



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:06 PM UTC

PDB ID : 2FA0 / pdb_00002fa0
Title : HMG-CoA synthase from Brassica juncea in complex with HMG-CoA and covalently bound to HMG-CoA
Authors : Pojer, F.; Ferrer, J.L.; Richard, S.B.; Noel, J.P.
Deposited on : 2005-12-06
Resolution : 2.49 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 types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (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

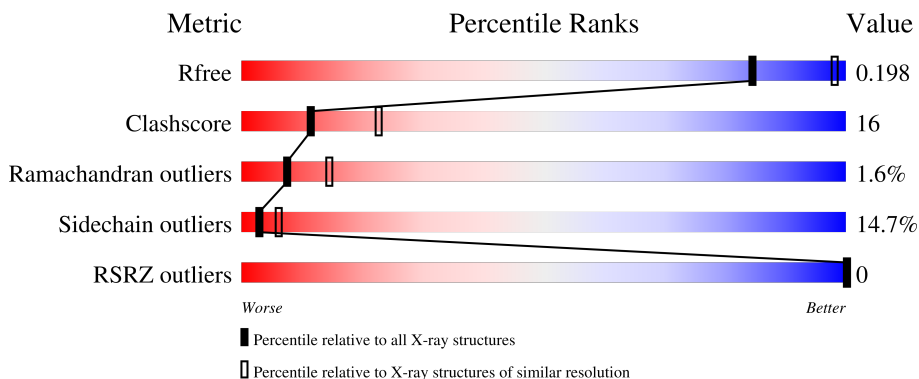
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.49 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	450	

2 Entry composition [i](#)

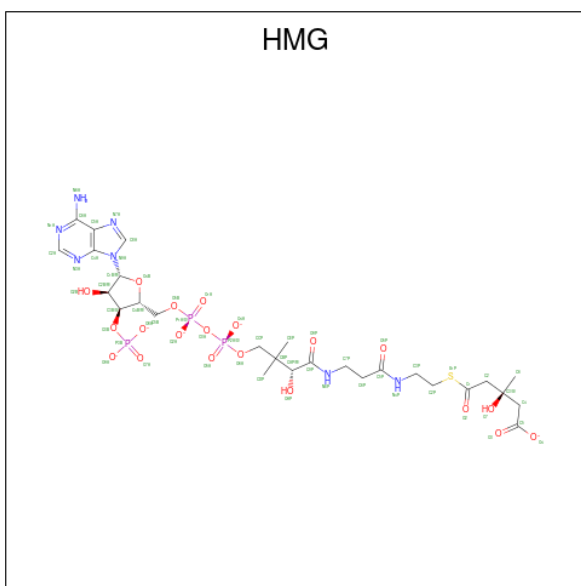
There are 3 unique types of molecules in this entry. The entry contains 3725 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 HMG-CoA synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	450	3510	2237	575	675	23	0	0	0

- Molecule 2 is 3-HYDROXY-3-METHYLGLUTARYL-COENZYME A (CCD ID: HMG) (formula: $C_{27}H_{39}N_7O_{20}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	A	1	116	54	14	40	6	2	0	1

- Molecule 3 is water.

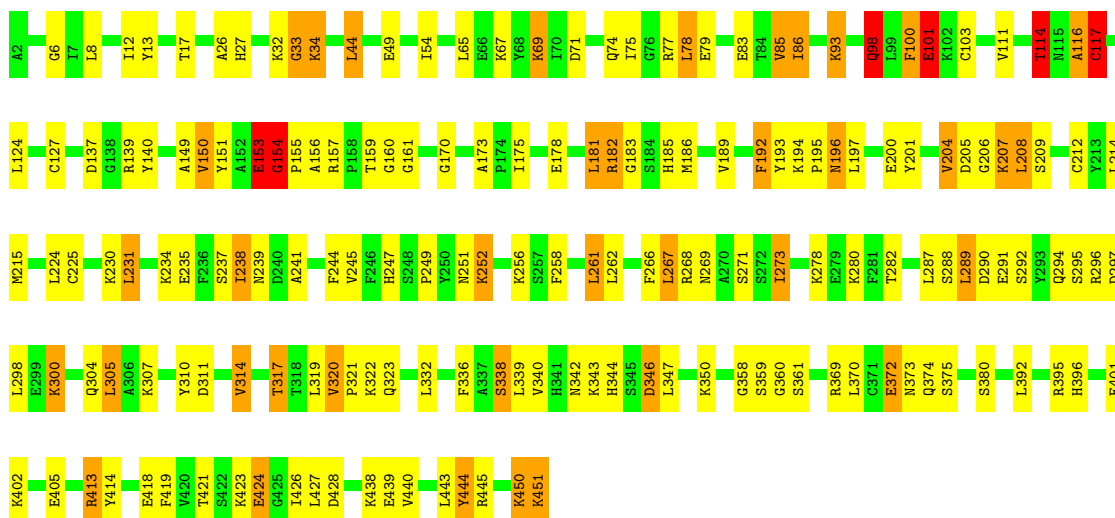
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	99	Total	O	0	0
			99	99		

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: HMG-CoA synthase

Chain A:  61% 29% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	61.33Å 61.33Å 411.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	68.52 – 2.49 68.52 – 2.49	Depositor EDS
% Data completeness (in resolution range)	97.7 (68.52-2.49) 97.7 (68.52-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.48Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.188 , 0.272 0.191 , 0.198	Depositor DCC
R_{free} test set	859 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtrriage
Anisotropy	0.228	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3725	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.31	6/3586 (0.2%)	1.44	39/4844 (0.8%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	ALA	C-N	26.69	1.73	1.33
1	A	85	VAL	CA-CB	7.04	1.62	1.54
1	A	124	LEU	CA-C	-5.84	1.45	1.52
1	A	155	PRO	CA-C	5.78	1.57	1.52
1	A	273	ILE	CA-CB	5.77	1.60	1.53
1	A	196	ASN	CG-OD1	5.49	1.33	1.23

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	ALA	CA-C-N	18.87	156.54	122.38
1	A	116	ALA	C-N-CA	18.87	156.54	122.38
1	A	116	ALA	O-C-N	-15.77	101.61	122.59
1	A	314	VAL	N-CA-C	8.42	119.05	111.81
1	A	117	CYS	N-CA-C	-8.16	103.03	113.16
1	A	114	THR	CB-CA-C	-7.85	94.37	109.37
1	A	201	TYR	N-CA-C	7.66	119.29	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ALA	CA-C-N	-7.31	112.27	119.87
1	A	173	ALA	C-N-CA	-7.31	112.27	119.87
1	A	85	VAL	N-CA-C	7.26	118.33	107.80
1	A	153	GLU	N-CA-C	6.88	125.44	110.80
1	A	34	LYS	N-CA-C	6.63	119.07	111.11
1	A	127	CYS	N-CA-C	6.42	118.35	111.36
1	A	439	GLU	N-CA-C	6.30	116.89	108.38
1	A	71	ASP	N-CA-C	-6.02	102.07	109.65
1	A	360	GLY	N-CA-C	6.01	123.05	114.64
1	A	54	ILE	CB-CA-C	-5.98	104.19	112.02
1	A	160	GLY	N-CA-C	-5.92	103.23	112.58
1	A	212	CYS	N-CA-C	-5.67	106.40	113.55
1	A	154	GLY	CA-C-N	-5.67	115.05	120.83
1	A	154	GLY	C-N-CA	-5.67	115.05	120.83
1	A	67	LYS	CA-C-N	-5.65	112.85	122.56
1	A	67	LYS	C-N-CA	-5.65	112.85	122.56
1	A	161	GLY	N-CA-C	-5.63	102.53	110.75
1	A	320	VAL	CA-C-N	-5.63	112.90	119.32
1	A	320	VAL	C-N-CA	-5.63	112.90	119.32
1	A	261	LEU	N-CA-C	-5.50	104.92	111.03
1	A	295	SER	N-CA-C	5.50	117.48	108.52
1	A	231	LEU	N-CA-C	5.39	117.59	111.02
1	A	423	LYS	N-CA-C	5.34	116.96	111.03
1	A	317	THR	N-CA-C	-5.34	106.44	113.17
1	A	150	VAL	CB-CA-C	5.31	120.00	111.29
1	A	33	GLY	N-CA-C	5.29	121.80	113.86
1	A	194	LYS	CA-C-N	5.28	124.98	119.64
1	A	194	LYS	C-N-CA	5.28	124.98	119.64
1	A	98	GLN	N-CA-C	5.17	117.59	111.33
1	A	204	VAL	CB-CA-C	-5.16	103.00	110.63
1	A	342	ASN	N-CA-C	5.15	116.71	111.14
1	A	86	ILE	N-CA-C	-5.12	107.41	111.81

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	114	THR	Peptide
1	A	117	CYS	Mainchain
1	A	149	ALA	Peptide
1	A	450	LYS	Peptide

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	3510	0	3456	115	0
2	A	116	0	77	18	0
3	A	99	0	0	12	0
All	All	3725	0	3533	115	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 16.

All (115) 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:A:116:ALA:C	1:A:117:CYS:N	1.73	1.45
1:A:249:PRO:HG3	2:A:500[A]:HMG:C2P	2.17	0.74
1:A:249:PRO:HG3	2:A:500[A]:HMG:H2P2	1.70	0.72
1:A:288:SER:O	1:A:291:GLU:HB2	1.88	0.72
1:A:249:PRO:CG	2:A:500[A]:HMG:H2P2	2.19	0.72
1:A:288:SER:HB3	1:A:291:GLU:OE1	1.90	0.71
1:A:116:ALA:HB1	2:A:500[A]:HMG:O3	1.91	0.70
1:A:33:GLY:HA3	2:A:500[A]:HMG:N1A	2.06	0.70
1:A:185:HIS:HD2	3:A:597:HOH:O	1.75	0.70
1:A:343:LYS:O	1:A:347:LEU:HG	1.91	0.69
1:A:287:LEU:HD22	1:A:291:GLU:HB3	1.76	0.67
1:A:114:THR:HG21	3:A:585:HOH:O	1.93	0.67
1:A:247:HIS:CD2	2:A:500[A]:HMG:O2	2.48	0.66
1:A:311:ASP:HB3	3:A:564:HOH:O	1.94	0.66
1:A:372:GLU:OE1	1:A:372:GLU:HA	1.94	0.66
1:A:49:GLU:OE2	1:A:413:ARG:HD2	1.96	0.66
1:A:358:GLY:HA2	2:A:500[A]:HMG:O3	1.97	0.64
1:A:424:GLU:H	1:A:424:GLU:CD	2.05	0.64
1:A:8:LEU:HD11	1:A:170:GLY:HA3	1.80	0.63
1:A:93:LYS:C	1:A:93:LYS:HD3	2.23	0.63
1:A:252:LYS:CE	2:A:500[A]:HMG:O8A	2.47	0.63
1:A:252:LYS:HE2	2:A:500[A]:HMG:O8A	1.99	0.63
1:A:450:LYS:O	1:A:451:LYS:HB3	1.99	0.62
1:A:200:GLU:HG3	1:A:414:TYR:CD1	2.34	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ASP:OD2	1:A:208:LEU:HD23	1.99	0.61
1:A:396:HIS:HB3	3:A:589:HOH:O	1.98	0.61
1:A:258:PHE:CD2	1:A:314:VAL:CG2	2.84	0.61
1:A:307:LYS:O	1:A:311:ASP:OD1	2.20	0.60
1:A:340:VAL:O	1:A:344:HIS:HB3	2.02	0.60
1:A:317:THR:HG22	1:A:339:LEU:HB2	1.83	0.59
1:A:116:ALA:O	1:A:117:CYS:N	2.32	0.58
1:A:373:ASN:CG	1:A:374:GLN:H	2.12	0.58
1:A:289:LEU:HA	1:A:292:SER:OG	2.03	0.57
1:A:237:SER:OG	1:A:239:ASN:HB2	2.05	0.57
1:A:74:GLN:O	1:A:139:ARG:HB3	2.06	0.56
1:A:49:GLU:HG2	1:A:421:THR:HG22	1.87	0.56
1:A:34:LYS:NZ	2:A:500[B]:HMG:O2A	2.34	0.56
1:A:247:HIS:NE2	2:A:500[A]:HMG:O2	2.39	0.56
1:A:339:LEU:HD12	1:A:339:LEU:O	2.05	0.56
1:A:402:LYS:HE2	3:A:598:HOH:O	2.05	0.56
1:A:193:TYR:CE1	1:A:195:PRO:HG3	2.42	0.55
1:A:114:THR:CG2	3:A:585:HOH:O	2.53	0.55
1:A:297:ASP:HA	1:A:300:LYS:HD3	1.87	0.55
1:A:323:GLN:OE1	1:A:395:ARG:NH1	2.40	0.55
1:A:451:LYS:HG2	1:A:451:LYS:O	2.07	0.54
1:A:78:LEU:C	1:A:78:LEU:HD23	2.33	0.54
1:A:209:SER:O	2:A:500[B]:HMG:H2P1	2.09	0.53
1:A:413:ARG:HD3	1:A:419:PHE:CE1	2.43	0.53
1:A:116:ALA:CB	2:A:500[A]:HMG:O3	2.56	0.52
1:A:266:PHE:HD1	1:A:267:LEU:HD23	1.75	0.52
1:A:183:GLY:HA2	3:A:555:HOH:O	2.10	0.52
1:A:238:ILE:CD1	1:A:258:PHE:CE2	2.93	0.51
1:A:320:VAL:HG23	1:A:338:SER:HB3	1.92	0.51
1:A:289:LEU:HG	1:A:290:ASP:H	1.75	0.51
1:A:401:GLU:HB2	3:A:535:HOH:O	2.09	0.51
1:A:182:ARG:NH2	3:A:501:HOH:O	2.44	0.51
1:A:392:LEU:HB3	1:A:395:ARG:HH21	1.76	0.51
1:A:156:ALA:O	1:A:159:THR:OG1	2.30	0.50
1:A:192:PHE:CD2	1:A:192:PHE:C	2.89	0.50
1:A:137:ASP:OD2	1:A:139:ARG:HD3	2.11	0.50
1:A:78:LEU:C	1:A:78:LEU:CD2	2.86	0.48
1:A:178:GLU:CG	1:A:369:ARG:HG2	2.44	0.48
1:A:443:LEU:HD12	1:A:445:ARG:HH11	1.78	0.47
1:A:100:PHE:CD1	1:A:100:PHE:N	2.82	0.47
1:A:69:LYS:HB2	1:A:69:LYS:HE2	1.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:VAL:O	1:A:321:PRO:C	2.55	0.47
1:A:401:GLU:OE1	1:A:401:GLU:HA	2.12	0.47
1:A:249:PRO:CG	2:A:500[A]:HMG:C2P	2.87	0.47
1:A:44:LEU:C	1:A:44:LEU:HD23	2.41	0.46
1:A:235:GLU:HG3	3:A:539:HOH:O	2.15	0.46
1:A:266:PHE:CD1	1:A:267:LEU:HD23	2.51	0.46
1:A:443:LEU:HD12	1:A:445:ARG:NH1	2.31	0.46
1:A:258:PHE:CD2	1:A:314:VAL:HG21	2.50	0.46
1:A:443:LEU:HB2	1:A:445:ARG:HD3	1.98	0.46
1:A:238:ILE:HD12	1:A:258:PHE:CZ	2.51	0.45
1:A:418:GLU:HA	1:A:440:VAL:O	2.16	0.45
1:A:98:GLN:O	1:A:101:GLU:HB2	2.16	0.45
1:A:296:ARG:O	1:A:297:ASP:C	2.58	0.45
1:A:200:GLU:HG3	1:A:414:TYR:CE1	2.50	0.45
1:A:225:CYS:SG	1:A:268:ARG:NH1	2.90	0.45
1:A:225:CYS:HB3	3:A:539:HOH:O	2.16	0.44
1:A:153:GLU:O	1:A:154:GLY:O	2.35	0.44
1:A:289:LEU:C	1:A:291:GLU:N	2.75	0.44
1:A:33:GLY:HA3	2:A:500[B]:HMG:N1A	2.33	0.43
1:A:77:ARG:HD3	1:A:79:GLU:OE2	2.19	0.43
1:A:249:PRO:HG3	2:A:500[A]:HMG:H2P1	1.97	0.43
1:A:310:TYR:CD1	1:A:314:VAL:HB	2.53	0.43
1:A:428:ASP:OD1	1:A:428:ASP:N	2.50	0.43
1:A:247:HIS:CD2	1:A:249:PRO:HD3	2.54	0.43
1:A:49:GLU:OE2	1:A:413:ARG:CD	2.66	0.43
1:A:249:PRO:CB	2:A:500[A]:HMG:H2P2	2.49	0.43
1:A:258:PHE:O	1:A:261:LEU:HB2	2.19	0.43
1:A:426:ILE:O	1:A:427:LEU:C	2.62	0.43
1:A:207:LYS:HZ3	1:A:208:LEU:HD13	1.83	0.42
1:A:241:ALA:HB3	1:A:244:PHE:CE1	2.55	0.42
1:A:151:TYR:HB2	1:A:157:ARG:HA	2.01	0.42
1:A:346:ASP:O	1:A:350:LYS:NZ	2.44	0.42
1:A:6:GLY:HA3	1:A:175:ILE:O	2.20	0.42
1:A:181:LEU:HD22	1:A:224:LEU:HG	1.99	0.42
1:A:224:LEU:HD23	1:A:224:LEU:C	2.45	0.41
1:A:359:SER:C	1:A:361:SER:HA	2.45	0.41
1:A:192:PHE:CD1	1:A:204:VAL:HG22	2.55	0.41
1:A:245:VAL:HG12	1:A:332:LEU:HD12	2.02	0.41
1:A:12:ILE:HG12	1:A:13:TYR:N	2.35	0.41
1:A:26:ALA:O	1:A:27:HIS:C	2.63	0.41
1:A:289:LEU:HD23	1:A:289:LEU:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLU:OE2	2:A:500[A]:HMG:O4	2.39	0.41
1:A:322:LYS:NZ	3:A:580:HOH:O	2.50	0.41
1:A:175:ILE:HD13	1:A:336:PHE:CZ	2.56	0.40
1:A:193:TYR:CD1	1:A:195:PRO:HG3	2.56	0.40
1:A:238:ILE:CD1	1:A:258:PHE:CZ	3.04	0.40
1:A:252:LYS:HG3	1:A:256:LYS:HE3	2.03	0.40
1:A:75:ILE:HD13	1:A:140:TYR:HB2	2.04	0.40
1:A:280:LYS:HD3	1:A:305:LEU:HD11	2.03	0.40
1:A:251:ASN:O	1:A:252:LYS:C	2.63	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	448/450 (100%)	407 (91%)	34 (8%)	7 (2%)	7 14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	153	GLU
1	A	444	TYR
1	A	154	GLY
1	A	206	GLY
1	A	32	LYS
1	A	101	GLU
1	A	111	VAL

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	381/381 (100%)	325 (85%)	56 (15%)	3 6

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	THR
1	A	44	LEU
1	A	65	LEU
1	A	69	LYS
1	A	78	LEU
1	A	85	VAL
1	A	86	ILE
1	A	93	LYS
1	A	98	GLN
1	A	100	PHE
1	A	101	GLU
1	A	103	CYS
1	A	114	THR
1	A	150	VAL
1	A	181	LEU
1	A	182	ARG
1	A	186	MET
1	A	189	VAL
1	A	192	PHE
1	A	196	ASN
1	A	197	LEU
1	A	207	LYS
1	A	208	LEU
1	A	214	LEU
1	A	215	MET
1	A	230	LYS
1	A	231	LEU
1	A	234	LYS
1	A	238	ILE
1	A	252	LYS

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Mol	Chain	Res	Type
1	A	262	LEU
1	A	267	LEU
1	A	269	ASN
1	A	271	SER
1	A	273	ILE
1	A	278	LYS
1	A	282	THR
1	A	289	LEU
1	A	294	GLN
1	A	298	LEU
1	A	300	LYS
1	A	304	GLN
1	A	305	LEU
1	A	319	LEU
1	A	338	SER
1	A	346	ASP
1	A	370	LEU
1	A	372	GLU
1	A	375	SER
1	A	380	SER
1	A	405	GLU
1	A	413	ARG
1	A	424	GLU
1	A	438	LYS
1	A	444	TYR
1	A	451	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	21	GLN
1	A	185	HIS
1	A	264	ASN
1	A	269	ASN
1	A	304	GLN

5.3.3 RNA

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](#)

2 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
2	HMG	A	500[B]	-	58,60,60	1.44	7 (12%)	81,90,90	2.56	19 (23%)
2	HMG	A	500[A]	1	58,60,60	1.63	9 (15%)	81,90,90	2.09	18 (22%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HMG	A	500[B]	-	-	26/60/77/77	0/3/3/3
2	HMG	A	500[A]	1	-	15/60/77/77	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500[A]	HMG	O3-C5	6.31	1.42	1.22
2	A	500[B]	HMG	P2A-O3A	5.44	1.65	1.59
2	A	500[B]	HMG	P1A-O3A	5.40	1.65	1.59
2	A	500[A]	HMG	P1A-O3A	4.69	1.64	1.59
2	A	500[A]	HMG	P2A-O3A	4.12	1.63	1.59
2	A	500[A]	HMG	C1-S1P	-3.73	1.67	1.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500[A]	HMG	O4-C5	3.34	1.41	1.30
2	A	500[B]	HMG	C1-S1P	-3.16	1.68	1.76
2	A	500[A]	HMG	C4-C3	-3.06	1.51	1.54
2	A	500[A]	HMG	C2-C3	-2.53	1.52	1.54
2	A	500[A]	HMG	C5A-N7A	-2.29	1.34	1.39
2	A	500[B]	HMG	C5A-N7A	-2.26	1.35	1.39
2	A	500[B]	HMG	C6-C3	2.26	1.55	1.52
2	A	500[B]	HMG	C8A-N7A	2.08	1.35	1.31
2	A	500[B]	HMG	O7-C3	2.06	1.47	1.44
2	A	500[A]	HMG	C4A-N9A	-2.06	1.33	1.37

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500[B]	HMG	C2-C1-S1P	12.02	128.73	113.56
2	A	500[B]	HMG	O2-C1-S1P	-7.02	113.75	122.68
2	A	500[A]	HMG	C3P-N4P-C5P	6.79	135.46	122.82
2	A	500[B]	HMG	C6-C3-C2	6.50	119.24	111.30
2	A	500[B]	HMG	C6-C3-C4	-6.33	103.56	111.30
2	A	500[B]	HMG	C7P-N8P-C9P	6.19	133.66	122.55
2	A	500[A]	HMG	N1A-C2A-N3A	-6.00	119.50	128.58
2	A	500[A]	HMG	C7P-N8P-C9P	5.87	133.09	122.55
2	A	500[B]	HMG	N1A-C2A-N3A	-5.16	120.76	128.58
2	A	500[B]	HMG	O6A-CCP-CBP	5.07	118.70	110.55
2	A	500[B]	HMG	C5A-C4A-N3A	-4.84	120.05	126.72
2	A	500[A]	HMG	N9A-C8A-N7A	-4.61	107.40	113.94
2	A	500[A]	HMG	C2-C1-S1P	4.51	119.25	113.56
2	A	500[A]	HMG	C2P-S1P-C1	4.44	114.97	101.84
2	A	500[A]	HMG	C5A-C4A-N3A	-4.21	120.91	126.72
2	A	500[A]	HMG	O4-C5-O3	-4.07	112.87	123.33
2	A	500[B]	HMG	O2-C1-C2	-4.03	117.48	123.66
2	A	500[A]	HMG	C4A-N9A-C8A	3.89	109.82	105.74
2	A	500[A]	HMG	C2A-N3A-C4A	3.61	120.65	111.83
2	A	500[B]	HMG	N9A-C8A-N7A	-3.56	108.89	113.94
2	A	500[B]	HMG	C2A-N3A-C4A	3.38	120.09	111.83
2	A	500[B]	HMG	N3A-C4A-N9A	3.29	132.77	127.17
2	A	500[A]	HMG	O2-C1-C2	-3.16	118.82	123.66
2	A	500[B]	HMG	CEP-CBP-CAP	3.10	114.05	108.77
2	A	500[A]	HMG	C5A-N7A-C8A	3.01	108.18	103.45
2	A	500[A]	HMG	N3A-C4A-N9A	2.97	132.22	127.17
2	A	500[B]	HMG	C5A-N7A-C8A	2.66	107.62	103.45
2	A	500[B]	HMG	C4A-N9A-C8A	2.63	108.50	105.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500[A]	HMG	C4A-C5A-N7A	-2.49	107.73	110.58
2	A	500[A]	HMG	O2A-P1A-O3A	2.47	113.95	107.27
2	A	500[B]	HMG	C4A-C5A-N7A	-2.43	107.81	110.58
2	A	500[B]	HMG	C2P-S1P-C1	2.25	108.51	101.84
2	A	500[A]	HMG	O3B-C3B-C4B	-2.16	102.41	110.03
2	A	500[A]	HMG	C2A-N1A-C6A	2.15	122.26	118.73
2	A	500[B]	HMG	O2A-P1A-O3A	2.10	112.94	107.27
2	A	500[B]	HMG	O7-C3-C4	2.09	115.59	109.05
2	A	500[A]	HMG	C4A-N9A-C1B	-2.01	121.92	126.63

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	500[A]	HMG	C5B-O5B-P1A-O1A
2	A	500[A]	HMG	C5B-O5B-P1A-O2A
2	A	500[A]	HMG	O4B-C4B-C5B-O5B
2	A	500[A]	HMG	C2P-C3P-N4P-C5P
2	A	500[A]	HMG	C2-C3-C4-C5
2	A	500[A]	HMG	O7-C3-C4-C5
2	A	500[A]	HMG	C6-C3-C4-C5
2	A	500[B]	HMG	C5B-O5B-P1A-O3A
2	A	500[B]	HMG	C5B-O5B-P1A-O1A
2	A	500[B]	HMG	C5B-O5B-P1A-O2A
2	A	500[B]	HMG	C9P-CAP-CBP-CEP
2	A	500[B]	HMG	N8P-C9P-CAP-OAP
2	A	500[B]	HMG	O2-C1-S1P-C2P
2	A	500[B]	HMG	C2-C1-S1P-C2P
2	A	500[B]	HMG	C1-C2-C3-O7
2	A	500[B]	HMG	C1-C2-C3-C6
2	A	500[B]	HMG	C1-C2-C3-C4
2	A	500[A]	HMG	C3B-C4B-C5B-O5B
2	A	500[B]	HMG	C2B-C1B-N9A-C4A
2	A	500[B]	HMG	C2B-C1B-N9A-C8A
2	A	500[B]	HMG	O9P-C9P-CAP-OAP
2	A	500[B]	HMG	O4B-C4B-C5B-O5B
2	A	500[A]	HMG	C6P-C7P-N8P-C9P
2	A	500[B]	HMG	OAP-CAP-CBP-CDP
2	A	500[B]	HMG	OAP-CAP-CBP-CEP
2	A	500[B]	HMG	C9P-CAP-CBP-CDP
2	A	500[B]	HMG	CEP-CBP-CCP-O6A
2	A	500[B]	HMG	C9P-CAP-CBP-CCP

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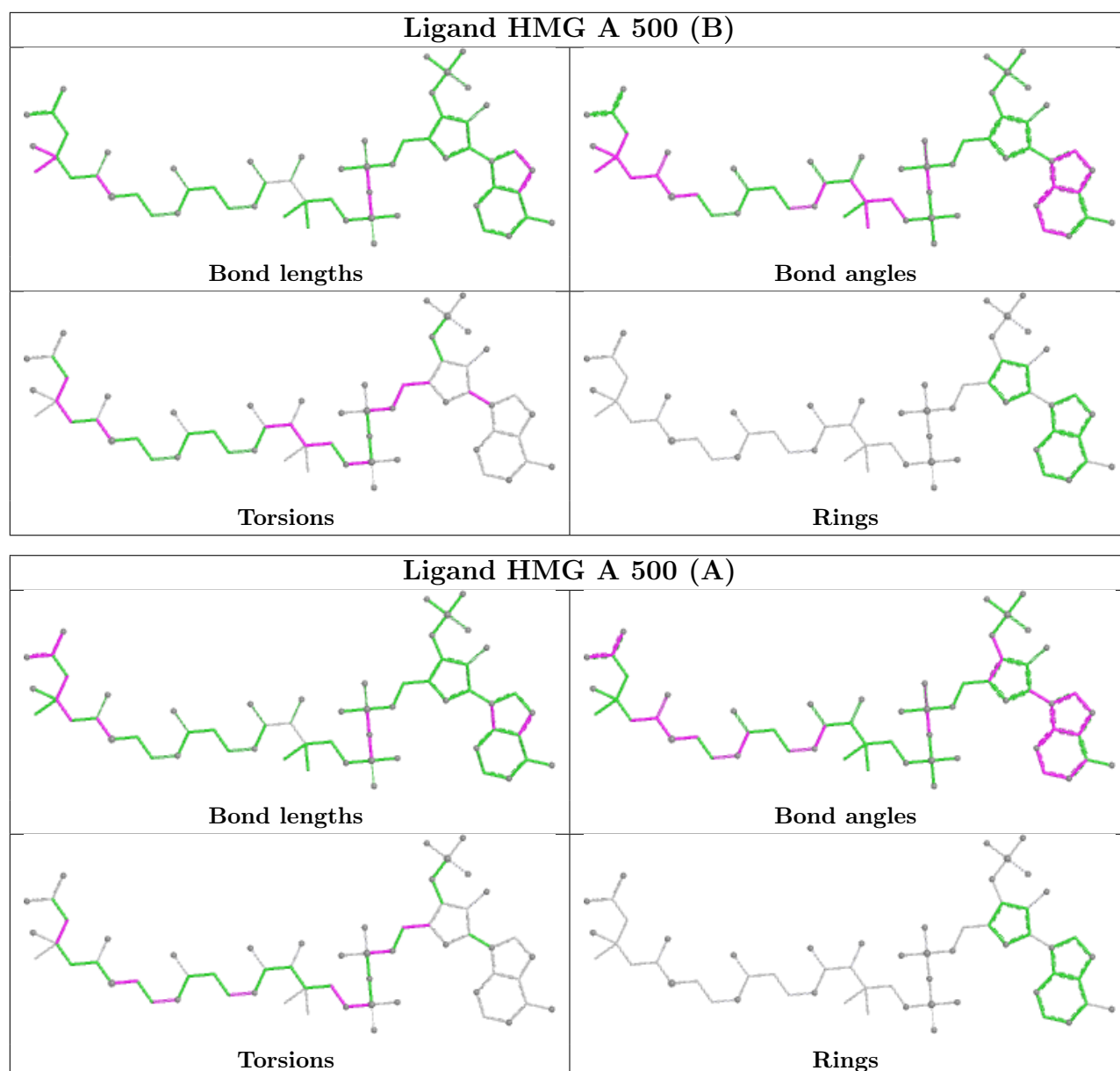
Mol	Chain	Res	Type	Atoms
2	A	500[A]	HMG	C3P-C2P-S1P-C1
2	A	500[A]	HMG	CCP-O6A-P2A-O5A
2	A	500[A]	HMG	CCP-O6A-P2A-O4A
2	A	500[A]	HMG	CCP-O6A-P2A-O3A
2	A	500[A]	HMG	C5B-O5B-P1A-O3A
2	A	500[B]	HMG	CCP-O6A-P2A-O5A
2	A	500[B]	HMG	CCP-O6A-P2A-O4A
2	A	500[B]	HMG	CCP-O6A-P2A-O3A
2	A	500[B]	HMG	C4B-C5B-O5B-P1A
2	A	500[B]	HMG	C2-C3-C4-C5
2	A	500[B]	HMG	C6-C3-C4-C5
2	A	500[A]	HMG	CBP-CCP-O6A-P2A
2	A	500[B]	HMG	CDP-CBP-CCP-O6A

There are no ring outliers.

2 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500[B]	HMG	3	0
2	A	500[A]	HMG	15	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	116:ALA	C	117:CYS	N	1.73

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, 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	450/450 (100%)	-0.34	0 100 100	2, 27, 55, 68	1 (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](#)

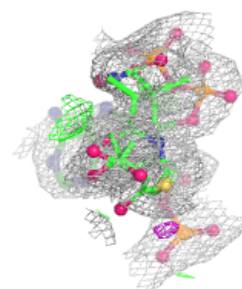
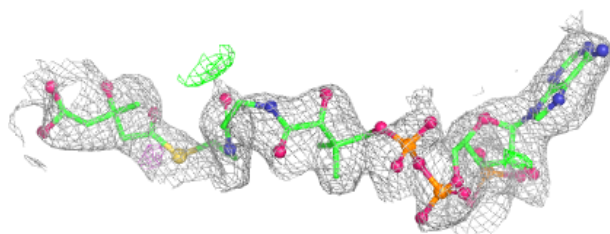
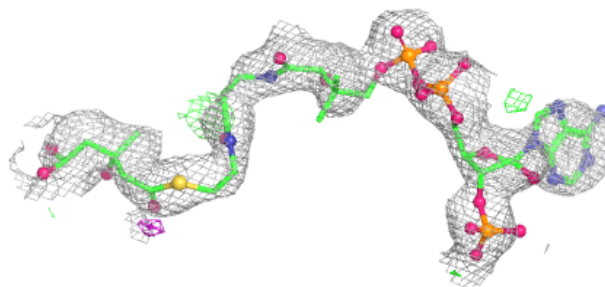
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	HMG	A	500[A]	58/58	0.93	0.09	16,28,37,38	58
2	HMG	A	500[B]	58/58	0.93	0.09	2,40,59,60	58

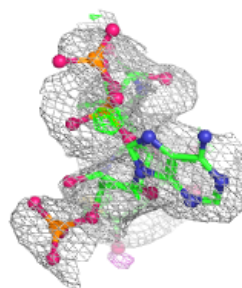
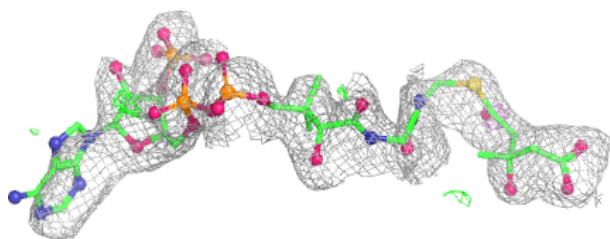
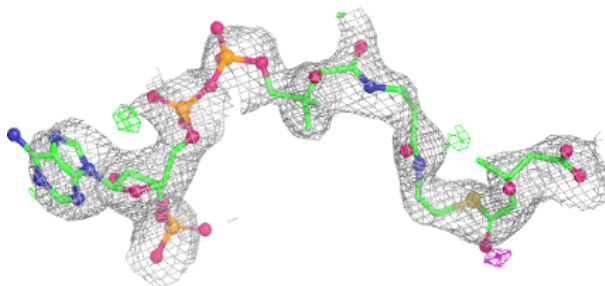
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HMG A 500 (A):

$2mF_o-DF_c$ (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 (B):**

$2mF_o-DF_c$ (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.