



Full wwPDB EM Validation Report ⓘ

Aug 6, 2026 – 05:17 PM JST

PDB ID : 8HBH / pdb_00008hbh
EMDB ID : EMD-34632
Title : Structure of human soluble guanylate cyclase in the NO-activated state at 3.1 angstrom
Authors : Chen, L.; Liu, R.
Deposited on : 2022-10-28
Resolution : 3.10 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 ⓘ 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:

EMDB validation analysis : 0.0.1.dev133
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

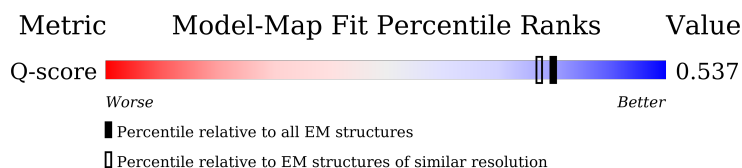
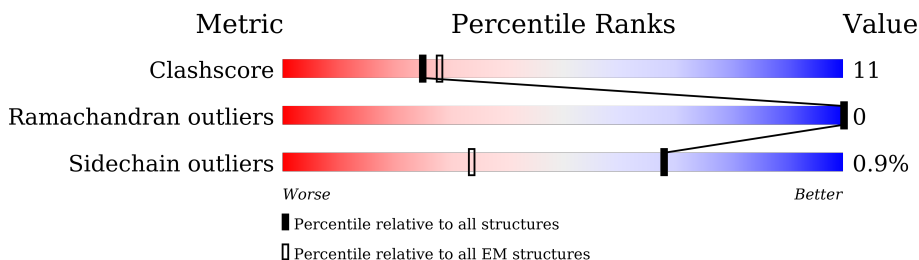
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.10 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14724 (2.60 - 3.60)

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 $\leq 5\%$. 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	690	7% 58% 17% 24%
2	B	619	71% 22% 7%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8502 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 Guanylate cyclase soluble subunit alpha-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	522	Total	C	N	O	S	0	0
			3922	2544	662	686	30		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	44	MET	VAL	variant	UNP Q02108
A	554	VAL	LEU	variant	UNP Q02108

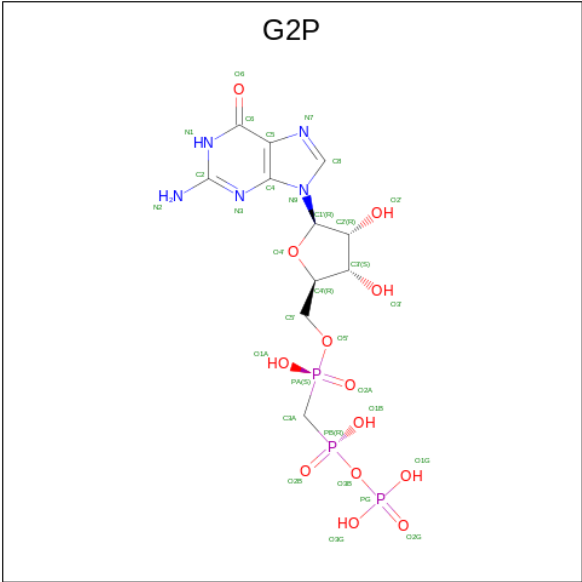
- Molecule 2 is a protein called Guanylate cyclase soluble subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	577	Total	C	N	O	S	3	0
			4501	2880	783	810	28		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

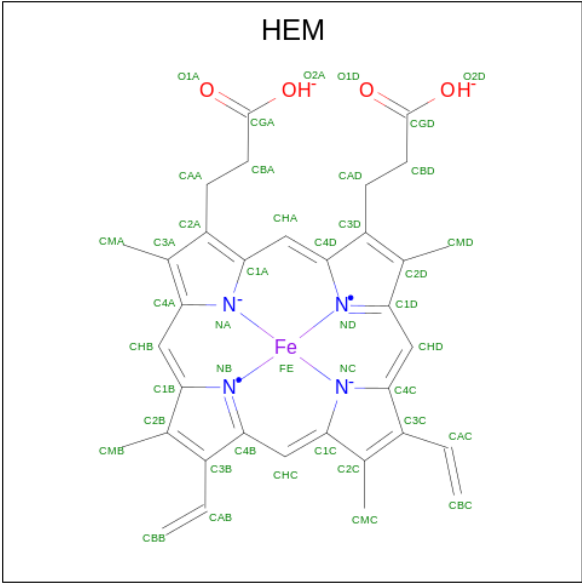
Mol	Chain	Residues	Atoms		AltConf
3	A	2	Total	Mg	0
			2	2	

- Molecule 4 is PHOSPHOMETHYLPHOSPHONIC ACID GUANYLATE ESTER (CCD ID: G2P) (formula: C₁₁H₁₈N₅O₁₃P₃).



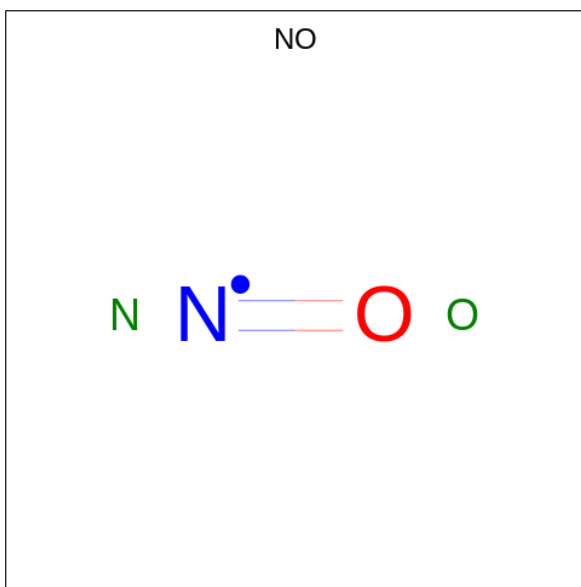
Mol	Chain	Residues	Atoms					AltConf
4	B	1	Total	C	N	O	P	0
			32	11	5	13	3	

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					AltConf
5	B	1	Total	C	Fe	N	O	0
			43	34	1	4	4	

- Molecule 6 is NITRIC OXIDE (CCD ID: NO) (formula: NO) (labeled as "Ligand of Interest" by depositor).

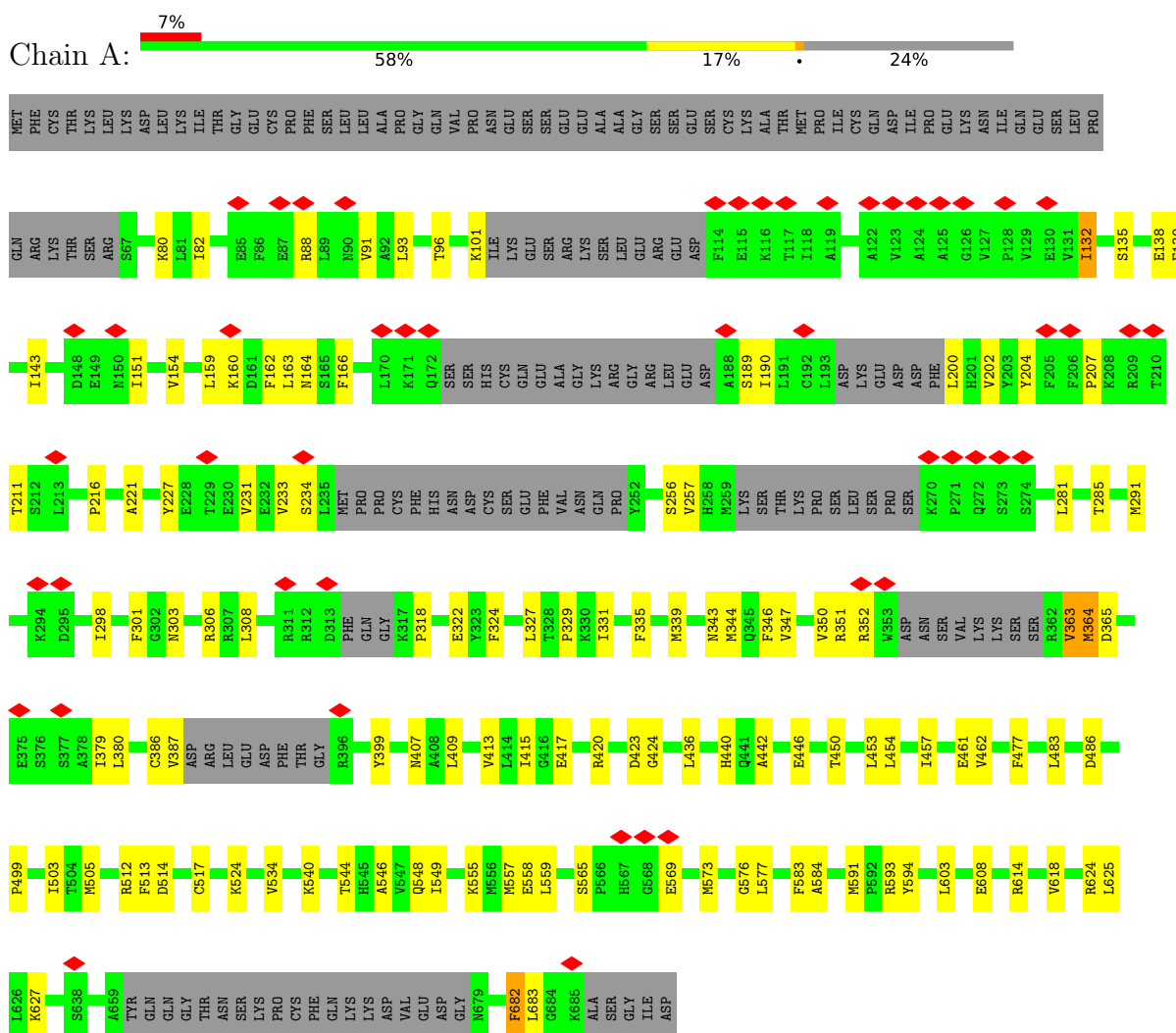


Mol	Chain	Residues	Atoms			AltConf
			Total	N	O	
6	B	1	2	1	1	0

3 Residue-property plots

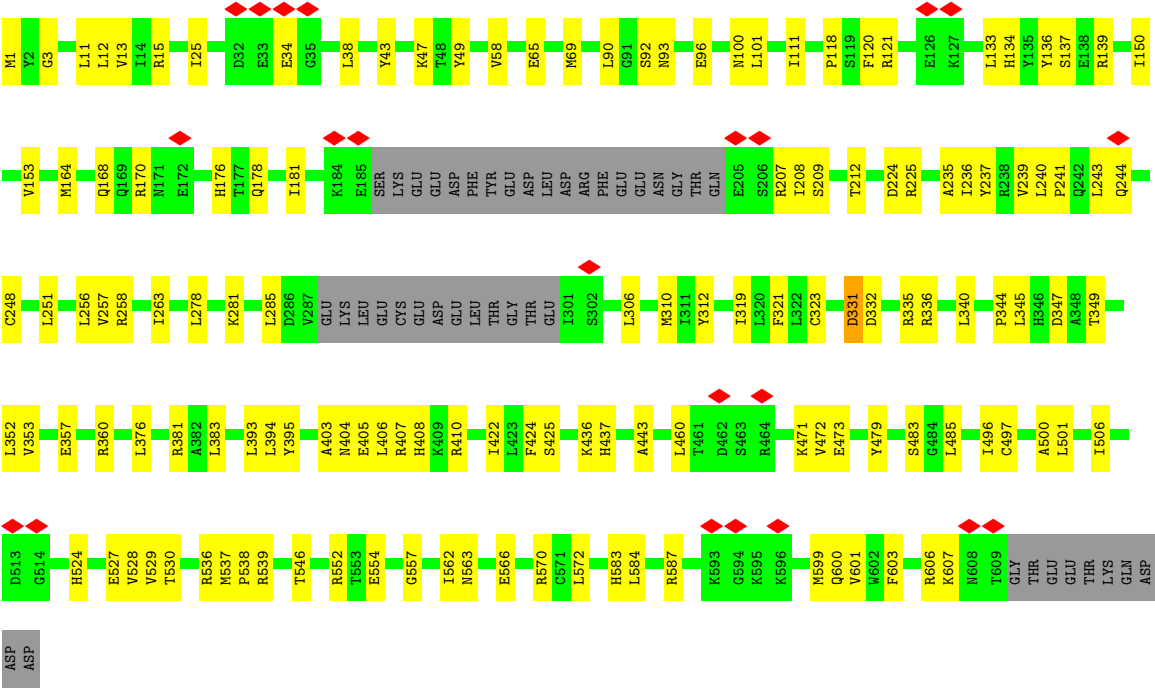
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Guanylate cyclase soluble subunit alpha-1



- Molecule 2: Guanylate cyclase soluble subunit beta-1





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	970628	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	11.848	Depositor
Minimum map value	-7.497	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.158	Depositor
Recommended contour level	1	Depositor
Map size ($\text{\AA}$)	267.52, 267.52, 267.52	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.045, 1.045, 1.045	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: NO, G2P, MG, HEM

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	0.31	0/3996	0.45	3/5409 (0.1%)
2	B	0.25	0/4602	0.40	0/6237
All	All	0.28	0/8598	0.42	3/11646 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	363	VAL	N-CA-C	8.44	118.52	110.42
1	A	682	PHE	N-CA-C	6.11	122.19	114.31
1	A	364	MET	N-CA-C	5.66	116.86	108.60

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	3922	0	3876	89	0
2	B	4501	0	4486	98	0
3	A	2	0	0	0	0
4	B	32	0	14	3	0
5	B	43	0	30	9	0
6	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8502	0	8406	178	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 11.

All (178) 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)
5:B:1002:HEM:HBC2	5:B:1002:HEM:HHD	1.32	1.04
1:A:163:LEU:HD12	1:A:190:ILE:HG21	1.69	0.74
2:B:583:HIS:HB2	2:B:607:LYS:HB2	1.72	0.72
1:A:577:LEU:HD12	1:A:618:VAL:HG22	1.74	0.69
2:B:251:LEU:O	2:B:281:LYS:NZ	2.25	0.69
1:A:555:LYS:NZ	1:A:558:GLU:OE1	2.26	0.67
2:B:65:GLU:O	2:B:69:MET:HG3	1.94	0.67
2:B:121:ARG:NH1	2:B:136:TYR:OH	2.28	0.67
2:B:137:SER:O	2:B:170:ARG:NH1	2.28	0.66
2:B:164:MET:HG2	2:B:181:ILE:HG12	1.77	0.66
1:A:387:VAL:HG11	1:A:413:VAL:HG22	1.79	0.65
1:A:135:SER:O	1:A:139:GLU:HG2	1.96	0.64
1:A:546:ALA:HB3	1:A:625:LEU:HD22	1.78	0.64
2:B:344:PRO:HG2	2:B:347:ASP:HB2	1.79	0.64
2:B:473:GLU:OE1	4:B:1001:G2P:N1	2.27	0.64
2:B:557:GLY:HA2	2:B:587:ARG:HH22	1.63	0.63
5:B:1002:HEM:HHD	5:B:1002:HEM:CBC	2.20	0.63
2:B:405:GLU:OE1	2:B:410:ARG:NH1	2.32	0.62
1:A:343:ASN:OD1	1:A:344:MET:N	2.33	0.61
1:A:204:TYR:HE2	1:A:207:PRO:HA	1.66	0.61
1:A:80:LYS:HD2	1:A:88:ARG:HH21	1.66	0.61
5:B:1002:HEM:HBC2	5:B:1002:HEM:CHD	2.12	0.60
2:B:424:PHE:HB3	2:B:554:GLU:HG2	1.84	0.60
1:A:190:ILE:HG23	1:A:202:VAL:HG13	1.83	0.60
2:B:471:LYS:HE3	2:B:479:TYR:CE2	2.37	0.60
1:A:442:ALA:O	1:A:446:GLU:HG2	2.03	0.58
1:A:350:VAL:CG2	1:A:364:MET:HG2	2.33	0.58
2:B:483:SER:HB2	2:B:496:ILE:HG12	1.86	0.58
2:B:118:PRO:O	2:B:258:ARG:NH2	2.37	0.58
1:A:138:GLU:HA	1:A:221:ALA:HB2	1.86	0.58
1:A:513:PHE:HE2	1:A:573:MET:HE1	1.66	0.58
2:B:257:VAL:HG12	2:B:258:ARG:HG3	1.86	0.57
2:B:90:LEU:O	2:B:100:ASN:ND2	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:483:LEU:HB3	1:A:577:LEU:HD23	1.86	0.56
1:A:204:TYR:CE2	1:A:207:PRO:HA	2.40	0.56
1:A:423:ASP:OD1	1:A:424:GLY:N	2.39	0.56
5:B:1002:HEM:CMB	5:B:1002:HEM:HBB2	2.36	0.56
2:B:524:HIS:HD2	2:B:546:THR:HG23	1.71	0.56
2:B:422:ILE:HD11	2:B:528:VAL:HG11	1.88	0.55
2:B:25:ILE:HG12	2:B:58:VAL:HG11	1.87	0.55
1:A:298:ILE:HD12	1:A:318:PRO:HB2	1.89	0.55
2:B:224:ASP:OD1	2:B:225:ARG:N	2.34	0.55
2:B:566:GLU:HB2	2:B:600:GLN:HB2	1.88	0.55
1:A:483:LEU:HD23	1:A:549:ILE:HG23	1.90	0.54
2:B:101:LEU:HD11	5:B:1002:HEM:HMC3	1.89	0.54
2:B:552:ARG:NH2	4:B:1001:G2P:O2A	2.42	0.53
1:A:82:ILE:O	1:A:211:THR:OG1	2.24	0.53
1:A:350:VAL:HG21	1:A:364:MET:HG2	1.90	0.53
2:B:381:ARG:HD2	2:B:381:ARG:O	2.09	0.53
1:A:327:LEU:HB2	1:A:347:VAL:HG23	1.91	0.53
1:A:499:PRO:HB2	2:B:529:VAL:HG23	1.90	0.53
1:A:216:PRO:HA	1:A:233:VAL:HG11	1.91	0.52
2:B:587:ARG:HH21	2:B:601:VAL:HG11	1.75	0.52
2:B:256:LEU:HD22	2:B:263:ILE:HB	1.90	0.52
1:A:234:SER:O	1:A:256:SER:HB3	2.09	0.52
2:B:168:GLN:HB2	2:B:178:GLN:HB2	1.91	0.52
1:A:415:ILE:HD12	2:B:111:ILE:HD11	1.93	0.51
2:B:331:ASP:O	2:B:335:ARG:HG3	2.10	0.51
2:B:34:GLU:OE2	2:B:47:LYS:NZ	2.42	0.51
2:B:239:VAL:HG12	2:B:285:LEU:HD22	1.92	0.51
2:B:256:LEU:HA	2:B:278:LEU:HA	1.91	0.51
2:B:552:ARG:HH22	4:B:1001:G2P:H4'	1.76	0.51
2:B:563:ASN:OD1	2:B:587:ARG:NH2	2.44	0.51
1:A:132:ILE:HD12	1:A:132:ILE:H	1.77	0.50
2:B:120:PHE:HD2	2:B:133:LEU:HD21	1.75	0.50
1:A:96:THR:HG1	1:A:135:SER:HG	1.57	0.50
1:A:291:MET:HA	1:A:379:ILE:O	2.11	0.50
1:A:91:VAL:HG12	1:A:143:ILE:HD11	1.93	0.49
1:A:454:LEU:HD13	2:B:393:LEU:HD23	1.93	0.49
2:B:566:GLU:OE2	2:B:570:ARG:NH2	2.44	0.49
1:A:524:LYS:HA	1:A:534:VAL:HG12	1.94	0.49
1:A:308:LEU:HD13	1:A:363:VAL:HG11	1.94	0.49
1:A:231:VAL:HG21	1:A:257:VAL:HG13	1.95	0.49
2:B:241:PRO:HA	2:B:244:GLN:HG3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:TYR:CE2	2:B:92:SER:HB3	2.48	0.49
2:B:353:VAL:O	2:B:357:GLU:HG2	2.13	0.49
1:A:513:PHE:CE2	1:A:573:MET:HE1	2.47	0.48
2:B:501:LEU:HD13	2:B:606:ARG:HG3	1.95	0.48
1:A:162:PHE:HB2	1:A:227:TYR:OH	2.12	0.48
1:A:512:ARG:HB3	1:A:559:LEU:HD13	1.95	0.48
2:B:111:ILE:O	2:B:360:ARG:NH2	2.40	0.48
1:A:591:MET:HE2	2:B:472:VAL:O	2.14	0.48
2:B:237:TYR:HE2	2:B:244:GLN:HA	1.78	0.48
1:A:343:ASN:O	2:B:345:LEU:HD13	2.14	0.47
2:B:136:TYR:CD1	2:B:176:HIS:HB3	2.49	0.47
1:A:160:LYS:O	1:A:164:ASN:ND2	2.48	0.47
1:A:483:LEU:HA	1:A:576:GLY:O	2.14	0.47
1:A:524:LYS:HE2	2:B:537:MET:HB3	1.96	0.47
2:B:12:LEU:HD11	2:B:69:MET:HB2	1.96	0.47
2:B:349:THR:HA	2:B:352:LEU:HB3	1.96	0.47
1:A:322:GLU:O	1:A:351:ARG:NH1	2.49	0.46
1:A:324:PHE:CD1	1:A:350:VAL:HG12	2.50	0.46
1:A:450:THR:HG23	2:B:394:LEU:HD22	1.97	0.46
1:A:624:ARG:HA	1:A:627:LYS:HD3	1.96	0.46
2:B:436:LYS:HD3	2:B:437:HIS:CD2	2.50	0.46
2:B:93:ASN:HB3	2:B:96:GLU:HG2	1.97	0.46
2:B:395:TYR:OH	2:B:404:ASN:OD1	2.24	0.46
1:A:303:ASN:O	1:A:306:ARG:HG2	2.15	0.46
2:B:120:PHE:CD2	2:B:133:LEU:HD21	2.50	0.46
2:B:235:ALA:O	2:B:239:VAL:HG23	2.16	0.46
1:A:281:LEU:O	1:A:285:THR:HG22	2.15	0.46
1:A:524:LYS:O	2:B:537:MET:HE3	2.16	0.46
1:A:151:ILE:O	1:A:154:VAL:HG12	2.16	0.45
1:A:486:ASP:HB2	1:A:608:GLU:OE2	2.16	0.45
2:B:497:CYS:SG	2:B:572:LEU:HD11	2.55	0.45
2:B:524:HIS:CE1	2:B:599:MET:HE1	2.51	0.45
1:A:163:LEU:HA	1:A:166:PHE:CD1	2.51	0.45
2:B:321:PHE:CE2	2:B:323:CYS:HB3	2.51	0.45
2:B:332:ASP:O	2:B:336:ARG:HD2	2.17	0.45
1:A:505:MET:CE	1:A:565:SER:HA	2.47	0.45
2:B:310:MET:HE2	2:B:321:PHE:CD1	2.51	0.45
1:A:512:ARG:HG2	1:A:512:ARG:NH1	2.33	0.44
2:B:13:VAL:HG11	2:B:25:ILE:HD12	1.99	0.44
2:B:437:HIS:CD2	2:B:437:HIS:N	2.85	0.44
2:B:485:LEU:HD21	2:B:530:THR:HG21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:PHE:HE2	1:A:584:ALA:HB3	1.82	0.44
2:B:164:MET:HE2	2:B:164:MET:HB3	1.63	0.44
1:A:517:CYS:CB	2:B:536:ARG:HH21	2.31	0.44
1:A:93:LEU:HA	1:A:96:THR:HG22	1.99	0.44
2:B:460:LEU:HD21	2:B:506:ILE:HD12	1.99	0.44
1:A:593:ARG:HG2	2:B:539:ARG:CZ	2.47	0.44
2:B:136:TYR:CE1	2:B:176:HIS:HB3	2.53	0.44
1:A:101:LYS:HE2	1:A:101:LYS:HB3	1.72	0.44
1:A:291:MET:HE1	2:B:207:ARG:HG3	2.00	0.44
2:B:3:GLY:HA3	2:B:38:LEU:O	2.17	0.43
1:A:557:MET:HE1	1:A:614:ARG:HA	1.99	0.43
2:B:237:TYR:CE2	2:B:244:GLN:HA	2.52	0.43
2:B:527:GLU:OE1	2:B:527:GLU:N	2.37	0.43
2:B:208:ILE:HD12	2:B:208:ILE:N	2.32	0.43
1:A:159:LEU:HD23	1:A:163:LEU:HD23	2.00	0.43
1:A:407:ASN:C	1:A:409:LEU:H	2.27	0.43
5:B:1002:HEM:HBB2	5:B:1002:HEM:HMB1	2.00	0.43
1:A:514:ASP:OD1	1:A:514:ASP:N	2.51	0.43
1:A:682:PHE:CE2	2:B:538:PRO:HG3	2.53	0.43
1:A:80:LYS:HD2	1:A:88:ARG:NH2	2.32	0.43
1:A:335:PHE:CE1	1:A:339:MET:HE2	2.53	0.43
2:B:406:LEU:HD21	2:B:538:PRO:HB3	2.00	0.43
2:B:150:ILE:HA	2:B:153:VAL:HG12	2.01	0.43
2:B:209:SER:HB3	2:B:212:THR:HG23	2.01	0.42
1:A:603:LEU:HD12	1:A:603:LEU:HA	1.90	0.42
1:A:417:GLU:OE1	1:A:420:ARG:NH2	2.53	0.42
2:B:134:HIS:HB3	2:B:176:HIS:HD2	1.85	0.42
2:B:584:LEU:HA	2:B:603:PHE:O	2.18	0.42
5:B:1002:HEM:CMB	5:B:1002:HEM:CBB	2.98	0.42
1:A:200:LEU:HB2	1:A:257:VAL:HB	2.01	0.42
1:A:365:ASP:HB3	1:A:386:CYS:SG	2.60	0.42
2:B:243:LEU:HD23	2:B:248:CYS:SG	2.60	0.42
2:B:408:HIS:O	2:B:410:ARG:N	2.52	0.42
2:B:306:LEU:HB3	2:B:323:CYS:HB2	2.01	0.42
1:A:132:ILE:HA	1:A:135:SER:OG	2.20	0.41
1:A:440:HIS:CE1	2:B:383:LEU:HD22	2.54	0.41
1:A:583:PHE:CE1	2:B:443:ALA:HB3	2.55	0.41
2:B:25:ILE:CG1	2:B:58:VAL:HG11	2.49	0.41
5:B:1002:HEM:HMB1	5:B:1002:HEM:CBB	2.50	0.41
2:B:500:ALA:HB1	2:B:562:ILE:HD13	2.01	0.41
1:A:436:LEU:HD23	2:B:376:LEU:HB3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:565:SER:OG	1:A:569:GLU:N	2.53	0.41
2:B:312:TYR:HD1	2:B:319:ILE:HG22	1.84	0.41
1:A:352:ARG:HB2	1:A:363:VAL:HB	2.03	0.41
1:A:461:GLU:HG2	1:A:462:VAL:N	2.36	0.41
2:B:403:ALA:O	2:B:407:ARG:HB2	2.20	0.41
1:A:308:LEU:HD22	1:A:364:MET:HE3	2.03	0.41
1:A:453:LEU:O	1:A:457:ILE:HG13	2.21	0.41
2:B:1:MET:O	2:B:43:TYR:N	2.45	0.41
2:B:11:LEU:O	2:B:15:ARG:HG3	2.21	0.41
2:B:101:LEU:HD11	5:B:1002:HEM:CMC	2.51	0.41
2:B:236:ILE:HG23	2:B:240:LEU:HD12	2.03	0.41
1:A:364:MET:HE2	1:A:364:MET:HB3	1.99	0.40
1:A:82:ILE:HD13	1:A:82:ILE:HA	1.87	0.40
1:A:329:PRO:HG2	1:A:346:PHE:CD2	2.56	0.40
1:A:291:MET:HG2	1:A:380:LEU:HA	2.02	0.40
1:A:512:ARG:HG2	1:A:512:ARG:HH11	1.87	0.40
1:A:544:THR:O	1:A:548:GLN:HG3	2.21	0.40
1:A:189:SER:O	1:A:190:ILE:HG13	2.20	0.40
1:A:462:VAL:HG21	1:A:594:TYR:CE1	2.56	0.40
2:B:49:TYR:OH	2:B:139:ARG:HB3	2.21	0.40
2:B:340:LEU:HD12	2:B:340:LEU:HA	1.92	0.40
2:B:471:LYS:HE3	2:B:479:TYR:HE2	1.84	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	502/690 (73%)	485 (97%)	17 (3%)	0	100	100
2	B	574/619 (93%)	557 (97%)	17 (3%)	0	100	100
All	All	1076/1309 (82%)	1042 (97%)	34 (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 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	404/613 (66%)	398 (98%)	6 (2%)	57	75
2	B	487/555 (88%)	485 (100%)	2 (0%)	84	86
All	All	891/1168 (76%)	883 (99%)	8 (1%)	68	80

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

Mol	Chain	Res	Type
1	A	132	ILE
1	A	301	PHE
1	A	331	ILE
1	A	503	ILE
1	A	540	LYS
1	A	683	LEU
2	B	331	ASP
2	B	425	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	143	GLN
2	B	231	GLN
2	B	260	HIS
2	B	437	HIS
2	B	581	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](#)

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
5	HEM	B	1002	6	50,50,50	1.73	11 (22%)	66,82,82	1.82	18 (27%)
4	G2P	B	1001	3	31,34,34	1.37	5 (16%)	46,54,54	1.81	7 (15%)
6	NO	B	1003	5	0,1,1	-	-	-		

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
5	HEM	B	1002	6	-	5/14/54/54	-
4	G2P	B	1001	3	-	4/19/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1002	HEM	FE-NB	4.86	2.09	1.94
5	B	1002	HEM	C1B-NB	-4.06	1.33	1.40
5	B	1002	HEM	FE-NC	3.79	2.08	1.95
5	B	1002	HEM	C4D-ND	-3.66	1.34	1.40
5	B	1002	HEM	C1C-C2C	-3.28	1.38	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1001	G2P	C5-C4	2.90	1.46	1.38
5	B	1002	HEM	C3C-C4C	-2.89	1.40	1.46
5	B	1002	HEM	FE-ND	-2.66	1.86	1.94
4	B	1001	G2P	C6-N1	-2.64	1.33	1.38
5	B	1002	HEM	C4B-NB	-2.60	1.33	1.38
4	B	1001	G2P	PA-O5'	2.45	1.61	1.57
4	B	1001	G2P	C5-N7	-2.26	1.34	1.39
5	B	1002	HEM	C1D-ND	-2.25	1.34	1.38
5	B	1002	HEM	O2A-CGA	-2.20	1.23	1.30
5	B	1002	HEM	C1C-NC	-2.07	1.35	1.39
4	B	1001	G2P	PB-O1B	2.02	1.61	1.56

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1001	G2P	C5-C4-N3	-5.81	119.04	128.46
5	B	1002	HEM	CHC-C4B-NB	5.08	129.94	124.42
5	B	1002	HEM	CHD-C1D-ND	4.83	129.66	124.42
4	B	1001	G2P	C2-N3-C4	4.69	120.66	112.30
4	B	1001	G2P	PB-O3B-PG	-4.34	117.33	132.62
4	B	1001	G2P	N9-C4-N3	4.29	134.55	125.94
5	B	1002	HEM	C1B-NB-C4B	3.38	108.57	105.07
5	B	1002	HEM	CHB-C1B-NB	3.34	128.49	124.37
5	B	1002	HEM	CHD-C1D-C2D	-3.29	119.84	124.98
4	B	1001	G2P	C6-C5-N7	3.25	136.29	130.25
5	B	1002	HEM	CHA-C4D-ND	3.04	128.12	124.37
5	B	1002	HEM	CAD-CBD-CGD	-2.92	107.32	113.60
5	B	1002	HEM	CHD-C4C-NC	2.89	127.56	124.44
5	B	1002	HEM	CAA-CBA-CGA	-2.82	107.53	113.60
5	B	1002	HEM	C4C-NC-C1C	2.68	107.97	105.35
4	B	1001	G2P	C4-C5-N7	-2.59	106.62	110.72
5	B	1002	HEM	O2D-CGD-O1D	-2.43	117.24	123.30
5	B	1002	HEM	C4A-NA-C1A	2.31	107.61	105.35
5	B	1002	HEM	C2A-C1A-NA	-2.22	107.66	110.15
5	B	1002	HEM	C1A-CHA-C4D	-2.18	121.18	126.34
5	B	1002	HEM	O2D-CGD-CBD	2.16	120.96	114.03
5	B	1002	HEM	CHB-C1B-C2B	-2.14	120.80	126.73
5	B	1002	HEM	C4C-CHD-C1D	-2.13	121.46	126.06
5	B	1002	HEM	O2A-CGA-CBA	2.11	120.79	114.03
4	B	1001	G2P	O6-C6-C5	-2.10	121.04	126.60

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1001	G2P	PB-C3A-PA-O2A
4	B	1001	G2P	PB-C3A-PA-O5'
5	B	1002	HEM	C3D-CAD-CBD-CGD
5	B	1002	HEM	C4D-C3D-CAD-CBD
5	B	1002	HEM	C2D-C3D-CAD-CBD
5	B	1002	HEM	C4C-C3C-CAC-CBC
4	B	1001	G2P	C4'-C5'-O5'-PA
5	B	1002	HEM	C2A-CAA-CBA-CGA
4	B	1001	G2P	O4'-C4'-C5'-O5'

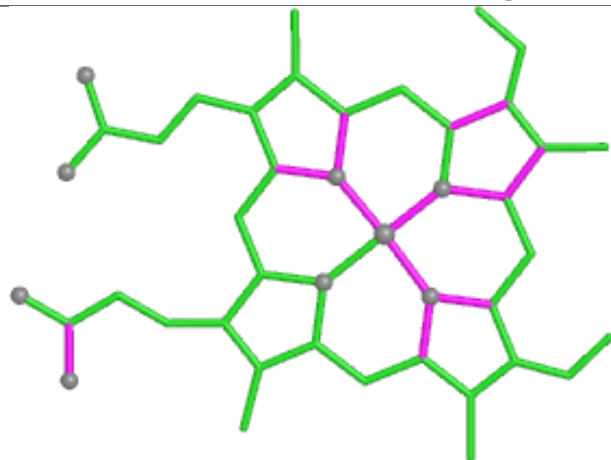
There are no ring outliers.

2 monomers are involved in 12 short contacts:

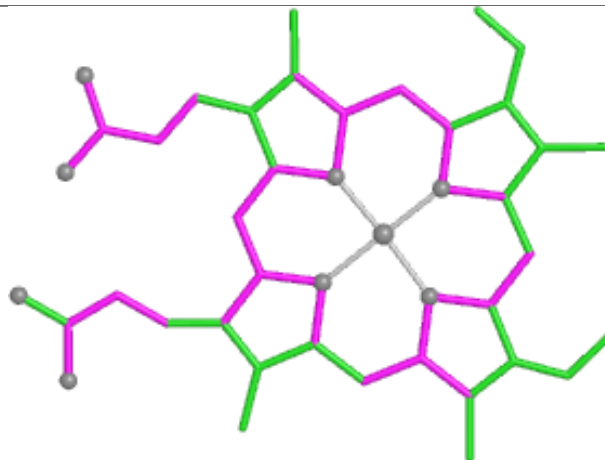
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1002	HEM	9	0
4	B	1001	G2P	3	0

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.

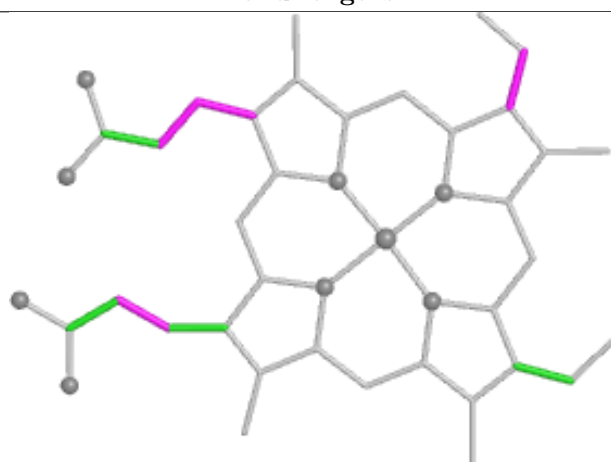
Ligand HEM B 1002



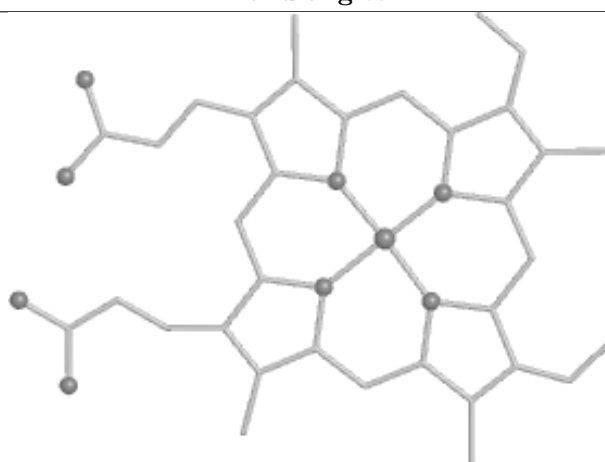
Bond lengths



Bond angles

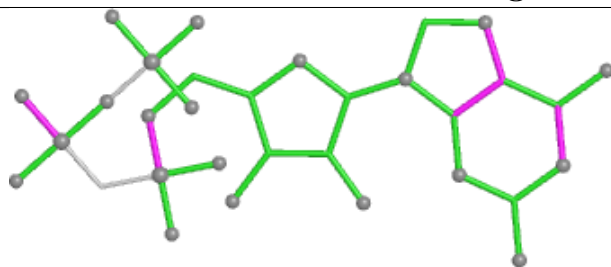


Torsions

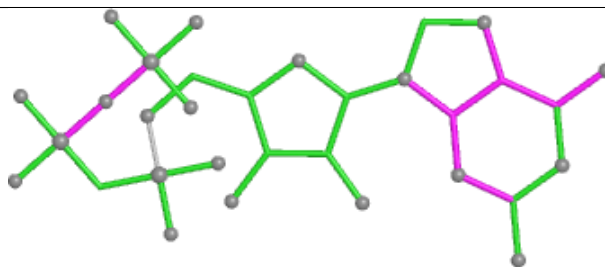


Rings

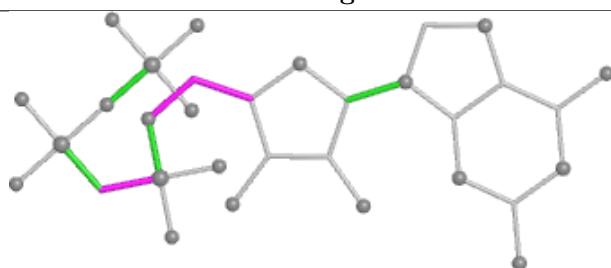
Ligand G2P B 1001



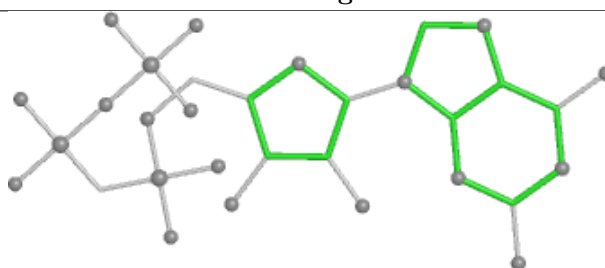
Bond lengths



Bond angles



Torsions



Rings

Ligand NO B 1003			
 Bond lengths		 Bond angles	
 Torsions		 Rings	

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.

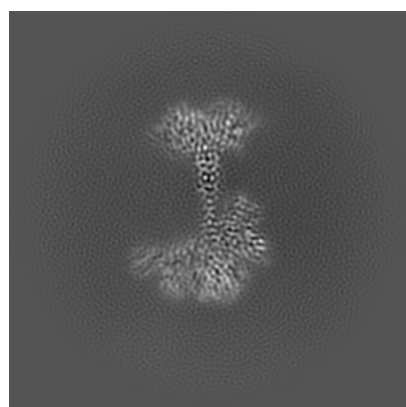
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34632. These allow visual inspection of the internal detail of the map and identification of artifacts.

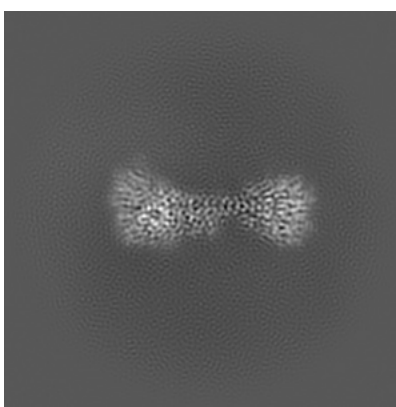
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

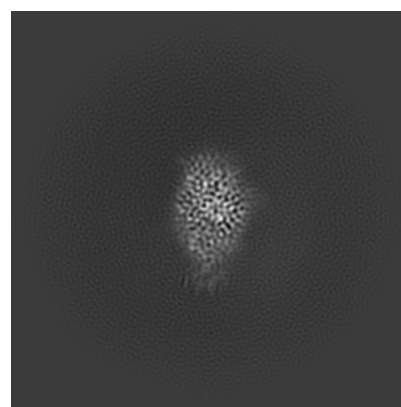
6.1.1 Primary map



X



Y

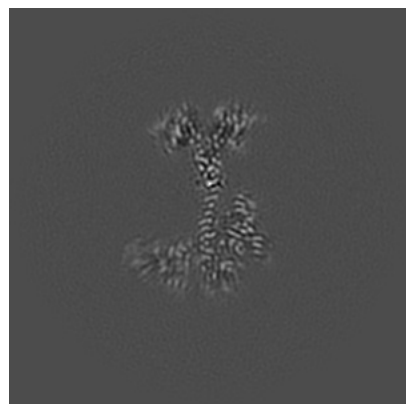


Z

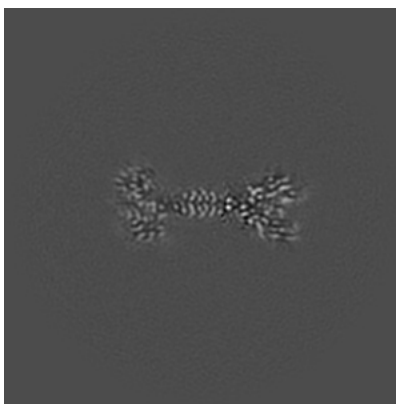
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

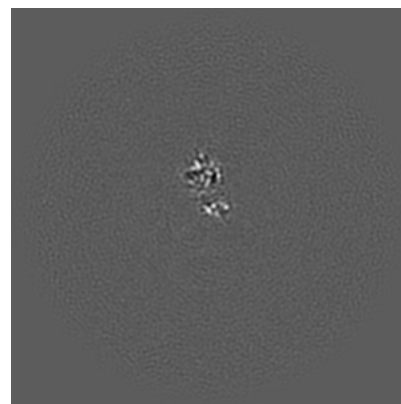
6.2.1 Primary map



X Index: 128



Y Index: 128

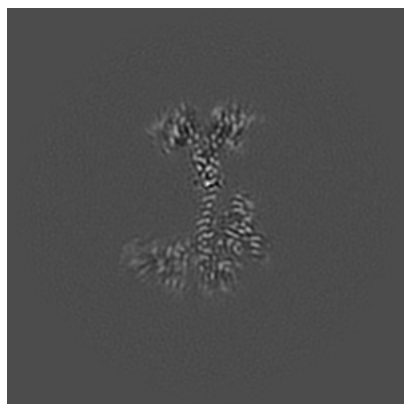


Z Index: 128

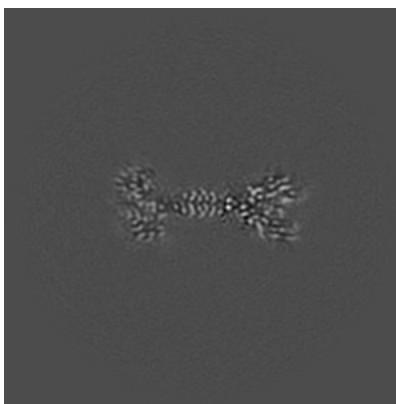
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

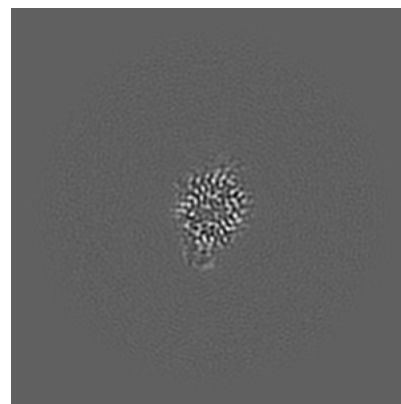
6.3.1 Primary map



X Index: 128



Y Index: 128

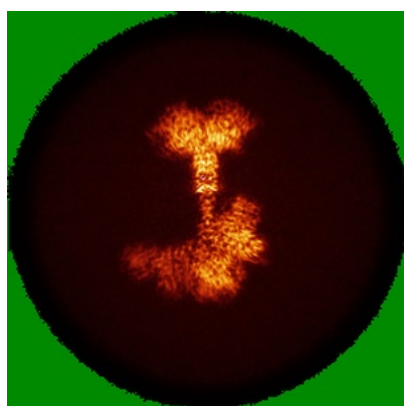


Z Index: 176

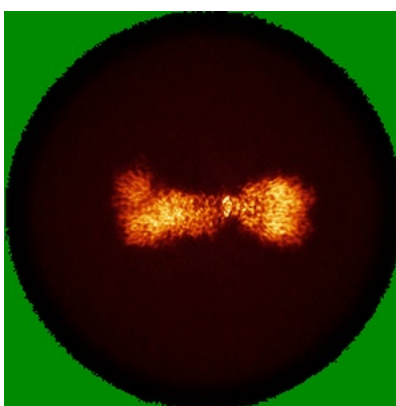
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

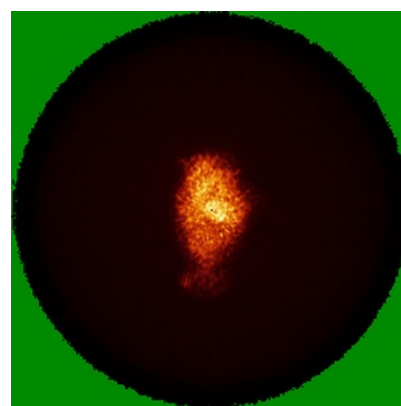
6.4.1 Primary map



X



Y

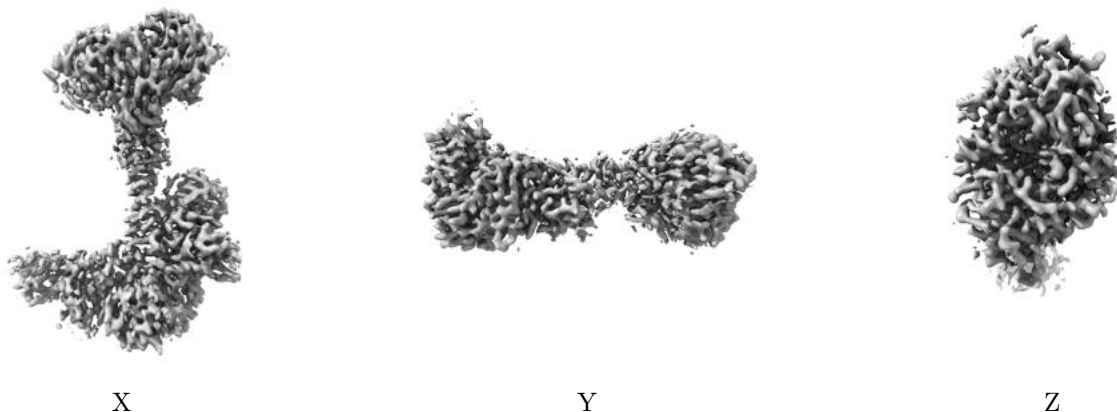


Z

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



The images above show the 3D surface view of the map at the recommended contour level 1.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

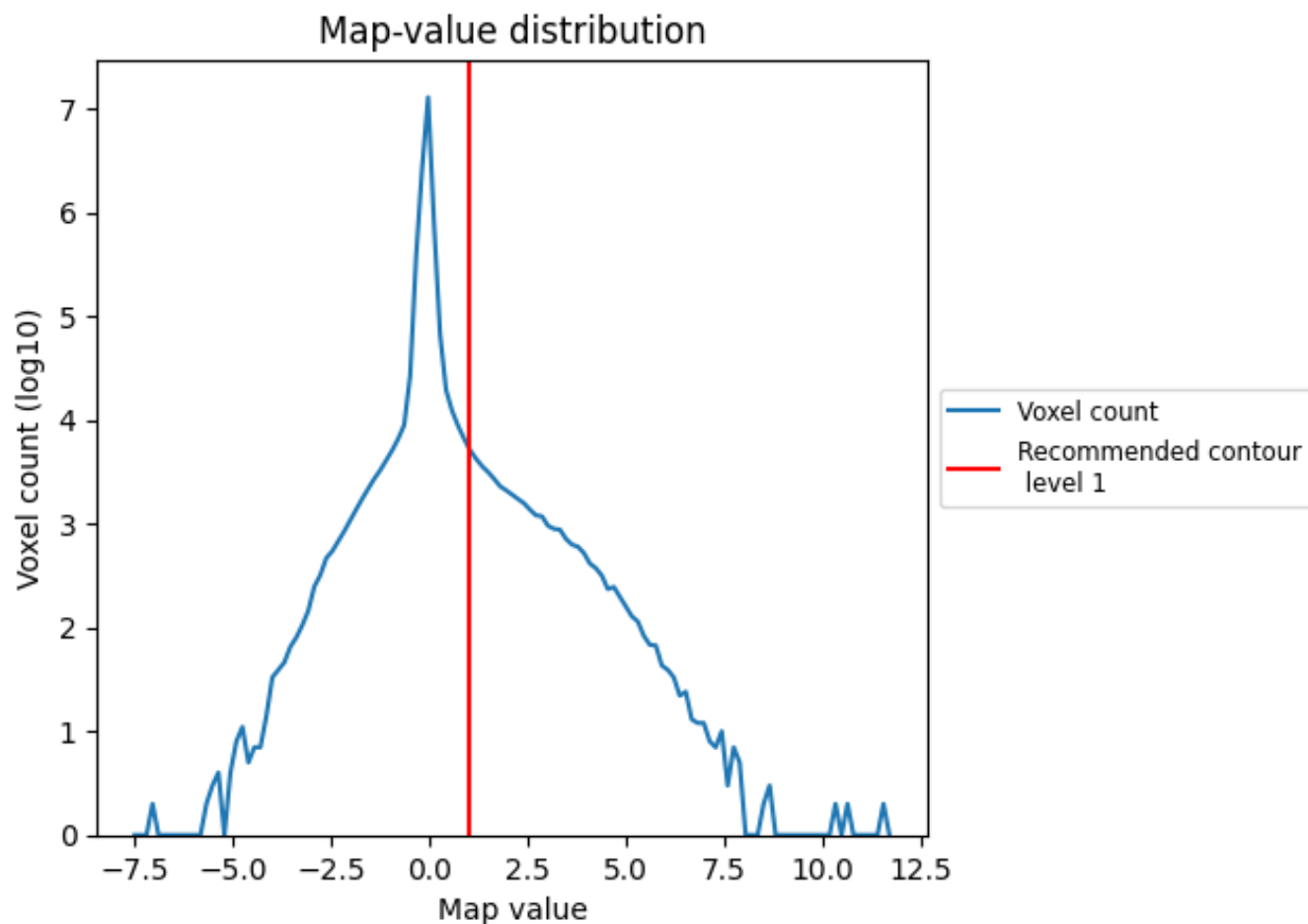
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

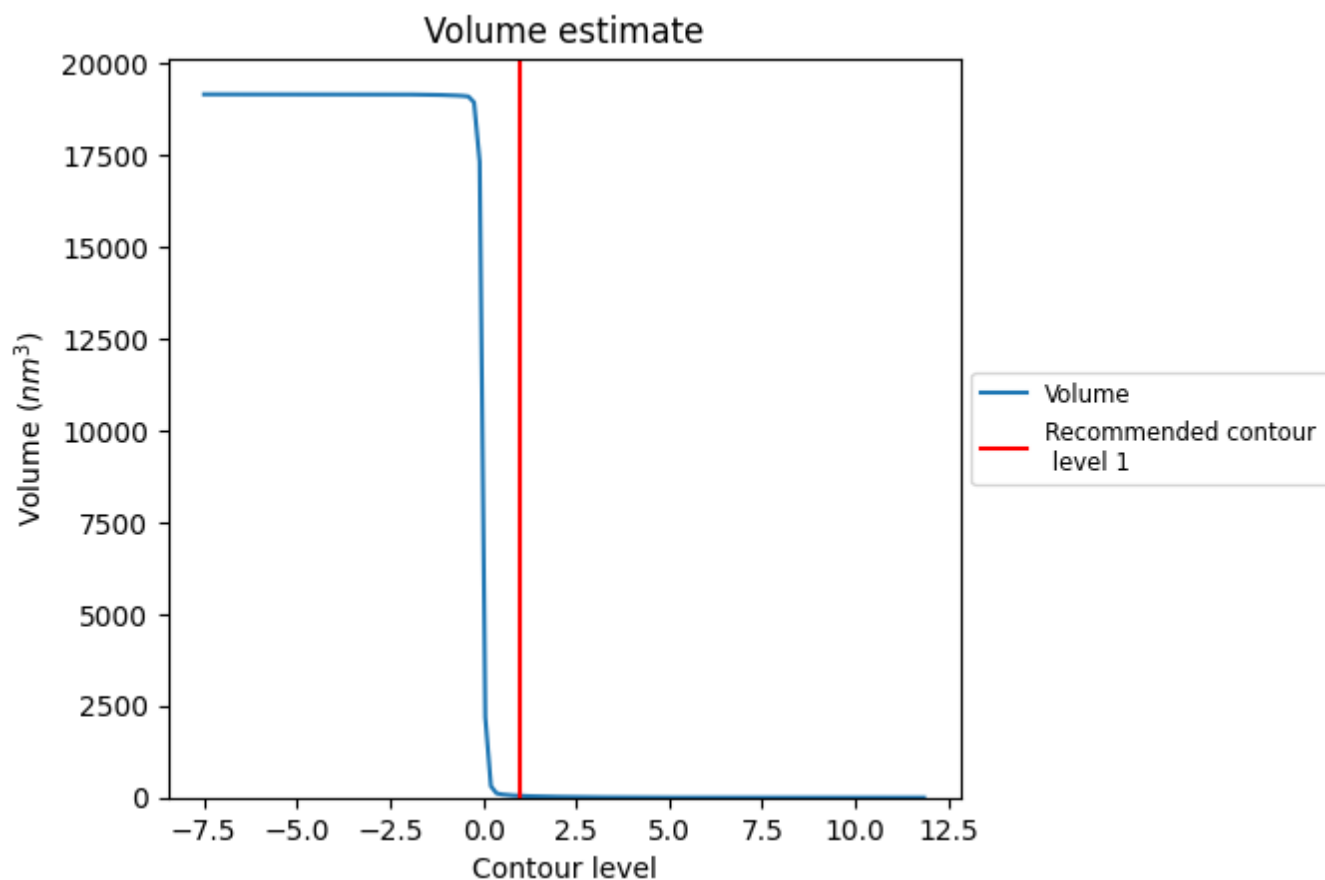
This section contains the results of statistical analysis of the map.

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.

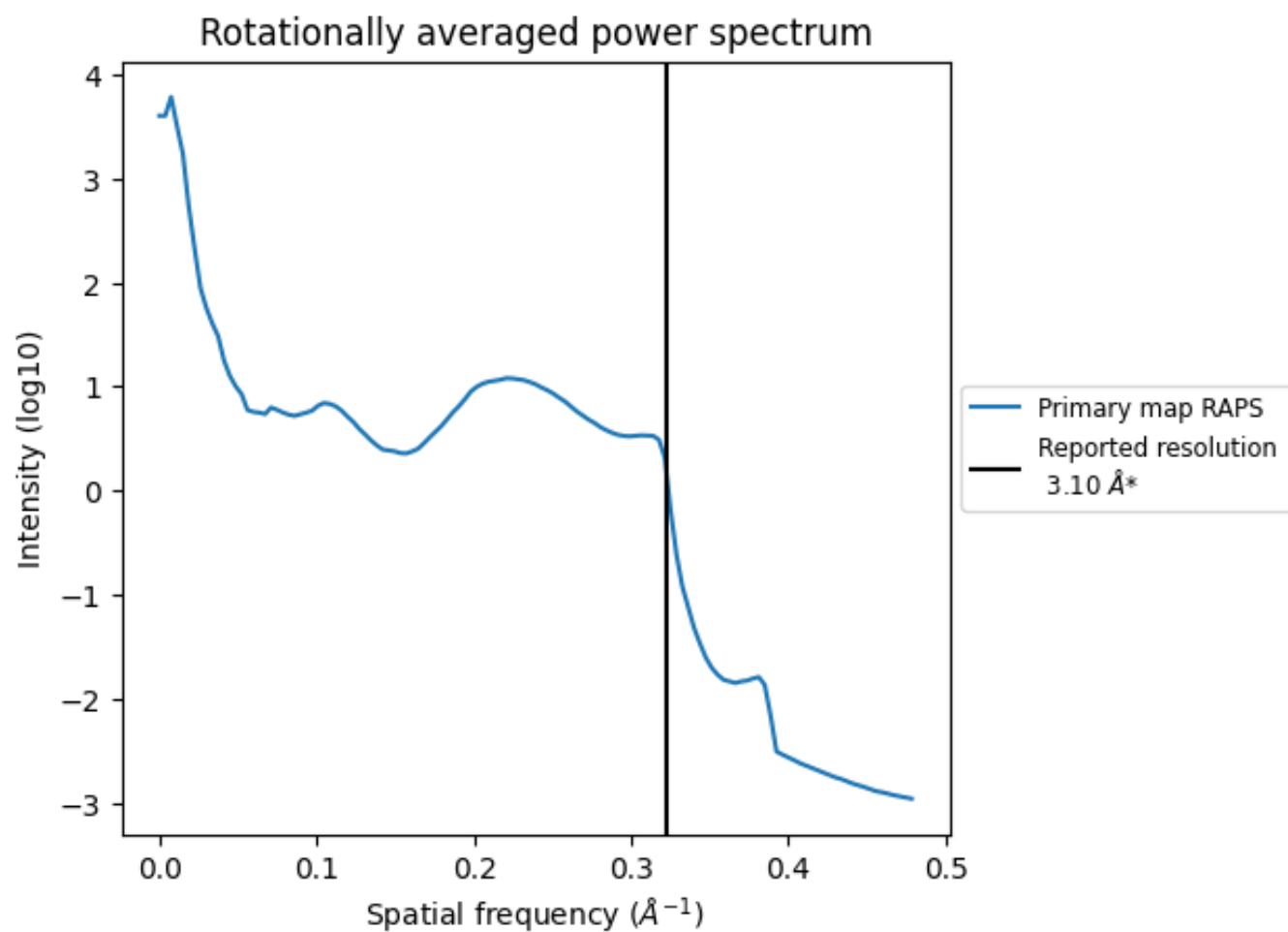
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 47 nm³; this corresponds to an approximate mass of 43 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 ⓘ



*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

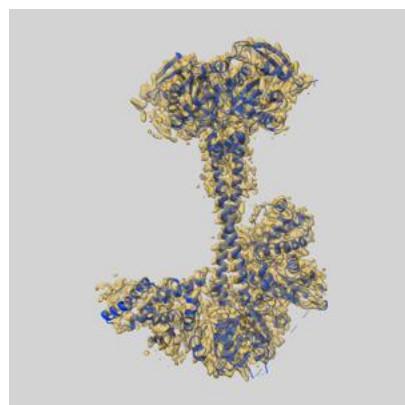
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

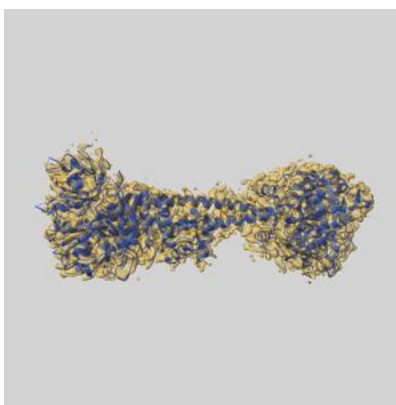
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-34632 and PDB model 8HBH. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

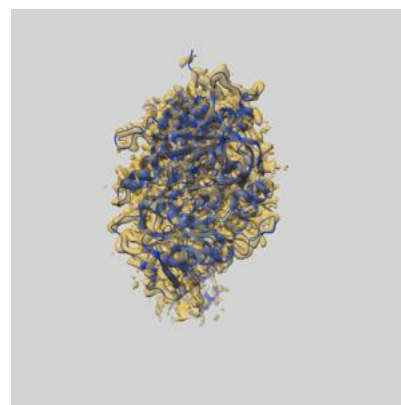
9.1 Map-model overlay [i](#)



X



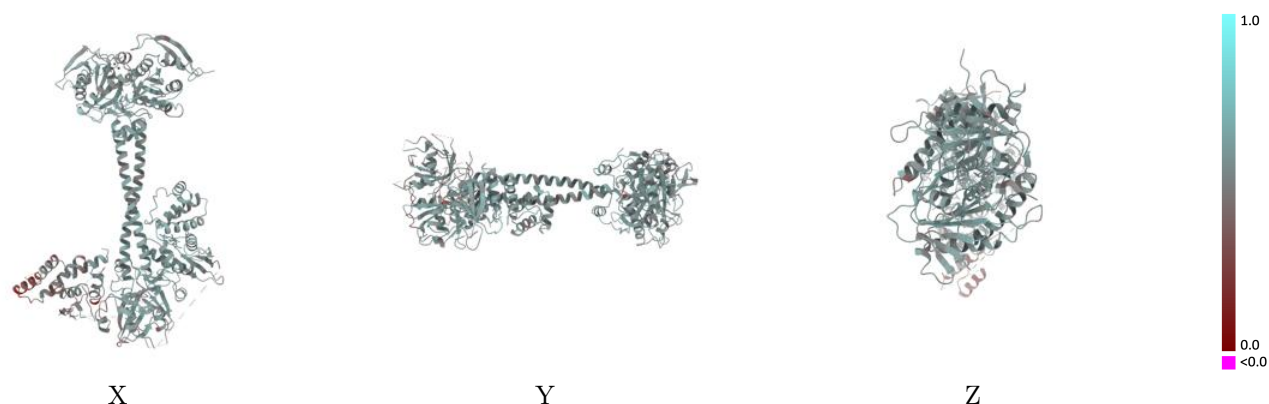
Y



Z

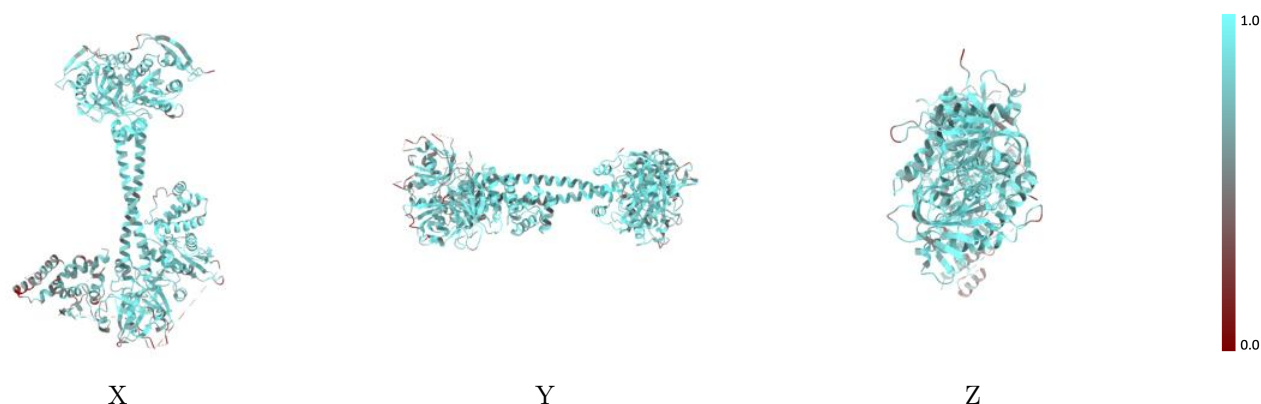
The images above show the 3D surface view of the map at the recommended contour level 1.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



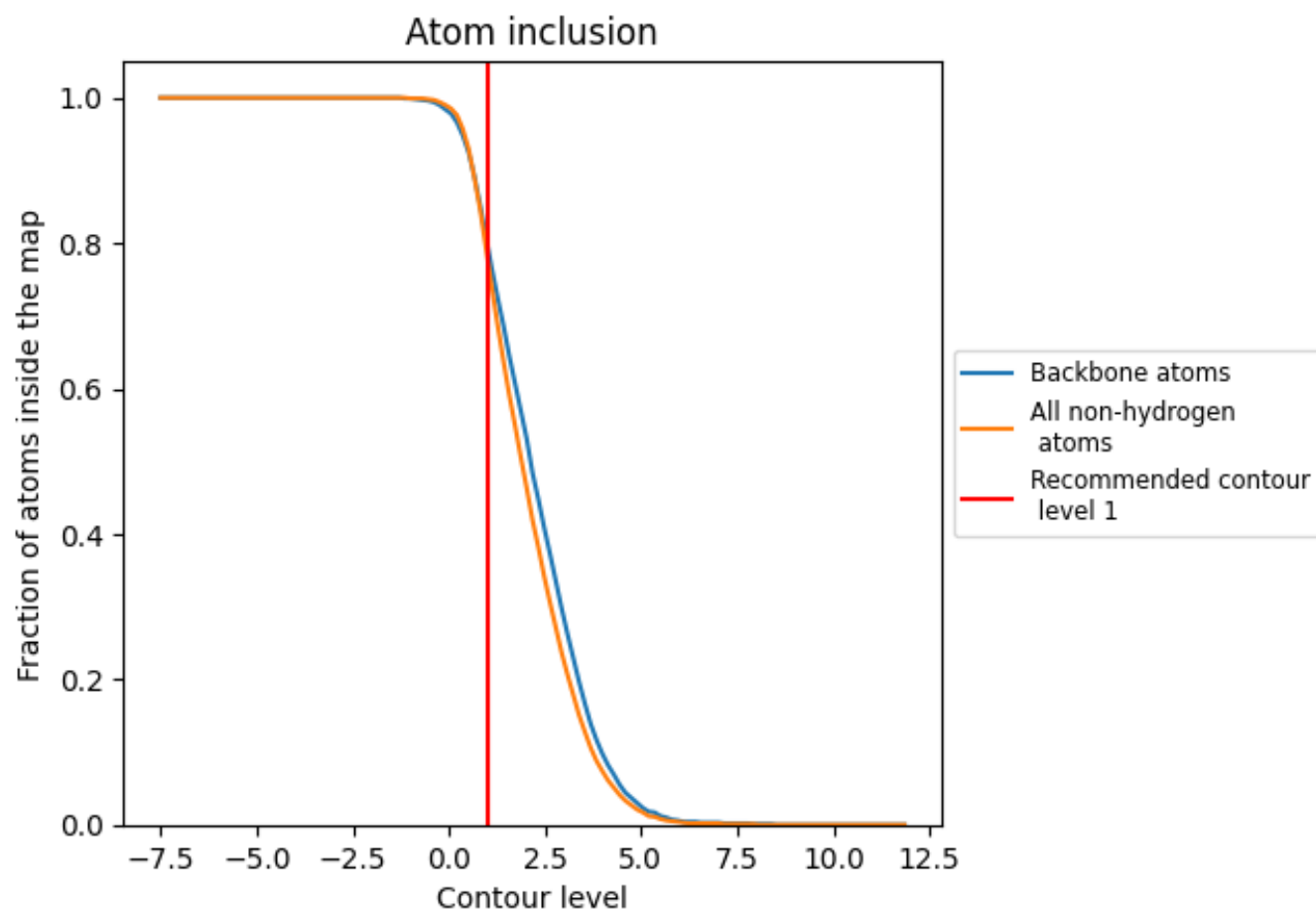
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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](#)



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 (1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 78% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7800	0.5370
A	0.7450	0.5180
B	0.8100	0.5530

