



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:54 PM UTC

PDB ID : 5WPK / pdb\_00005wpk  
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](mailto: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 types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

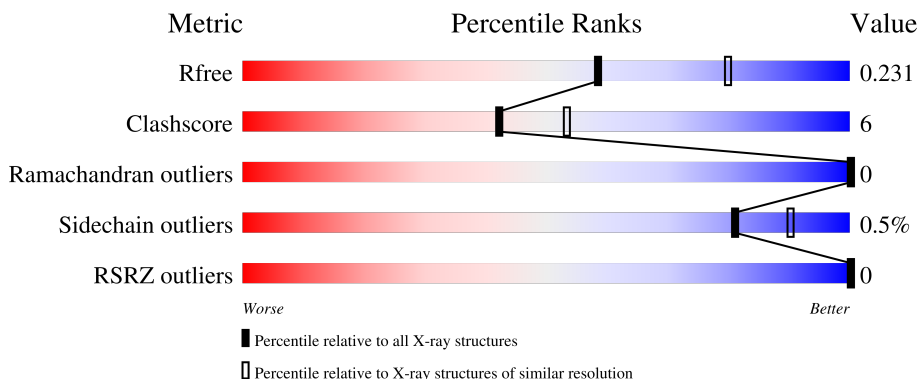
# 1 Overall quality at a glance

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))
$R_{free}$	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	 75% 14% 11%
1	B	426	 86% 13% .

## 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.

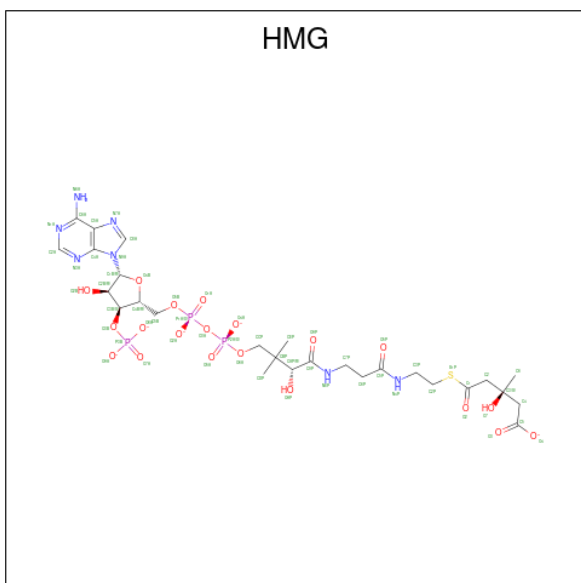
- Molecule 1 is a protein called 3-hydroxy-3-methylglutaryl coenzyme A reductase.

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

There are 6 discrepancies between the modelled and reference sequences:

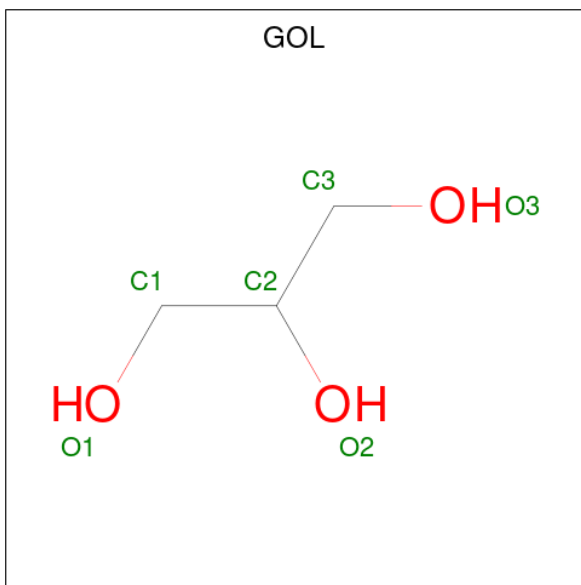
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A0D6J7E8
A	0	HIS	-	expression tag	UNP A0A0D6J7E8
A	355	GLU	VAL	conflict	UNP A0A0D6J7E8
B	-1	GLY	-	expression tag	UNP A0A0D6J7E8
B	0	HIS	-	expression tag	UNP A0A0D6J7E8
B	355	GLU	VAL	conflict	UNP A0A0D6J7E8

- Molecule 2 is 3-HYDROXY-3-METHYLGLUTARYL-COENZYME A (CCD ID: HMG) (formula: C<sub>27</sub>H<sub>39</sub>N<sub>7</sub>O<sub>20</sub>P<sub>3</sub>S).



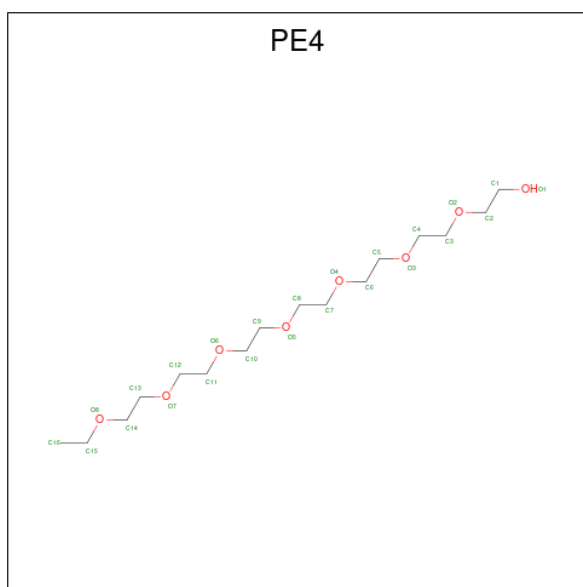
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	A	1	Total	C	N	O	P	S	0	0
			58	27	7	20	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			58	27	7	20	3	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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 (CCD ID: PE4) (formula: C<sub>16</sub>H<sub>34</sub>O<sub>8</sub>).



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

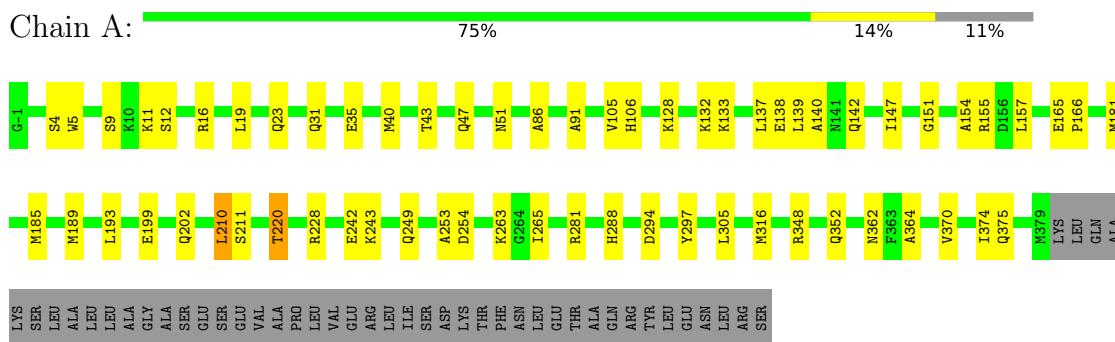
- Molecule 5 is water.

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

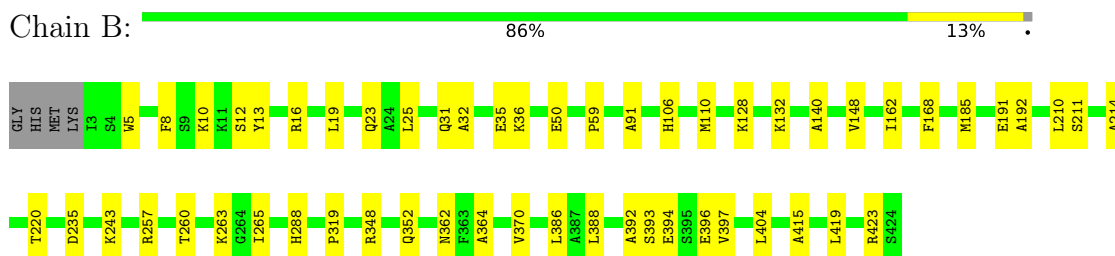
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron 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 dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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: 3-hydroxy-3-methylglutaryl coenzyme A reductase



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



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	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
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.88 (at 2.30Å)	Xtrriage
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 (3.62%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.9	Xtrriage
Anisotropy	0.502	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 32.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.216 for l,-k,h	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	6617	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage'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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.22	1/2922 (0.0%)	0.38	0/3961
1	B	0.14	0/3221	0.36	0/4371
All	All	0.18	1/6143 (0.0%)	0.37	0/8332

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	210	LEU	C-O	-6.73	1.15	1.23

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	2872	0	2864	45	0
1	B	3162	0	3120	41	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	187	0	0	1	1
All	All	6617	0	6127	77	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 (77) 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:110[A]:MET:HE3	1:B:210:LEU:HD11	1.76	0.67
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:A:11:LYS:HZ2	1:A:19:LEU:HD12	1.66	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:253:ALA:C	1:B:162:ILE:HD11	2.30	0.57
1:A:265:ILE:HG12	1:A:362:ASN:HB2	1.86	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:B:148:VAL:HG11	1:B:185:MET:HE1	1.87	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:B:263:LYS:NZ	5:B:607:HOH:O	2.26	0.47
1:A:9:SER:OG	5:A:604:HOH:O	2.20	0.47
1:A:12:SER:O	1:A:16:ARG:HG3	2.14	0.47
1:B:140:ALA:HA	3:B:504:GOL:O1	2.14	0.47
1:B:348:ARG:O	1:B:352:GLN:HG3	2.14	0.47
1:A:185:MET:HE3	1:A:185:MET:HB2	1.77	0.47
1:B:5:TRP:CZ2	1:B:23:GLN:HG3	2.50	0.47
1:B:162:ILE:HG23	1:B:168:PHE:HB2	1.96	0.47
1:A:242:GLU:HA	1:A:305:LEU:HD11	1.96	0.46
1:A:40:MET:HE1	1:B:59:PRO:HD3	1.97	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:13:TYR:CE2	4:B:501:PE4:H62	2.51	0.45
1:B:243:LYS:HB3	1:B:370:VAL:HB	1.98	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:HD13	1:B:386:LEU:HA	1.89	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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 [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	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 [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	297/344 (86%)	295 (99%)	2 (1%)	76	87
1	B	322/344 (94%)	321 (100%)	1 (0%)	86	93
All	All	619/688 (90%)	616 (100%)	3 (0%)	81	90

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	210	LEU
1	A	220	THR
1	B	235	ASP

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	109	GLN
1	A	113	GLN
1	A	121	ASN
1	A	298	GLN
1	A	344	ASN
1	B	109	GLN
1	B	126	GLN
1	B	344	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

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	501	-	15,15,23	0.54	0	14,14,22	0.31	0
2	HMG	B	502	-	58,60,60	2.66	18 (31%)	81,90,90	1.93	21 (25%)
4	PE4	B	505	-	15,15,23	0.54	0	14,14,22	0.20	0
3	GOL	B	503	-	5,5,5	0.37	0	5,5,5	0.29	0
2	HMG	A	500	-	58,60,60	2.63	18 (31%)	81,90,90	1.90	16 (19%)
3	GOL	A	501	-	5,5,5	0.38	0	5,5,5	0.33	0
3	GOL	B	504	-	5,5,5	0.38	0	5,5,5	0.82	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	501	-	-	11/13/13/21	-
2	HMG	B	502	-	-	10/60/77/77	0/3/3/3
4	PE4	B	505	-	-	7/13/13/21	-
3	GOL	B	503	-	-	4/4/4/4	-
2	HMG	A	500	-	-	13/60/77/77	0/3/3/3
3	GOL	A	501	-	-	2/4/4/4	-
3	GOL	B	504	-	-	0/4/4/4	-

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	HMG	O4B-C1B	8.95	1.62	1.42
2	A	500	HMG	O4B-C1B	8.87	1.62	1.42
2	A	500	HMG	C2B-C1B	-7.13	1.31	1.53
2	B	502	HMG	C2B-C1B	-7.05	1.31	1.53
2	B	502	HMG	O4B-C4B	-6.09	1.31	1.45
2	A	500	HMG	O4B-C4B	-6.08	1.31	1.45
2	B	502	HMG	P2A-O3A	5.63	1.65	1.59
2	B	502	HMG	P1A-O3A	5.60	1.65	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HMG	P1A-O3A	5.60	1.65	1.59
2	B	502	HMG	C9P-N8P	5.40	1.46	1.33
2	A	500	HMG	P2A-O3A	5.29	1.65	1.59
2	B	502	HMG	C5P-N4P	5.22	1.45	1.33
2	A	500	HMG	C9P-N8P	5.15	1.45	1.33
2	A	500	HMG	C5P-N4P	5.14	1.45	1.33
2	B	502	HMG	C6A-N6A	4.48	1.45	1.34
2	A	500	HMG	C6A-N6A	4.42	1.45	1.34
2	B	502	HMG	P3B-O3B	3.64	1.65	1.59
2	A	500	HMG	P3B-O3B	3.23	1.65	1.59
2	A	500	HMG	O7-C3	-3.03	1.40	1.44
2	B	502	HMG	O7-C3	-2.79	1.40	1.44
2	A	500	HMG	O2B-C2B	2.78	1.49	1.43
2	B	502	HMG	O2B-C2B	2.77	1.49	1.43
2	B	502	HMG	C5A-C4A	-2.77	1.34	1.39
2	A	500	HMG	C5A-C4A	-2.75	1.34	1.39
2	A	500	HMG	C1-S1P	2.56	1.82	1.76
2	B	502	HMG	C1-S1P	2.43	1.82	1.76
2	A	500	HMG	C5A-N7A	-2.43	1.34	1.39
2	A	500	HMG	O3B-C3B	-2.43	1.35	1.44
2	B	502	HMG	C5A-N7A	-2.42	1.34	1.39
2	B	502	HMG	O9P-C9P	-2.37	1.18	1.23
2	A	500	HMG	O9P-C9P	-2.35	1.18	1.23
2	B	502	HMG	O3B-C3B	-2.29	1.36	1.44
2	B	502	HMG	O5P-C5P	-2.29	1.18	1.23
2	A	500	HMG	O5P-C5P	-2.25	1.18	1.23
2	A	500	HMG	C8A-N9A	-2.11	1.34	1.37
2	B	502	HMG	C8A-N9A	-2.10	1.34	1.37

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HMG	C2-C1-S1P	5.98	121.11	113.56
2	A	500	HMG	N1A-C2A-N3A	-5.64	120.05	128.58
2	B	502	HMG	N1A-C2A-N3A	-5.45	120.33	128.58
2	B	502	HMG	C2-C1-S1P	5.33	120.28	113.56
2	A	500	HMG	C5A-C4A-N3A	-5.11	119.68	126.72
2	B	502	HMG	C5A-C4A-N3A	-4.74	120.19	126.72
2	B	502	HMG	N9A-C8A-N7A	-4.44	107.64	113.94
2	A	500	HMG	N9A-C8A-N7A	-4.29	107.85	113.94
2	A	500	HMG	C1B-N9A-C8A	-4.16	117.87	127.09
2	A	500	HMG	N6A-C6A-N1A	-3.97	109.54	118.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	HMG	N6A-C6A-N1A	-3.96	109.57	118.38
2	B	502	HMG	C6P-C5P-N4P	3.83	123.32	116.34
2	B	502	HMG	C1B-N9A-C8A	-3.62	119.05	127.09
2	A	500	HMG	C2A-N3A-C4A	3.49	120.36	111.83
2	B	502	HMG	C2A-N3A-C4A	3.26	119.81	111.83
2	A	500	HMG	N3A-C4A-N9A	3.16	132.54	127.17
2	B	502	HMG	C7P-C6P-C5P	-3.15	107.15	112.39
2	A	500	HMG	C4A-N9A-C1B	3.08	133.83	126.63
2	B	502	HMG	C5A-N7A-C8A	3.05	108.25	103.45
2	A	500	HMG	C5A-N7A-C8A	2.99	108.15	103.45
2	B	502	HMG	C5A-C6A-N6A	2.97	130.63	123.29
2	A	500	HMG	O2-C1-S1P	-2.93	118.95	122.68
2	A	500	HMG	C5A-C6A-N6A	2.82	130.27	123.29
2	B	502	HMG	C2P-S1P-C1	2.80	110.11	101.84
2	B	502	HMG	N3A-C4A-N9A	2.76	131.85	127.17
2	B	502	HMG	O5P-C5P-N4P	-2.75	117.64	123.03
2	B	502	HMG	CAP-C9P-N8P	2.56	121.34	116.48
2	B	502	HMG	C4A-N9A-C8A	2.55	108.42	105.74
2	B	502	HMG	C4A-N9A-C1B	2.52	132.53	126.63
2	B	502	HMG	C4A-C5A-N7A	-2.48	107.75	110.58
2	B	502	HMG	O2-C1-C2	-2.47	119.87	123.66
2	A	500	HMG	C4A-N9A-C8A	2.44	108.30	105.74
2	A	500	HMG	O2-C1-C2	-2.44	119.92	123.66
2	A	500	HMG	C2P-C3P-N4P	-2.34	107.53	112.41
2	A	500	HMG	C4A-C5A-N7A	-2.33	107.92	110.58
2	B	502	HMG	O2-C1-S1P	-2.23	119.85	122.68
2	B	502	HMG	O4-C5-C4	2.10	120.99	114.35

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500	HMG	CCP-O6A-P2A-O5A
2	A	500	HMG	CCP-O6A-P2A-O3A
2	A	500	HMG	C5B-O5B-P1A-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
2	B	502	HMG	C2-C1-S1P-C2P

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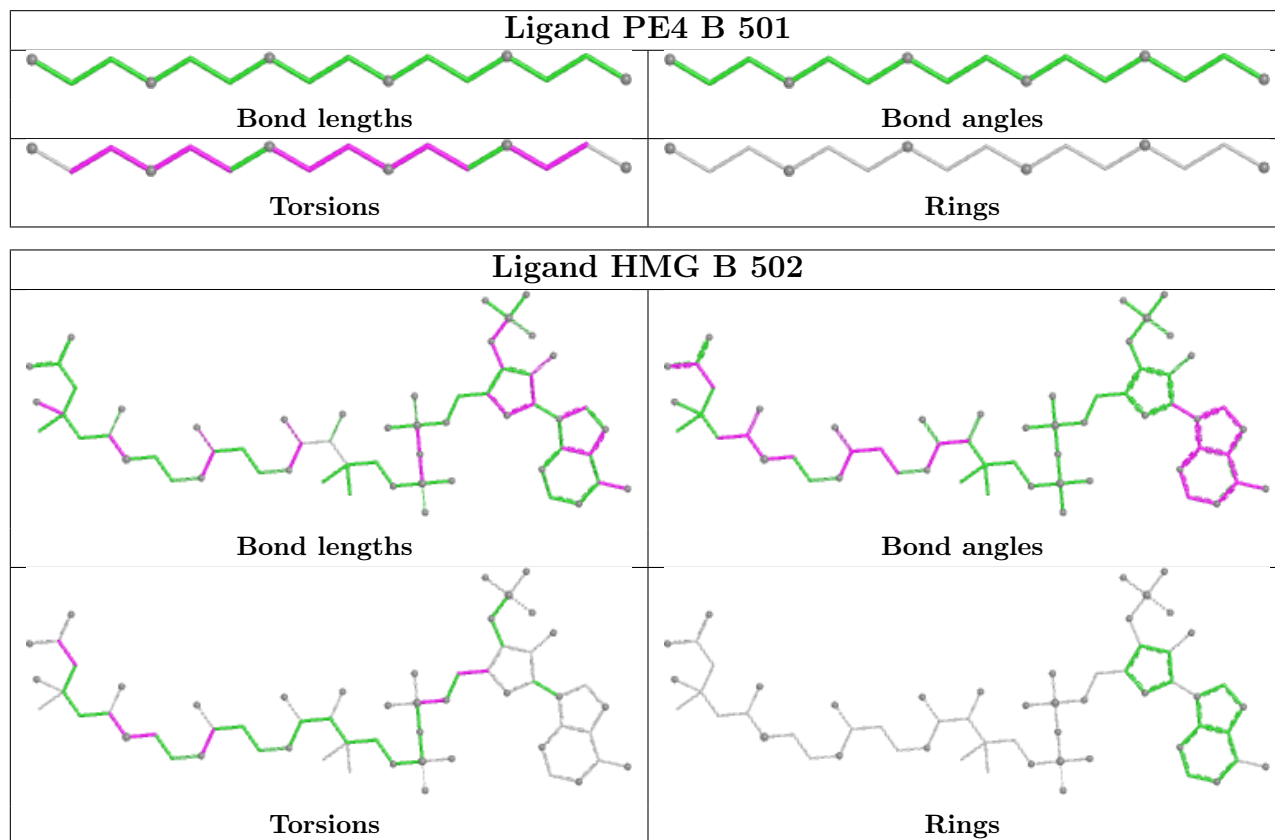
Mol	Chain	Res	Type	Atoms
3	A	501	GOL	O1-C1-C2-C3
3	B	503	GOL	O1-C1-C2-C3
3	B	503	GOL	O2-C2-C3-O3
2	B	502	HMG	C6P-C5P-N4P-C3P
4	B	501	PE4	C9-C10-O6-C11
4	B	501	PE4	O4-C7-C8-O5
2	B	502	HMG	O5P-C5P-N4P-C3P
4	B	501	PE4	O6-C10-C9-O5
4	B	505	PE4	O3-C5-C6-O4
4	B	505	PE4	O4-C7-C8-O5
4	B	501	PE4	O6-C11-C12-O7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

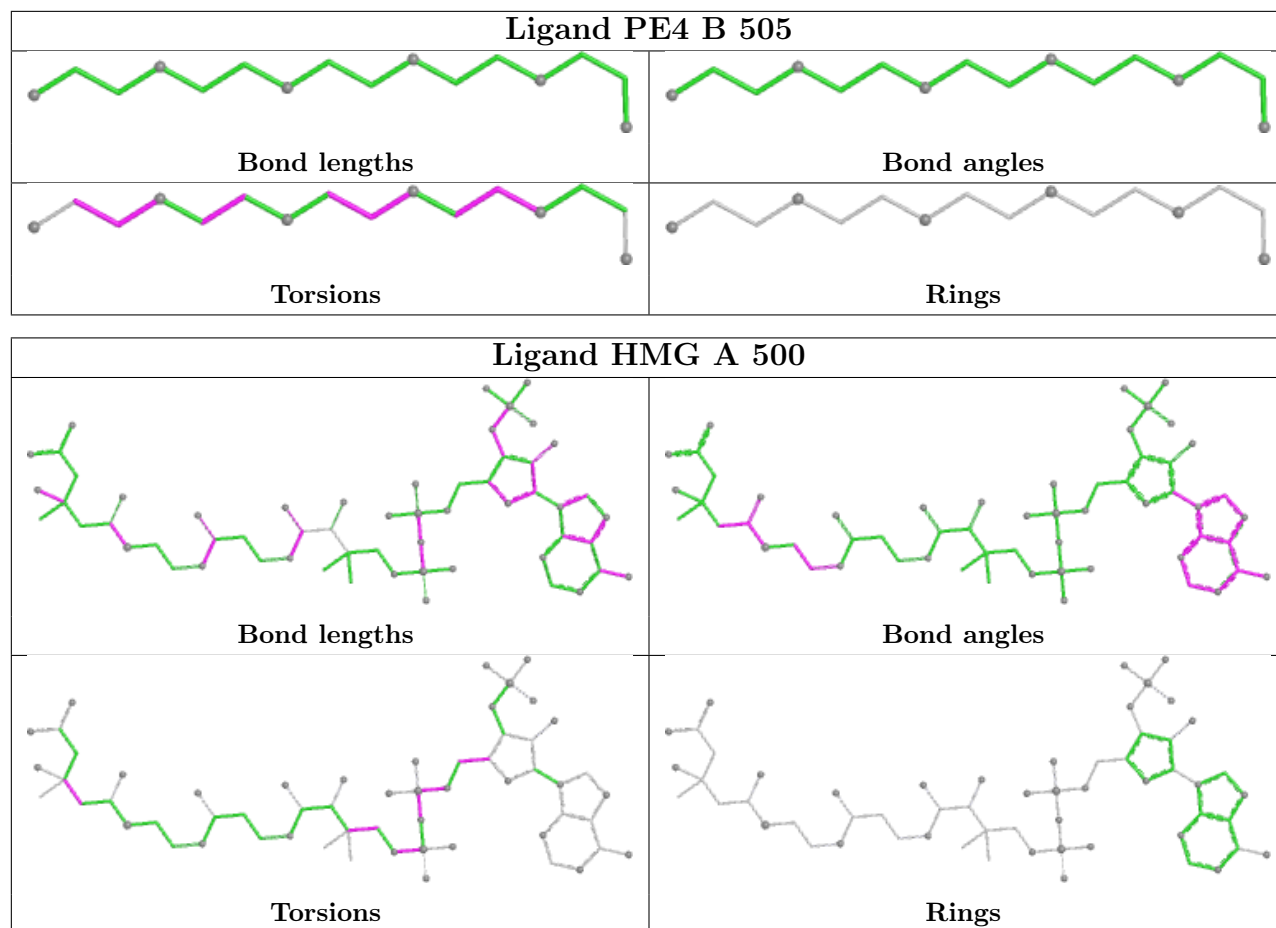
There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	501	PE4	1	0
2	B	502	HMG	2	0
4	B	505	PE4	1	0
3	B	503	GOL	2	0
2	A	500	HMG	1	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	381/426 (89%)	-1.62	0 <a href="#">100</a> <a href="#">100</a>	12, 24, 47, 82	1 (0%)
1	B	422/426 (99%)	-1.55	0 <a href="#">100</a> <a href="#">100</a>	13, 26, 59, 79	3 (0%)
All	All	803/852 (94%)	-1.58	0 <a href="#">100</a> <a href="#">100</a>	12, 25, 57, 82	4 (0%)

There are no RSRZ outliers to report.

### 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](#)

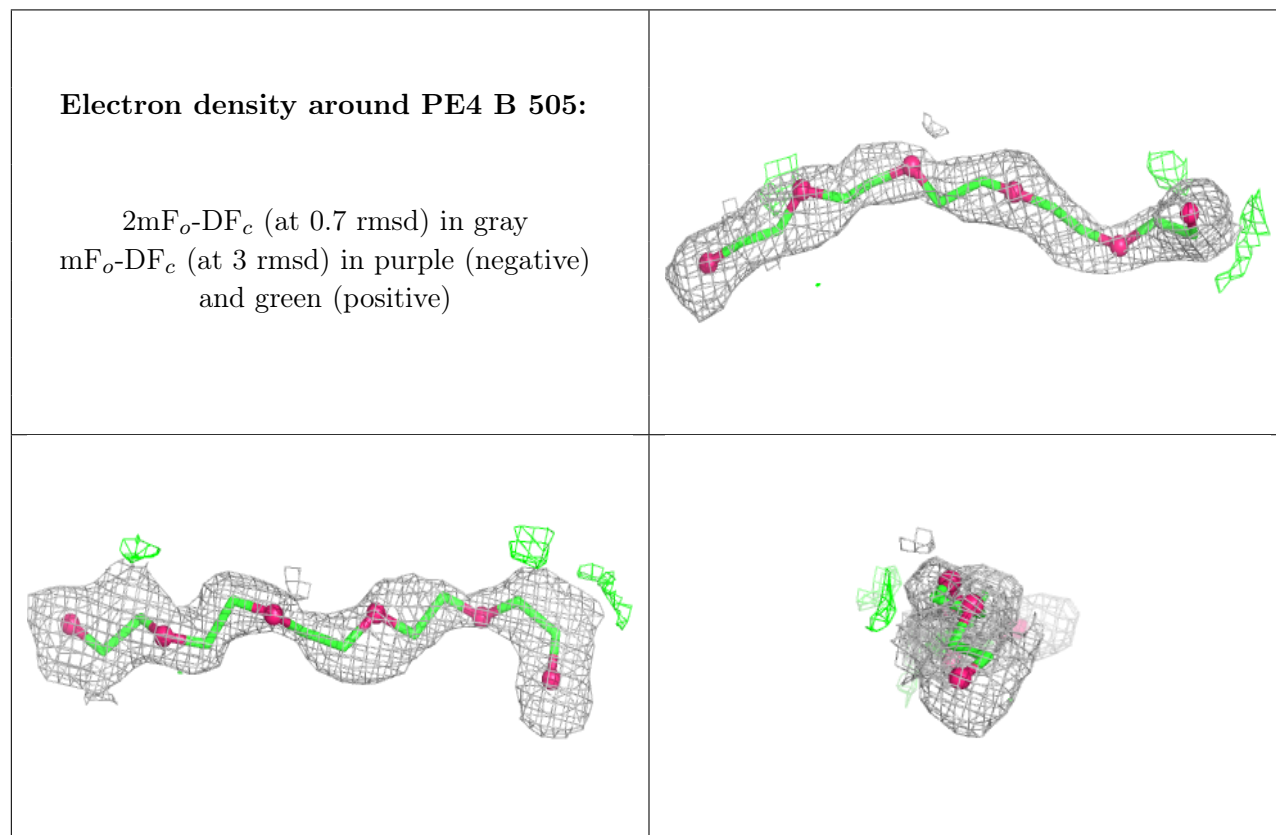
There are no oligosaccharides in this entry.

### 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, 95<sup>th</sup> 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.

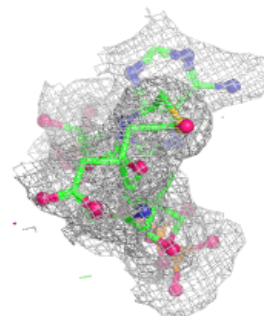
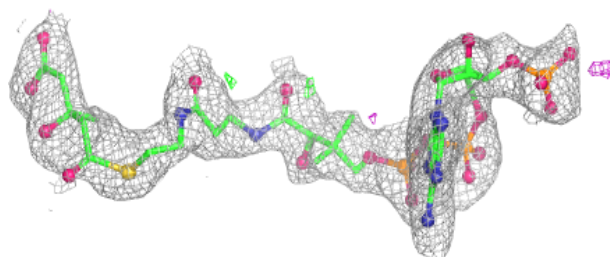
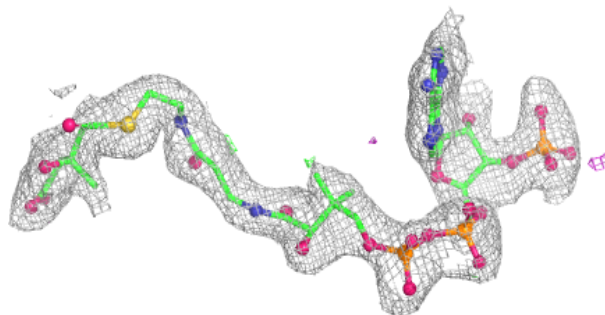
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	PE4	B	505	16/24	0.98	0.04	31,38,44,46	0
2	HMG	B	502	58/58	0.99	0.03	19,38,54,64	0
3	GOL	A	501	6/6	0.99	0.03	27,34,42,45	0
3	GOL	B	503	6/6	0.99	0.04	26,35,36,37	0
3	GOL	B	504	6/6	0.99	0.03	45,51,54,55	0
4	PE4	B	501	16/24	0.99	0.03	30,40,53,56	0
2	HMG	A	500	58/58	0.99	0.02	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.

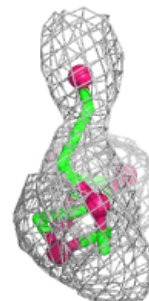
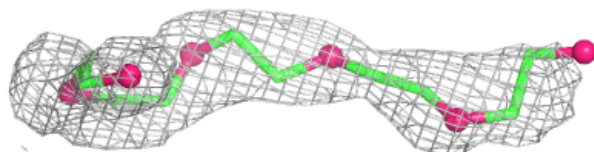
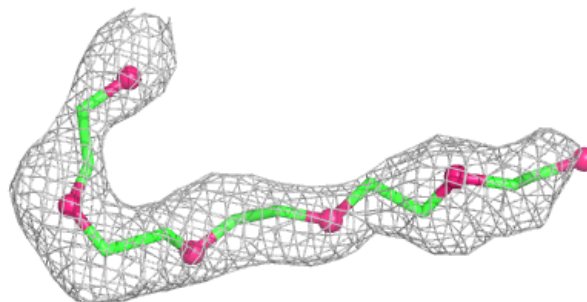


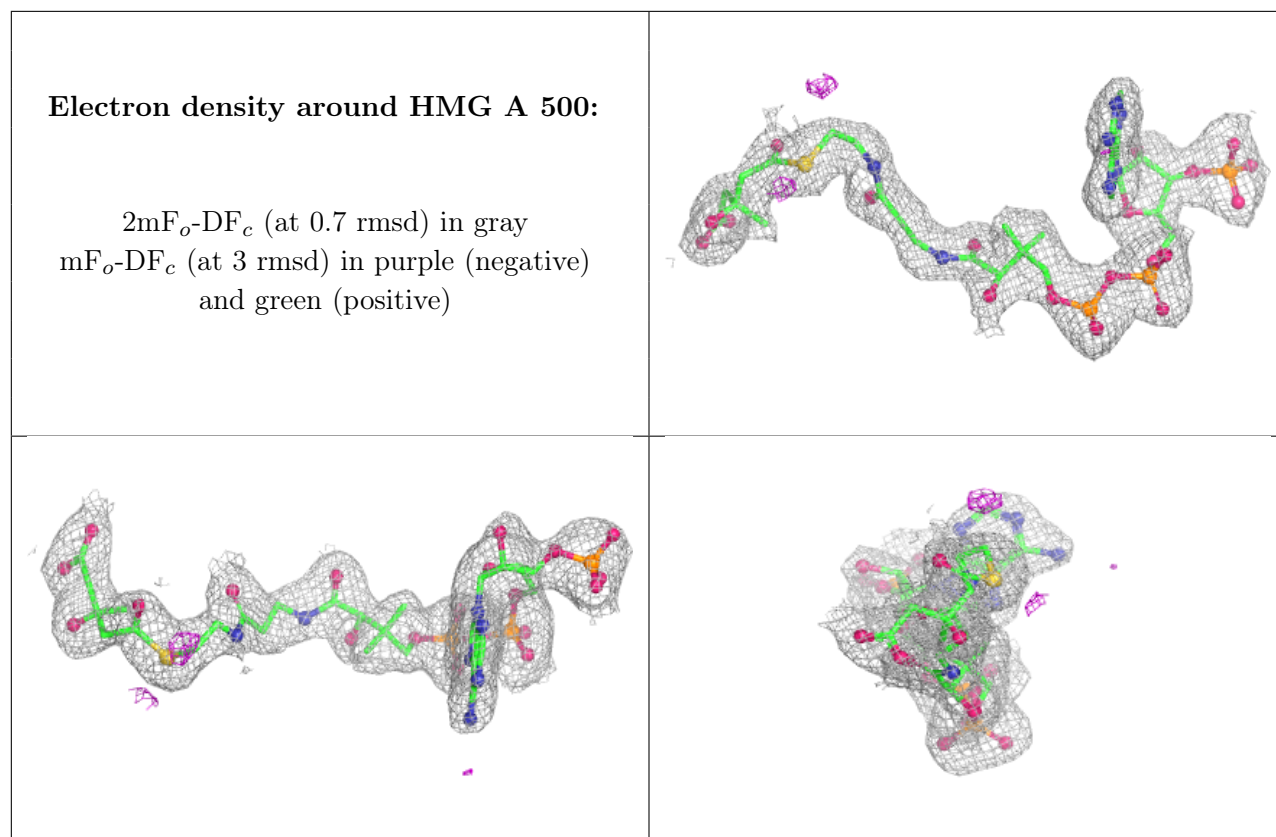
**Electron density around HMG B 502:**

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

**Electron density around 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)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.