



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 01:47 PM UTC

PDB ID : 7FIH / pdb_00007fih
EMDB ID : EMD-31597
Title : luteinizing hormone/choriogonadotropin receptor(S277I)-chorionic gonadotropin-Gs-Org43553 complex
Authors : Duan, J.; Xu, P.; Cheng, X.; Mao, C.; Croll, T.; He, X.; Shi, J.; Luan, X.; Yin, W.; You, E.; Liu, Q.; Zhang, S.; Jiang, H.; Zhang, Y.; Jiang, Y.; Xu, H.E.
Deposited on : 2021-07-31
Resolution : 3.20 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.dev132
Mogul : 2022.3.0, CSD as543be (2022)
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) : 2.49

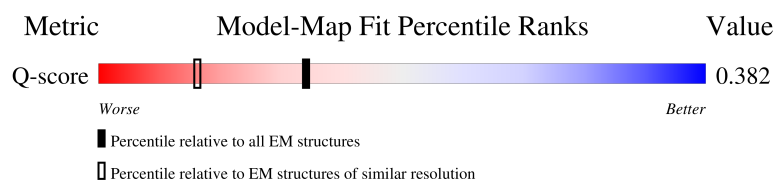
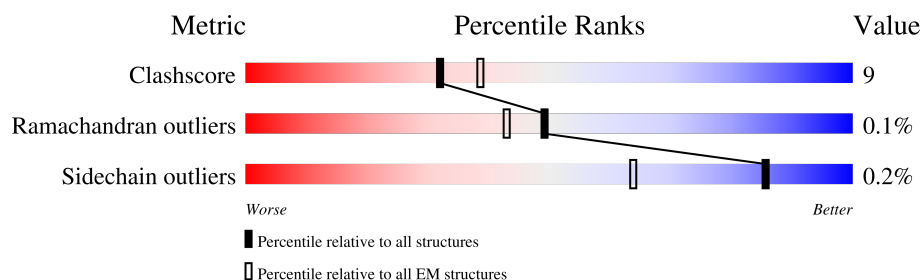
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.20 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.



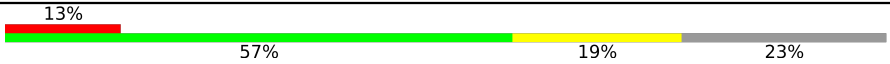
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15020 (2.70 - 3.70)

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	361	9% 52% 14% 34%
2	B	350	8% 76% 20% .
3	G	71	18% 70% 10% 20%
4	R	697	9% 66% 13% 21%

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Mol	Chain	Length	Quality of chain
5	X	115	
6	Y	165	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
7	55Z	R	801	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 10853 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 Engineered Guanine nucleotide-binding protein G(s) subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	238	Total	C	N	O	S	0	0
			1972	1245	354	364	9		

- Molecule 2 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	338	Total	C	N	O	S	0	0
			2601	1604	467	509	21		

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-9	HIS	-	expression tag	UNP P54311
B	-8	HIS	-	expression tag	UNP P54311
B	-7	HIS	-	expression tag	UNP P54311
B	-6	HIS	-	expression tag	UNP P54311
B	-5	HIS	-	expression tag	UNP P54311
B	-4	HIS	-	expression tag	UNP P54311
B	-3	GLY	-	expression tag	UNP P54311
B	-2	SER	-	expression tag	UNP P54311
B	-1	LEU	-	expression tag	UNP P54311
B	0	LEU	-	expression tag	UNP P54311
B	1	GLN	-	expression tag	UNP P54311

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	57	Total	C	N	O	S	0	0
			436	273	77	83	3		

- Molecule 4 is a protein called Lutropin-choriogonadotropic hormone receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	R	550	Total	C	N	O	S	0	0
			4314	2814	699	771	30		

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	11	ASP	-	expression tag	UNP P22888
R	12	TYR	-	expression tag	UNP P22888
R	13	LYS	-	expression tag	UNP P22888
R	14	ASP	-	expression tag	UNP P22888
R	15	ASP	-	expression tag	UNP P22888
R	16	ASP	-	expression tag	UNP P22888
R	17	ASP	-	expression tag	UNP P22888
R	18	VAL	-	expression tag	UNP P22888
R	19	GLU	-	expression tag	UNP P22888
R	20	ASN	-	expression tag	UNP P22888
R	21	LEU	-	expression tag	UNP P22888
R	22	TYR	-	expression tag	UNP P22888
R	23	PHE	-	expression tag	UNP P22888
R	24	GLN	-	expression tag	UNP P22888
R	25	GLY	-	expression tag	UNP P22888
R	26	ALA	-	expression tag	UNP P22888
R	27	SER	-	expression tag	UNP P22888
R	277	ILE	SER	engineered mutation	UNP P22888
R	700	HIS	-	expression tag	UNP P22888
R	701	HIS	-	expression tag	UNP P22888
R	702	HIS	-	expression tag	UNP P22888
R	703	HIS	-	expression tag	UNP P22888
R	704	HIS	-	expression tag	UNP P22888
R	705	HIS	-	expression tag	UNP P22888
R	706	HIS	-	expression tag	UNP P22888
R	707	HIS	-	expression tag	UNP P22888

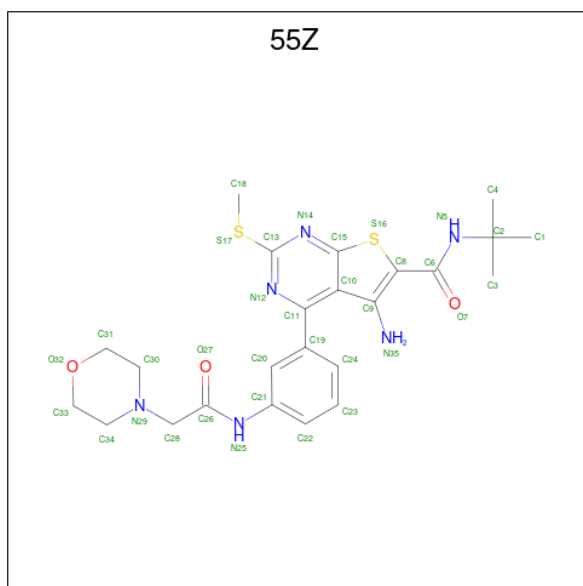
- Molecule 5 is a protein called Glycoprotein hormones alpha chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	X	88	Total	C	N	O	S	0	0
			679	420	118	128	13		

- Molecule 6 is a protein called Choriogonadotropin subunit beta 3.

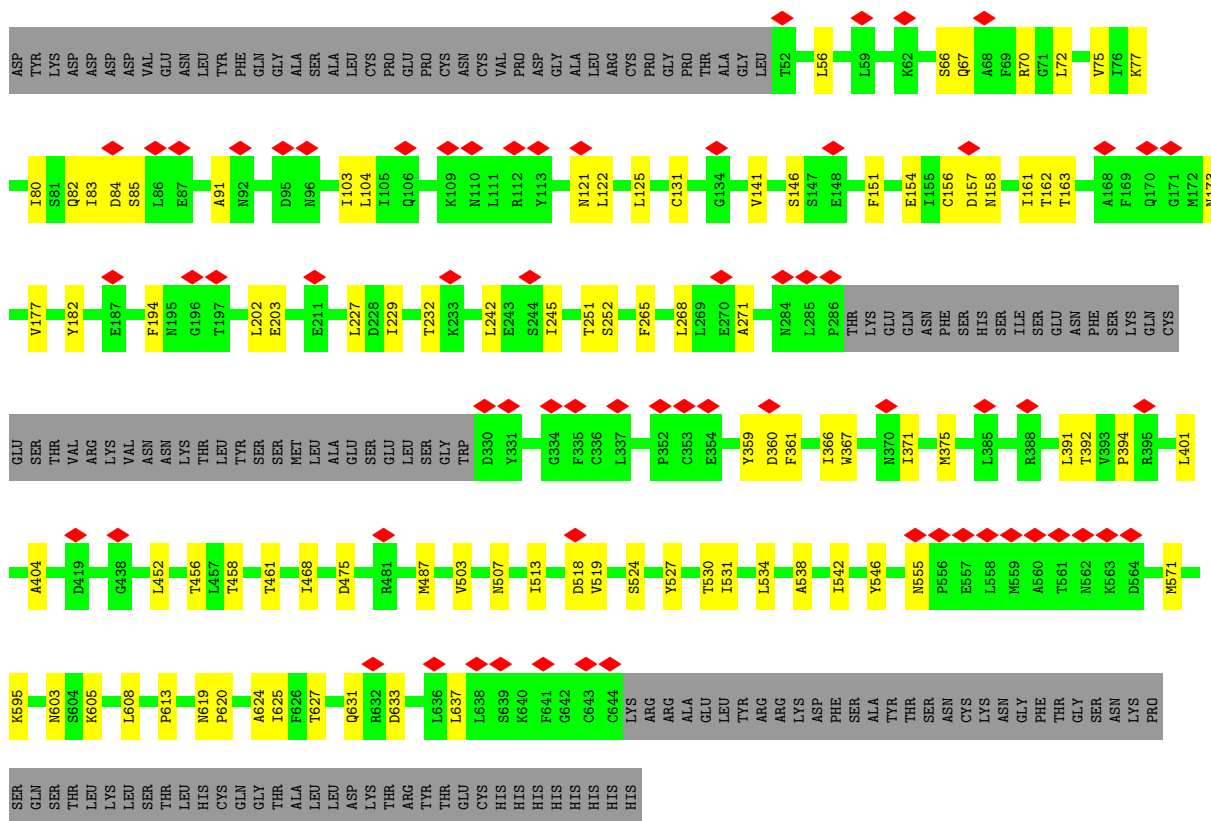
Mol	Chain	Residues	Atoms					AltConf	Trace
6	Y	110	Total	C	N	O	S	0	0
			816	505	151	147	13		

- Molecule 7 is 5-azanyl-N-tert-butyl-2-methylsulfanyl-4-[3-(2-morpholin-4-ylethanoylamino) phenyl]thieno[2,3-d]pyrimidine-6-carboxamide (CCD ID: 55Z) (formula: C₂₄H₃₀N₆O₃S₂).

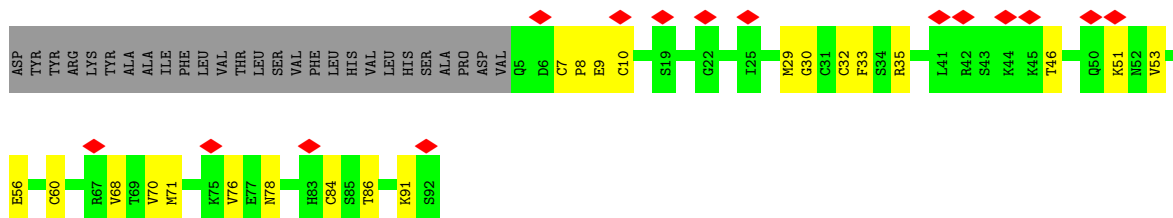


Mol	Chain	Residues	Atoms					AltConf
7	R	1	Total	C	N	O	S	0
			35	24	6	3	2	

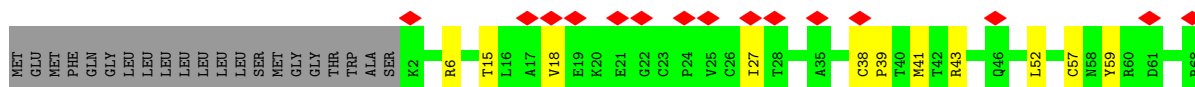
- Molecule 4: Lutropin-choriogonadotropic hormone receptor

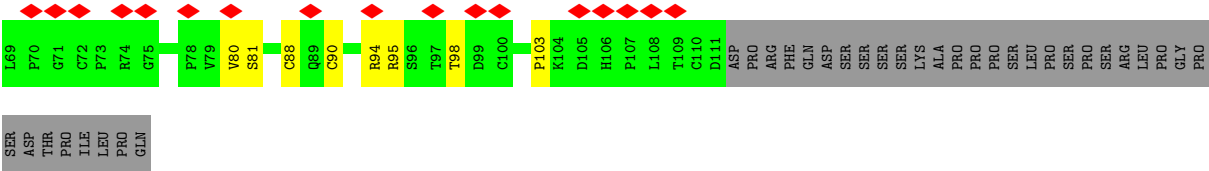


- Molecule 5: Glycoprotein hormones alpha chain



- Molecule 6: Choriogonadotropin subunit beta 3





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	413532	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 ($e^-/\text{\AA}^2$)	70	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.254	Depositor
Minimum map value	-0.001	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.040	Depositor
Recommended contour level	0.358	Depositor
Map size ($\text{\AA}$)	278.272, 278.272, 278.272	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.087, 1.087, 1.087	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 55Z

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	0.21	0/2011	0.53	2/2706 (0.1%)
2	B	0.14	0/2648	0.36	0/3589
3	G	0.15	0/442	0.29	0/597
4	R	0.18	0/4410	0.41	0/5986
5	X	0.21	0/695	0.58	0/938
6	Y	0.18	0/832	0.42	0/1138
All	All	0.18	0/11038	0.43	2/14954 (0.0%)

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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	255	ARG	N-CA-C	-11.43	98.44	112.38
1	A	254	ASN	CA-C-O	-5.54	112.58	120.51

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	252	ASP	Mainchain

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	1972	0	1950	38	0
2	B	2601	0	2505	49	0
3	G	436	0	448	7	0
4	R	4314	0	4388	70	0
5	X	679	0	647	19	0
6	Y	816	0	817	17	0
7	R	35	0	0	3	0
All	All	10853	0	10755	185	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 9.

All (185) 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:233:LYS:NZ	2:B:188:MET:SD	2.40	0.94
2:B:114:CYS:SG	2:B:122:SER:OG	2.29	0.91
4:R:268:LEU:HD21	4:R:271:ALA:HB2	1.53	0.90
2:B:74:SER:OG	2:B:76:ASP:OD1	1.96	0.82
4:R:371:ILE:HG22	4:R:375:MET:HE2	1.62	0.79
4:R:507:ASN:ND2	4:R:518:ASP:OD2	2.16	0.78
4:R:56:LEU:O	4:R:83:ILE:HD11	1.84	0.78
2:B:27:ASP:OD1	2:B:28:ALA:N	2.21	0.73
1:A:254:ASN:C	1:A:256:LEU:N	2.44	0.72
5:X:60:CYS:HA	5:X:84:CYS:HA	1.72	0.70
2:B:333:ASP:O	2:B:334:SER:OG	2.08	0.70
2:B:71:VAL:N	2:B:105:TYR:OH	2.25	0.70
1:A:230:GLU:O	1:A:234:TRP:NE1	2.25	0.69
1:A:253:TYR:HB3	1:A:301:TYR:CE2	2.29	0.68
4:R:392:THR:HG22	4:R:394:PRO:HD2	1.75	0.68
2:B:149:CYS:O	2:B:150:ARG:NH1	2.28	0.67
1:A:260:LEU:HD21	1:A:334:GLU:HG3	1.77	0.66
4:R:546:TYR:HD1	4:R:571:MET:HE3	1.60	0.66
2:B:54:HIS:ND1	2:B:80:ILE:HD12	2.10	0.66
4:R:595:LYS:HE3	4:R:595:LYS:HA	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:R:625:ILE:O	4:R:631:GLN:NE2	2.29	0.65
4:R:518:ASP:O	4:R:524:SER:OG	2.11	0.65
1:A:361:ASN:O	1:A:365:ILE:HD12	1.97	0.65
2:B:283:ARG:HG3	3:G:51:LEU:HD21	1.78	0.65
4:R:519:VAL:O	4:R:519:VAL:HG12	1.96	0.65
4:R:80:ILE:HG23	4:R:83:ILE:HD12	1.79	0.65
1:A:253:TYR:HB3	1:A:301:TYR:CD2	2.32	0.64
4:R:371:ILE:CG2	4:R:375:MET:HE2	2.28	0.63
4:R:194:PHE:HE2	4:R:202:LEU:HD22	1.65	0.62
1:A:260:LEU:HD21	1:A:334:GLU:CG	2.30	0.61
4:R:154:GLU:OE2	5:X:91:LYS:NZ	2.31	0.61
5:X:60:CYS:HB3	5:X:84:CYS:HB3	1.81	0.61
1:A:60:MET:HE1	1:A:361:ASN:HB2	1.84	0.59
2:B:173:THR:HG22	2:B:175:GLN:HG2	1.85	0.59
1:A:59:GLN:O	1:A:62:ILE:HG22	2.02	0.59
4:R:146:SER:OG	4:R:173:ASN:OD1	2.18	0.58
1:A:253:TYR:CZ	1:A:298:ILE:HG12	2.39	0.58
2:B:283:ARG:NH1	2:B:298:ASP:OD1	2.37	0.57
2:B:210:LEU:HD22	2:B:255:LEU:HD22	1.87	0.57
4:R:82:GLN:HA	5:X:46:THR:HG21	1.86	0.57
4:R:503:VAL:O	4:R:503:VAL:HG12	2.03	0.57
2:B:283:ARG:NH2	2:B:298:ASP:OD2	2.37	0.57
5:X:10:CYS:HA	5:X:29:MET:O	2.05	0.57
5:X:7:CYS:O	6:Y:6:ARG:NE	2.35	0.56
4:R:157:ASP:OD1	6:Y:95:ARG:NH1	2.38	0.56
4:R:229:ILE:HG22	4:R:229:ILE:O	2.05	0.55
1:A:254:ASN:C	1:A:256:LEU:H	2.11	0.54
2:B:71:VAL:HG23	2:B:105:TYR:CE1	2.42	0.54
1:A:317:GLU:N	1:A:320:GLU:OE2	2.41	0.54
1:A:374:GLN:O	1:A:378:LEU:HD13	2.08	0.54
2:B:267:ASP:OD1	2:B:268:ASN:N	2.42	0.53
4:R:80:ILE:CG2	4:R:83:ILE:HD12	2.38	0.53
2:B:86:THR:OG1	2:B:88:ASN:OD1	2.26	0.53
1:A:209:GLU:OE1	1:A:209:GLU:HA	2.09	0.53
6:Y:18:VAL:HG23	6:Y:27:ILE:HB	1.90	0.53
1:A:224:VAL:HG13	1:A:224:VAL:O	2.09	0.52
4:R:633:ASP:O	4:R:637:LEU:HD23	2.09	0.52
1:A:11:THR:C	1:A:12:LEU:HD22	2.35	0.52
2:B:34:THR:HG21	2:B:300:LEU:HB3	1.92	0.52
4:R:268:LEU:C	4:R:268:LEU:HD23	2.35	0.51
4:R:359:TYR:OH	4:R:603:ASN:OD1	2.13	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:ASP:O	1:A:20:VAL:HG23	2.10	0.51
4:R:151:PHE:HD2	4:R:177:VAL:HG23	1.75	0.51
2:B:18:ILE:HD11	3:G:23:ALA:HA	1.93	0.51
1:A:269:ASN:OD1	1:A:270:ARG:N	2.43	0.51
4:R:83:ILE:HG22	4:R:83:ILE:O	2.11	0.51
4:R:404:ALA:HB2	4:R:452:LEU:HD23	1.93	0.51
4:R:251:THR:O	4:R:252:SER:OG	2.22	0.51
2:B:95:LEU:HD22	2:B:100:VAL:HG21	1.92	0.50
1:A:285:ASP:O	1:A:289:GLU:OE1	2.28	0.50
4:R:182:TYR:OH	4:R:203:GLU:OE1	2.17	0.50
4:R:131:CYS:HA	4:R:156:CYS:HB3	1.93	0.50
1:A:306:ALA:HB1	1:A:307:ARG:NH2	2.27	0.50
5:X:68:VAL:HG23	5:X:76:VAL:HB	1.94	0.49
4:R:72:LEU:HD13	4:R:75:VAL:HG11	1.94	0.49
6:Y:43:ARG:HH21	6:Y:52:LEU:HD21	1.77	0.49
2:B:272:GLY:N	2:B:290:ASP:OD2	2.46	0.49
4:R:242:LEU:O	4:R:245:ILE:HG22	2.12	0.49
4:R:141:VAL:O	4:R:141:VAL:HG12	2.12	0.48
4:R:513:ILE:HD12	4:R:608:LEU:HD21	1.95	0.48
7:R:801:55Z:O7	7:R:801:55Z:N35	2.45	0.48
5:X:53:VAL:HG21	6:Y:59:TYR:HE2	1.79	0.48
4:R:530:THR:O	4:R:534:LEU:HD13	2.13	0.48
2:B:254:ASP:OD1	2:B:255:LEU:N	2.46	0.48
4:R:141:VAL:O	4:R:141:VAL:CG1	2.61	0.48
5:X:10:CYS:HA	5:X:30:GLY:HA3	1.96	0.48
6:Y:15:THR:O	6:Y:15:THR:HG23	2.14	0.47
4:R:162:THR:HG23	4:R:163:THR:HG22	1.96	0.47
1:A:56:ILE:O	1:A:60:MET:HG2	2.13	0.47
2:B:263:THR:O	2:B:263:THR:HG23	2.14	0.47
2:B:158:VAL:O	2:B:158:VAL:HG13	2.15	0.47
6:Y:41:MET:HG2	6:Y:52:LEU:HD13	1.96	0.47
4:R:608:LEU:O	4:R:613:PRO:HD3	2.16	0.46
4:R:122:LEU:HD12	4:R:125:LEU:HD12	1.96	0.46
4:R:66:SER:OG	4:R:91:ALA:O	2.33	0.46
1:A:346:ARG:HB2	1:A:346:ARG:NH1	2.31	0.46
2:B:286:LEU:HD22	2:B:327:VAL:HG21	1.98	0.46
4:R:487:MET:HA	4:R:487:MET:HE2	1.98	0.46
2:B:44:GLN:HA	2:B:44:GLN:OE1	2.16	0.45
2:B:101:MET:HE1	2:B:117:LEU:HD21	1.98	0.45
5:X:35:ARG:HG2	5:X:56:GLU:HB3	1.98	0.45
4:R:103:ILE:HD12	4:R:125:LEU:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:THR:HG21	1:A:356:SER:HB3	1.98	0.45
2:B:105:TYR:O	2:B:151:PHE:HB2	2.16	0.45
2:B:247:ASP:OD1	2:B:249:THR:OG1	2.30	0.45
4:R:103:ILE:C	4:R:104:LEU:HD12	2.42	0.45
4:R:84:ASP:N	4:R:84:ASP:OD1	2.50	0.45
5:X:70:VAL:HG12	5:X:71:MET:N	2.31	0.45
4:R:391:LEU:H	4:R:391:LEU:HD23	1.82	0.45
4:R:366:ILE:HG23	4:R:367:TRP:N	2.31	0.45
4:R:67:GLN:OE1	4:R:70:ARG:NH2	2.50	0.45
1:A:253:TYR:HB3	1:A:301:TYR:HE2	1.79	0.44
1:A:378:LEU:HD12	4:R:468:ILE:HD12	1.97	0.44
4:R:151:PHE:CD2	4:R:177:VAL:HG23	2.51	0.44
4:R:232:THR:HG22	4:R:232:THR:O	2.17	0.44
2:B:248:ALA:HB1	2:B:269:ILE:HG22	1.99	0.44
4:R:624:ALA:O	4:R:627:THR:HG22	2.17	0.44
6:Y:94:ARG:O	6:Y:98:THR:N	2.39	0.44
2:B:133:VAL:O	2:B:133:VAL:HG13	2.17	0.44
6:Y:57:CYS:HA	6:Y:90:CYS:HA	2.00	0.44
2:B:35:ASN:OD1	2:B:36:ASN:N	2.50	0.44
2:B:86:THR:O	2:B:88:ASN:ND2	2.50	0.44
3:G:15:LEU:O	3:G:19:LEU:HD13	2.18	0.44
4:R:619:ASN:N	4:R:620:PRO:HD2	2.33	0.44
4:R:360:ASP:OD1	4:R:361:PHE:N	2.50	0.43
4:R:519:VAL:O	4:R:519:VAL:CG1	2.65	0.43
5:X:9:GLU:OE2	6:Y:6:ARG:NH2	2.49	0.43
2:B:154:ASP:OD1	2:B:154:ASP:C	2.60	0.43
4:R:458:THR:CG2	4:R:538:ALA:HB1	2.47	0.43
6:Y:80:VAL:HG22	6:Y:81:SER:N	2.34	0.43
6:Y:38:CYS:HB2	6:Y:39:PRO:HD2	2.00	0.43
4:R:475:ASP:OD1	4:R:475:ASP:C	2.61	0.43
6:Y:59:TYR:HA	6:Y:88:CYS:SG	2.59	0.43
1:A:321:ASP:O	1:A:324:VAL:HG12	2.19	0.43
2:B:121:CYS:HB2	2:B:146:LEU:HD13	2.00	0.43
5:X:56:GLU:OE1	6:Y:103:PRO:HA	2.19	0.43
2:B:6:GLN:O	2:B:9:GLN:NE2	2.52	0.43
4:R:229:ILE:HG22	4:R:232:THR:HB	2.01	0.43
1:A:371:ASP:OD1	1:A:375:ARG:NE	2.52	0.43
2:B:13:GLN:HA	2:B:13:GLN:OE1	2.19	0.42
2:B:61:MET:HG3	2:B:317:CYS:SG	2.59	0.42
2:B:73:ALA:HB1	2:B:100:VAL:HG11	2.01	0.42
3:G:13:ARG:HA	3:G:16:VAL:HG12	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:ASP:OD2	2:B:173:THR:OG1	2.35	0.42
4:R:401:LEU:HB2	4:R:456:THR:HG21	2.01	0.42
1:A:371:ASP:OD2	4:R:555:ASN:ND2	2.47	0.42
1:A:375:ARG:HB3	1:A:379:ARG:NH1	2.34	0.42
2:B:283:ARG:CG	3:G:51:LEU:HD21	2.49	0.42
4:R:202:LEU:HD23	4:R:227:LEU:HD13	2.00	0.42
4:R:268:LEU:HD21	4:R:271:ALA:CB	2.37	0.42
4:R:527:TYR:O	4:R:531:ILE:HG12	2.20	0.42
1:A:282:ASN:OD1	1:A:283:LYS:N	2.52	0.42
2:B:264:TYR:OH	2:B:299:ALA:O	2.25	0.42
4:R:461:THR:HG21	4:R:542:ILE:HA	2.01	0.42
4:R:620:PRO:O	4:R:625:ILE:N	2.33	0.42
5:X:51:LYS:NZ	6:Y:95:ARG:O	2.38	0.42
1:A:11:THR:O	1:A:12:LEU:HD22	2.20	0.42
2:B:118:ASP:N	2:B:118:ASP:OD1	2.51	0.42
6:Y:80:VAL:HG22	6:Y:81:SER:H	1.85	0.42
1:A:346:ARG:HB2	1:A:346:ARG:HH11	1.84	0.41
5:X:32:CYS:HB3	5:X:33:PHE:H	1.71	0.41
2:B:71:VAL:HG23	2:B:105:TYR:OH	2.19	0.41
3:G:10:ALA:O	3:G:13:ARG:HG2	2.20	0.41
4:R:121:ASN:O	4:R:122:LEU:HD23	2.21	0.41
4:R:503:VAL:O	4:R:503:VAL:CG1	2.69	0.41
7:R:801:55Z:O27	7:R:801:55Z:C34	2.68	0.41
3:G:15:LEU:HD23	3:G:19:LEU:HD13	2.01	0.41
4:R:77:LYS:HZ1	5:X:86:THR:HG21	1.85	0.41
2:B:150:ARG:HB3	2:B:192:LEU:HD13	2.03	0.41
2:B:75:GLN:HG2	2:B:99:TRP:HE3	1.85	0.41
4:R:77:LYS:HZ1	5:X:86:THR:CG2	2.34	0.41
2:B:291:ASP:OD1	2:B:292:PHE:N	2.54	0.41
1:A:27:GLU:OE1	2:B:89:LYS:NZ	2.51	0.41
1:A:54:SER:O	1:A:57:VAL:HG22	2.21	0.41
1:A:249:ASP:OD1	1:A:251:SER:N	2.48	0.40
2:B:90:VAL:HG12	2:B:91:HIS:ND1	2.36	0.40
7:R:801:55Z:O27	7:R:801:55Z:C20	2.68	0.40
1:A:253:TYR:HD1	1:A:253:TYR:HA	1.59	0.40
4:R:83:ILE:O	4:R:85:SER:N	2.54	0.40
4:R:265:PHE:CD1	4:R:268:LEU:HD12	2.56	0.40
1:A:337:ARG:HG3	1:A:337:ARG:HH21	1.86	0.40
2:B:183:HIS:CE1	2:B:203:ALA:CB	3.05	0.40
4:R:158:ASN:HB2	4:R:161:ILE:HD12	2.04	0.40
5:X:8:PRO:HA	6:Y:6:ARG:NE	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:X:78:ASN:OD1	5:X:78:ASN:C	2.65	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	234/361 (65%)	228 (97%)	5 (2%)	1 (0%)	30	62
2	B	336/350 (96%)	327 (97%)	9 (3%)	0	100	100
3	G	55/71 (78%)	55 (100%)	0	0	100	100
4	R	546/697 (78%)	527 (96%)	19 (4%)	0	100	100
5	X	86/115 (75%)	81 (94%)	5 (6%)	0	100	100
6	Y	108/165 (66%)	101 (94%)	7 (6%)	0	100	100
All	All	1365/1759 (78%)	1319 (97%)	45 (3%)	1 (0%)	49	79

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	254	ASN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	215/315 (68%)	214 (100%)	1 (0%)	81	85
2	B	281/292 (96%)	281 (100%)	0	100	100
3	G	46/58 (79%)	46 (100%)	0	100	100
4	R	481/621 (78%)	480 (100%)	1 (0%)	87	89
5	X	80/104 (77%)	80 (100%)	0	100	100
6	Y	94/145 (65%)	94 (100%)	0	100	100
All	All	1197/1535 (78%)	1195 (100%)	2 (0%)	85	89

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	254	ASN
4	R	605	LYS

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

Mol	Chain	Res	Type
1	A	268	ASN
1	A	380	GLN
4	R	107	ASN
4	R	190	GLN
4	R	562	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

1 ligand is 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
7	55Z	R	801	-	38,38,38	3.55	22 (57%)	51,55,55	4.74	28 (54%)

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
7	55Z	R	801	-	-	5/23/31/31	0/4/4/4

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	R	801	55Z	C10-C9	-7.36	1.38	1.45
7	R	801	55Z	C9-C8	-6.94	1.27	1.39
7	R	801	55Z	O7-C6	-6.40	1.11	1.23
7	R	801	55Z	C34-C33	-5.35	1.30	1.50
7	R	801	55Z	C19-C11	5.21	1.54	1.49
7	R	801	55Z	C20-C21	-5.10	1.31	1.39
7	R	801	55Z	C26-N25	4.96	1.46	1.35
7	R	801	55Z	C6-N5	4.88	1.40	1.34
7	R	801	55Z	C30-C31	-4.84	1.32	1.50
7	R	801	55Z	C15-N14	-4.76	1.27	1.33
7	R	801	55Z	C13-N14	-4.30	1.27	1.34
7	R	801	55Z	C34-N29	-4.26	1.35	1.46
7	R	801	55Z	C2-N5	-4.24	1.42	1.48
7	R	801	55Z	C30-N29	-3.94	1.36	1.46
7	R	801	55Z	C8-S16	-3.83	1.67	1.74
7	R	801	55Z	C13-N12	-3.65	1.28	1.34
7	R	801	55Z	O27-C26	-3.59	1.16	1.23
7	R	801	55Z	C15-S16	3.02	1.79	1.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	R	801	55Z	C9-N35	2.76	1.40	1.35
7	R	801	55Z	C6-C8	-2.57	1.44	1.49
7	R	801	55Z	C23-C24	-2.12	1.35	1.38
7	R	801	55Z	C21-N25	2.02	1.45	1.41

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	R	801	55Z	C18-S17-C13	23.25	119.70	102.25
7	R	801	55Z	C28-N29-C34	9.22	125.48	111.14
7	R	801	55Z	C10-C9-C8	8.97	118.50	112.09
7	R	801	55Z	C28-N29-C30	6.83	121.77	111.14
7	R	801	55Z	C2-N5-C6	-6.39	119.77	125.96
7	R	801	55Z	S16-C15-N14	5.87	130.53	121.69
7	R	801	55Z	C8-C9-N35	-5.49	116.78	124.01
7	R	801	55Z	C20-C19-C11	-5.24	110.60	120.09
7	R	801	55Z	C24-C19-C11	5.03	128.78	120.62
7	R	801	55Z	C10-C15-N14	-4.70	122.47	126.44
7	R	801	55Z	C28-C26-N25	4.70	122.93	114.23
7	R	801	55Z	C10-C15-S16	-4.38	107.00	112.40
7	R	801	55Z	C31-C30-N29	4.08	116.32	110.12
7	R	801	55Z	C23-C24-C19	-3.80	116.21	120.54
7	R	801	55Z	C22-C21-C20	-3.75	115.12	119.66
7	R	801	55Z	O32-C31-C30	3.55	119.42	111.77
7	R	801	55Z	O32-C33-C34	3.53	119.38	111.77
7	R	801	55Z	C8-C6-N5	2.94	119.91	116.06
7	R	801	55Z	C10-C11-C19	2.90	127.94	123.53
7	R	801	55Z	C10-C11-N12	-2.76	118.15	122.06
7	R	801	55Z	N14-C13-N12	-2.76	121.97	127.00
7	R	801	55Z	O27-C26-N25	-2.73	118.75	123.64
7	R	801	55Z	C23-C22-C21	2.67	122.81	119.73
7	R	801	55Z	C33-C34-N29	2.61	114.08	110.12
7	R	801	55Z	C13-N14-C15	2.56	120.50	116.62
7	R	801	55Z	O7-C6-C8	-2.54	115.65	120.15
7	R	801	55Z	C26-C28-N29	2.27	118.70	113.41
7	R	801	55Z	C22-C21-N25	2.21	127.83	120.41

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	R	801	55Z	N12-C13-S17-C18

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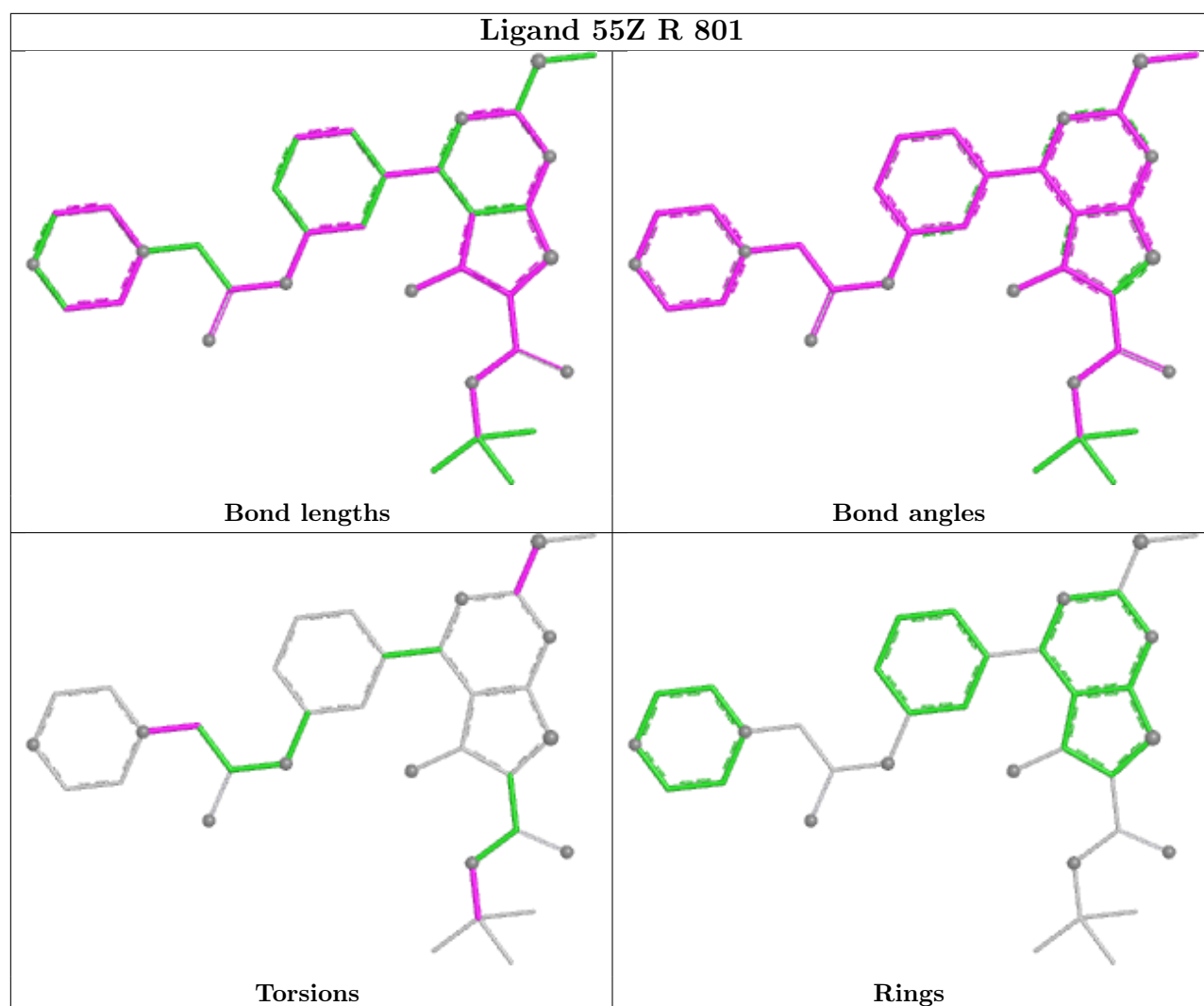
Mol	Chain	Res	Type	Atoms
7	R	801	55Z	N14-C13-S17-C18
7	R	801	55Z	C26-C28-N29-C34
7	R	801	55Z	C1-C2-N5-C6
7	R	801	55Z	C4-C2-N5-C6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	R	801	55Z	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.



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.

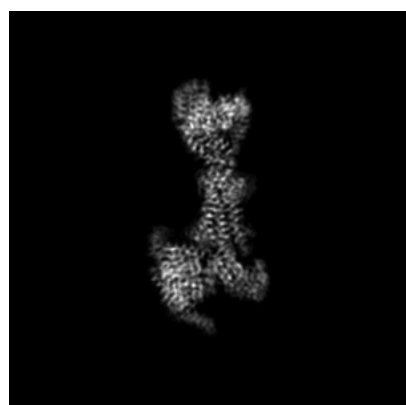
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31597. 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](#)

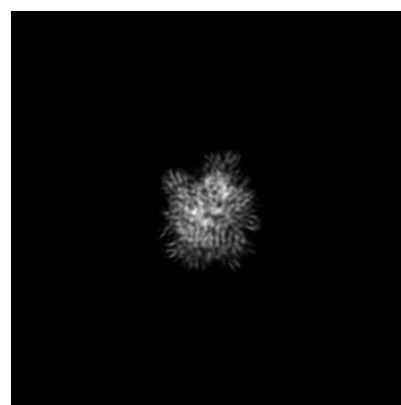
6.1.1 Primary map



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: 136



Y Index: 136

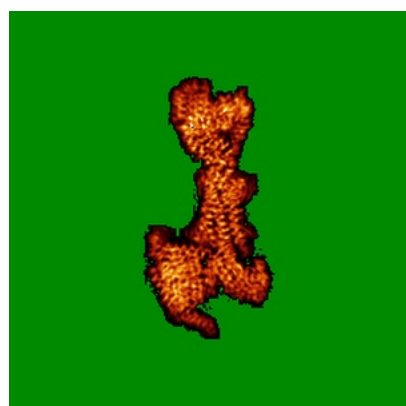


Z Index: 89

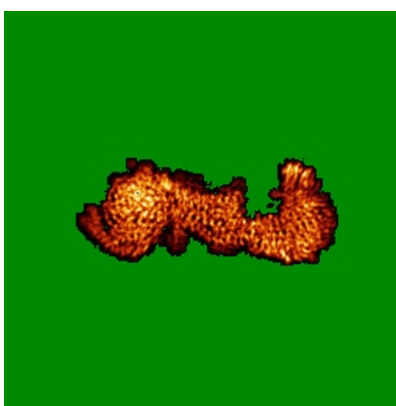
The images above show the largest variance slices of the map in three orthogonal directions.

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

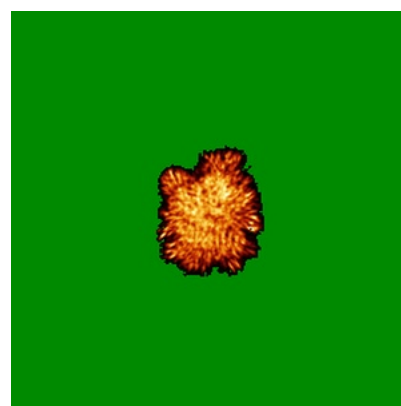
6.4.1 Primary map



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

This section was not generated.

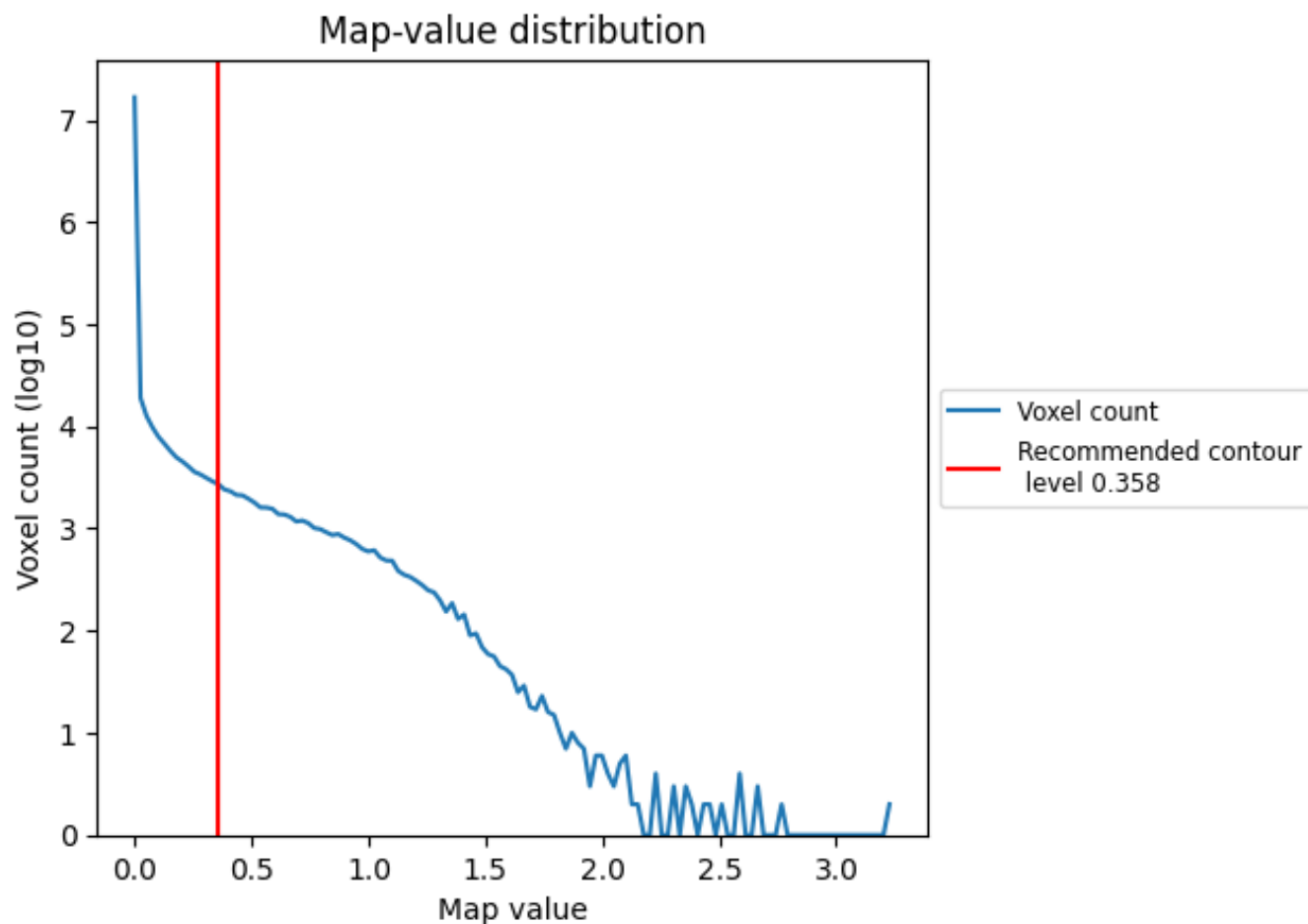
6.6 Mask visualisation

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

7 Map analysis [i](#)

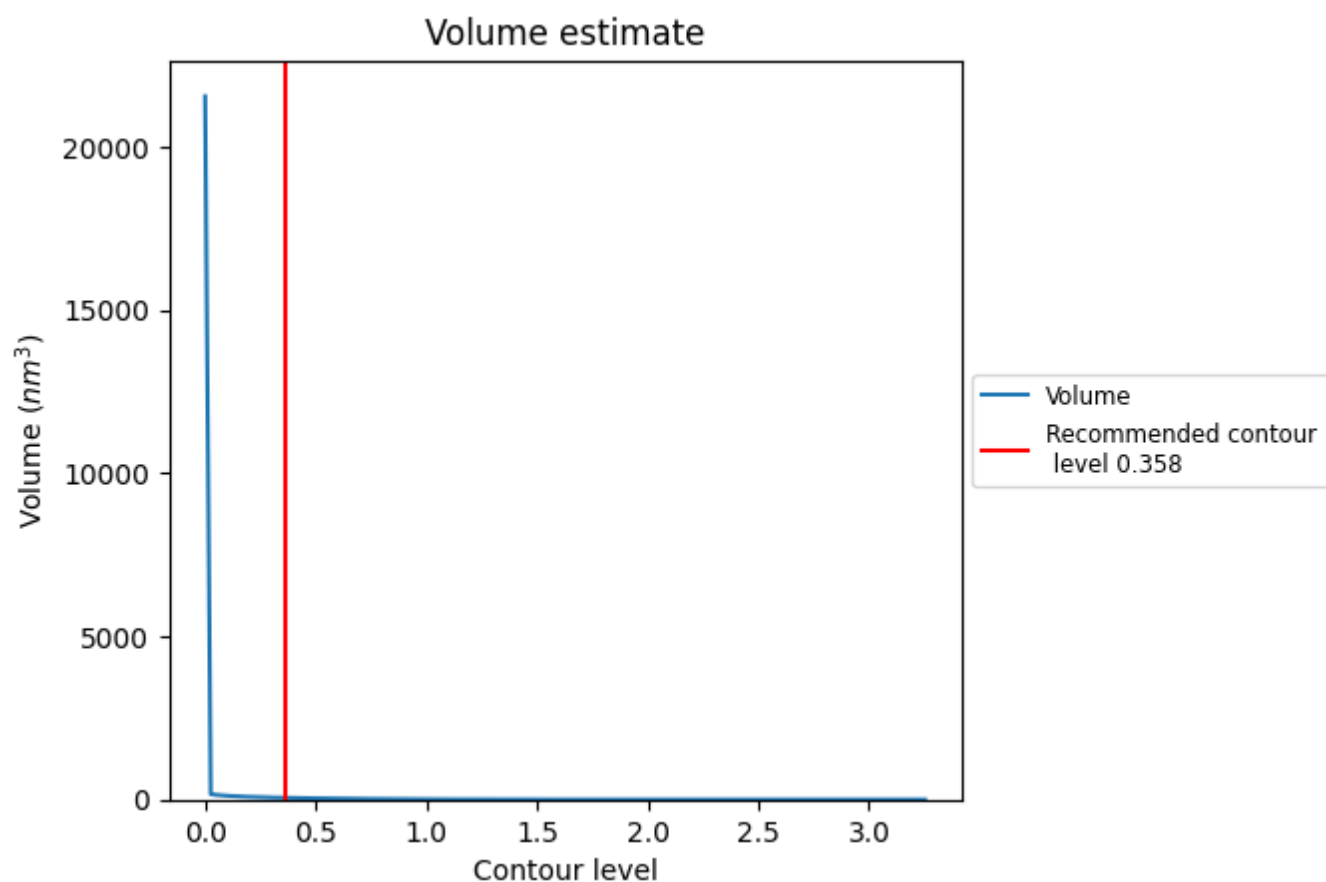
This section contains the results of statistical analysis of the map.

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.

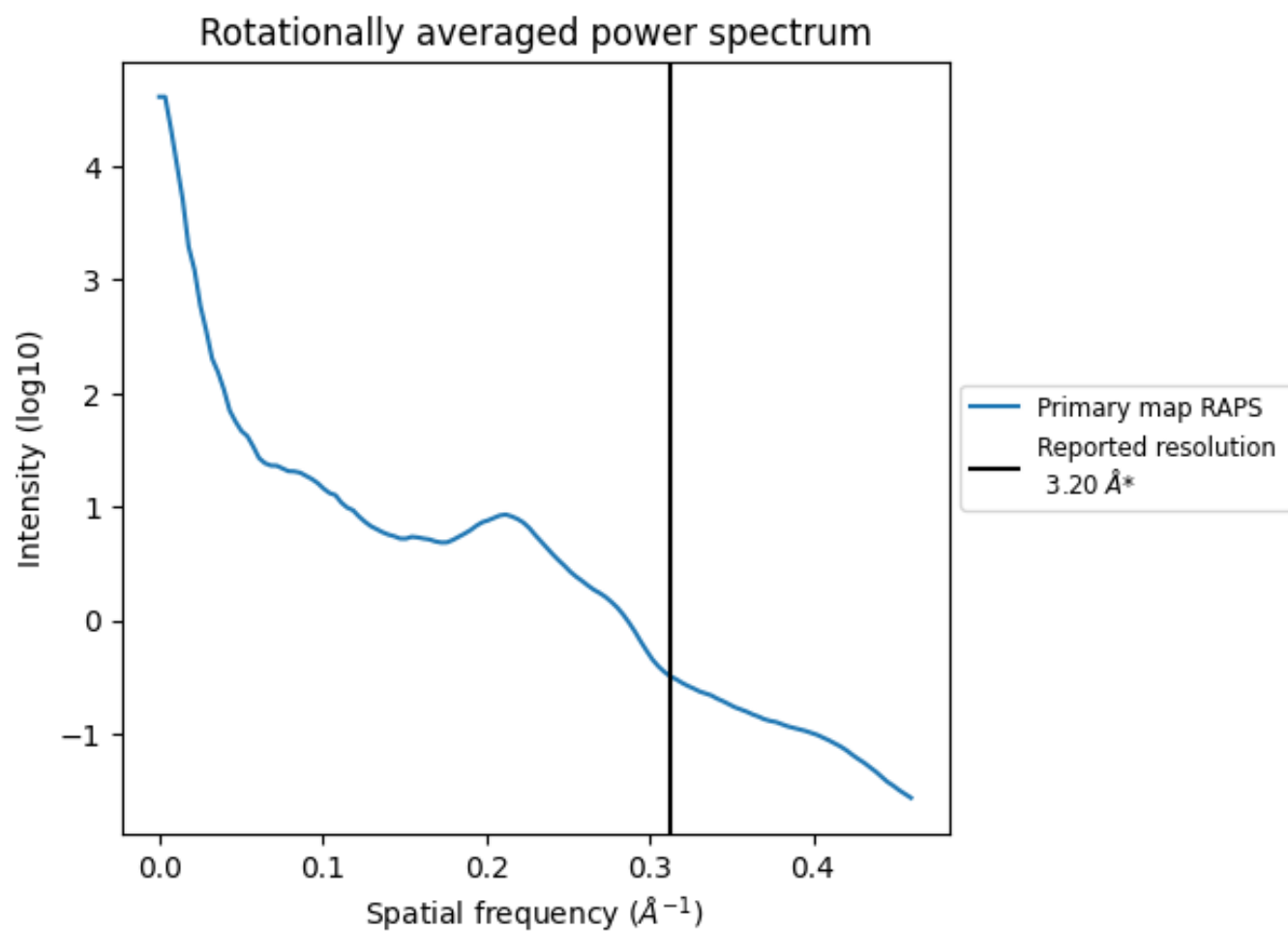
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 54 nm^3 ; this corresponds to an approximate mass of 48 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 0.312 $\text{\AA}^{-1}$

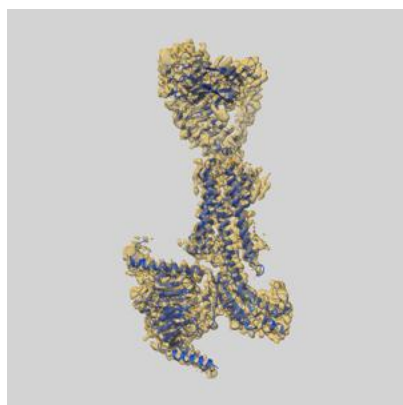
8 Fourier-Shell correlation ⓘ

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

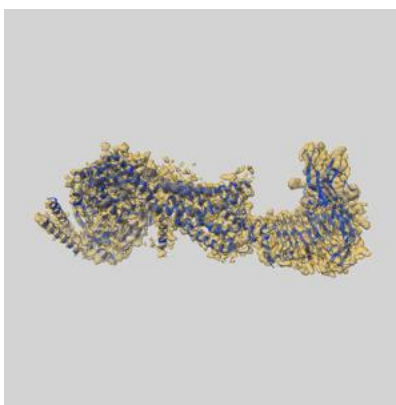
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-31597 and PDB model 7FIH. Per-residue inclusion information can be found in section [3](#) on page [7](#).

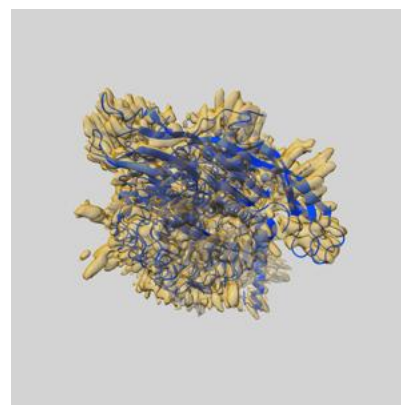
9.1 Map-model overlay [i](#)



X



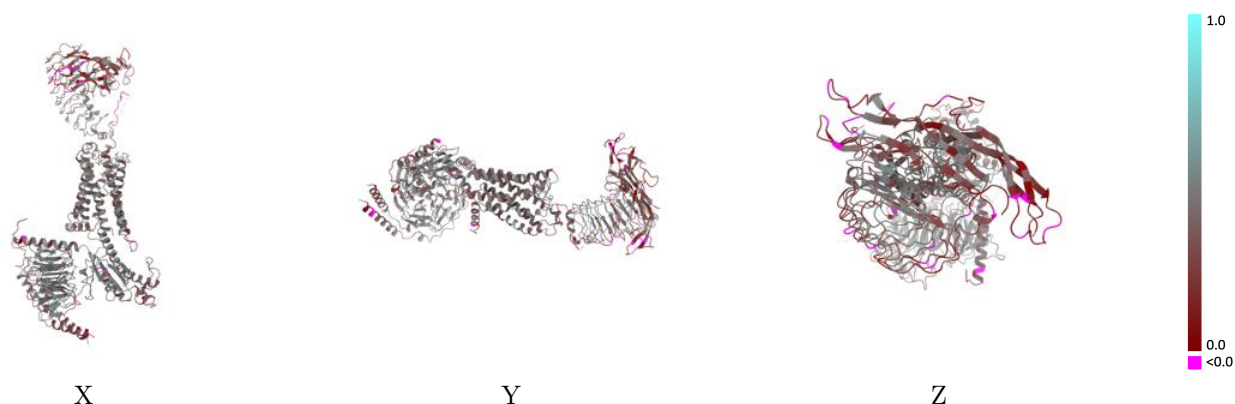
Y



Z

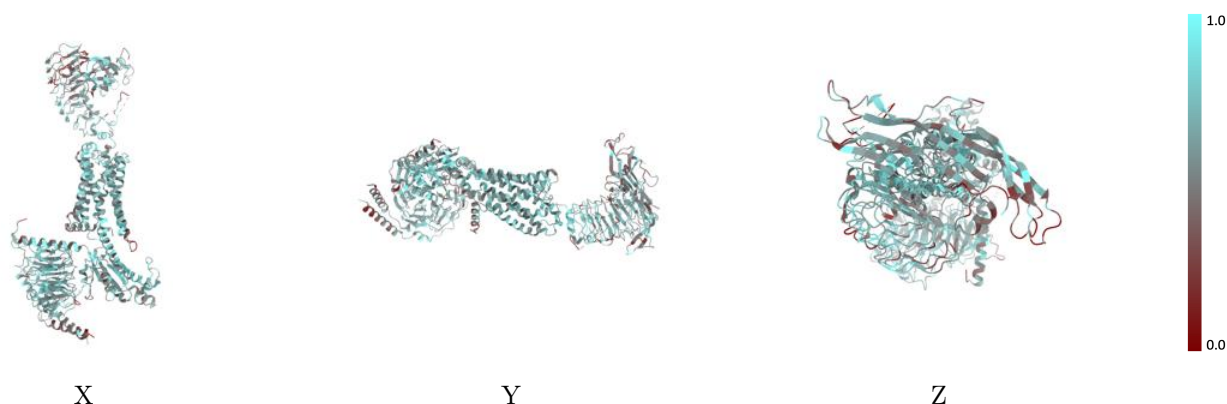
The images above show the 3D surface view of the map at the recommended contour level 0.358 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



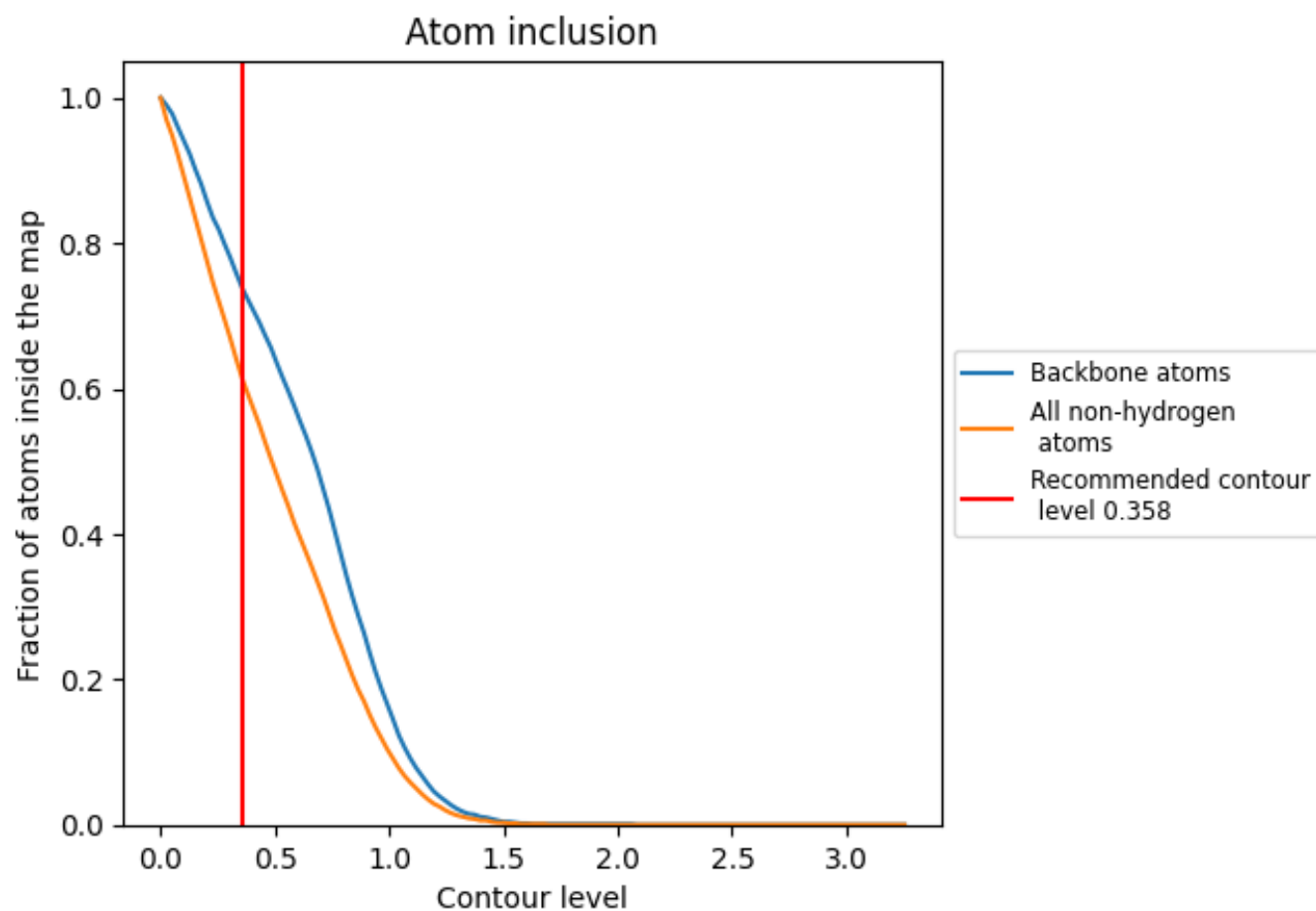
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.358).

9.4 Atom inclusion ⓘ



At the recommended contour level, 74% of all backbone atoms, 61% 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 (0.358) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6120	0.3820
A	0.6370	0.4100
B	0.6350	0.4330
G	0.5620	0.3450
R	0.6240	0.3870
X	0.5290	0.2640
Y	0.5170	0.2420

1.0

0.0

<0.0