



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 08:37 AM UTC

PDB ID : 6LW1 / pdb_00006lw1
EMDB ID : EMD-30002
Title : Cryo-EM structure of TLR7/Cpd-7 (DSR-139970) complex in open form
Authors : Zhang, Z.; Ohto, U.; Shimizu, T.
Deposited on : 2020-02-06
Resolution : 2.80 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

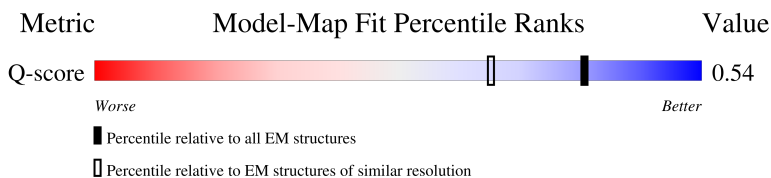
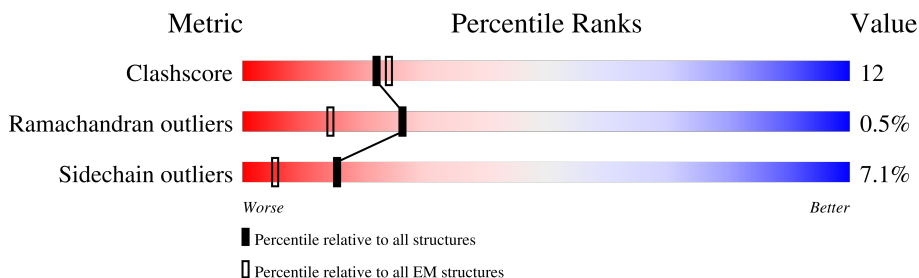
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.80 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	11806 (2.30 - 3.30)

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	823	
1	B	823	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12412 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 Toll-like receptor 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	763	Total	C	N	O	S	0	0
			6170	3951	1051	1138	30		
1	B	763	Total	C	N	O	S	0	0
			6170	3951	1051	1138	30		

There are 40 discrepancies between the modelled and reference sequences:

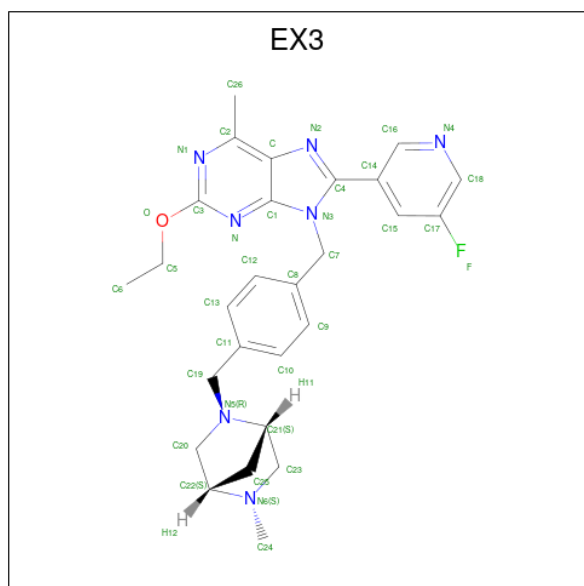
Chain	Residue	Modelled	Actual	Comment	Reference
A	23	ARG	-	expression tag	UNP B3Y653
A	24	SER	-	expression tag	UNP B3Y653
A	25	PRO	-	expression tag	UNP B3Y653
A	26	TRP	-	expression tag	UNP B3Y653
A	167	GLN	ASN	engineered mutation	UNP B3Y653
A	389	GLN	ASN	engineered mutation	UNP B3Y653
A	440	LEU	SER	engineered mutation	UNP B3Y653
A	441	VAL	GLU	engineered mutation	UNP B3Y653
A	442	PRO	VAL	engineered mutation	UNP B3Y653
A	443	ARG	GLY	engineered mutation	UNP B3Y653
A	444	GLY	PHE	engineered mutation	UNP B3Y653
A	445	SER	CYS	engineered mutation	UNP B3Y653
A	488	GLN	ASN	engineered mutation	UNP B3Y653
A	799	GLN	ASN	engineered mutation	UNP B3Y653
A	840	GLU	-	expression tag	UNP B3Y653
A	841	PHE	-	expression tag	UNP B3Y653
A	842	LEU	-	expression tag	UNP B3Y653
A	843	VAL	-	expression tag	UNP B3Y653
A	844	PRO	-	expression tag	UNP B3Y653
A	845	ARG	-	expression tag	UNP B3Y653
B	23	ARG	-	expression tag	UNP B3Y653
B	24	SER	-	expression tag	UNP B3Y653
B	25	PRO	-	expression tag	UNP B3Y653
B	26	TRP	-	expression tag	UNP B3Y653
B	167	GLN	ASN	engineered mutation	UNP B3Y653
B	389	GLN	ASN	engineered mutation	UNP B3Y653

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Chain	Residue	Modelled	Actual	Comment	Reference
B	440	LEU	SER	engineered mutation	UNP B3Y653
B	441	VAL	GLU	engineered mutation	UNP B3Y653
B	442	PRO	VAL	engineered mutation	UNP B3Y653
B	443	ARG	GLY	engineered mutation	UNP B3Y653
B	444	GLY	PHE	engineered mutation	UNP B3Y653
B	445	SER	CYS	engineered mutation	UNP B3Y653
B	488	GLN	ASN	engineered mutation	UNP B3Y653
B	799	GLN	ASN	engineered mutation	UNP B3Y653
B	840	GLU	-	expression tag	UNP B3Y653
B	841	PHE	-	expression tag	UNP B3Y653
B	842	LEU	-	expression tag	UNP B3Y653
B	843	VAL	-	expression tag	UNP B3Y653
B	844	PRO	-	expression tag	UNP B3Y653
B	845	ARG	-	expression tag	UNP B3Y653

- Molecule 2 is 2-ethoxy-8-(5-fluoranylpyridin-3-yl)-6-methyl-9-[[4-[[[(1S,4S)-5-methyl-2,5-diazabicyclo[2.2.1]heptan-2-yl]methyl]phenyl]methyl]purine (CCD ID: EX3) (formula: C₂₇H₃₀FN₇O) (labeled as "Ligand of Interest" by depositor).

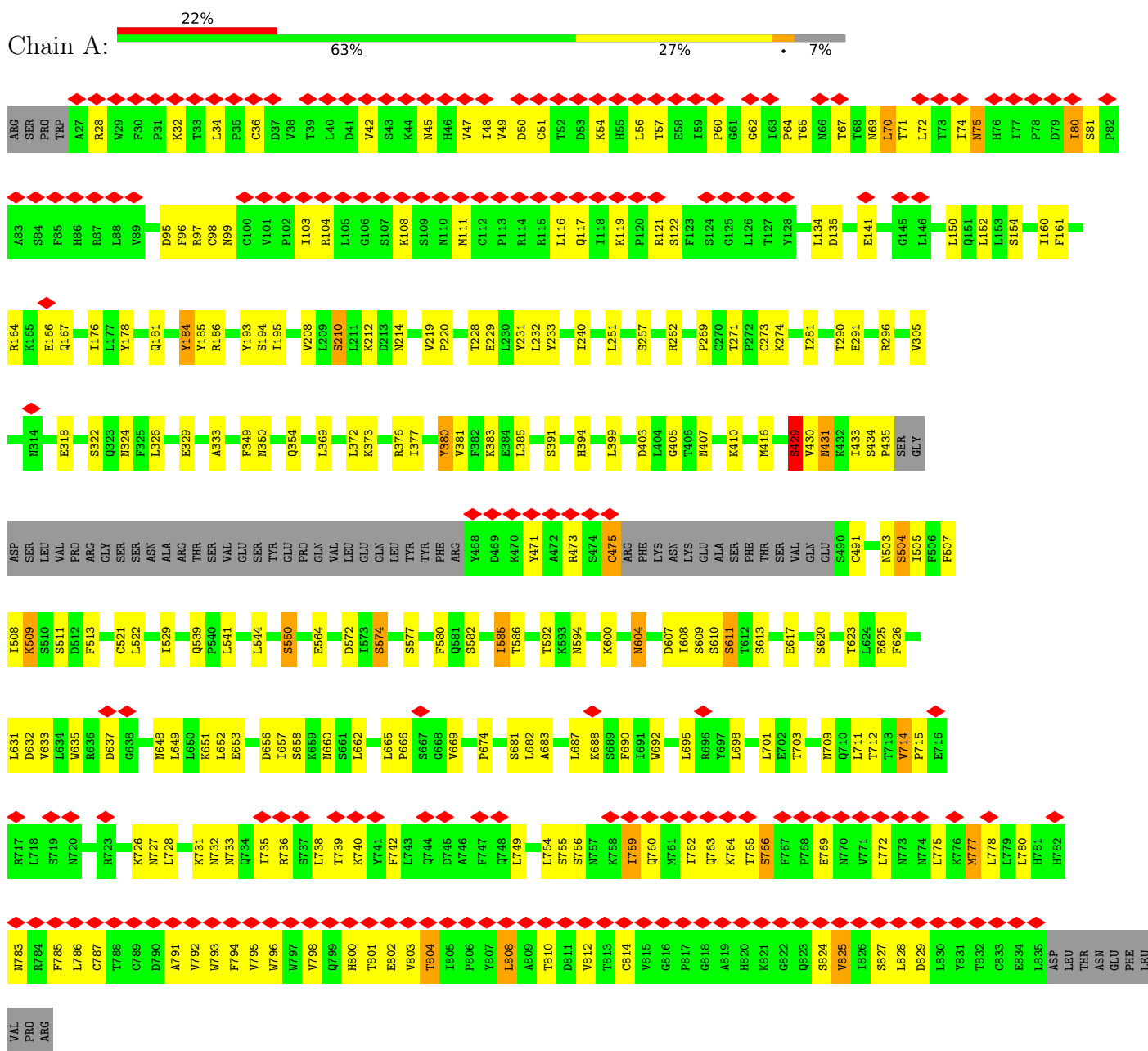


Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	F	N	O	0
			36	27	1	7	1	
2	B	1	Total	C	F	N	O	0
			36	27	1	7	1	

3 Residue-property plots

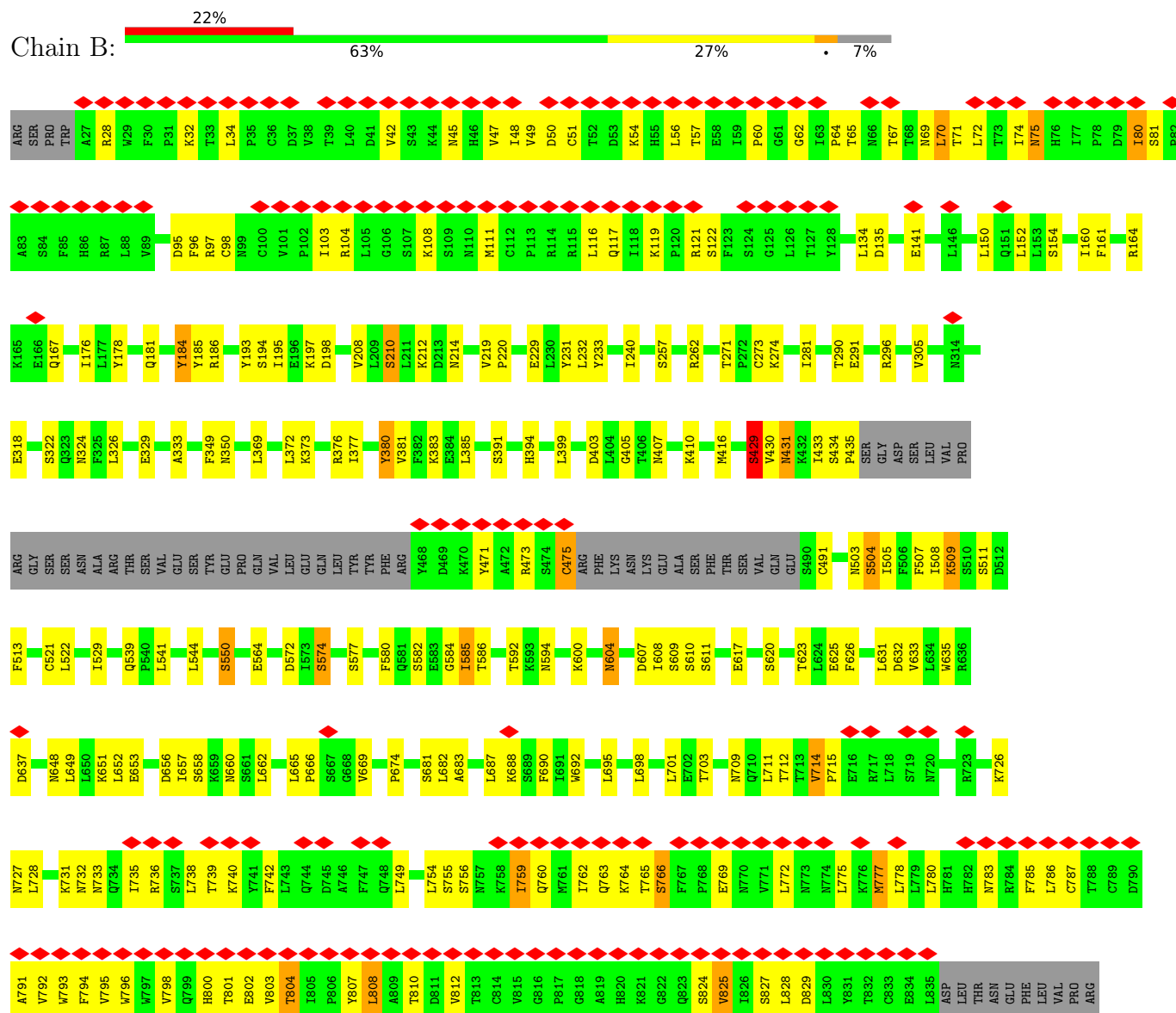
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Toll-like receptor 7



● Molecule 1: Toll-like receptor 7

Chain B:



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C2	Depositor
Number of particles used	301212	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	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	0.242	Depositor
Minimum map value	-0.172	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.028	Depositor
Map size ($\text{\AA}$)	219.12, 219.12, 219.12	wwPDB
Map dimensions	220, 220, 220	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.996, 0.996, 0.996	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EX3

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	0.36	0/6298	0.45	0/8534
1	B	0.36	0/6298	0.45	0/8534
All	All	0.36	0/12596	0.45	0/17068

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	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	380	TYR	Peptide
1	A	429	SER	Peptide
1	B	380	TYR	Peptide
1	B	429	SER	Peptide

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	6170	0	6233	145	0
1	B	6170	0	6233	145	0
2	A	36	0	0	0	0
2	B	36	0	0	0	0
All	All	12412	0	12466	286	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 12.

All (286) 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:429:SER:O	1:A:431:ASN:N	2.12	0.82
1:B:429:SER:O	1:B:431:ASN:N	2.12	0.82
1:B:796:TRP:O	1:B:800:HIS:HB2	1.80	0.82
1:A:796:TRP:O	1:A:800:HIS:HB2	1.80	0.80
1:B:742:PHE:H	1:B:766:SER:HB2	1.48	0.78
1:A:48:ILE:HG23	1:A:69:ASN:HB2	1.66	0.78
1:A:742:PHE:H	1:A:766:SER:HB2	1.48	0.78
1:A:262:ARG:NH1	1:A:324:ASN:O	2.18	0.77
1:B:262:ARG:NH1	1:B:324:ASN:O	2.18	0.77
1:B:48:ILE:HG23	1:B:69:ASN:HB2	1.66	0.75
1:B:711:LEU:HB2	1:B:733:ASN:HD21	1.51	0.75
1:A:429:SER:O	1:A:431:ASN:ND2	2.16	0.74
1:A:711:LEU:HB2	1:A:733:ASN:HD21	1.51	0.74
1:B:609:SER:HB2	1:B:633:VAL:HG21	1.69	0.74
1:B:652:LEU:HD23	1:B:674:PRO:HG3	1.71	0.73
1:A:609:SER:HB2	1:A:633:VAL:HG21	1.70	0.73
1:A:652:LEU:HD23	1:A:674:PRO:HG3	1.71	0.72
1:B:798:VAL:HG11	1:B:825:VAL:HG12	1.71	0.72
1:A:798:VAL:HG11	1:A:825:VAL:HG12	1.71	0.72
1:B:711:LEU:HB2	1:B:733:ASN:ND2	2.05	0.72
1:A:711:LEU:HB2	1:A:733:ASN:ND2	2.05	0.71
1:B:433:ILE:HB	1:B:503:ASN:HD22	1.56	0.69
1:B:429:SER:O	1:B:431:ASN:ND2	2.16	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:ILE:HB	1:A:503:ASN:HD22	1.56	0.69
1:B:74:ILE:O	1:B:75:ASN:ND2	2.26	0.67
1:B:103:ILE:HD12	1:B:108:LYS:HE2	1.76	0.66
1:B:754:LEU:HB2	1:B:780:LEU:HD23	1.77	0.66
1:B:47:VAL:HB	1:B:67:THR:HA	1.78	0.66
1:A:103:ILE:HD12	1:A:108:LYS:HE2	1.76	0.65
1:A:47:VAL:HB	1:A:67:THR:HA	1.78	0.65
1:A:74:ILE:O	1:A:75:ASN:ND2	2.26	0.65
1:A:754:LEU:HB2	1:A:780:LEU:HD23	1.77	0.64
1:A:71:THR:HA	1:A:95:ASP:HB3	1.79	0.64
1:B:71:THR:HA	1:B:95:ASP:HB3	1.79	0.64
1:A:435:PRO:HD3	1:A:505:ILE:HA	1.80	0.63
1:B:435:PRO:HD3	1:B:505:ILE:HA	1.80	0.62
1:A:385:LEU:HG	1:A:416:MET:HE1	1.81	0.62
1:A:572:ASP:OD1	1:A:574:SER:OG	2.17	0.62
1:B:385:LEU:HG	1:B:416:MET:HE1	1.81	0.62
1:B:405:GLY:O	1:B:407:ASN:ND2	2.31	0.62
1:A:405:GLY:O	1:A:407:ASN:ND2	2.31	0.62
1:B:572:ASP:OD1	1:B:574:SER:OG	2.17	0.61
1:B:824:SER:HB2	1:B:827:SER:HB2	1.83	0.60
1:A:824:SER:HB2	1:A:827:SER:HB2	1.83	0.60
1:A:521:CYS:SG	1:A:522:LEU:N	2.75	0.59
1:B:635:TRP:HB3	1:B:666:PRO:HG3	1.84	0.59
1:A:98:CYS:N	1:A:135:ASP:O	2.34	0.59
1:A:32:LYS:NZ	1:A:62:GLY:O	2.36	0.59
1:A:28:ARG:NH1	1:A:801:THR:O	2.33	0.59
1:B:521:CYS:SG	1:B:522:LEU:N	2.75	0.59
1:A:471:TYR:OH	1:A:473:ARG:NH2	2.36	0.58
1:A:625:GLU:HA	1:A:656:ASP:HB3	1.85	0.58
1:A:329:GLU:O	1:A:333:ALA:N	2.36	0.58
1:A:635:TRP:HB3	1:A:666:PRO:HG3	1.84	0.58
1:B:625:GLU:HA	1:B:656:ASP:HB3	1.85	0.58
1:A:141:GLU:HA	1:A:160:ILE:HG23	1.86	0.57
1:A:504:SER:O	1:A:504:SER:OG	2.16	0.57
1:A:801:THR:HG23	1:A:803:VAL:H	1.69	0.57
1:B:329:GLU:O	1:B:333:ALA:N	2.36	0.57
1:B:32:LYS:NZ	1:B:62:GLY:O	2.36	0.57
1:A:769:GLU:HA	1:A:772:LEU:HB2	1.87	0.57
1:B:98:CYS:N	1:B:135:ASP:O	2.34	0.57
1:B:740:LYS:HE3	1:B:763:GLN:HE21	1.70	0.57
1:B:764:LYS:HB2	1:B:793:TRP:CD2	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:703:THR:HA	1:A:727:ASN:HB2	1.86	0.57
1:B:471:TYR:OH	1:B:473:ARG:NH2	2.36	0.57
1:B:801:THR:HG23	1:B:803:VAL:H	1.69	0.57
1:B:54:LYS:HB3	1:B:56:LEU:HD13	1.87	0.56
1:A:764:LYS:HB2	1:A:793:TRP:CD2	2.40	0.56
1:B:504:SER:O	1:B:504:SER:OG	2.16	0.56
1:B:703:THR:HA	1:B:727:ASN:HB2	1.86	0.56
1:B:656:ASP:HA	1:B:681:SER:HB3	1.88	0.56
1:A:775:LEU:HD22	1:A:778:LEU:HD12	1.88	0.56
1:A:740:LYS:HE3	1:A:763:GLN:HE21	1.70	0.56
1:B:28:ARG:HD2	1:B:802:GLU:HG2	1.87	0.56
1:B:141:GLU:HA	1:B:160:ILE:HG23	1.86	0.56
1:B:769:GLU:HA	1:B:772:LEU:HB2	1.87	0.56
1:A:28:ARG:HD2	1:A:802:GLU:HG2	1.87	0.56
1:A:541:LEU:HD13	1:A:544:LEU:HD13	1.88	0.56
1:A:164:ARG:H	1:A:167:GLN:HE21	1.54	0.56
1:B:775:LEU:HD22	1:B:778:LEU:HD12	1.88	0.55
1:B:164:ARG:H	1:B:167:GLN:HE21	1.54	0.55
1:A:54:LYS:HB3	1:A:56:LEU:HD13	1.87	0.55
1:B:682:LEU:O	1:B:709:ASN:ND2	2.40	0.55
1:B:508:ILE:HG23	1:B:513:PHE:HE2	1.72	0.55
1:B:791:ALA:HB1	1:B:794:PHE:HB3	1.89	0.55
1:A:508:ILE:HG23	1:A:513:PHE:HE2	1.72	0.54
1:A:682:LEU:O	1:A:709:ASN:ND2	2.40	0.54
1:B:541:LEU:HD13	1:B:544:LEU:HD13	1.88	0.54
1:A:656:ASP:HA	1:A:681:SER:HB3	1.88	0.54
1:A:759:ILE:HD11	1:A:783:ASN:ND2	2.22	0.54
1:B:121:ARG:HG3	1:B:121:ARG:HH11	1.73	0.54
1:B:759:ILE:HD11	1:B:783:ASN:ND2	2.22	0.54
1:A:431:ASN:H	1:A:431:ASN:HD22	1.56	0.54
1:B:431:ASN:HD22	1:B:431:ASN:H	1.56	0.53
1:A:121:ARG:HG3	1:A:121:ARG:HH11	1.73	0.53
1:A:791:ALA:HB1	1:A:794:PHE:HB3	1.89	0.53
1:B:28:ARG:NH1	1:B:801:THR:O	2.33	0.53
1:A:658:SER:HB2	1:A:683:ALA:O	2.09	0.52
1:B:662:LEU:HD22	1:B:665:LEU:HD23	1.91	0.52
1:A:764:LYS:HB2	1:A:793:TRP:CE2	2.45	0.52
1:A:808:LEU:HD23	1:A:812:VAL:HG11	1.91	0.52
1:B:582:SER:HB2	1:B:585:ILE:HD11	1.91	0.52
1:A:296:ARG:NH1	1:A:318:GLU:OE2	2.42	0.52
1:A:662:LEU:HD22	1:A:665:LEU:HD23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:ARG:NH1	1:B:318:GLU:OE2	2.42	0.52
1:B:233:TYR:HB2	1:B:257:SER:O	2.10	0.52
1:B:764:LYS:HB2	1:B:793:TRP:CE2	2.45	0.52
1:B:808:LEU:HD23	1:B:812:VAL:HG11	1.91	0.52
1:B:80:ILE:HG22	1:B:81:SER:H	1.75	0.52
1:B:161:PHE:HA	1:B:193:TYR:CE1	2.45	0.51
1:B:219:VAL:HG13	1:B:232:LEU:HD11	1.91	0.51
1:A:80:ILE:HG22	1:A:81:SER:H	1.75	0.51
1:A:582:SER:HB2	1:A:585:ILE:HD11	1.91	0.51
1:B:732:ASN:HA	1:B:756:SER:HB2	1.92	0.51
1:A:233:TYR:HB2	1:A:257:SER:O	2.10	0.51
1:B:658:SER:HB2	1:B:683:ALA:O	2.09	0.51
1:A:161:PHE:HA	1:A:193:TYR:CE1	2.45	0.51
1:A:732:ASN:HA	1:A:756:SER:HB2	1.92	0.51
1:A:96:PHE:HB3	1:A:134:LEU:HD23	1.93	0.50
1:A:632:ASP:OD2	1:A:633:VAL:HG13	2.12	0.50
1:B:96:PHE:HB3	1:B:134:LEU:HD23	1.93	0.50
1:A:434:SER:OG	1:B:410:LYS:NZ	2.24	0.50
1:A:219:VAL:HG13	1:A:232:LEU:HD11	1.91	0.50
1:A:762:ILE:HG13	1:A:785:PHE:HE1	1.76	0.50
1:B:97:ARG:HG2	1:B:135:ASP:HB2	1.93	0.49
1:B:690:PHE:HB3	1:B:711:LEU:HD21	1.94	0.49
1:A:56:LEU:HD23	1:A:60:PRO:HD3	1.94	0.49
1:A:349:PHE:C	1:A:350:ASN:HD22	2.21	0.49
1:B:762:ILE:HG13	1:B:785:PHE:HE1	1.77	0.49
1:B:632:ASP:OD2	1:B:633:VAL:HG13	2.11	0.49
1:A:690:PHE:HB3	1:A:711:LEU:HD21	1.94	0.49
1:B:305:VAL:HG23	1:B:326:LEU:HD21	1.95	0.49
1:B:349:PHE:C	1:B:350:ASN:HD22	2.21	0.49
1:A:97:ARG:HG2	1:A:135:ASP:HB2	1.93	0.49
1:A:305:VAL:HG23	1:A:326:LEU:HD21	1.95	0.49
1:B:772:LEU:HA	1:B:775:LEU:HD13	1.95	0.48
1:B:376:ARG:HA	1:B:403:ASP:HB3	1.95	0.48
1:A:376:ARG:HA	1:A:403:ASP:HB3	1.95	0.48
1:A:649:LEU:O	1:A:674:PRO:HB3	2.14	0.48
1:B:714:VAL:HG13	1:B:735:ILE:HD13	1.96	0.48
1:B:600:LYS:HG3	1:B:623:THR:HB	1.96	0.48
1:A:54:LYS:H	1:A:75:ASN:HD21	1.62	0.48
1:A:48:ILE:HG22	1:A:50:ASP:OD2	2.14	0.48
1:A:269:PRO:HG3	1:B:584:GLY:C	2.39	0.47
1:A:772:LEU:HA	1:A:775:LEU:HD13	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:714:VAL:HG13	1:A:735:ILE:HD13	1.96	0.47
1:B:56:LEU:HD23	1:B:60:PRO:HD3	1.94	0.47
1:B:574:SER:HB3	1:B:604:ASN:HD22	1.79	0.47
1:B:97:ARG:HH21	1:B:475:CYS:HA	1.79	0.47
1:B:649:LEU:O	1:B:674:PRO:HB3	2.14	0.47
1:A:240:ILE:HG13	1:A:281:ILE:HG12	1.97	0.47
1:A:97:ARG:HH21	1:A:475:CYS:HA	1.79	0.47
1:A:600:LYS:HG3	1:A:623:THR:HB	1.96	0.47
1:B:48:ILE:HG22	1:B:50:ASP:OD2	2.14	0.47
1:B:208:VAL:HG13	1:B:229:GLU:HB2	1.97	0.47
1:A:290:THR:OG1	1:A:291:GLU:OE1	2.27	0.47
1:B:383:LYS:HE3	1:B:383:LYS:HB2	1.69	0.47
1:B:769:GLU:OE1	1:B:801:THR:OG1	2.32	0.46
1:A:208:VAL:HG13	1:A:229:GLU:HB2	1.97	0.46
1:B:54:LYS:H	1:B:75:ASN:HD21	1.62	0.46
1:B:240:ILE:HG13	1:B:281:ILE:HG12	1.97	0.46
1:B:508:ILE:HD11	1:B:529:ILE:HD13	1.98	0.46
1:A:574:SER:HB3	1:A:604:ASN:HD22	1.79	0.46
1:A:769:GLU:OE1	1:A:801:THR:OG1	2.32	0.46
1:A:326:LEU:HB2	1:A:350:ASN:OD1	2.16	0.46
1:A:410:LYS:NZ	1:B:434:SER:OG	2.23	0.46
1:B:210:SER:O	1:B:210:SER:OG	2.31	0.46
1:A:154:SER:HA	1:A:178:TYR:HB2	1.98	0.45
1:A:380:TYR:O	1:A:380:TYR:CD1	2.70	0.45
1:B:380:TYR:CD1	1:B:380:TYR:O	2.70	0.45
1:A:692:TRP:CE2	1:A:715:PRO:HD3	2.51	0.45
1:B:692:TRP:CE2	1:B:715:PRO:HD3	2.51	0.45
1:A:119:LYS:O	1:A:122:SER:HB2	2.17	0.45
1:A:787:CYS:HB2	1:A:814:CYS:HB3	1.82	0.45
1:B:64:PRO:O	1:B:67:THR:HG22	2.16	0.45
1:A:508:ILE:HD11	1:A:529:ILE:HD13	1.98	0.45
1:B:690:PHE:HZ	1:B:695:LEU:HD21	1.82	0.45
1:B:726:LYS:O	1:B:727:ASN:ND2	2.50	0.45
1:A:539:GLN:HA	1:A:564:GLU:HG3	1.99	0.45
1:B:326:LEU:HB2	1:B:350:ASN:OD1	2.16	0.45
1:A:152:LEU:HD12	1:A:176:ILE:HD12	1.99	0.45
1:B:119:LYS:O	1:B:122:SER:HB2	2.17	0.45
1:B:290:THR:OG1	1:B:291:GLU:OE1	2.27	0.45
1:A:690:PHE:HZ	1:A:695:LEU:HD21	1.82	0.45
1:A:759:ILE:H	1:A:759:ILE:HG13	1.66	0.45
1:A:64:PRO:O	1:A:67:THR:HG22	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:726:LYS:O	1:A:727:ASN:ND2	2.50	0.44
1:B:210:SER:HA	1:B:231:TYR:HB2	1.99	0.44
1:B:736:ARG:C	1:B:759:ILE:HA	2.42	0.44
1:B:791:ALA:O	1:B:795:VAL:HG23	2.17	0.44
1:A:391:SER:HA	1:A:394:HIS:ND1	2.33	0.44
1:B:391:SER:HA	1:B:394:HIS:ND1	2.33	0.44
1:A:736:ARG:C	1:A:759:ILE:HA	2.42	0.44
1:B:152:LEU:HD12	1:B:176:ILE:HD12	1.99	0.44
1:B:154:SER:HA	1:B:178:TYR:HB2	1.98	0.44
1:B:577:SER:OG	1:B:607:ASP:HB2	2.18	0.44
1:A:383:LYS:HB2	1:A:383:LYS:HE3	1.69	0.44
1:B:738:LEU:HD21	1:B:754:LEU:HD13	2.00	0.44
1:A:574:SER:HB3	1:A:604:ASN:ND2	2.33	0.43
1:B:539:GLN:HA	1:B:564:GLU:HG3	1.99	0.43
1:A:507:PHE:HE1	1:A:509:LYS:HG3	1.83	0.43
1:A:577:SER:OG	1:A:607:ASP:HB2	2.18	0.43
1:B:195:ILE:HD13	1:B:220:PRO:HB3	2.00	0.43
1:B:574:SER:HB3	1:B:604:ASN:ND2	2.33	0.43
1:A:210:SER:HA	1:A:231:TYR:HB2	1.99	0.43
1:A:791:ALA:O	1:A:795:VAL:HG23	2.17	0.43
1:B:631:LEU:HD12	1:B:660:ASN:HD22	1.83	0.43
1:A:34:LEU:HD23	1:A:34:LEU:HA	1.82	0.43
1:A:193:TYR:HB2	1:A:214:ASN:O	2.19	0.43
1:A:373:LYS:O	1:A:399:LEU:HD12	2.19	0.43
1:B:507:PHE:HE1	1:B:509:LYS:HG3	1.83	0.43
1:B:808:LEU:HD23	1:B:808:LEU:HA	1.87	0.43
1:B:121:ARG:HG3	1:B:121:ARG:NH1	2.34	0.43
1:B:632:ASP:OD2	1:B:633:VAL:N	2.52	0.43
1:A:166:GLU:OE2	1:A:166:GLU:N	2.46	0.42
1:A:354:GLN:HA	1:B:507:PHE:CE1	2.54	0.42
1:A:772:LEU:HB3	1:A:803:VAL:HG21	2.01	0.42
1:B:70:LEU:HD22	1:B:71:THR:H	1.84	0.42
1:B:373:LYS:O	1:B:399:LEU:HD12	2.19	0.42
1:B:760:GLN:HB3	1:B:786:LEU:CD2	2.50	0.42
1:A:121:ARG:HG3	1:A:121:ARG:NH1	2.34	0.42
1:A:195:ILE:HD13	1:A:220:PRO:HB3	2.00	0.42
1:B:617:GLU:HG2	1:B:648:ASN:HB3	2.02	0.42
1:B:626:PHE:HB3	1:B:657:ILE:HG22	2.02	0.42
1:B:687:LEU:HD12	1:B:690:PHE:HB2	2.01	0.42
1:A:184:TYR:CZ	1:A:186:ARG:HB2	2.55	0.42
1:A:369:LEU:HB3	1:A:372:LEU:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:369:LEU:HB3	1:B:372:LEU:HB2	2.01	0.42
1:B:772:LEU:HB3	1:B:803:VAL:HG21	2.01	0.42
1:A:70:LEU:HD22	1:A:71:THR:H	1.84	0.42
1:A:738:LEU:HD21	1:A:754:LEU:HD13	2.00	0.42
1:B:651:LYS:O	1:B:653:GLU:HG3	2.20	0.42
1:A:550:SER:HA	1:A:574:SER:O	2.19	0.42
1:A:736:ARG:O	1:A:759:ILE:HA	2.20	0.42
1:B:184:TYR:CZ	1:B:186:ARG:HB2	2.55	0.42
1:B:193:TYR:HB2	1:B:214:ASN:O	2.19	0.42
1:A:631:LEU:HD12	1:A:660:ASN:HD22	1.83	0.42
1:A:687:LEU:HD12	1:A:690:PHE:HB2	2.01	0.42
1:A:760:GLN:HB3	1:A:786:LEU:CD2	2.50	0.42
1:A:828:LEU:HD23	1:A:828:LEU:HA	1.86	0.42
1:B:181:GLN:HG3	1:B:212:LYS:HE3	2.01	0.42
1:B:429:SER:C	1:B:431:ASN:H	2.15	0.42
1:A:181:GLN:HG3	1:A:212:LYS:HE3	2.01	0.42
1:A:651:LYS:O	1:A:653:GLU:HG3	2.20	0.42
1:A:688:LYS:HD3	1:A:688:LYS:HA	1.85	0.42
1:B:550:SER:HA	1:B:574:SER:O	2.19	0.42
1:B:70:LEU:HD22	1:B:71:THR:N	2.35	0.41
1:A:632:ASP:OD2	1:A:633:VAL:N	2.52	0.41
1:A:777:MET:HA	1:A:804:THR:HG22	2.02	0.41
1:B:736:ARG:O	1:B:759:ILE:HA	2.20	0.41
1:A:273:CYS:SG	1:A:274:LYS:N	2.93	0.41
1:A:51:CYS:HB2	1:A:72:LEU:HD23	2.03	0.41
1:A:611:SER:OG	1:A:613:SER:O	2.35	0.41
1:A:617:GLU:HG2	1:A:648:ASN:HB3	2.02	0.41
1:A:626:PHE:HB3	1:A:657:ILE:HG22	2.02	0.41
1:B:763:GLN:H	1:B:763:GLN:CD	2.25	0.41
1:A:580:PHE:CD1	1:A:608:ILE:HG12	2.56	0.41
1:A:70:LEU:HD22	1:A:71:THR:N	2.35	0.41
1:A:431:ASN:ND2	1:A:431:ASN:H	2.19	0.41
1:B:580:PHE:CD1	1:B:608:ILE:HG12	2.56	0.41
1:B:807:TYR:O	1:B:812:VAL:HG12	2.21	0.41
1:A:99:ASN:HD22	1:A:99:ASN:HA	1.62	0.41
1:B:273:CYS:SG	1:B:274:LYS:N	2.93	0.41
1:B:698:LEU:HD13	1:B:701:LEU:HD22	2.03	0.41
1:B:787:CYS:HB3	1:B:828:LEU:HD11	2.02	0.41
1:A:104:ARG:HE	1:A:186:ARG:HA	1.86	0.41
1:A:698:LEU:HD13	1:A:701:LEU:HD22	2.03	0.41
1:B:104:ARG:HE	1:B:186:ARG:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:TYR:CE1	1:A:186:ARG:HG2	2.56	0.40
1:B:380:TYR:O	1:B:380:TYR:CG	2.74	0.40
1:A:635:TRP:CH2	1:A:669:VAL:HG21	2.56	0.40
1:B:51:CYS:HB2	1:B:72:LEU:HD23	2.02	0.40
1:B:150:LEU:HD12	1:B:150:LEU:HA	1.93	0.40
1:A:36:CYS:HB3	1:A:50:ASP:O	2.22	0.40
1:A:150:LEU:HD12	1:A:150:LEU:HA	1.93	0.40
1:A:228:THR:HA	1:A:251:LEU:HA	2.02	0.40
1:A:759:ILE:HD11	1:A:783:ASN:CG	2.47	0.40
1:B:185:TYR:CE1	1:B:186:ARG:HG2	2.56	0.40
1:B:50:ASP:HA	1:B:71:THR:HB	2.04	0.40
1:B:635:TRP:CH2	1:B:669:VAL:HG21	2.56	0.40
1:B:34:LEU:HD23	1:B:34:LEU:HA	1.82	0.40
1:B:197:LYS:HG3	1:B:198:ASP:OD1	2.22	0.40
1:B:688:LYS:HA	1:B:688:LYS:HD3	1.85	0.40
1:B:777:MET:HA	1:B:804:THR:HG22	2.02	0.40
1:B:787:CYS:SG	1:B:828:LEU:HD11	2.61	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	757/823 (92%)	675 (89%)	78 (10%)	4 (0%)	24	55
1	B	757/823 (92%)	675 (89%)	78 (10%)	4 (0%)	24	55
All	All	1514/1646 (92%)	1350 (89%)	156 (10%)	8 (0%)	26	55

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	430	VAL

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Mol	Chain	Res	Type
1	B	430	VAL
1	A	429	SER
1	B	429	SER
1	A	431	ASN
1	B	431	ASN
1	A	381	VAL
1	B	381	VAL

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	718/774 (93%)	667 (93%)	51 (7%)	13	39
1	B	718/774 (93%)	667 (93%)	51 (7%)	13	39
All	All	1436/1548 (93%)	1334 (93%)	102 (7%)	15	39

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	VAL
1	A	45	ASN
1	A	49	VAL
1	A	57	THR
1	A	65	THR
1	A	70	LEU
1	A	75	ASN
1	A	80	ILE
1	A	111	MET
1	A	116	LEU
1	A	117	GLN
1	A	184	TYR
1	A	194	SER
1	A	210	SER
1	A	271	THR
1	A	322	SER

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Mol	Chain	Res	Type
1	A	377	ILE
1	A	429	SER
1	A	475	CYS
1	A	491	CYS
1	A	504	SER
1	A	509	LYS
1	A	511	SER
1	A	550	SER
1	A	574	SER
1	A	585	ILE
1	A	586	THR
1	A	592	THR
1	A	594	ASN
1	A	604	ASN
1	A	610	SER
1	A	611	SER
1	A	620	SER
1	A	637	ASP
1	A	712	THR
1	A	714	VAL
1	A	728	LEU
1	A	731	LYS
1	A	739	THR
1	A	749	LEU
1	A	755	SER
1	A	759	ILE
1	A	765	THR
1	A	766	SER
1	A	777	MET
1	A	792	VAL
1	A	804	THR
1	A	808	LEU
1	A	810	THR
1	A	825	VAL
1	A	829	ASP
1	B	42	VAL
1	B	45	ASN
1	B	49	VAL
1	B	57	THR
1	B	65	THR
1	B	70	LEU
1	B	75	ASN

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Mol	Chain	Res	Type
1	B	80	ILE
1	B	111	MET
1	B	116	LEU
1	B	117	GLN
1	B	184	TYR
1	B	194	SER
1	B	210	SER
1	B	271	THR
1	B	322	SER
1	B	377	ILE
1	B	429	SER
1	B	475	CYS
1	B	491	CYS
1	B	504	SER
1	B	509	LYS
1	B	511	SER
1	B	550	SER
1	B	574	SER
1	B	585	ILE
1	B	586	THR
1	B	592	THR
1	B	594	ASN
1	B	604	ASN
1	B	610	SER
1	B	611	SER
1	B	620	SER
1	B	637	ASP
1	B	712	THR
1	B	714	VAL
1	B	728	LEU
1	B	731	LYS
1	B	739	THR
1	B	749	LEU
1	B	755	SER
1	B	759	ILE
1	B	765	THR
1	B	766	SER
1	B	777	MET
1	B	792	VAL
1	B	804	THR
1	B	808	LEU
1	B	810	THR

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Mol	Chain	Res	Type
1	B	825	VAL
1	B	829	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (58) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	ASN
1	A	75	ASN
1	A	76	HIS
1	A	110	ASN
1	A	159	ASN
1	A	173	ASN
1	A	181	GLN
1	A	241	GLN
1	A	252	GLN
1	A	276	ASN
1	A	280	GLN
1	A	300	ASN
1	A	314	ASN
1	A	350	ASN
1	A	364	GLN
1	A	389	GLN
1	A	398	ASN
1	A	407	ASN
1	A	503	ASN
1	A	594	ASN
1	A	604	ASN
1	A	708	HIS
1	A	720	ASN
1	A	727	ASN
1	A	732	ASN
1	A	733	ASN
1	A	744	GLN
1	A	760	GLN
1	A	773	ASN
1	A	781	HIS
1	B	66	ASN
1	B	75	ASN
1	B	76	HIS
1	B	99	ASN
1	B	110	ASN
1	B	159	ASN

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Mol	Chain	Res	Type
1	B	181	GLN
1	B	241	GLN
1	B	252	GLN
1	B	276	ASN
1	B	280	GLN
1	B	300	ASN
1	B	314	ASN
1	B	350	ASN
1	B	364	GLN
1	B	389	GLN
1	B	398	ASN
1	B	503	ASN
1	B	515	HIS
1	B	708	HIS
1	B	720	ASN
1	B	727	ASN
1	B	732	ASN
1	B	733	ASN
1	B	744	GLN
1	B	760	GLN
1	B	773	ASN
1	B	781	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EX3	B	1001	-	41,41,41	1.91	8 (19%)	47,60,60	2.53	18 (38%)
2	EX3	A	1001	-	41,41,41	1.91	8 (19%)	47,60,60	2.53	18 (38%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EX3	B	1001	-	-	0/15/35/35	0/7/6/6
2	EX3	A	1001	-	-	0/15/35/35	0/7/6/6

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	EX3	C22-N6	-5.84	1.40	1.48
2	A	1001	EX3	C22-N6	-5.82	1.40	1.48
2	A	1001	EX3	C25-C22	-5.64	1.45	1.53
2	B	1001	EX3	C25-C22	-5.64	1.45	1.53
2	A	1001	EX3	C4-N3	-3.98	1.34	1.38
2	B	1001	EX3	C4-N3	-3.98	1.34	1.38
2	A	1001	EX3	C-N2	-3.68	1.32	1.39
2	B	1001	EX3	C-N2	-3.66	1.32	1.39
2	B	1001	EX3	C-C1	3.66	1.45	1.39
2	A	1001	EX3	C-C1	3.63	1.45	1.39
2	A	1001	EX3	C1-N3	-3.32	1.31	1.39
2	B	1001	EX3	C1-N3	-3.28	1.32	1.39
2	A	1001	EX3	C20-C22	-2.99	1.46	1.52
2	B	1001	EX3	C20-C22	-2.96	1.46	1.52
2	A	1001	EX3	C-C2	2.71	1.46	1.39
2	B	1001	EX3	C-C2	2.70	1.46	1.39

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	EX3	C-C1-N	-7.74	119.02	127.18

Continued on next page...

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	EX3	C-C1-N	-7.72	119.05	127.18
2	A	1001	EX3	C25-C22-C20	6.34	111.58	101.91
2	B	1001	EX3	C25-C22-C20	6.32	111.56	101.91
2	A	1001	EX3	C25-C22-N6	5.71	115.23	104.00
2	B	1001	EX3	C25-C22-N6	5.70	115.22	104.00
2	B	1001	EX3	N-C1-N3	5.34	134.63	125.67
2	A	1001	EX3	N-C1-N3	5.33	134.62	125.67
2	B	1001	EX3	C1-C-N2	-4.15	106.67	110.91
2	A	1001	EX3	C1-C-N2	-4.12	106.70	110.91
2	A	1001	EX3	C22-C25-C21	-3.66	75.22	95.10
2	B	1001	EX3	C22-C25-C21	-3.66	75.24	95.10
2	B	1001	EX3	C-N2-C4	3.56	107.76	104.29
2	A	1001	EX3	C-N2-C4	3.55	107.74	104.29
2	B	1001	EX3	C8-C7-N3	-3.34	108.03	113.46
2	A	1001	EX3	C8-C7-N3	-3.30	108.10	113.46
2	A	1001	EX3	C26-C2-N1	3.24	121.56	116.53
2	B	1001	EX3	C26-C2-N1	3.22	121.52	116.53
2	B	1001	EX3	C24-N6-C23	-2.74	108.81	112.48
2	B	1001	EX3	C18-N4-C16	2.73	121.22	117.51
2	A	1001	EX3	C24-N6-C23	-2.73	108.83	112.48
2	A	1001	EX3	C18-N4-C16	2.71	121.19	117.51
2	A	1001	EX3	C16-C14-C4	2.61	126.65	120.18
2	B	1001	EX3	C16-C14-C4	2.59	126.61	120.18
2	A	1001	EX3	C15-C14-C4	-2.58	114.00	120.25
2	B	1001	EX3	C15-C14-C4	-2.57	114.02	120.25
2	A	1001	EX3	N1-C3-N	-2.35	123.37	127.65
2	B	1001	EX3	N1-C3-N	-2.35	123.39	127.65
2	A	1001	EX3	N3-C4-N2	-2.30	110.66	112.96
2	B	1001	EX3	N3-C4-N2	-2.28	110.68	112.96
2	A	1001	EX3	C26-C2-C	-2.22	120.83	123.12
2	B	1001	EX3	C26-C2-C	-2.21	120.85	123.12
2	A	1001	EX3	C14-C4-N3	2.21	127.25	124.24
2	B	1001	EX3	C14-C4-N3	2.21	127.25	124.24
2	B	1001	EX3	C14-C16-N4	-2.19	120.25	123.50
2	A	1001	EX3	C14-C16-N4	-2.16	120.29	123.50

There are no chirality outliers.

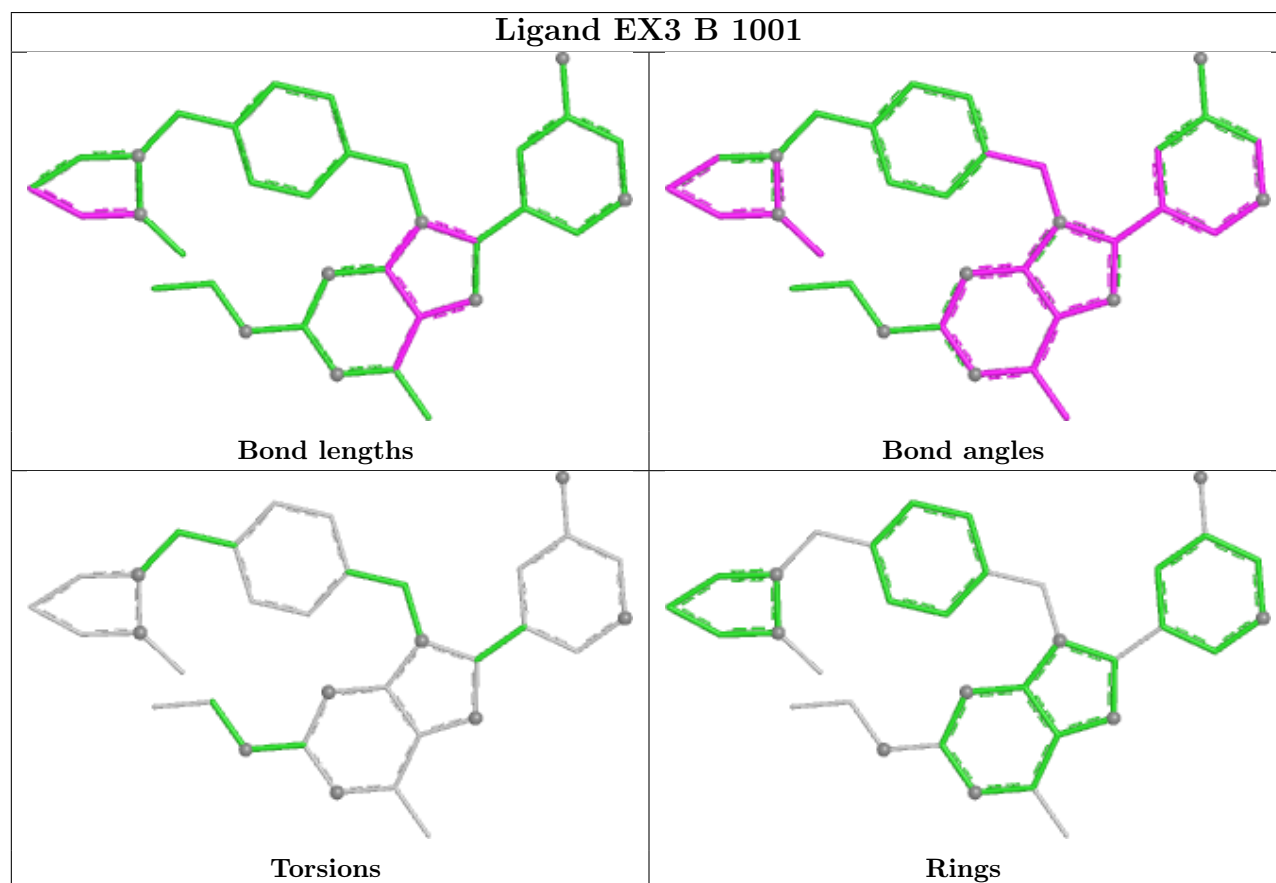
There are no torsion outliers.

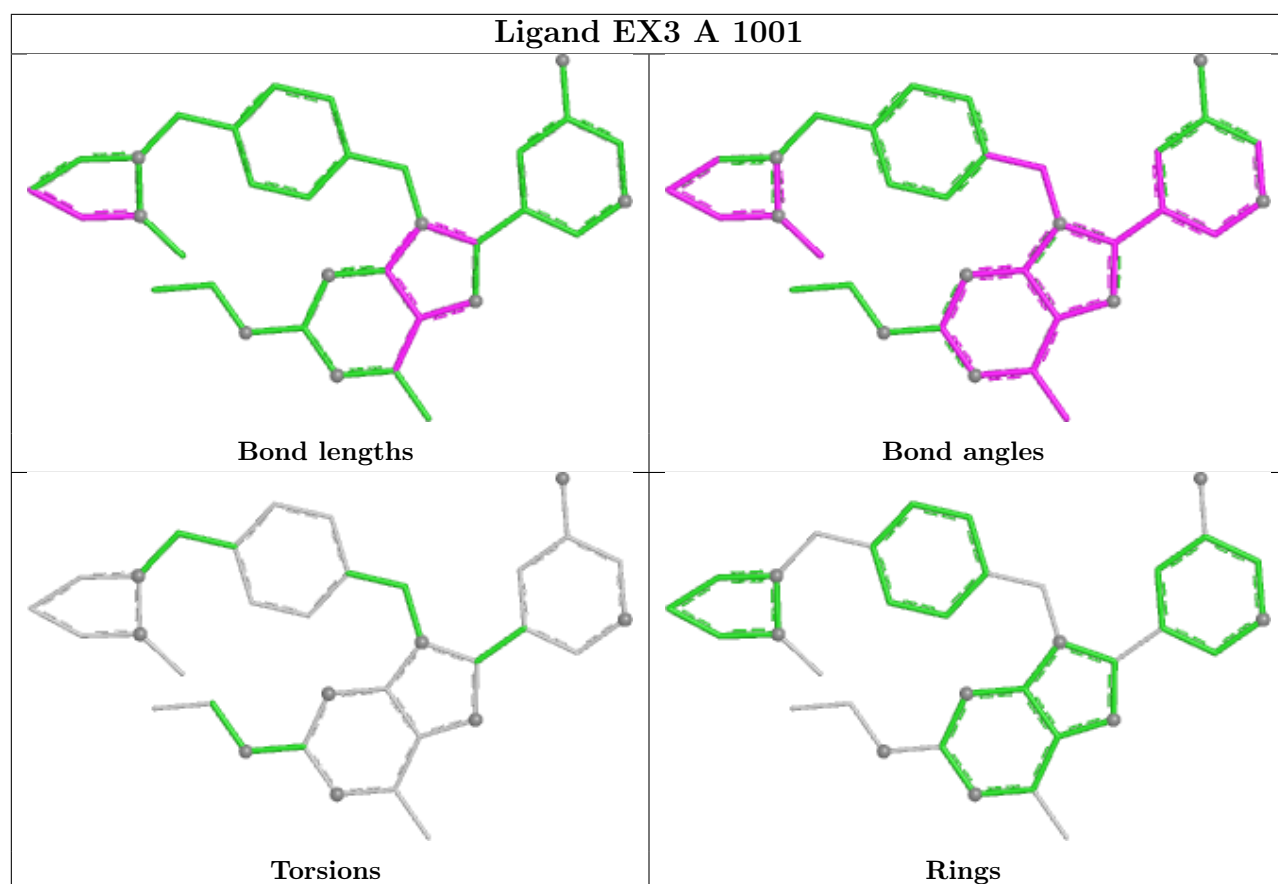
There are no ring outliers.

No monomer is involved in short contacts.

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.

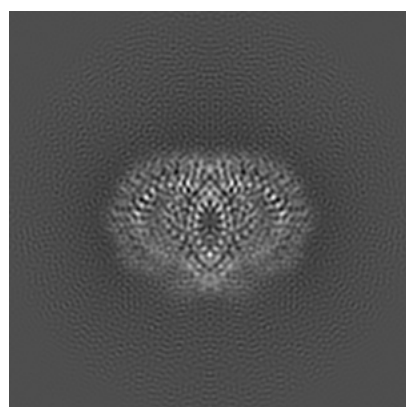
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-30002. These allow visual inspection of the internal detail of the map and identification of artifacts.

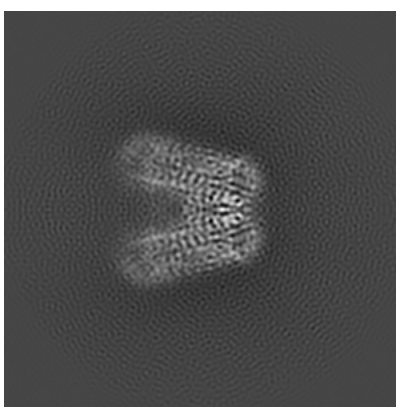
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

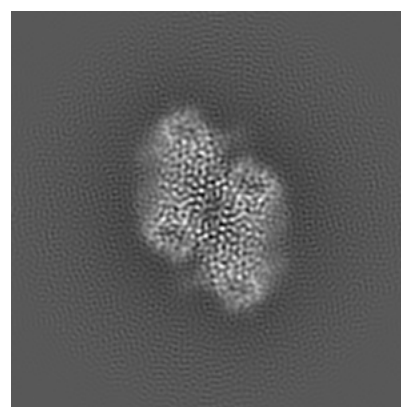
6.1.1 Primary map



X



Y

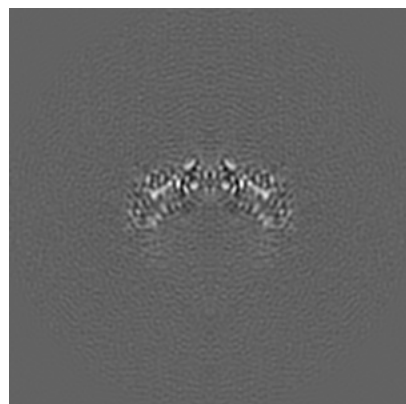


Z

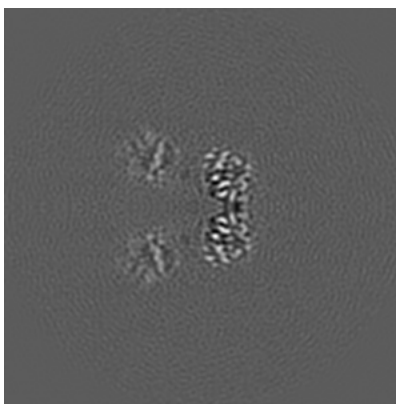
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

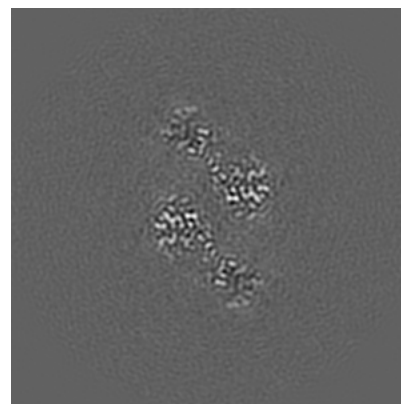
6.2.1 Primary map



X Index: 110



Y Index: 110

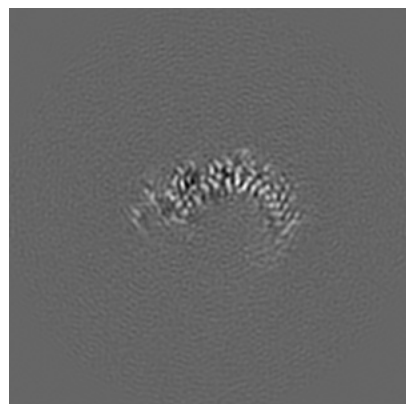


Z Index: 110

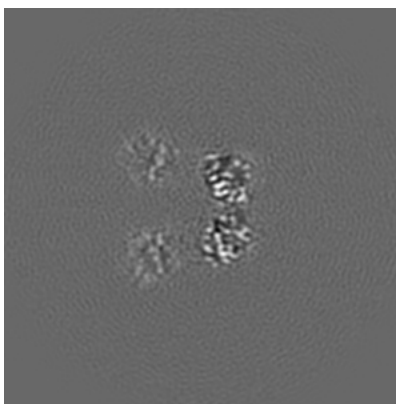
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

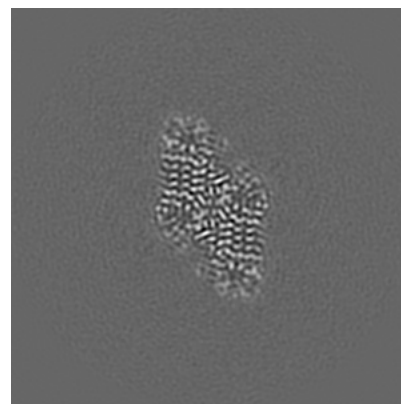
6.3.1 Primary map



X Index: 105



Y Index: 109

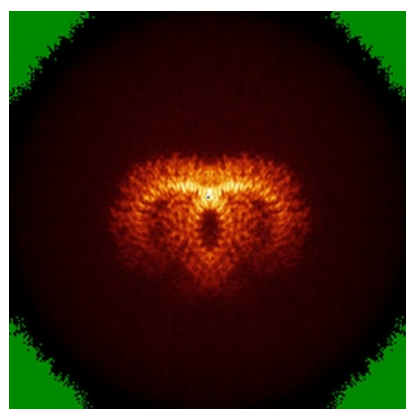


Z Index: 122

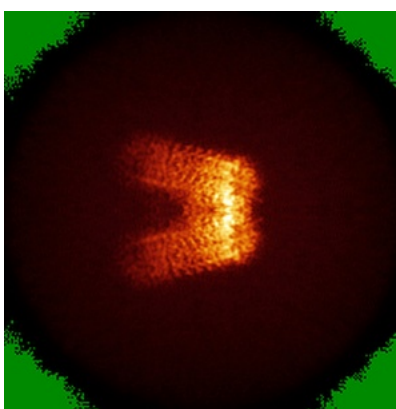
The images above show the largest variance slices of the map in three orthogonal directions.

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

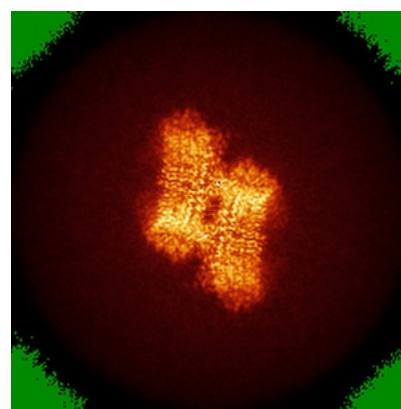
6.4.1 Primary map



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.

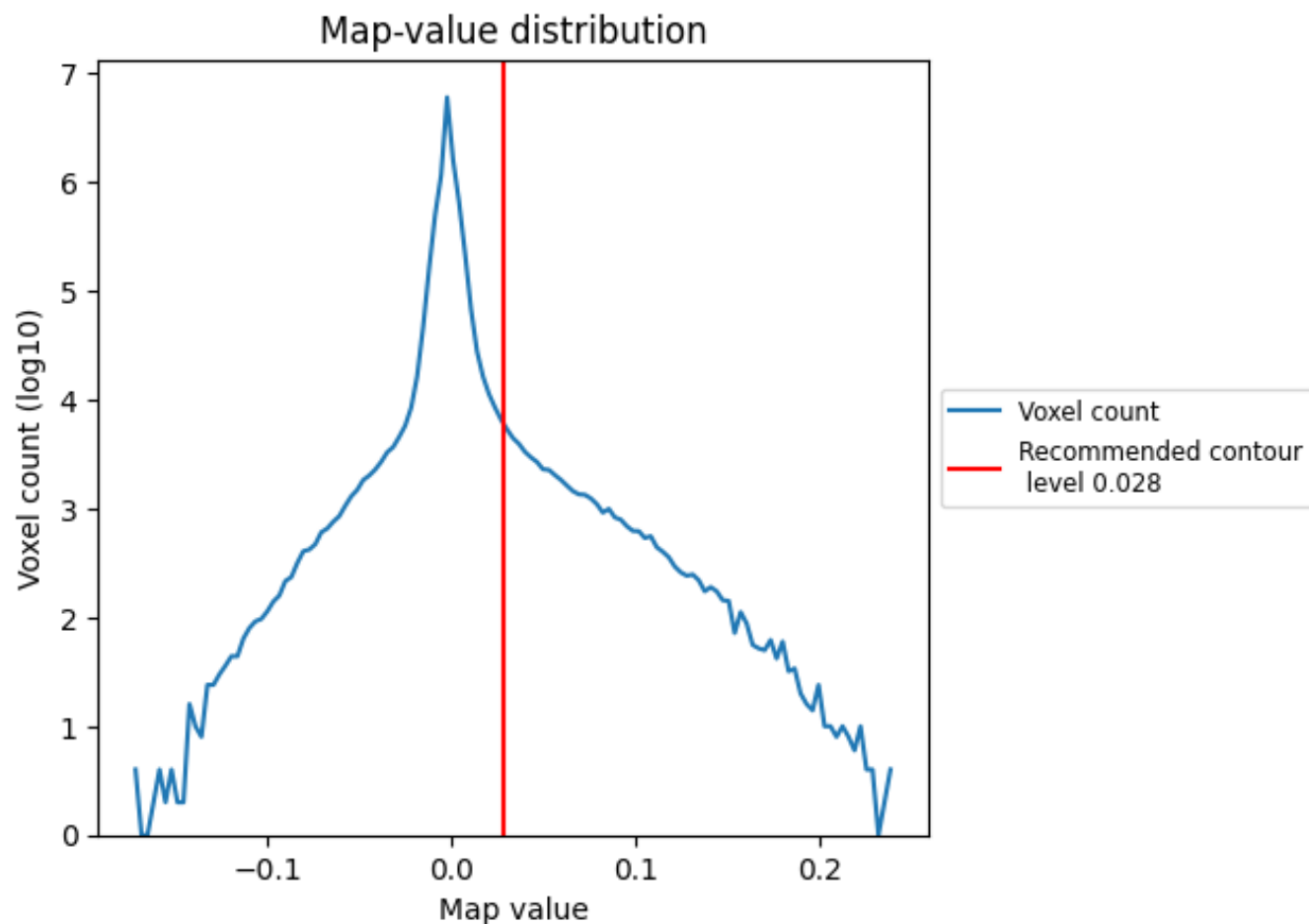
6.6 Mask visualisation

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

7 Map analysis [i](#)

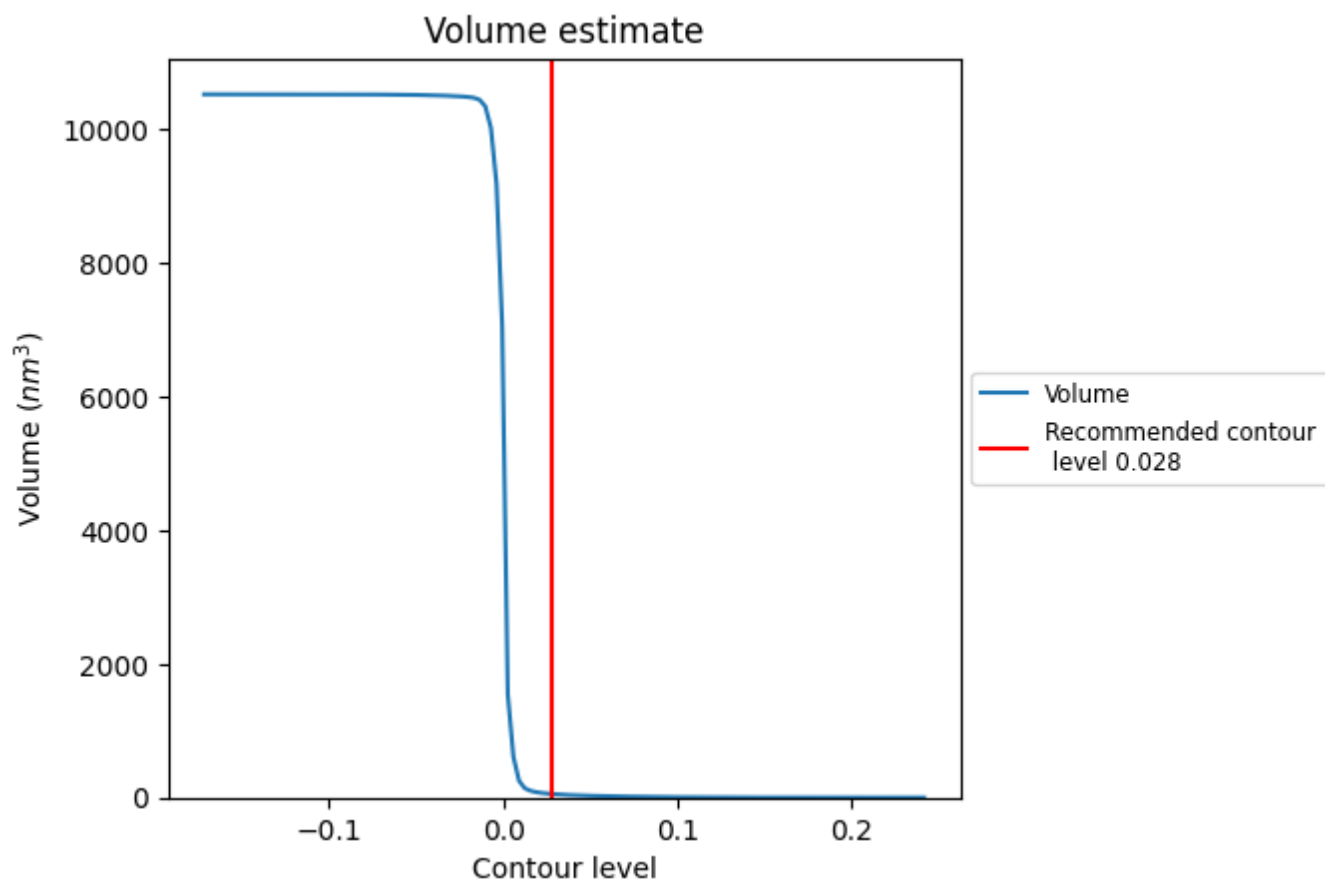
This section contains the results of statistical analysis of the map.

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.

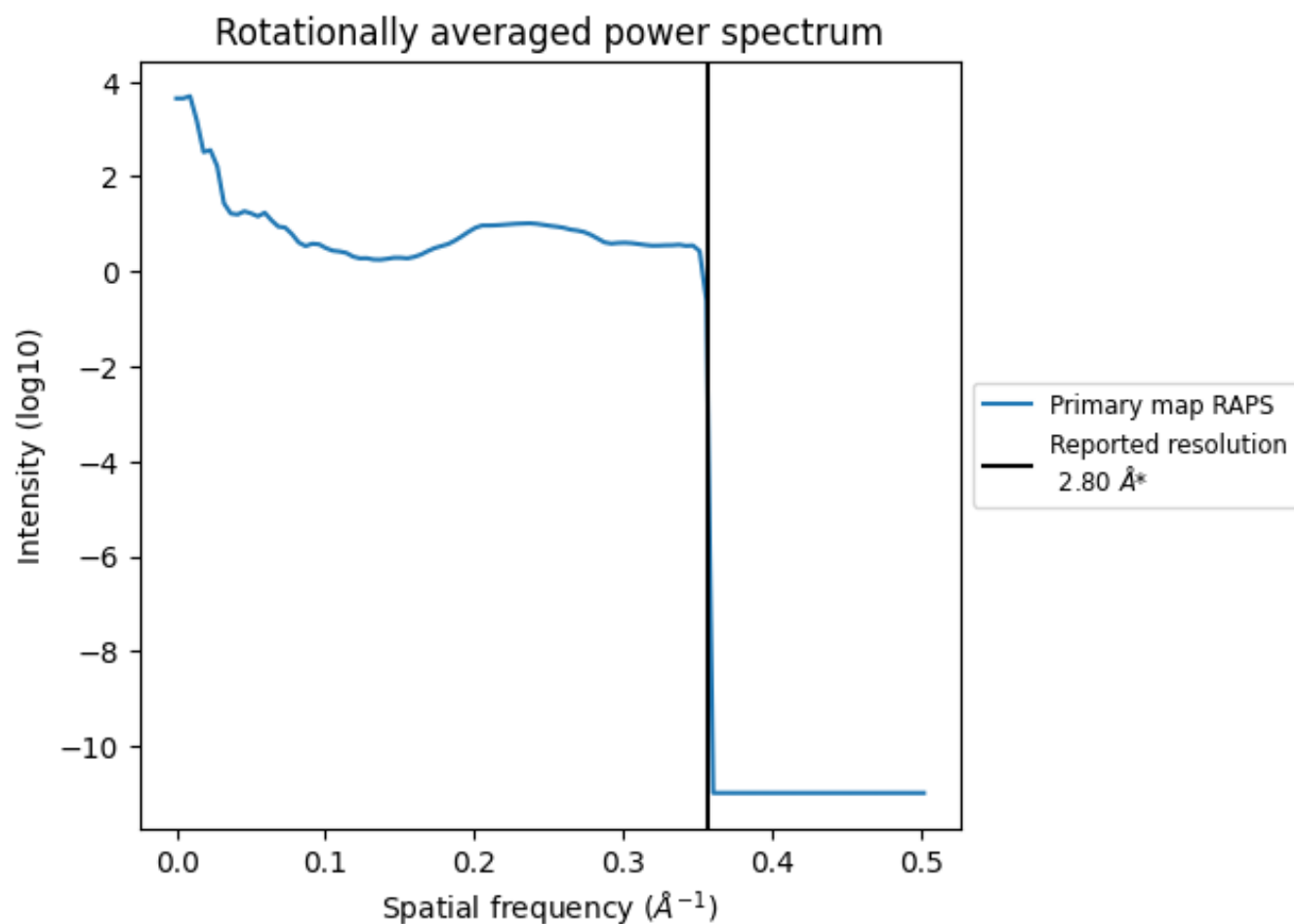
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 54 nm³; 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 0.357 $\text{\AA}^{-1}$

8 Fourier-Shell correlation ⓘ

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

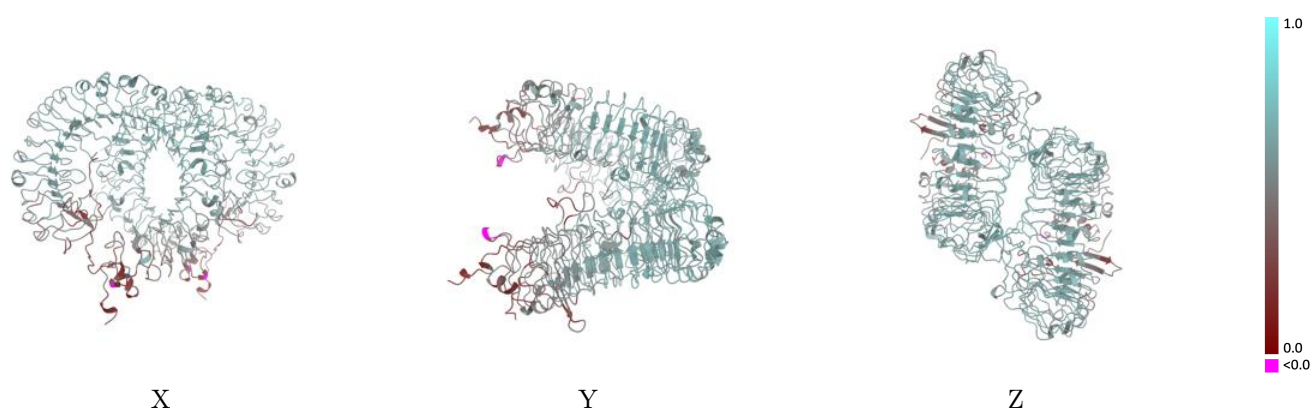
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-30002 and PDB model 6LW1. Per-residue inclusion information can be found in section 3 on page 5.

9.1 Map-model overlay [i](#)

This section was not generated.

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

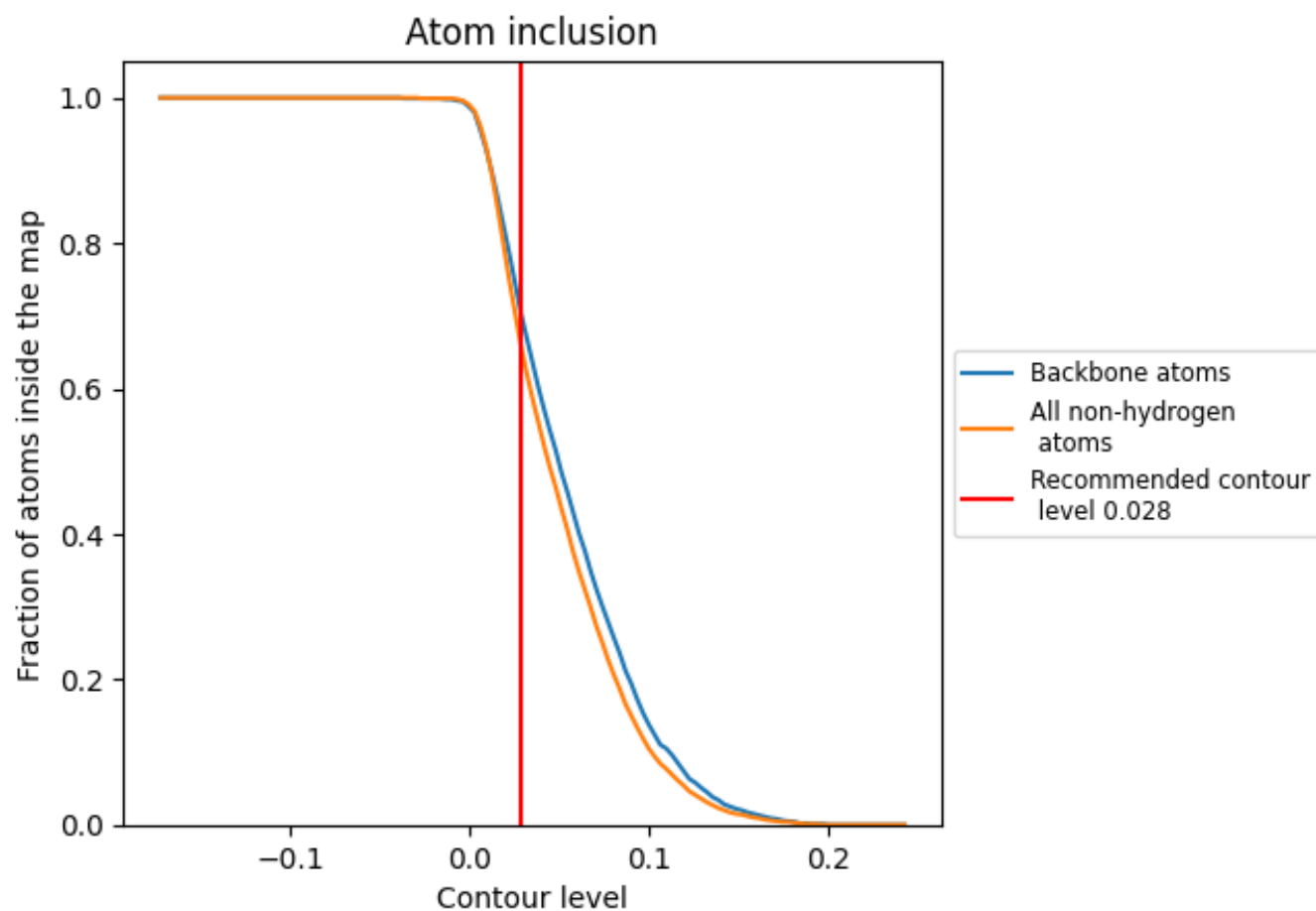


The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 67% 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 (0.028) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6670	0.5400
A	0.6680	0.5410
B	0.6660	0.5400

