



Full wwPDB EM Validation Report ⓘ

Apr 5, 2026 – 04:26 PM UTC

PDB ID : 9S3Z / pdb_00009s3z
EMDB ID : EMD-54556
Title : Cerebellar GluA1/4 LBD tetramer (focused refinement)
Authors : Sengupta, N.; Scrutton, A.; Greger, I.H.; Krieger, J.M.
Deposited on : 2025-07-25
Resolution : 3.55 Å (reported)
Based on initial models : ., 7OCE

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

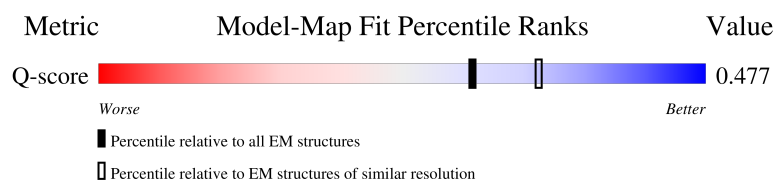
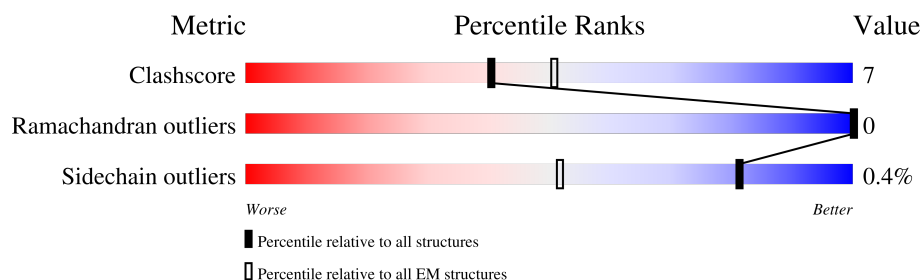
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.55 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	12819 (3.05 - 4.05)

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	888	 25% 5% 70%
1	C	888	 24% 6% 70%
2	B	863	 26% 6% 68%
2	D	863	 27% 5% 68%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8355 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 Glutamate receptor 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	268	Total	C	N	O	S	0	0
			2054	1310	347	384	13		
1	C	268	Total	C	N	O	S	0	0
			2051	1309	345	384	13		

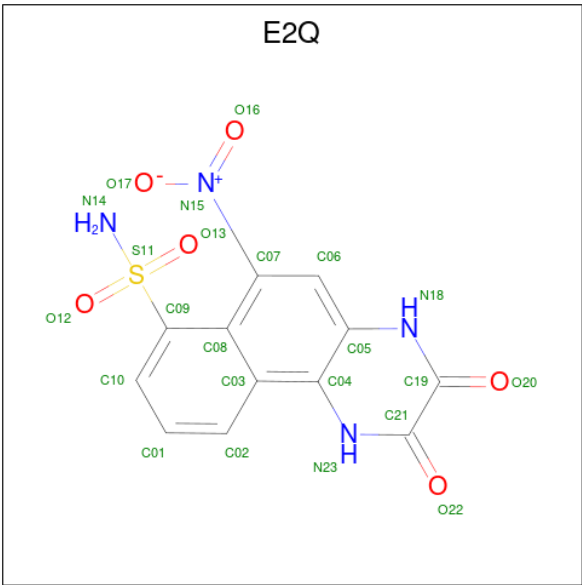
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	740	ASN	GLY	conflict	UNP A0A286ZS63
A	750	ASN	SER	conflict	UNP A0A286ZS63
A	754	LEU	VAL	conflict	UNP A0A286ZS63
A	760	ASN	SER	conflict	UNP A0A286ZS63
C	740	ASN	GLY	conflict	UNP A0A286ZS63
C	750	ASN	SER	conflict	UNP A0A286ZS63
C	754	LEU	VAL	conflict	UNP A0A286ZS63
C	760	ASN	SER	conflict	UNP A0A286ZS63

- Molecule 2 is a protein called Glutamate receptor 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	276	Total	C	N	O	S	0	0
			2081	1335	346	387	13		
2	D	275	Total	C	N	O	S	0	0
			2077	1333	345	386	13		

- Molecule 3 is 6-nitro-2,3-bis(oxidanylidene)-1,4-dihydrobenzo[f]quinoxaline-7-sulfonamide (CCD ID: E2Q) (formula: C₁₂H₈N₄O₆S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C	N	O	S	0
			23	12	4	6	1	
3	B	1	Total	C	N	O	S	0
			23	12	4	6	1	
3	C	1	Total	C	N	O	S	0
			23	12	4	6	1	
3	D	1	Total	C	N	O	S	0
			23	12	4	6	1	

- Molecule 1: Glutamate receptor 1

Chain C:



ASN	ALA	GLY	ASP	CYS	LEU	ALA	ASN	PRO	ALA	VAL	PRO	TRP	GLY	GLN	GLY	ILE	ASP	ILE	GLN	ARG	ALA	ALA	LEU	GLN	GLN	VAL	ARG	PHE	GLY	LEU	THR	GLY	ASN	GLN	VAL	GLN	PHE	ASN	GLU	GLY	LYS	GLY	ARG	ARG	THR	ASN	TYR	THR	LEU	HIS	VAL	ILE	GLU	MET	LYS	HIS	ASP	GLY	ILE	ARG
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[illegible][illegible]

GLJ GLY ARG ASP GLN THR THR SER SER GLN ASN GLJ PHE GLY ILE ILE PHE ASN SER SER TRP TRP PHE LEU LEU PHE GLN GLN GLY CYS ASP ASP SER SER PRO ARG ARG ILE VAL VAL GLY GLY VAL TRP TRP PHE PHE THR LEU ILE ILE ILE SER SER TYR

[illegible]

L699	L700	E701	S702	L703	R711	D715	N722	L730	A731	A745	K748	L749	Q752	K759	W763	G767	E768	S771	LYS	ASP	SER	GLY	SER	GLY	LYS	ASP	LYS	THR	SER	ALA	LEU	SER	SER	LEU	SER	SER	ASN	VAL	ALA	GLY	VAL	PHE	TYR	THR	ILE	LEU	ILE	GLY	GLY	LEU	LYS
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ALA	MET	LEU	VAL	ALA	LEU	ILE	GLU	PHE	CYS	TYR	LYS	SER	ARG	SER	GLU	LYS	ARG	MET	GLY	PHE	CYS	LEU	ILE	PRO	GLN	GLN	SER	ILE	ASN	GLU	ALA	ALA	ILE	ARG	THR	SER	SER	THR	LEU	PRO	ARG	ASN	SER	GLY	ALA	GLY	GLY	GLY	GLY	GLY	ASN	GLY	ARG	VAL
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SER HIS ASP PHE PRO LYS SER MET GLN SER ILE PRO CYS MET SER HIS SER THR GLY MET PRO LEU GLY ALA THR GLY LEU

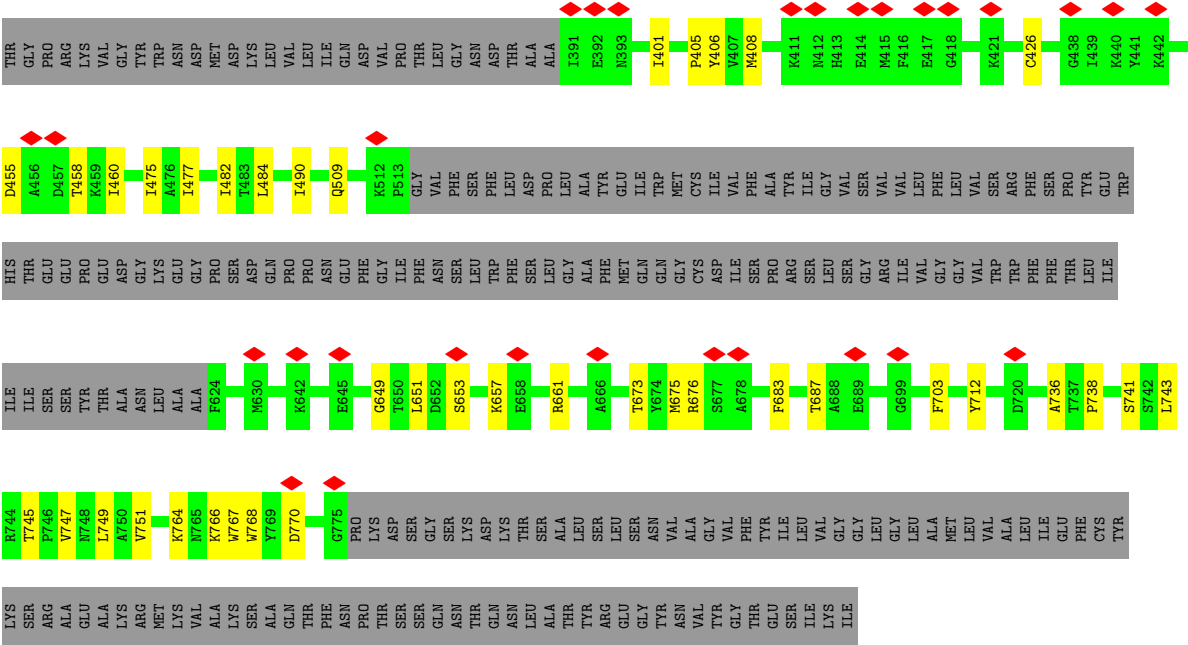
- Molecule 2: Glutamate receptor 4

Chain B:



ALA	PHE	PRO	SER	SER	VAL	GLN	GLY	LEU	PHE	ILE	ASN	THR	ASP	GLN	GLY	THR	ALA	ALA	PHE	LEU	ALA	ILE	PHE	LEU	HIS	ASN	THR	SER	PRO	PRO	ALA	ALA	SER	GLU	GLU	PRO	PHE	ASN	LEU	VAL	THR	PRO	HIS	VAL	ASP	ASN	ASN	ILE	GLU	THR	ALA	ASN	SER	PHE	ALA	ALA	VAL	THR
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	329869	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.710	Depositor
Minimum map value	-2.321	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.045	Depositor
Recommended contour level	0.48	Depositor
Map size (Å)	363.44, 363.44, 363.44	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.82600003, 0.82600003, 0.82600003	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: E2Q

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	0.12	0/2093	0.31	0/2832
1	C	0.11	0/2090	0.30	0/2828
2	B	0.10	0/2122	0.29	0/2871
2	D	0.12	0/2118	0.32	0/2866
All	All	0.11	0/8423	0.31	0/11397

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	2054	0	2015	31	0
1	C	2051	0	2008	33	0
2	B	2081	0	2042	29	0
2	D	2077	0	2039	24	0
3	A	23	0	0	0	0
3	B	23	0	0	1	0
3	C	23	0	0	0	0
3	D	23	0	0	0	0
All	All	8355	0	8104	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 7.

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)
2:B:643:GLN:HE22	2:B:646:ILE:HB	1.52	0.75
2:D:651:LEU:HD23	2:D:653:SER:H	1.51	0.74
1:A:425:ALA:HB2	1:A:472:VAL:HG21	1.71	0.73
1:C:425:ALA:HB2	1:C:472:VAL:HG21	1.70	0.71
1:A:462:GLU:HG2	1:A:467:ARG:HB2	1.78	0.66
2:D:405:PRO:HA	2:D:408:MET:HE3	1.79	0.65
1:C:645:THR:HG22	1:C:699:LEU:HB2	1.79	0.64
2:B:651:LEU:HD23	2:B:653:SER:H	1.64	0.62
2:D:673:THR:HA	2:D:676:ARG:HH21	1.64	0.62
1:C:494:LEU:HD12	1:C:495:GLY:H	1.64	0.62
1:A:496:ILE:HD13	1:A:651:THR:HG23	1.81	0.61
1:C:494:LEU:HD11	1:C:701:GLU:HB3	1.84	0.60
1:A:638:GLN:HE22	1:A:641:ILE:H	1.47	0.60
1:A:745:ALA:O	1:A:749:LEU:HG	2.02	0.59
2:D:649:GLY:HA2	2:D:675:MET:HE3	1.85	0.58
2:D:747:VAL:O	2:D:751:VAL:HG23	2.04	0.58
1:C:462:GLU:HG2	1:C:467:ARG:HB2	1.86	0.57
1:C:495:GLY:HA3	1:C:722:ASN:HB3	1.85	0.57
2:B:747:VAL:O	2:B:751:VAL:HG23	2.05	0.57
1:A:667:TRP:HA	1:A:670:MET:HG3	1.87	0.57
1:C:745:ALA:O	1:C:749:LEU:HG	2.04	0.57
1:C:629:ILE:HG21	1:C:635:LEU:HD13	1.87	0.57
1:A:664:GLU:O	1:A:668:THR:HG23	2.05	0.56
1:C:664:GLU:O	1:C:668:THR:HG23	2.06	0.56
1:A:645:THR:HG22	1:A:699:LEU:HB2	1.88	0.55
2:B:405:PRO:HA	2:B:408:MET:HE3	1.88	0.55
1:C:424:LEU:HD21	1:C:730:ILE:HD11	1.89	0.54
1:A:759:LYS:O	1:A:763:TRP:HB2	2.08	0.54
2:B:766:LYS:HA	2:B:770:ASP:HB2	1.90	0.53
1:A:629:ILE:HG21	1:A:635:LEU:HD13	1.91	0.52
1:C:397:LEU:HD12	1:C:440:ILE:HD12	1.93	0.51
1:A:744:LEU:O	1:A:748:LYS:HG3	2.09	0.51
2:B:715:GLN:HE21	2:B:769:TYR:HD1	1.58	0.51
2:D:475:ILE:HG22	2:D:477:ILE:HG13	1.93	0.50
2:B:490:ILE:HD12	2:B:736:ALA:HB1	1.93	0.49
1:A:748:LYS:HG2	2:D:484:LEU:HD22	1.95	0.49
2:D:766:LYS:HA	2:D:770:ASP:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:452:ASP:OD1	1:C:453:THR:HG23	2.13	0.49
1:C:667:TRP:HA	1:C:670:MET:HG3	1.94	0.49
1:C:681:THR:HG23	1:C:684:GLU:H	1.77	0.49
2:B:407:VAL:HG22	2:B:427:VAL:HG23	1.94	0.48
2:B:738:PRO:HG2	2:B:741:SER:HB2	1.94	0.48
1:A:447:GLY:HA3	1:A:460:VAL:HB	1.96	0.48
2:B:406:TYR:HB3	2:B:426:CYS:SG	2.54	0.48
1:C:748:LYS:O	1:C:752:GLN:HG3	2.14	0.47
2:B:405:PRO:HB3	2:B:712:TYR:CE1	2.50	0.47
1:A:424:LEU:HD21	1:A:730:ILE:HD11	1.96	0.47
2:B:407:VAL:HG13	2:B:422:TYR:HB3	1.96	0.47
1:A:690:ARG:HG2	1:A:715:ASP:OD2	2.15	0.47
2:B:437:ILE:HB	2:B:439:ILE:HG22	1.97	0.47
1:C:711:ARG:HH21	1:C:767:GLY:HA2	1.80	0.46
2:D:455:ASP:HB3	2:D:458:THR:HG22	1.96	0.46
1:A:638:GLN:HE22	1:A:641:ILE:N	2.13	0.46
1:C:683:GLU:HA	1:C:686:MET:HE2	1.97	0.46
2:D:406:TYR:HB3	2:D:426:CYS:SG	2.55	0.46
2:B:478:ALA:HB1	2:B:479:PRO:HD2	1.97	0.46
2:B:703:PHE:CD2	2:B:705:LEU:HD23	2.51	0.46
2:D:405:PRO:HB3	2:D:712:TYR:CE1	2.50	0.46
1:C:638:GLN:HE22	1:C:641:ILE:H	1.63	0.46
1:C:679:VAL:HG11	1:C:685:GLY:N	2.31	0.46
2:D:675:MET:HE2	2:D:683:PHE:HE1	1.80	0.46
2:D:738:PRO:HG2	2:D:741:SER:HB2	1.98	0.45
2:B:497:MET:HE3	2:B:708:THR:HG21	1.98	0.45
1:C:401:TYR:HB3	1:C:421:CYS:SG	2.56	0.45
1:A:644:GLY:HA2	1:A:670:MET:HE1	1.98	0.45
2:B:710:ASN:OD1	2:B:723:LYS:HB2	2.16	0.45
1:A:638:GLN:NE2	1:A:641:ILE:H	2.13	0.45
1:C:759:LYS:O	1:C:763:TRP:HB2	2.17	0.45
2:D:745:THR:O	2:D:749:LEU:HD23	2.17	0.45
2:B:435:LYS:HD3	2:B:435:LYS:C	2.42	0.44
1:A:401:TYR:CD2	1:A:474:PRO:HG3	2.53	0.44
1:A:678:PHE:N	1:A:678:PHE:CD1	2.85	0.44
1:A:405:LYS:HG2	1:A:418:GLU:OE1	2.17	0.44
1:A:489:LYS:HE2	2:D:482:ILE:HD13	1.99	0.44
1:C:392:ILE:HG23	1:C:439:GLU:HG3	1.99	0.44
1:C:748:LYS:HE3	1:C:748:LYS:HB3	1.73	0.44
2:B:745:THR:O	2:B:749:LEU:HD23	2.18	0.44
1:C:494:LEU:HD21	1:C:701:GLU:OE1	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:397:VAL:HB	2:B:473:ALA:HA	1.99	0.44
2:B:400:THR:HG21	2:B:407:VAL:HG21	1.99	0.44
2:D:490:ILE:HD12	2:D:736:ALA:HB1	2.00	0.44
1:A:460:VAL:HG13	1:A:485:ILE:HD13	1.99	0.43
2:B:673:THR:HA	2:B:676:ARG:HH21	1.83	0.43
2:D:687:THR:HG23	2:D:703:PHE:HE2	1.83	0.43
1:A:656:ARG:HG3	1:A:667:TRP:CZ2	2.52	0.43
2:B:475:ILE:HG22	2:B:477:ILE:HG13	2.01	0.43
1:C:678:PHE:CD1	1:C:678:PHE:N	2.87	0.43
2:B:703:PHE:HD2	2:B:705:LEU:HD23	1.83	0.43
1:A:643:TYR:CE1	1:A:666:MET:HG2	2.53	0.43
1:C:460:VAL:O	1:C:464:VAL:HG23	2.19	0.43
1:C:686:MET:HE2	1:C:686:MET:HB3	1.86	0.43
2:D:509:GLN:N	2:D:509:GLN:OE1	2.52	0.43
1:A:401:TYR:HB3	1:A:421:CYS:SG	2.58	0.42
2:B:430:ALA:HA	2:B:477:ILE:HD13	2.00	0.42
2:B:766:LYS:HE2	2:B:771:LYS:NZ	2.34	0.42
1:C:494:LEU:HD13	1:C:703:THR:HG23	2.00	0.42
2:B:640:LEU:HG	2:B:648:TYR:CD2	2.54	0.42
1:A:495:GLY:HA3	1:A:722:ASN:HB3	2.01	0.42
1:A:400:PRO:HG3	1:A:707:TYR:CE1	2.55	0.42
1:C:447:GLY:HA3	1:C:460:VAL:HB	2.01	0.42
1:C:485:ILE:HD12	1:C:731:ALA:HB1	2.00	0.42
2:D:657:LYS:HE2	2:D:661:ARG:HH21	1.85	0.42
2:B:709:MET:HG2	3:B:901:E2Q:O12	2.20	0.42
2:D:475:ILE:HD11	2:D:743:LEU:HD13	2.01	0.41
2:D:458:THR:HG23	2:D:460:ILE:H	1.85	0.41
1:A:748:LYS:HB3	1:A:748:LYS:HE3	1.70	0.41
1:C:680:ARG:H	1:C:680:ARG:HG2	1.53	0.41
1:C:396:ILE:HG21	1:C:446:TYR:CE1	2.55	0.41
1:C:396:ILE:HG21	1:C:446:TYR:HE1	1.85	0.41
1:A:680:ARG:HB3	1:A:680:ARG:CZ	2.51	0.41
1:A:460:VAL:O	1:A:464:VAL:HG23	2.20	0.41
2:B:649:GLY:HA2	2:B:675:MET:HE3	2.03	0.40
2:D:651:LEU:HD23	2:D:653:SER:N	2.27	0.40
2:D:766:LYS:HE3	2:D:767:TRP:NE1	2.37	0.40
2:D:764:LYS:O	2:D:768:TRP:HB2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all 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	264/888 (30%)	257 (97%)	7 (3%)	0	100	100
1	C	264/888 (30%)	261 (99%)	3 (1%)	0	100	100
2	B	272/863 (32%)	270 (99%)	2 (1%)	0	100	100
2	D	271/863 (31%)	263 (97%)	8 (3%)	0	100	100
All	All	1071/3502 (31%)	1051 (98%)	20 (2%)	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	210/757 (28%)	210 (100%)	0	100	100
1	C	209/757 (28%)	209 (100%)	0	100	100
2	B	211/740 (28%)	209 (99%)	2 (1%)	70	76
2	D	211/740 (28%)	210 (100%)	1 (0%)	81	80
All	All	841/2994 (28%)	838 (100%)	3 (0%)	81	80

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

Mol	Chain	Res	Type
2	B	394	ARG
2	B	401	ILE
2	D	401	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	643	GLN
1	C	710	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	$\# Z > 2$
3	E2Q	B	901	-	25,25,25	2.86	10 (40%)	32,39,39	4.07	12 (37%)
3	E2Q	C	901	-	25,25,25	2.87	10 (40%)	32,39,39	4.07	12 (37%)
3	E2Q	D	901	-	25,25,25	2.88	10 (40%)	32,39,39	4.06	12 (37%)
3	E2Q	A	901	-	25,25,25	2.84	10 (40%)	32,39,39	4.08	13 (40%)

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
3	E2Q	B	901	-	-	2/8/10/10	0/3/3/3
3	E2Q	C	901	-	-	6/8/10/10	0/3/3/3
3	E2Q	D	901	-	-	4/8/10/10	0/3/3/3
3	E2Q	A	901	-	-	6/8/10/10	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	901	E2Q	C05-N18	5.95	1.50	1.39
3	D	901	E2Q	C05-N18	5.87	1.49	1.39
3	A	901	E2Q	C05-N18	5.78	1.49	1.39
3	B	901	E2Q	C19-N18	5.76	1.43	1.35
3	D	901	E2Q	C19-N18	5.76	1.43	1.35
3	C	901	E2Q	C21-N23	5.75	1.43	1.35
3	A	901	E2Q	C19-N18	5.71	1.43	1.35
3	C	901	E2Q	C05-N18	5.63	1.49	1.39
3	B	901	E2Q	S11-N14	5.56	1.71	1.60
3	D	901	E2Q	S11-N14	5.55	1.71	1.60
3	C	901	E2Q	C19-N18	5.54	1.43	1.35
3	C	901	E2Q	S11-N14	5.50	1.71	1.60
3	D	901	E2Q	C21-N23	5.46	1.43	1.35
3	A	901	E2Q	S11-N14	5.40	1.71	1.60
3	A	901	E2Q	C21-N23	5.34	1.43	1.35
3	B	901	E2Q	C21-N23	5.17	1.43	1.35
3	C	901	E2Q	C04-N23	5.16	1.49	1.39
3	D	901	E2Q	C04-N23	4.81	1.49	1.39
3	A	901	E2Q	C04-N23	4.79	1.49	1.39
3	B	901	E2Q	C04-N23	4.68	1.48	1.39
3	C	901	E2Q	C09-S11	3.19	1.82	1.77
3	D	901	E2Q	C09-S11	3.02	1.81	1.77
3	B	901	E2Q	C09-S11	2.99	1.81	1.77
3	A	901	E2Q	C09-S11	2.94	1.81	1.77
3	D	901	E2Q	O16-N15	-2.89	1.17	1.22
3	B	901	E2Q	O16-N15	-2.84	1.17	1.22
3	A	901	E2Q	O16-N15	-2.82	1.17	1.22
3	C	901	E2Q	O16-N15	-2.81	1.18	1.22
3	B	901	E2Q	O22-C21	-2.67	1.18	1.23
3	D	901	E2Q	O22-C21	-2.63	1.18	1.23
3	A	901	E2Q	O22-C21	-2.58	1.18	1.23
3	C	901	E2Q	O20-C19	-2.58	1.18	1.23
3	A	901	E2Q	O20-C19	-2.49	1.18	1.23
3	B	901	E2Q	C03-C08	-2.42	1.38	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	901	E2Q	O20-C19	-2.38	1.19	1.23
3	A	901	E2Q	C03-C08	-2.34	1.38	1.42
3	D	901	E2Q	C03-C08	-2.33	1.38	1.42
3	B	901	E2Q	O20-C19	-2.26	1.19	1.23
3	C	901	E2Q	O22-C21	-2.21	1.19	1.23
3	C	901	E2Q	C03-C08	-2.14	1.38	1.42

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	E2Q	C08-C09-S11	12.24	136.96	121.74
3	C	901	E2Q	C08-C09-S11	12.15	136.84	121.74
3	D	901	E2Q	C08-C09-S11	12.08	136.76	121.74
3	B	901	E2Q	C08-C09-S11	12.06	136.73	121.74
3	B	901	E2Q	O13-S11-O12	-11.60	100.95	118.80
3	D	901	E2Q	O13-S11-O12	-11.52	101.08	118.80
3	C	901	E2Q	O13-S11-O12	-11.39	101.28	118.80
3	A	901	E2Q	O13-S11-O12	-11.28	101.44	118.80
3	A	901	E2Q	C10-C09-S11	-10.81	102.78	117.53
3	D	901	E2Q	C10-C09-S11	-10.63	103.03	117.53
3	B	901	E2Q	C10-C09-S11	-10.58	103.10	117.53
3	C	901	E2Q	C10-C09-S11	-10.48	103.23	117.53
3	C	901	E2Q	C09-C08-C03	6.54	120.37	116.78
3	D	901	E2Q	C09-C08-C03	5.81	119.96	116.78
3	B	901	E2Q	C09-C08-C03	5.80	119.96	116.78
3	A	901	E2Q	C09-C08-C03	5.34	119.71	116.78
3	B	901	E2Q	C19-C21-N23	5.08	120.37	116.82
3	D	901	E2Q	C19-C21-N23	4.65	120.07	116.82
3	C	901	E2Q	C21-C19-N18	4.57	120.02	116.82
3	A	901	E2Q	C19-C21-N23	4.51	119.97	116.82
3	A	901	E2Q	C21-C19-N18	4.07	119.66	116.82
3	D	901	E2Q	C21-C19-N18	3.64	119.36	116.82
3	A	901	E2Q	C05-N18-C19	-3.55	119.85	124.82
3	C	901	E2Q	C19-C21-N23	3.48	119.25	116.82
3	D	901	E2Q	C05-N18-C19	-3.44	120.01	124.82
3	B	901	E2Q	C05-N18-C19	-3.42	120.03	124.82
3	C	901	E2Q	C05-N18-C19	-3.36	120.12	124.82
3	B	901	E2Q	C21-C19-N18	3.20	119.06	116.82
3	B	901	E2Q	O12-S11-N14	3.18	111.93	107.35
3	D	901	E2Q	O12-S11-N14	3.16	111.90	107.35
3	B	901	E2Q	O13-S11-N14	3.03	111.72	107.35
3	A	901	E2Q	O12-S11-N14	3.00	111.67	107.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	E2Q	O12-S11-C09	2.99	111.60	107.26
3	D	901	E2Q	O13-S11-N14	2.95	111.60	107.35
3	C	901	E2Q	O13-S11-C09	2.90	111.47	107.26
3	A	901	E2Q	O13-S11-C09	2.89	111.45	107.26
3	C	901	E2Q	O12-S11-C09	2.89	111.45	107.26
3	A	901	E2Q	O13-S11-N14	2.85	111.46	107.35
3	D	901	E2Q	O12-S11-C09	2.82	111.35	107.26
3	B	901	E2Q	O12-S11-C09	2.73	111.22	107.26
3	A	901	E2Q	C02-C03-C04	-2.66	119.91	124.70
3	C	901	E2Q	O12-S11-N14	2.58	111.06	107.35
3	C	901	E2Q	O13-S11-N14	2.55	111.03	107.35
3	B	901	E2Q	C02-C03-C04	-2.44	120.31	124.70
3	D	901	E2Q	C02-C03-C04	-2.43	120.32	124.70
3	C	901	E2Q	C02-C03-C04	-2.32	120.53	124.70
3	A	901	E2Q	C04-N23-C21	-2.11	119.93	124.80
3	B	901	E2Q	C04-N23-C21	-2.04	120.07	124.80
3	D	901	E2Q	C04-N23-C21	-2.02	120.13	124.80

There are no chirality outliers.

All (18) torsion outliers are listed below:

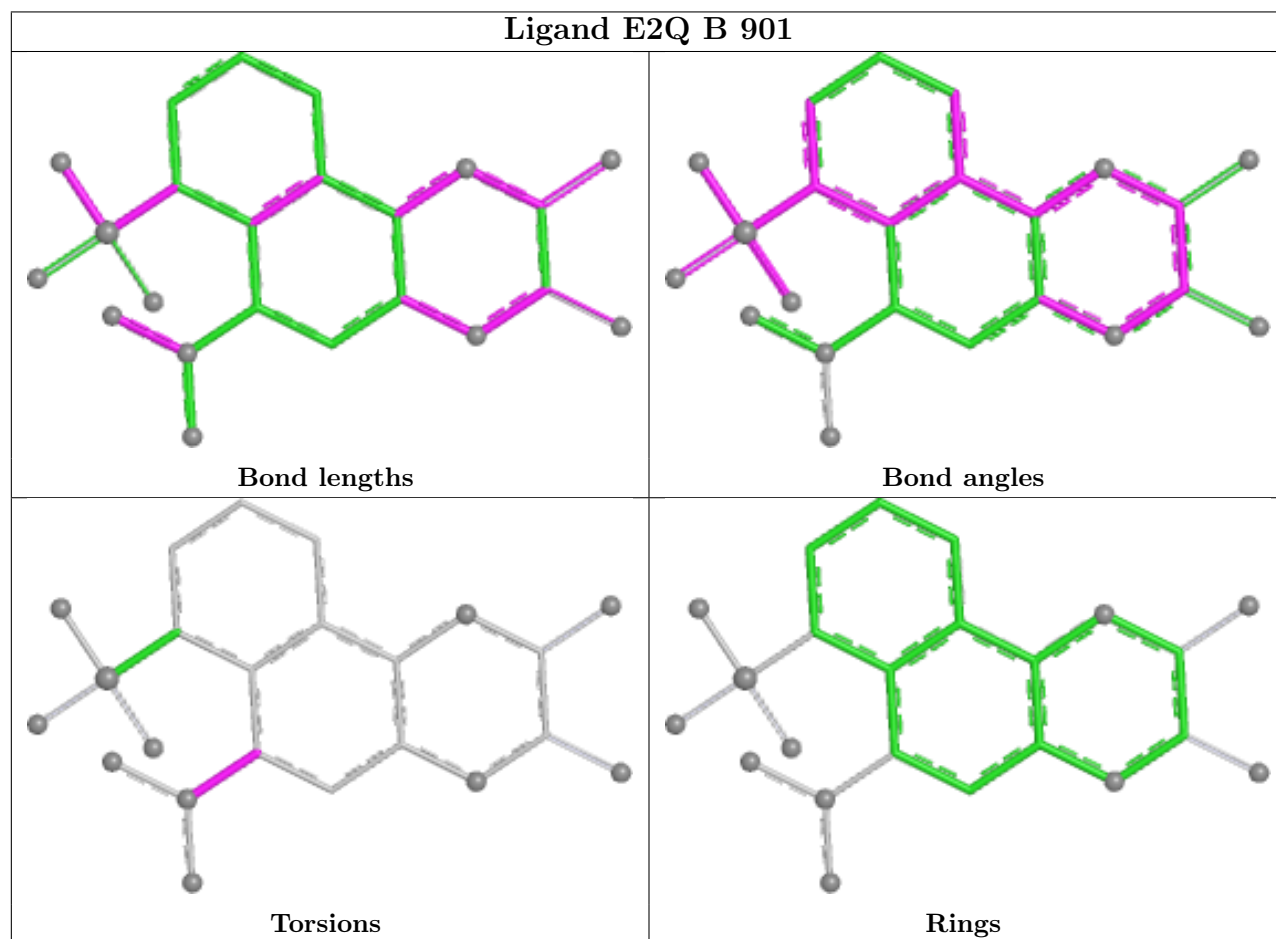
Mol	Chain	Res	Type	Atoms
3	A	901	E2Q	C08-C09-S11-N14
3	A	901	E2Q	C08-C09-S11-O12
3	B	901	E2Q	C06-C07-N15-O16
3	B	901	E2Q	C08-C07-N15-O16
3	C	901	E2Q	C08-C09-S11-O13
3	D	901	E2Q	C06-C07-N15-O16
3	D	901	E2Q	C08-C07-N15-O16
3	A	901	E2Q	C10-C09-S11-N14
3	D	901	E2Q	C08-C09-S11-O12
3	A	901	E2Q	C10-C09-S11-O12
3	A	901	E2Q	C06-C07-N15-O16
3	C	901	E2Q	C06-C07-N15-O16
3	C	901	E2Q	C10-C09-S11-O13
3	A	901	E2Q	C08-C07-N15-O16
3	C	901	E2Q	C08-C07-N15-O16
3	C	901	E2Q	C10-C09-S11-N14
3	C	901	E2Q	C08-C09-S11-N14
3	D	901	E2Q	C10-C09-S11-O12

There are no ring outliers.

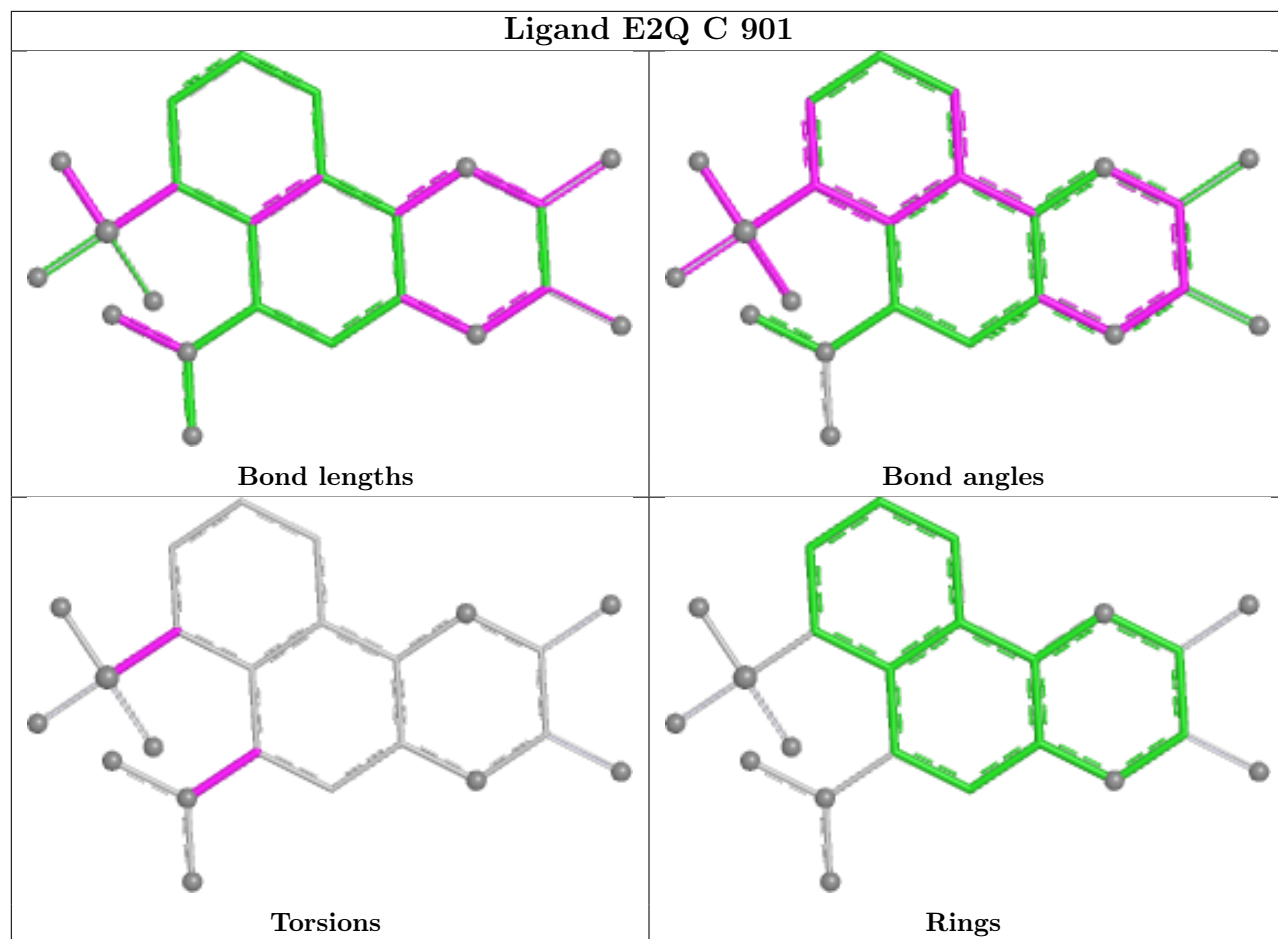
1 monomer is involved in 1 short contact:

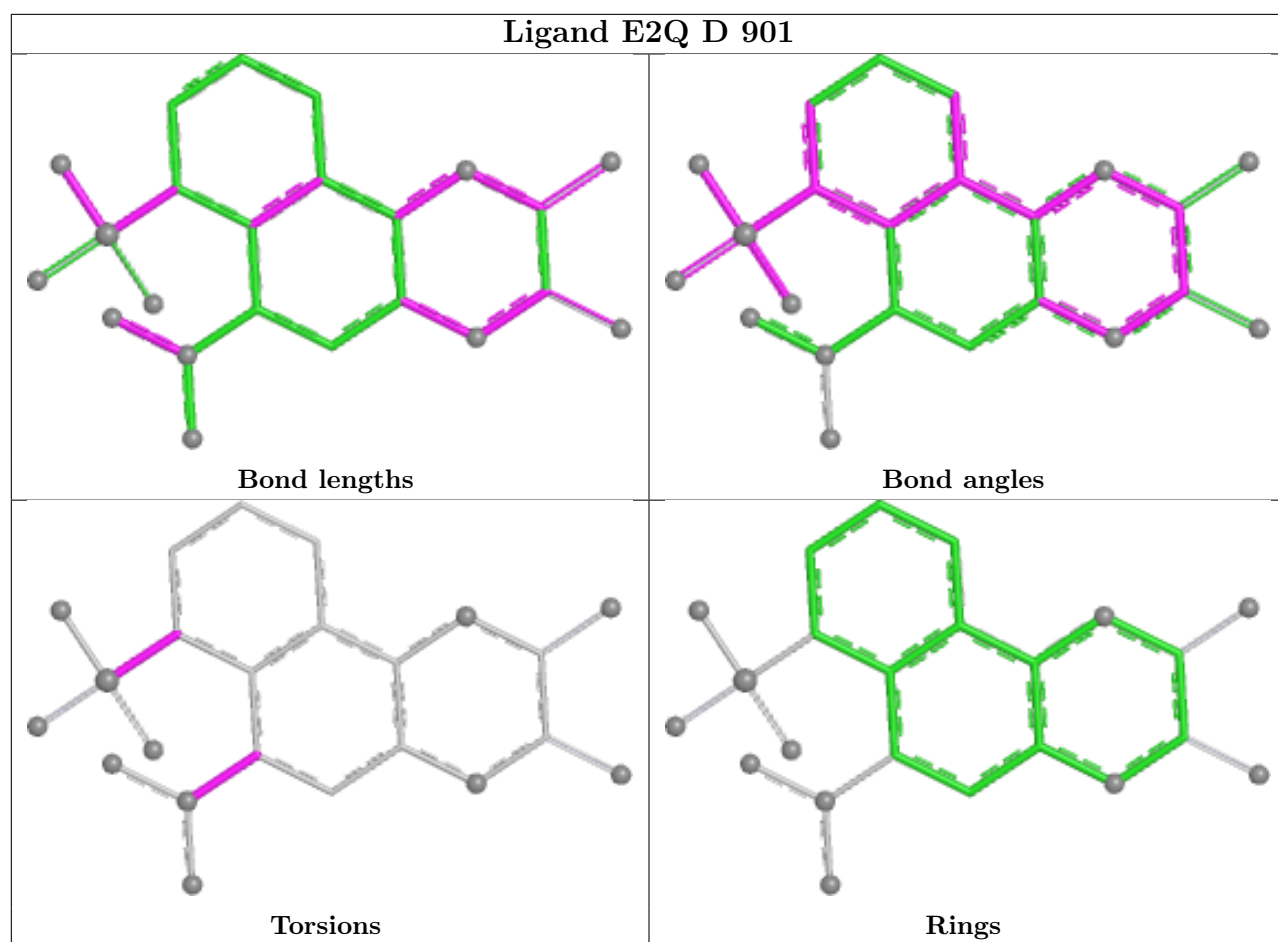
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	901	E2Q	1	0

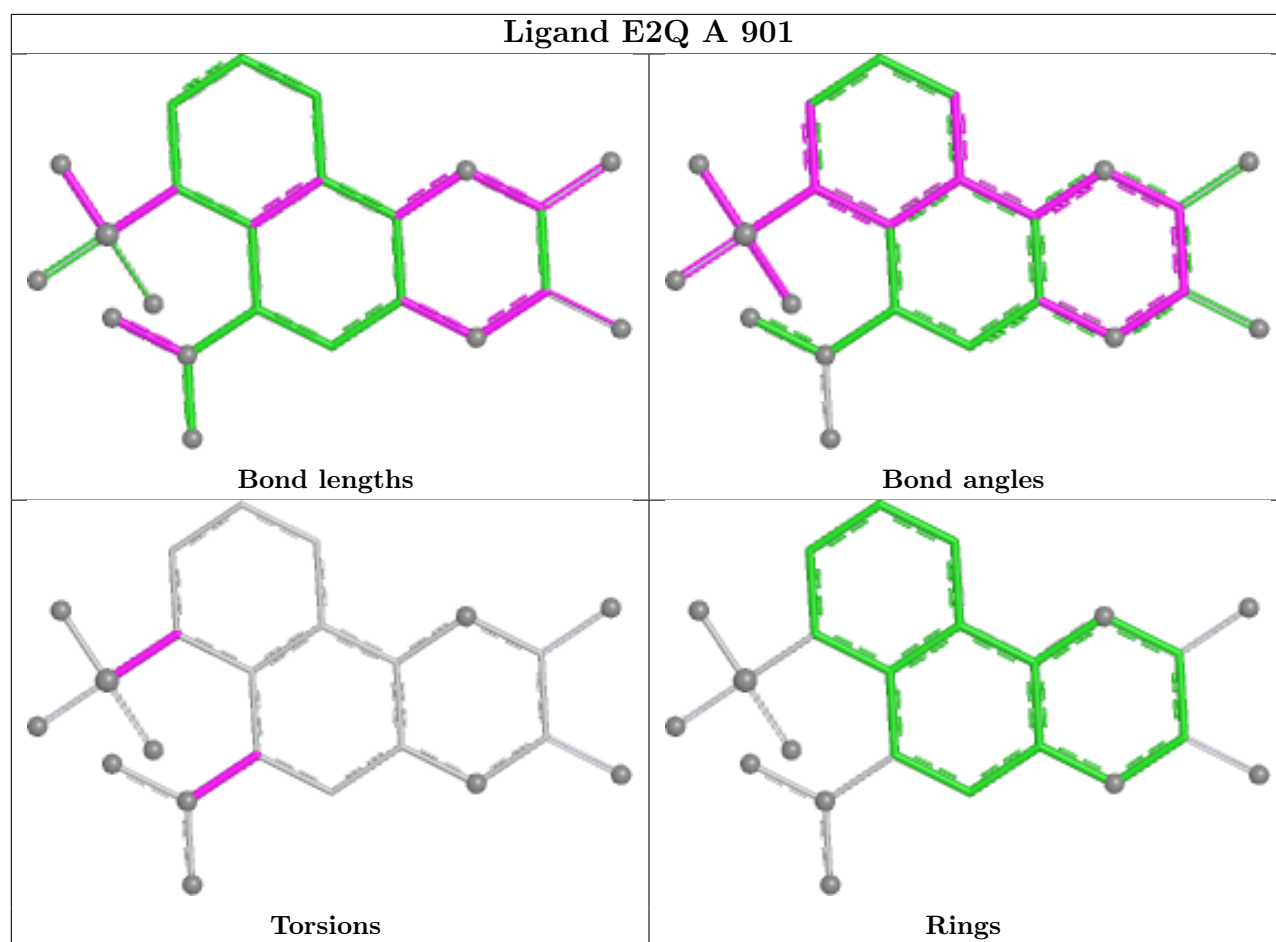
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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 E2Q C 901







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.

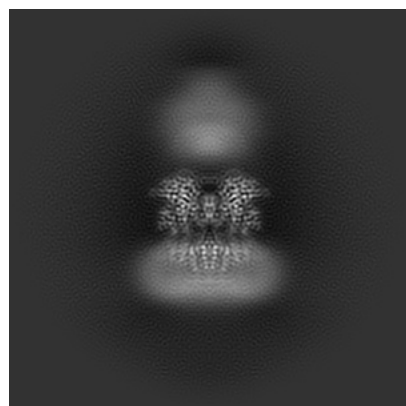
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-54556. These allow visual inspection of the internal detail of the map and identification of artifacts.

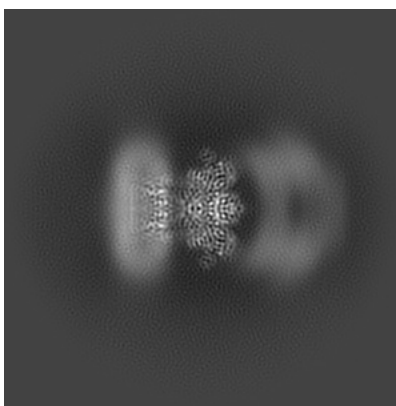
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

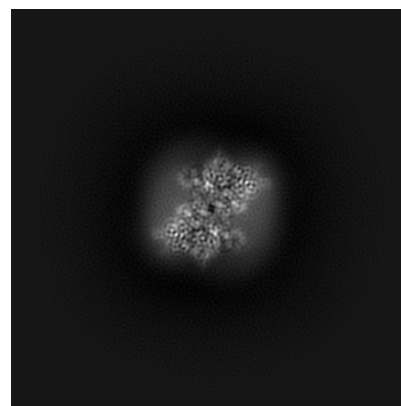
6.1.1 Primary map



X

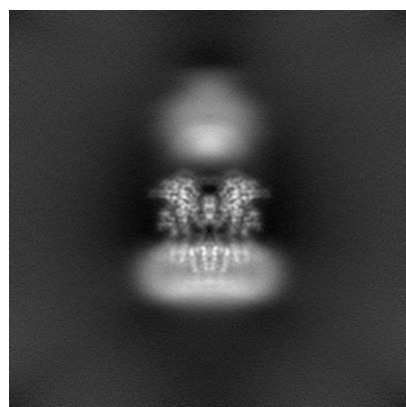


Y

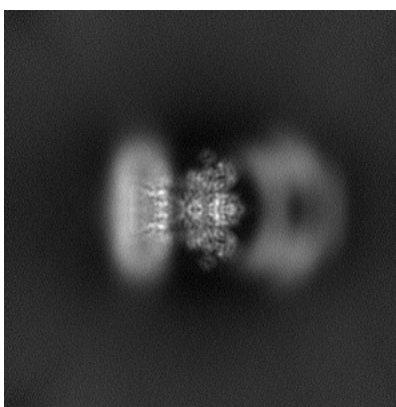


Z

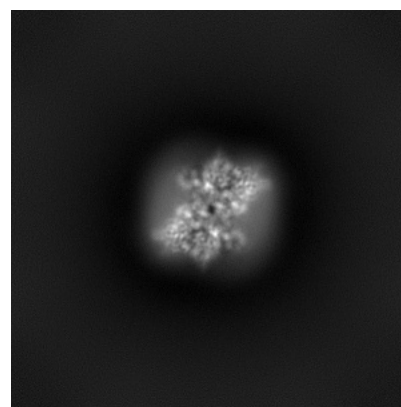
6.1.2 Raw map



X



Y

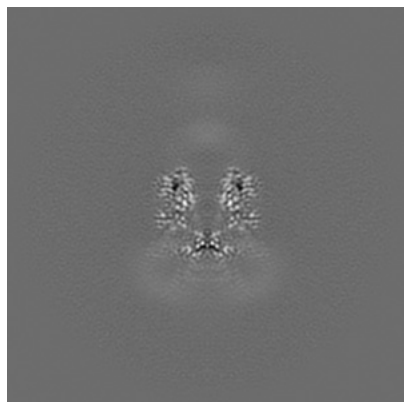


Z

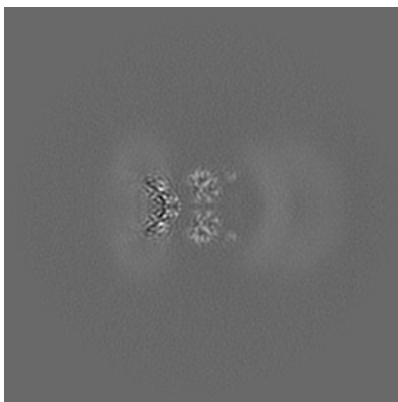
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

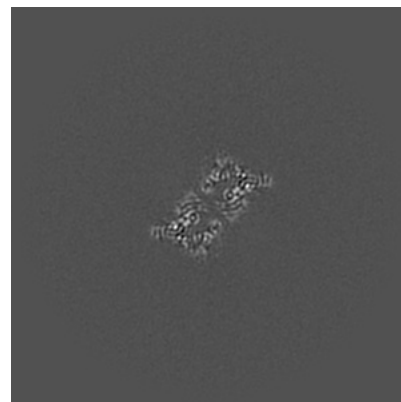
6.2.1 Primary map



X Index: 220

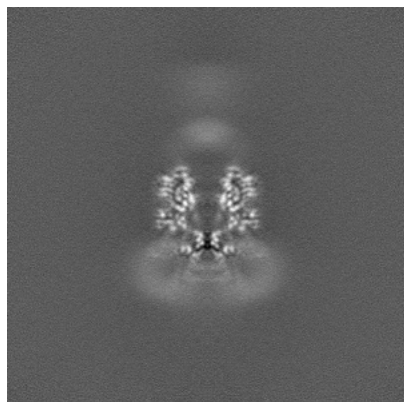


Y Index: 220

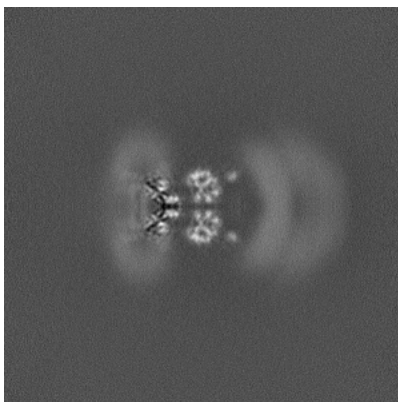


Z Index: 220

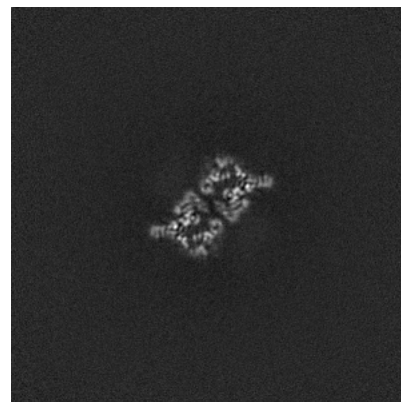
6.2.2 Raw map



X Index: 220



Y Index: 220

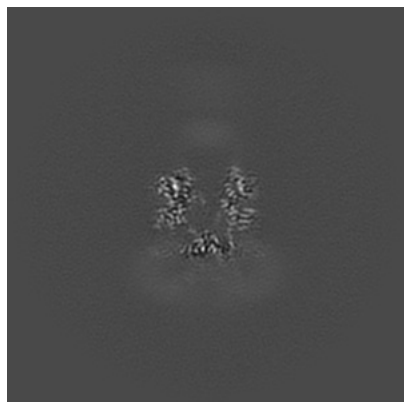


Z Index: 220

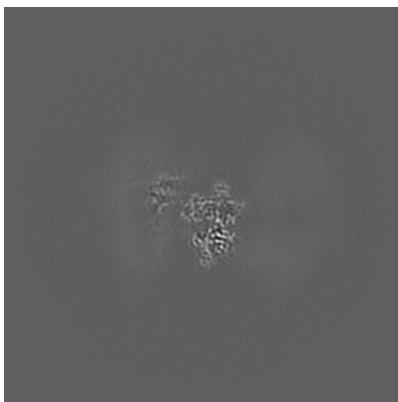
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

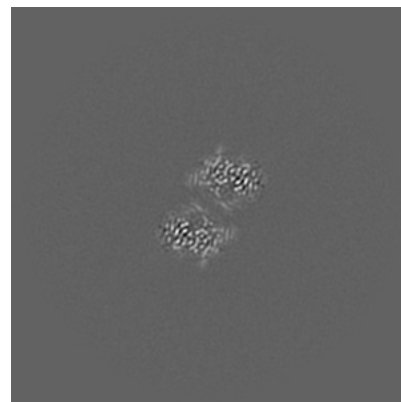
6.3.1 Primary map



X Index: 217

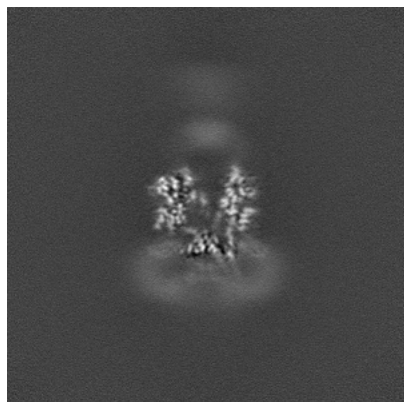


Y Index: 192

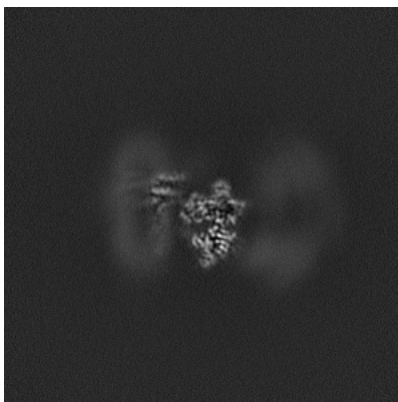


Z Index: 234

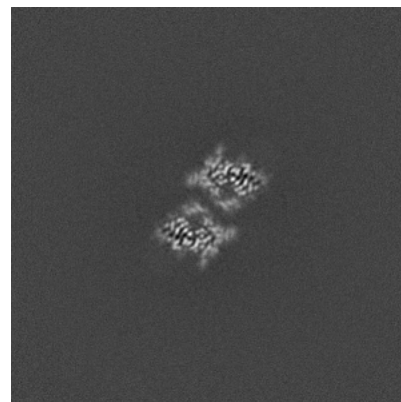
6.3.2 Raw map



X Index: 216



Y Index: 191

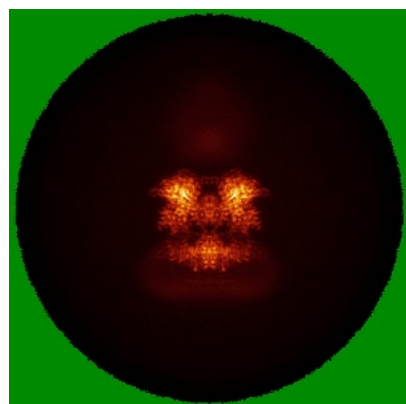


Z Index: 233

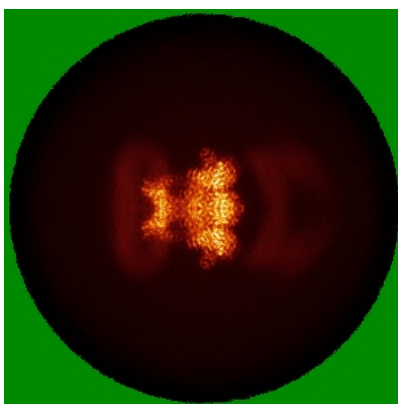
The images above show the largest variance slices of the map in three orthogonal directions.

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

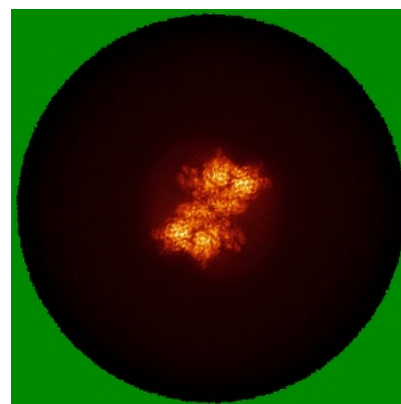
6.4.1 Primary map



X

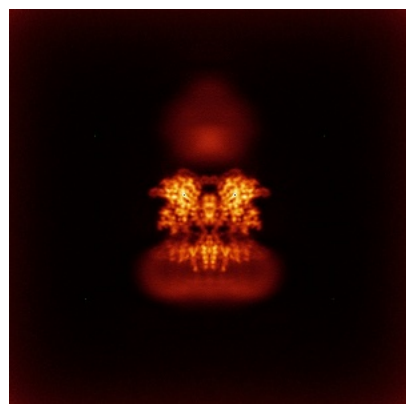


Y

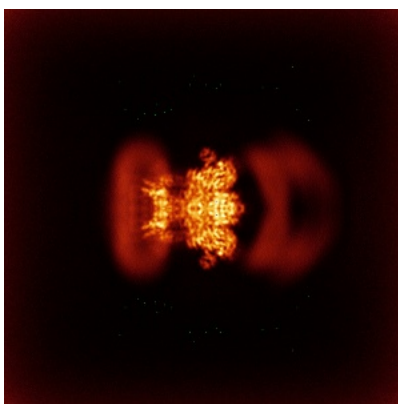


Z

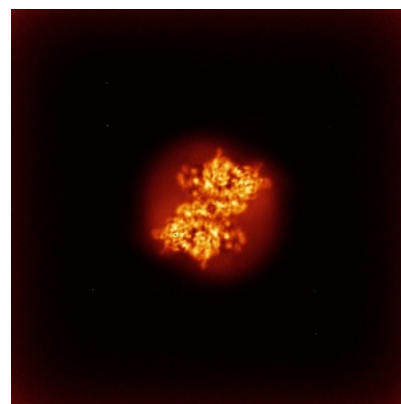
6.4.2 Raw map



X



Y

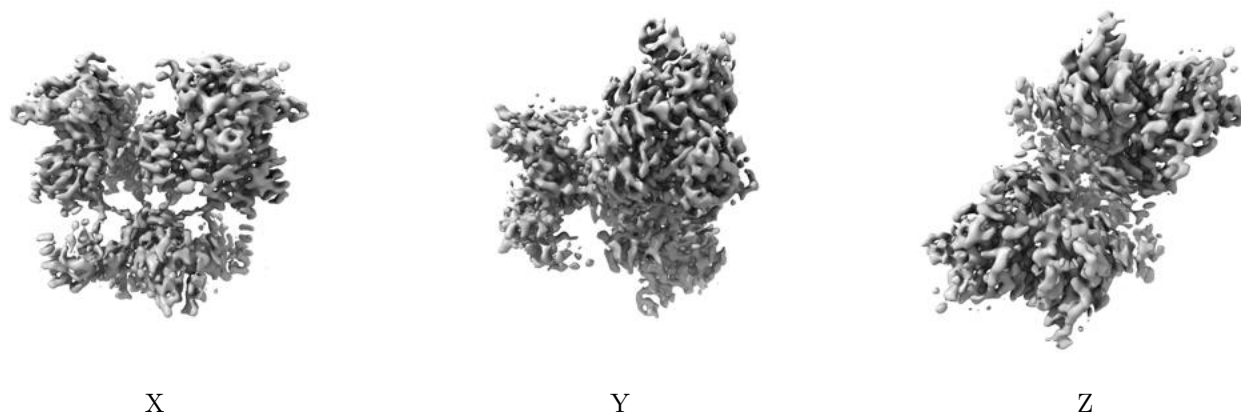


Z

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.

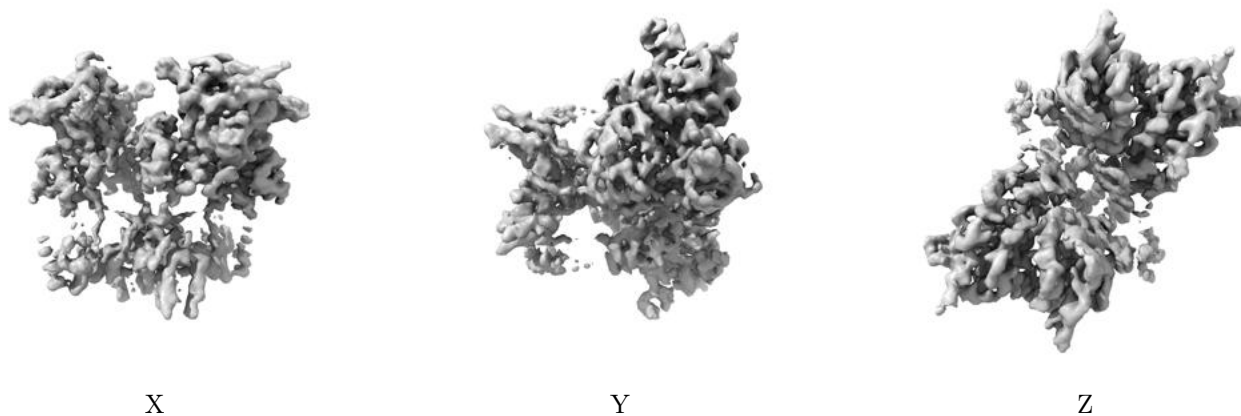
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.48. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

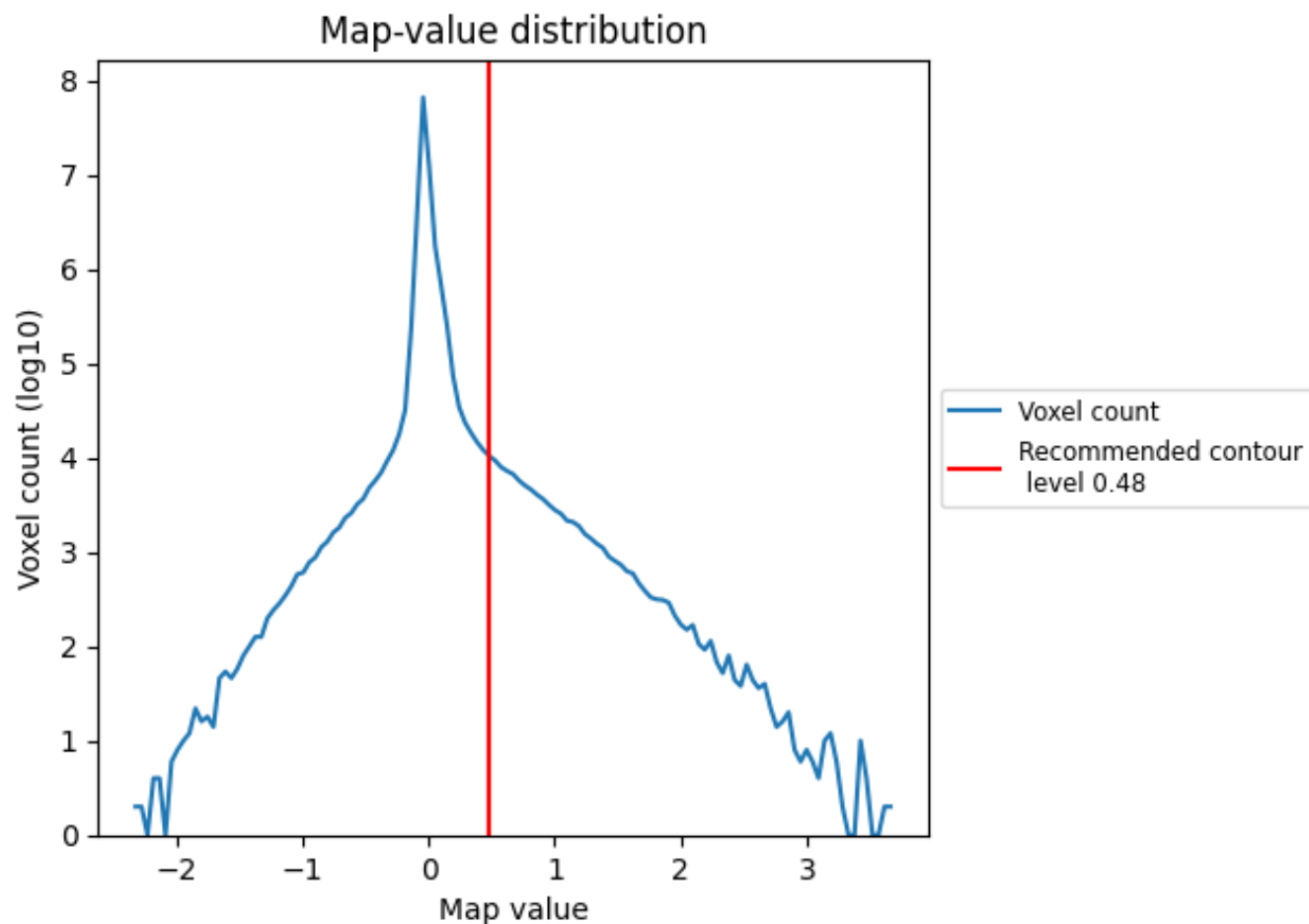
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

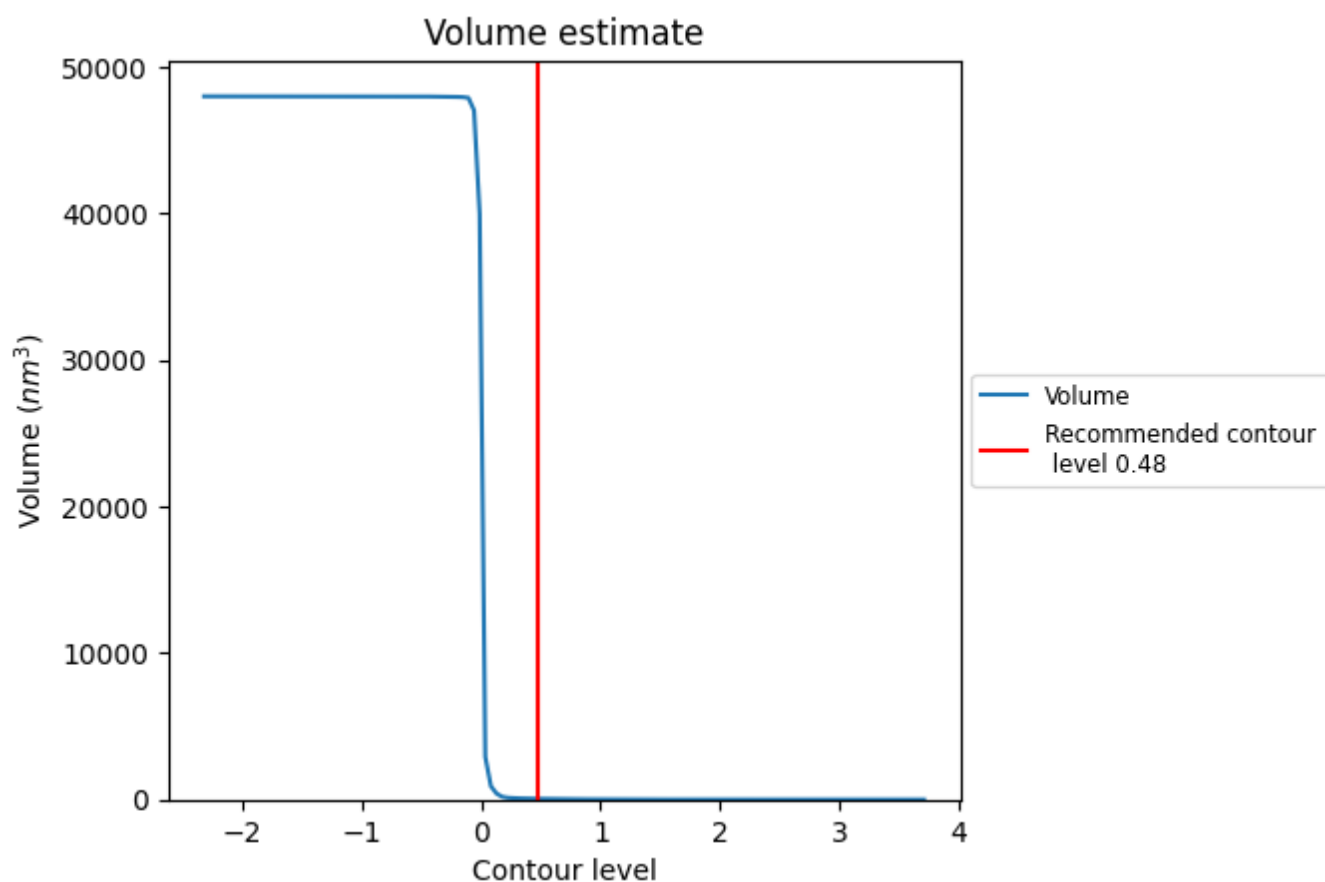
This section contains the results of statistical analysis of the map.

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.

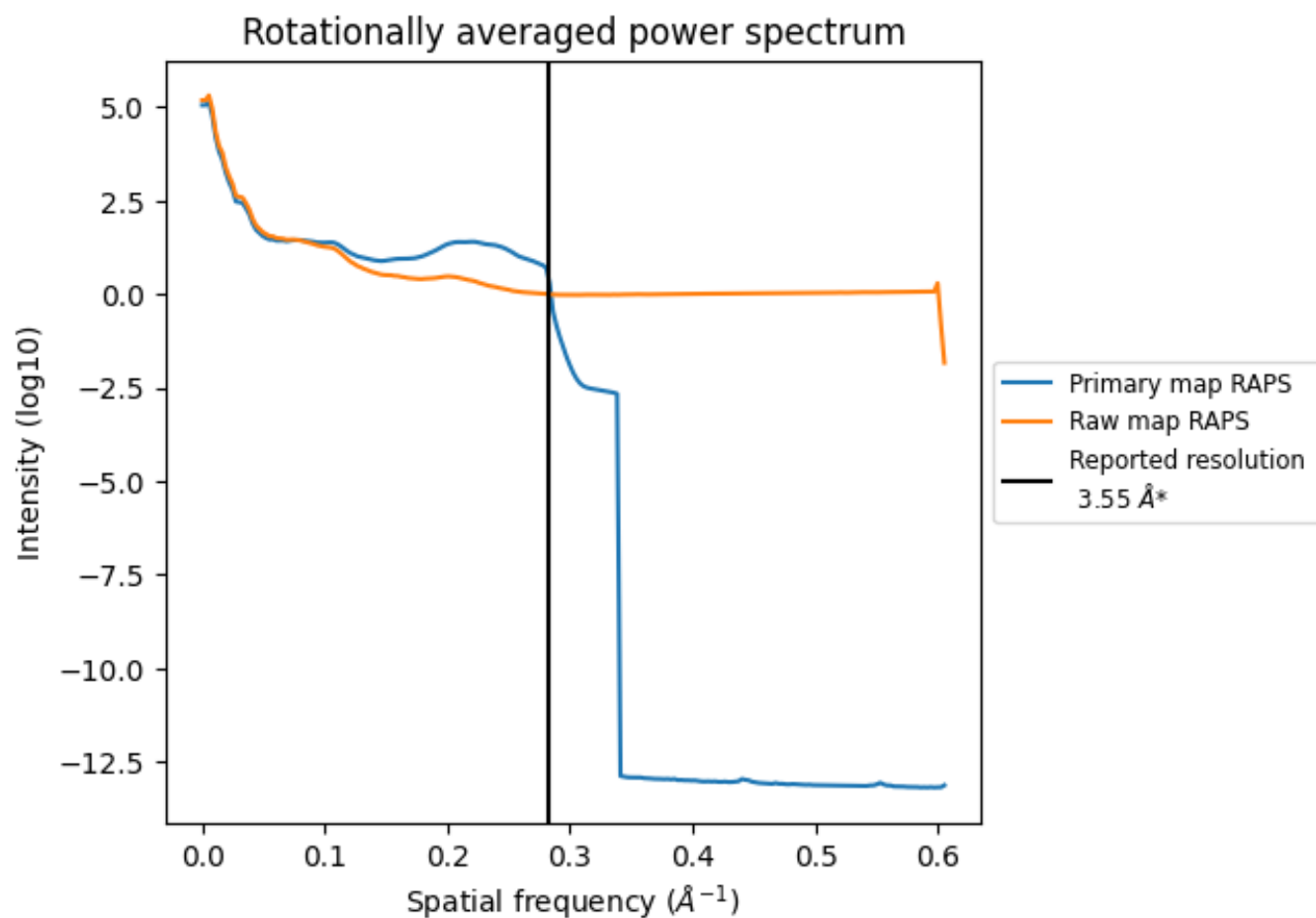
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 52 nm³; this corresponds to an approximate mass of 47 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 ⓘ

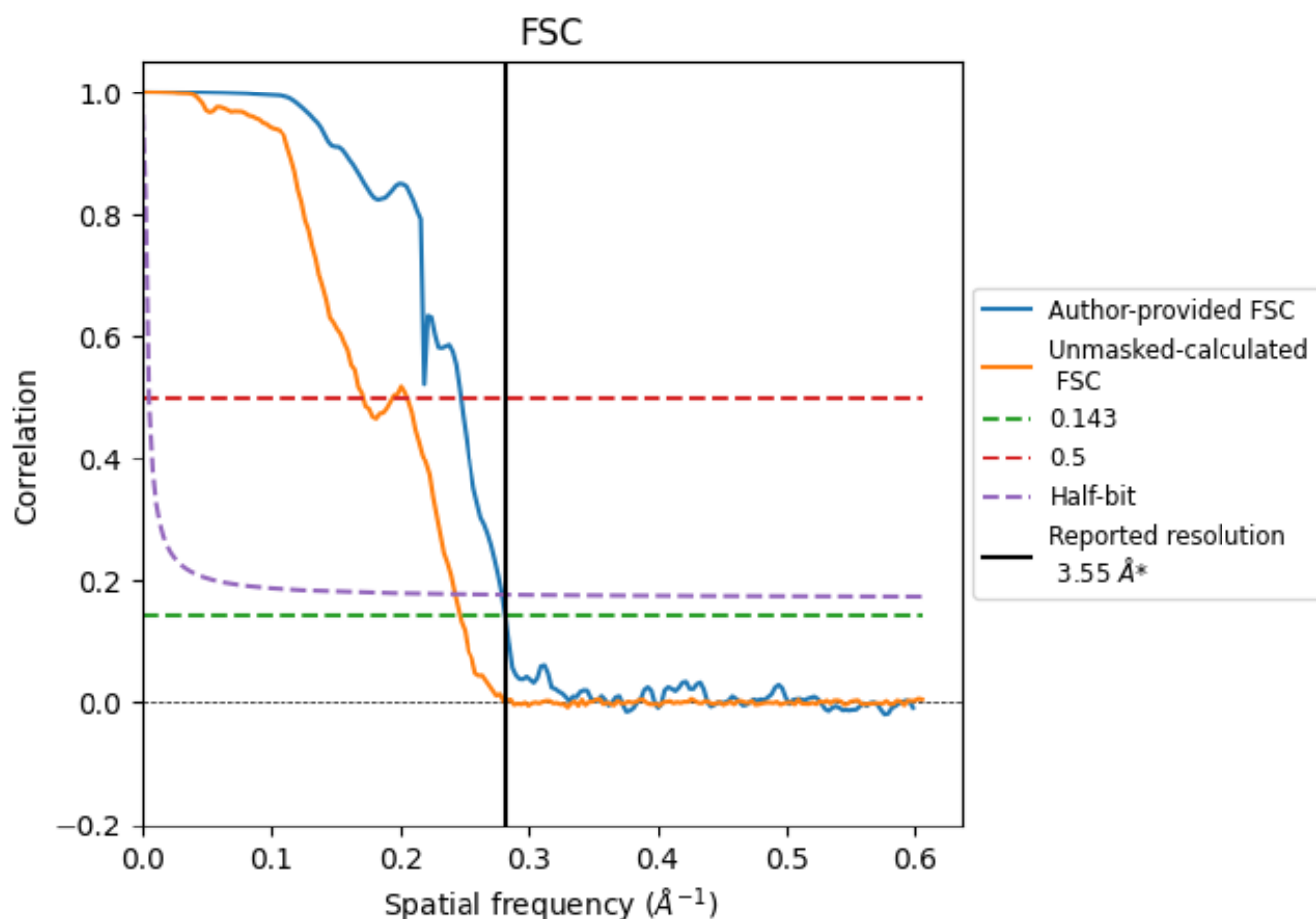


*Reported resolution corresponds to spatial frequency of 0.282 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.282 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

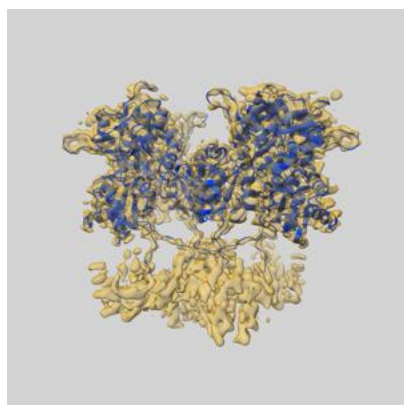
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.55	-	-
Author-provided FSC curve	3.55	4.05	3.58
Unmasked-calculated*	4.06	5.83	4.12

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.06 differs from the reported value 3.55 by more than 10 %

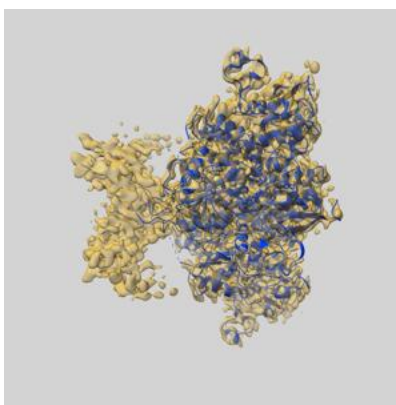
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-54556 and PDB model 9S3Z. Per-residue inclusion information can be found in section 3 on page 5.

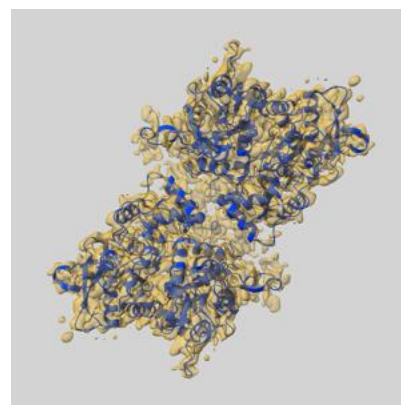
9.1 Map-model overlay [i](#)



X



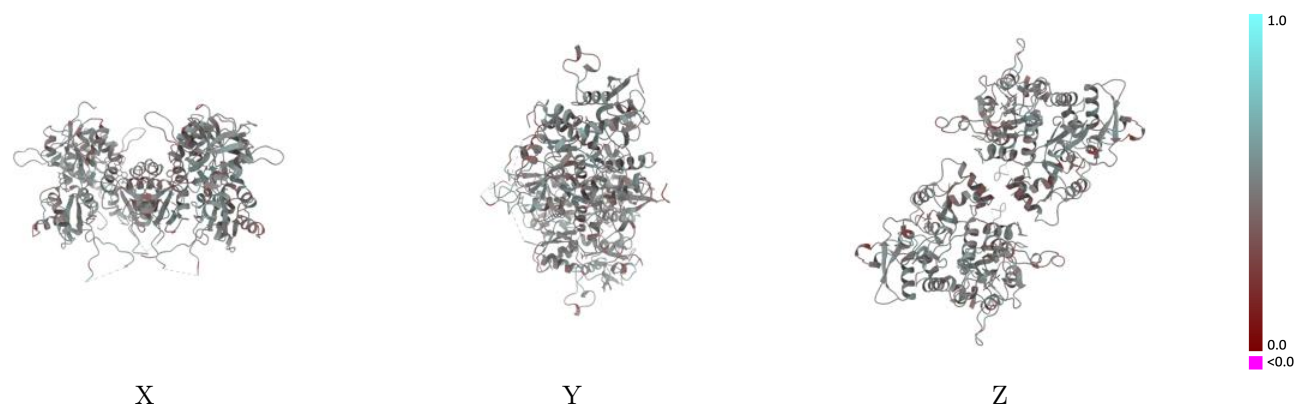
Y



Z

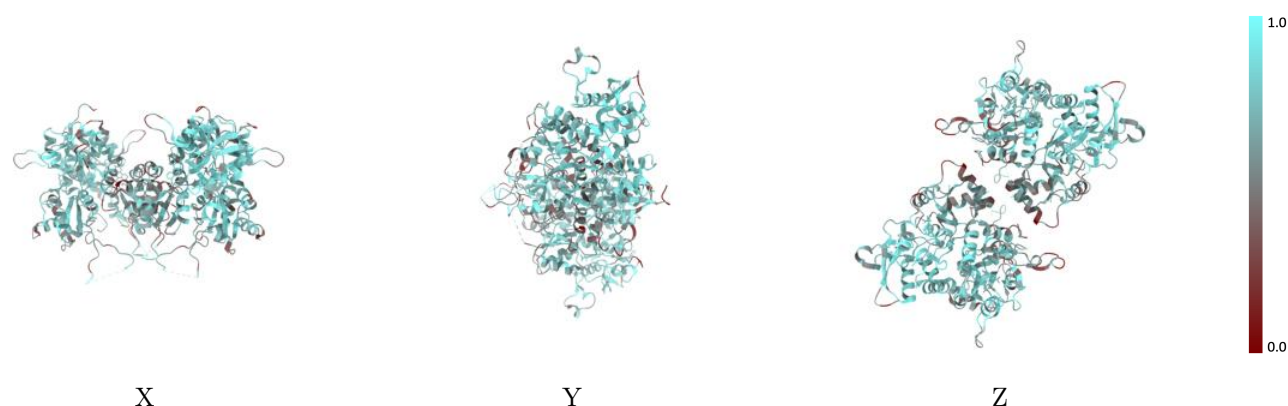
The images above show the 3D surface view of the map at the recommended contour level 0.48 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



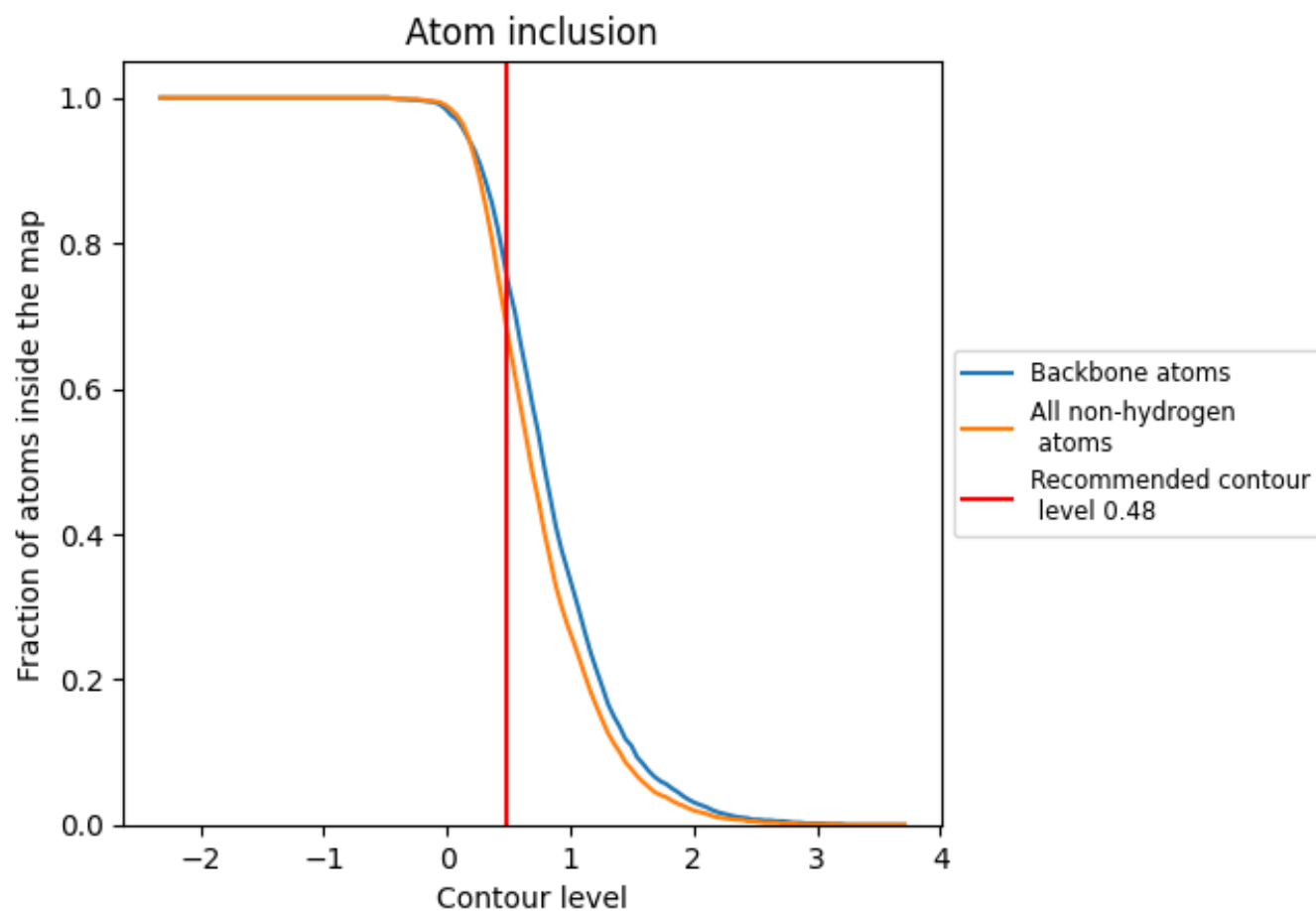
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.48).

9.4 Atom inclusion [i](#)



At the recommended contour level, 76% of all backbone atoms, 69% 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 (0.48) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6890	0.4770
A	0.6980	0.4790
B	0.6830	0.4770
C	0.6960	0.4790
D	0.6810	0.4720

