:A:138:GLU:O	1:A:142:GLN:HG3	2.07	0.54
1:A:43:THR:O	1:A:47:GLN:NE2	2.41	0.54
1:A:254:ASP:HB2	1:B:162:ILE:HD12	1.90	0.53
1:A:31:GLN:O	1:A:35:GLU:HG3	2.09	0.52
1:A:288:HIS:HB3	1:B:288:HIS:HB3	1.90	0.52
1:A:228:ARG:NH2	5:A:626:HOH:O	2.43	0.52
1:B:393:SER:O	1:B:397:VAL:HG23	2.09	0.52
1:A:374:ILE:HD12	2:A:500:HMG:H62	1.92	0.52
1:A:253:ALA:C	1:B:162:ILE:HD11	2.30	0.52
1:A:133:LYS:HG3	1:A:157:LEU:HB3	1.92	0.51
1:B:31:GLN:O	1:B:35:GLU:HG3	2.10	0.51
1:B:257:ARG:NH2	2:B:502:HMG:O3	2.43	0.51
1:A:4:SER:O	5:A:602:HOH:O	2.19	0.51
1:B:265:ILE:HG12	1:B:362:ASN:HB2	1.92	0.50
1:A:243:LYS:HB3	1:A:370:VAL:CG1	2.42	0.49
1:A:86:ALA:HB2	1:B:50:GLU:HG2	1.95	0.49
1:A:243:LYS:HB3	1:A:370:VAL:HG13	1.95	0.48
1:A:263:LYS:NZ	5:A:611:HOH:O	2.35	0.48
1:B:91:ALA:HB2	1:B:364:ALA:HB2	1.94	0.48
1:B:392:ALA:HB1	1:B:396:GLU:HG3	1.94	0.48
1:B:128:LYS:O	1:B:132:LYS:HG3	2.14	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:SER:O	1:A:16:ARG:HG3	2.14	0.47
1:A:9:SER:OG	5:A:604:HOH:O	2.20	0.47
1:B:348:ARG:O	1:B:352:GLN:HG3	2.14	0.47
1:B:140:ALA:HA	3:B:504:GOL:O1	2.14	0.47
1:B:162:ILE:HG23	1:B:168:PHE:HB2	1.96	0.47
1:B:5:TRP:CZ2	1:B:23:GLN:HG3	2.50	0.47
1:A:242:GLU:HA	1:A:305:LEU:HD11	1.96	0.46
1:A:294:ASP:O	5:A:603:HOH:O	2.20	0.46
1:B:12:SER:O	1:B:16:ARG:HG3	2.15	0.46
1:A:375:GLN:NE2	1:B:191:GLU:HG2	2.30	0.46
1:B:214:ALA:O	1:B:319:PRO:HB3	2.16	0.45
1:B:243:LYS:HB3	1:B:370:VAL:HB	1.98	0.45
1:B:13:TYR:CE2	4:B:501:PE4:H62	2.51	0.45
1:B:210:LEU:HD12	1:B:211:SER:H	1.82	0.45
1:A:348:ARG:O	1:A:352:GLN:HG3	2.17	0.45
1:B:32:ALA:O	1:B:36:LYS:HG3	2.16	0.45
1:B:192:ALA:CB	3:B:504:GOL:H11	2.46	0.44
1:A:11:LYS:NZ	1:A:19:LEU:HD12	2.31	0.44
1:B:419:LEU:HD11	1:B:423:ARG:CZ	2.47	0.44
1:A:137:LEU:HD22	1:A:154:ALA:HB1	2.00	0.44
1:A:105:VAL:HG23	1:A:352:GLN:HE21	1.83	0.43
1:A:211:SER:HA	1:B:260:THR:OG1	2.18	0.43
1:B:8:PHE:CD1	1:B:19:LEU:HD23	2.54	0.43
1:B:404:LEU:HD11	1:B:415:ALA:HB2	2.00	0.43
1:A:147:ILE:HG12	1:A:185:MET:SD	2.59	0.43
1:A:5:TRP:CZ2	1:A:23:GLN:HG3	2.54	0.42
1:B:386:LEU:HA	1:B:386:LEU:HD13	1.89	0.42
1:B:263:LYS:NZ	5:B:607:HOH:O	2.26	0.42
1:A:165:GLU:HA	1:A:166:PRO:HA	1.82	0.42
1:B:10:LYS:HA	1:B:388:LEU:HD21	2.00	0.42
1:A:220:THR:HA	1:A:316:MET:O	2.20	0.41
1:A:140:ALA:O	1:A:189:MET:HG2	2.20	0.41
1:B:106:HIS:HE1	3:B:503:GOL:O1	2.02	0.41
1:B:162:ILE:HD13	1:B:162:ILE:HG21	1.64	0.40
1:A:281:ARG:HB2	1:B:263:LYS:HE2	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:802:HOH:O	5:B:621:HOH:O[1_554]	2.17	0.03

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	380/426 (89%)	368 (97%)	12 (3%)	0	100	100
1	B	423/426 (99%)	412 (97%)	11 (3%)	0	100	100
All	All	803/852 (94%)	780 (97%)	23 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	297/344 (86%)	293 (99%)	4 (1%)	69	82
1	B	322/344 (94%)	319 (99%)	3 (1%)	78	89
All	All	619/688 (90%)	612 (99%)	7 (1%)	73	86

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	181	MET
1	A	210	LEU
1	A	220	THR
1	A	281	ARG
1	B	235	ASP
1	B	281	ARG
1	B	403	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	187	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PE4	B	505	-	15,15,23	0.53	0	14,14,22	0.20	0
3	GOL	B	503	-	5,5,5	0.37	0	5,5,5	0.26	0
2	HMG	A	500	-	49,60,60	3.65	15 (30%)	59,90,90	1.87	9 (15%)
4	PE4	B	501	-	15,15,23	0.53	0	14,14,22	0.31	0
2	HMG	B	502	-	49,60,60	3.66	15 (30%)	59,90,90	1.96	13 (22%)
3	GOL	A	501	-	5,5,5	0.37	0	5,5,5	0.29	0
3	GOL	B	504	-	5,5,5	0.37	0	5,5,5	0.76	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
4	PE4	B	505	-	-	7/13/13/21	-
3	GOL	B	503	-	-	4/4/4/4	-
2	HMG	A	500	-	-	13/54/77/77	0/3/3/3
4	PE4	B	501	-	-	11/13/13/21	-
2	HMG	B	502	-	-	10/54/77/77	0/3/3/3
3	GOL	A	501	-	-	2/4/4/4	-
3	GOL	B	504	-	-	0/4/4/4	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	HMG	O4B-C1B	15.49	1.62	1.41
2	A	500	HMG	O4B-C1B	15.34	1.62	1.41
2	A	500	HMG	C2B-C1B	-14.95	1.31	1.53
2	B	502	HMG	C2B-C1B	-14.78	1.31	1.53
2	B	502	HMG	O4B-C4B	-6.06	1.31	1.45
2	A	500	HMG	O4B-C4B	-6.05	1.31	1.45
2	B	502	HMG	C9P-N8P	5.81	1.46	1.33
2	A	500	HMG	C9P-N8P	5.53	1.45	1.33
2	B	502	HMG	C5P-N4P	5.45	1.45	1.33
2	A	500	HMG	C5P-N4P	5.36	1.45	1.33
2	B	502	HMG	P3B-O3B	3.49	1.65	1.59
2	B	502	HMG	C6A-N6A	3.18	1.45	1.34
2	A	500	HMG	C6A-N6A	3.14	1.45	1.34
2	A	500	HMG	P3B-O3B	3.12	1.65	1.59
2	A	500	HMG	O7-C3	-2.93	1.40	1.44
2	A	500	HMG	O2B-C2B	2.92	1.49	1.43
2	B	502	HMG	O2B-C2B	2.91	1.49	1.43
2	B	502	HMG	O7-C3	-2.69	1.40	1.44
2	B	502	HMG	C5A-C4A	-2.56	1.34	1.40
2	A	500	HMG	C5A-C4A	-2.55	1.34	1.40
2	A	500	HMG	C1-S1P	2.54	1.82	1.76
2	B	502	HMG	C1-S1P	2.42	1.82	1.76
2	A	500	HMG	O3B-C3B	-2.28	1.35	1.44
2	B	502	HMG	O9P-C9P	-2.28	1.18	1.23
2	A	500	HMG	O9P-C9P	-2.27	1.18	1.23
2	A	500	HMG	C2A-N3A	2.25	1.35	1.32
2	B	502	HMG	O5P-C5P	-2.23	1.18	1.23
2	A	500	HMG	O5P-C5P	-2.20	1.18	1.23
2	B	502	HMG	O3B-C3B	-2.15	1.36	1.44

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	HMG	C2A-N3A	2.06	1.35	1.32

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	HMG	C5A-C6A-N6A	6.77	130.63	120.35
2	A	500	HMG	C5A-C6A-N6A	6.52	130.27	120.35
2	A	500	HMG	C2-C1-S1P	6.04	121.11	113.63
2	A	500	HMG	N3A-C2A-N1A	-5.52	120.05	128.68
2	B	502	HMG	C2-C1-S1P	5.38	120.28	113.63
2	B	502	HMG	N3A-C2A-N1A	-5.34	120.33	128.68
2	A	500	HMG	N6A-C6A-N1A	-4.35	109.54	118.57
2	B	502	HMG	N6A-C6A-N1A	-4.34	109.57	118.57
2	B	502	HMG	C6P-C5P-N4P	4.10	123.32	116.42
2	A	500	HMG	C1B-N9A-C4A	4.09	133.83	126.64
2	B	502	HMG	C1B-N9A-C4A	3.35	132.53	126.64
2	B	502	HMG	C7P-C6P-C5P	-3.13	107.15	112.36
2	B	502	HMG	O5P-C5P-N4P	-2.85	117.64	123.01
2	B	502	HMG	P1A-O3A-P2A	-2.84	123.07	132.83
2	A	500	HMG	O2-C1-S1P	-2.82	118.95	122.61
2	B	502	HMG	C2P-S1P-C1	2.64	110.11	101.87
2	B	502	HMG	CAP-C9P-N8P	2.39	121.34	116.58
2	A	500	HMG	P1A-O3A-P2A	-2.36	124.73	132.83
2	B	502	HMG	O2-C1-C2	-2.36	119.87	123.64
2	A	500	HMG	C2P-C3P-N4P	-2.33	107.53	112.42
2	A	500	HMG	O2-C1-C2	-2.32	119.92	123.64
2	B	502	HMG	O2-C1-S1P	-2.13	119.85	122.61

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	503	GOL	O1-C1-C2-C3
3	B	503	GOL	O2-C2-C3-O3
2	A	500	HMG	CCP-O6A-P2A-O5A
2	A	500	HMG	CCP-O6A-P2A-O3A
2	A	500	HMG	C5B-O5B-P1A-O1A
2	A	500	HMG	C5B-O5B-P1A-O2A
2	B	502	HMG	C5B-O5B-P1A-O3A
2	B	502	HMG	C5B-O5B-P1A-O1A
2	B	502	HMG	C3P-C2P-S1P-C1
2	B	502	HMG	O2-C1-S1P-C2P

Continued on next page...

Continued from previous page...

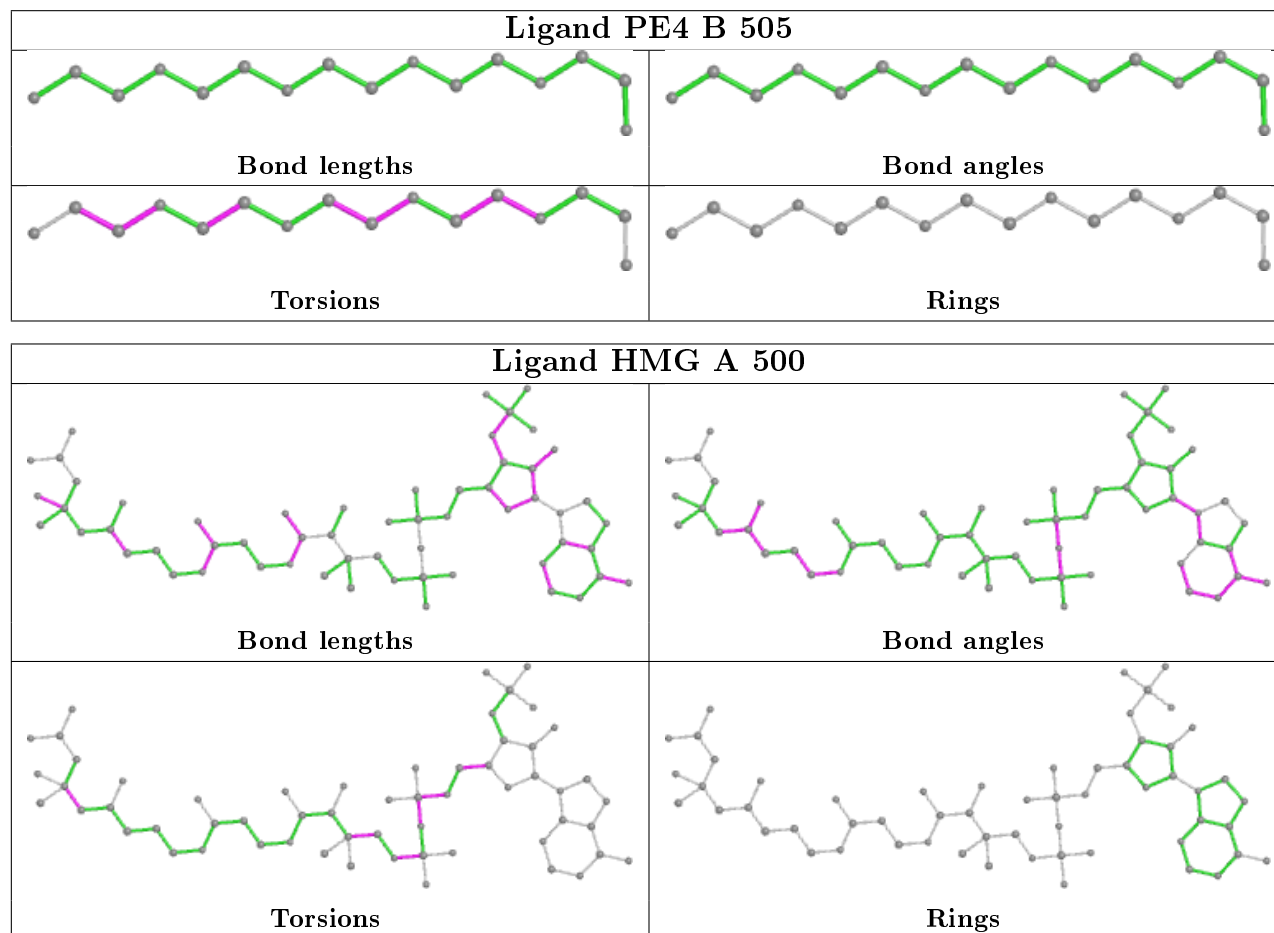
Mol	Chain	Res	Type	Atoms
2	B	502	HMG	C2-C1-S1P-C2P
3	A	501	GOL	O1-C1-C2-C3
2	B	502	HMG	C6P-C5P-N4P-C3P
4	B	501	PE4	C9-C10-O6-C11
4	B	501	PE4	O4-C7-C8-O5
2	B	502	HMG	O5P-C5P-N4P-C3P
4	B	501	PE4	O6-C10-C9-O5
4	B	505	PE4	O4-C7-C8-O5
4	B	501	PE4	O6-C11-C12-O7
4	B	505	PE4	O3-C5-C6-O4
3	B	503	GOL	C1-C2-C3-O3
3	B	503	GOL	O1-C1-C2-O2
3	A	501	GOL	O1-C1-C2-O2
4	B	505	PE4	O6-C10-C9-O5
2	A	500	HMG	C3B-C4B-C5B-O5B
4	B	501	PE4	O2-C3-C4-O3
4	B	501	PE4	C7-C8-O5-C9
4	B	505	PE4	C10-C9-O5-C8
4	B	505	PE4	C4-C3-O2-C2
4	B	501	PE4	C12-C11-O6-C10
4	B	501	PE4	C3-C4-O3-C5
4	B	505	PE4	C6-C5-O3-C4
2	A	500	HMG	CAP-CBP-CCP-O6A
4	B	501	PE4	C5-C6-O4-C7
2	A	500	HMG	O4B-C4B-C5B-O5B
2	A	500	HMG	CEP-CBP-CCP-O6A
4	B	501	PE4	O3-C5-C6-O4
2	A	500	HMG	P2A-O3A-P1A-O1A
4	B	501	PE4	C8-C7-O4-C6
2	A	500	HMG	CDP-CBP-CCP-O6A
4	B	505	PE4	O2-C3-C4-O3
2	B	502	HMG	C3B-C4B-C5B-O5B
2	A	500	HMG	C5B-O5B-P1A-O3A
2	A	500	HMG	P2A-O3A-P1A-O2A
2	A	500	HMG	C1-C2-C3-C4
2	B	502	HMG	CCP-O6A-P2A-O4A
2	B	502	HMG	S1P-C2P-C3P-N4P

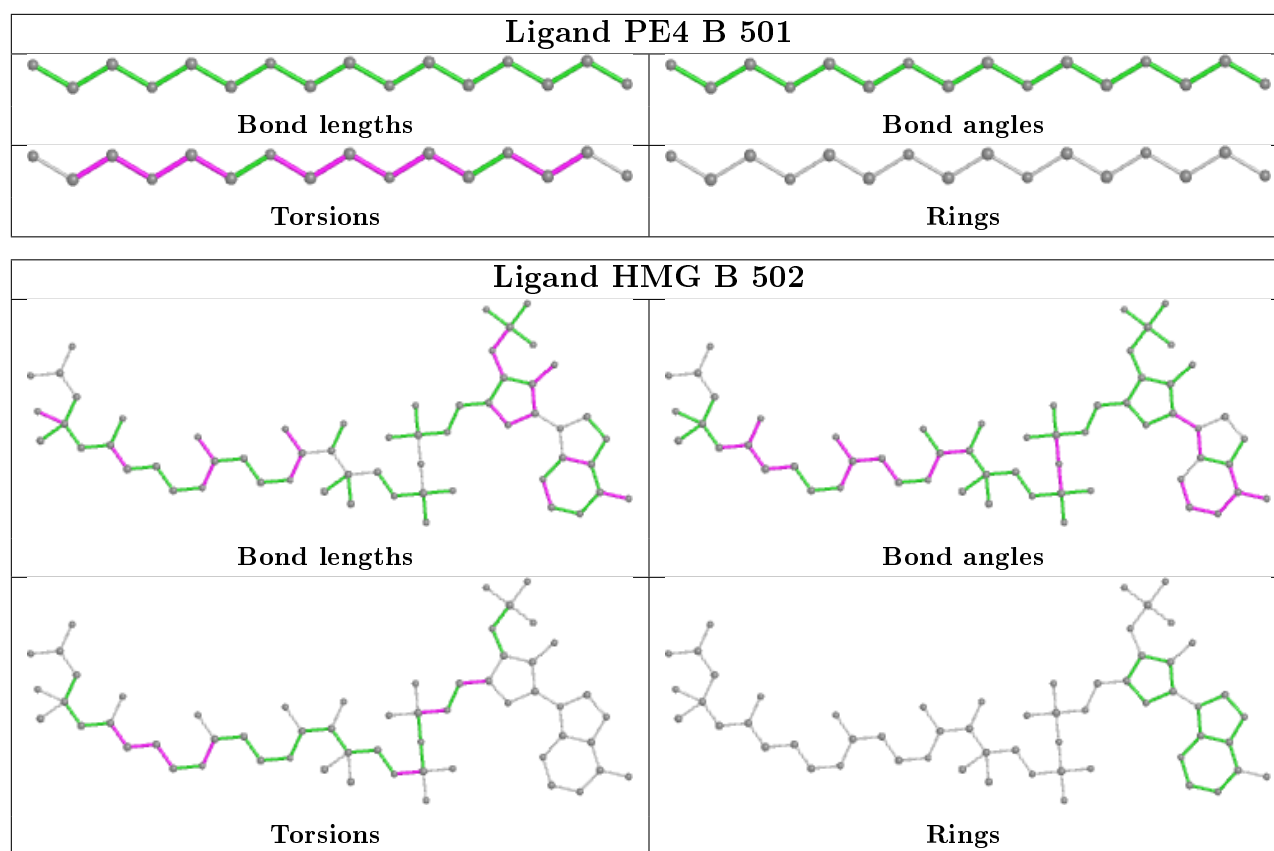
There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	505	PE4	1	0
3	B	503	GOL	2	0
2	A	500	HMG	1	0
4	B	501	PE4	1	0
2	B	502	HMG	2	0
3	A	501	GOL	1	0
3	B	504	GOL	2	0

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





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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	381/426 (89%)	-0.26	7 (1%) 68 74	12, 24, 47, 82	0
1	B	422/426 (99%)	-0.01	21 (4%) 28 35	13, 26, 59, 79	0
All	All	803/852 (94%)	-0.13	28 (3%) 44 51	12, 25, 57, 82	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	147	ILE	7.0
1	A	379	MET	5.5
1	A	374	ILE	4.6
1	B	181	MET	4.5
1	B	164	GLY	4.2
1	B	182	GLY	4.2
1	B	146	SER	4.0
1	A	377	GLY	3.7
1	B	143	ALA	3.5
1	B	177	THR	3.2
1	B	418	TYR	3.0
1	B	386	LEU	3.0
1	B	403	ARG	3.0
1	B	399	PRO	2.9
1	B	148	VAL	2.8
1	B	180	ALA	2.7
1	B	192	ALA	2.7
1	B	233	GLN	2.7
1	B	145	PRO	2.6
1	B	130	ALA	2.6
1	B	232	ARG	2.4
1	A	0	HIS	2.2
1	B	235	ASP	2.2
1	A	375	GLN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	378	HIS	2.1
1	A	1	MET	2.1
1	B	141	ASN	2.1
1	B	231	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

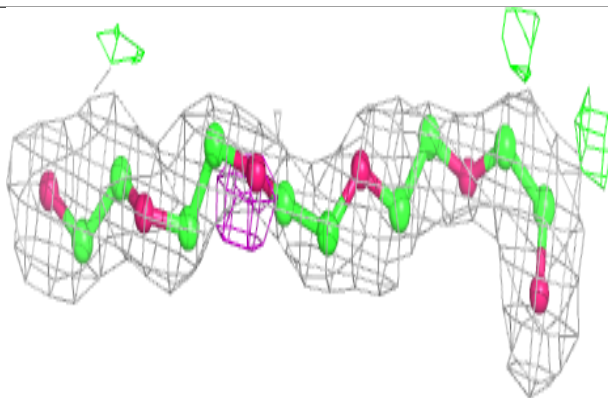
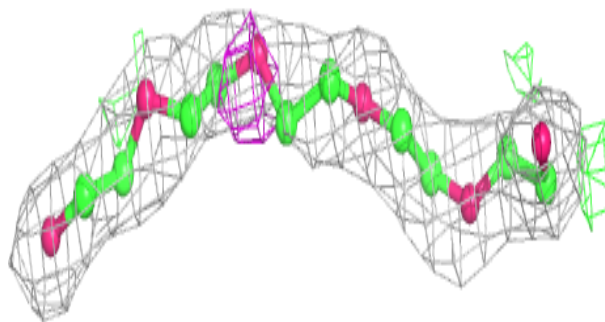
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	B	504	6/6	0.67	0.22	45,51,54,55	0
3	GOL	B	503	6/6	0.87	0.19	26,35,36,37	0
4	PE4	B	505	16/24	0.87	0.27	31,38,44,46	0
4	PE4	B	501	16/24	0.88	0.26	30,40,53,56	0
3	GOL	A	501	6/6	0.91	0.14	27,34,42,45	0
2	HMG	B	502	58/58	0.93	0.16	19,38,54,64	0
2	HMG	A	500	58/58	0.96	0.12	18,31,44,54	0

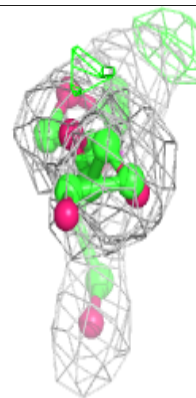
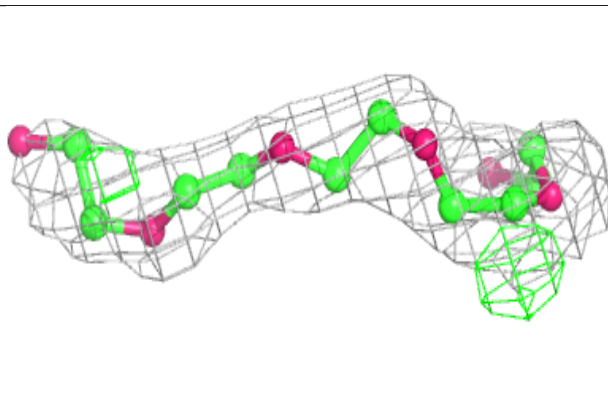
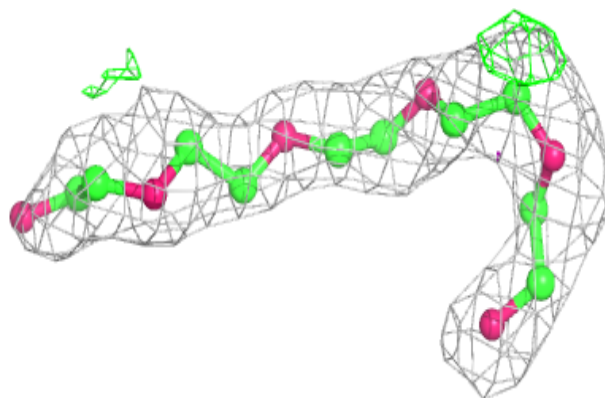
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around PE4 B 505:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

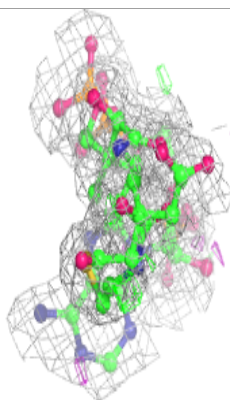
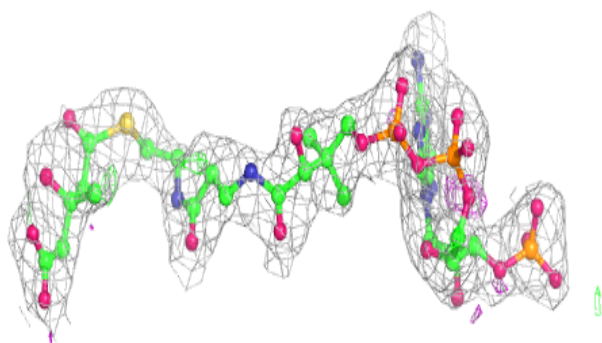
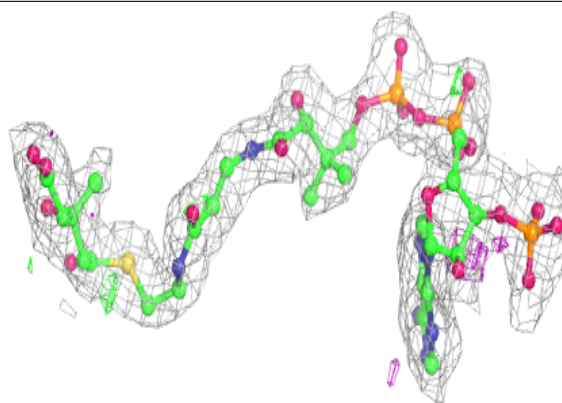
**Electron density around PE4 B 501:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

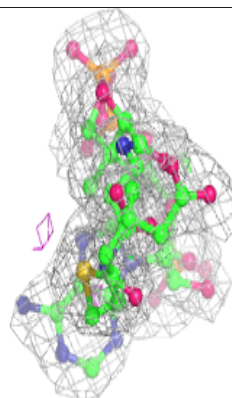
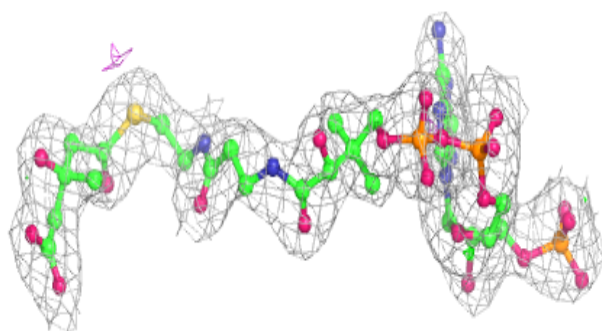
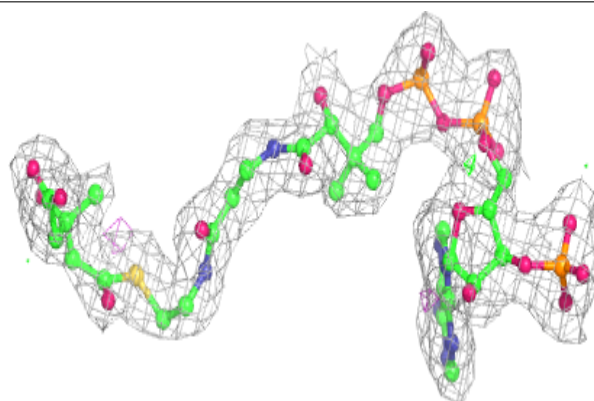


Electron density around HMG B 502:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around HMG A 500:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.