



Full wwPDB EM Validation Report ⓘ

Nov 19, 2022 – 08:54 am GMT

PDB ID : 5LY6
EMDB ID : EMD-4118
Title : CryoEM structure of the membrane pore complex of Pneumolysin at 4.5Å
Authors : van Pee, K.; Neuhaus, A.; D'Imprima, E.; Mills, D.J.; Kuehlbrandt, W.; Yildiz, O.
Deposited on : 2016-09-24
Resolution : 4.50 Å(reported)
Based on initial model : 5AOD

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.dev43
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

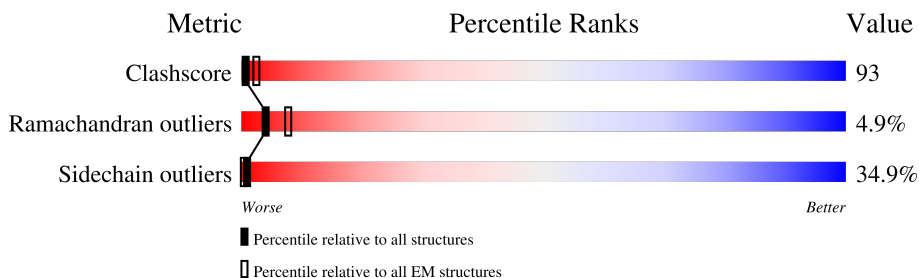
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.50 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	B	471	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3725 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 Pneumolysin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	471	3725	2350	633	735	7	0	0

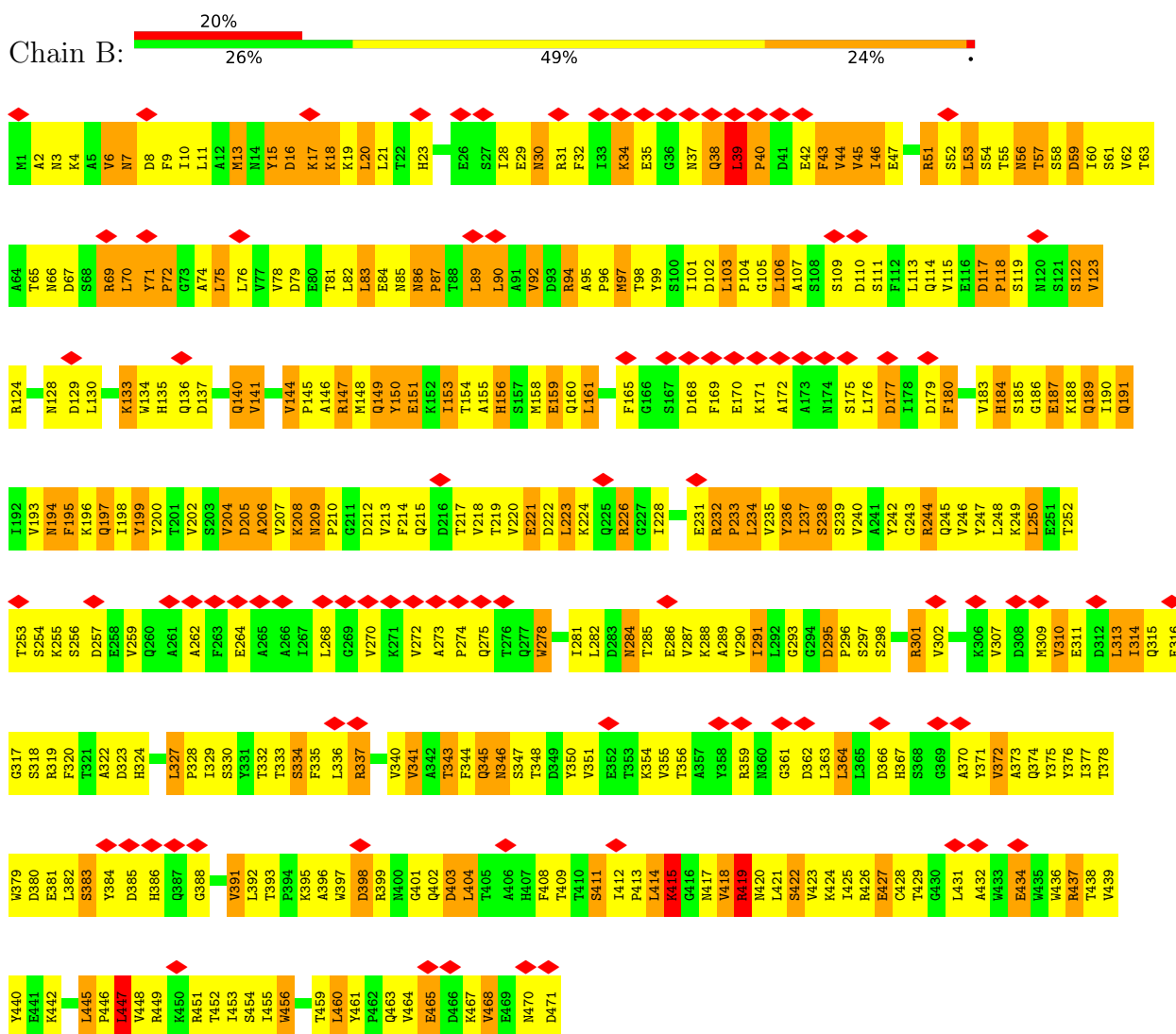
There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	172	ALA	THR	conflict	UNP Q04IN8
B	173	ALA	GLY	conflict	UNP Q04IN8
B	182	ALA	SER	conflict	UNP Q04IN8
B	260	GLN	GLU	conflict	UNP Q04IN8
B	266	ALA	LEU	conflict	UNP Q04IN8
B	268	LEU	LYS	conflict	UNP Q04IN8
B	277	GLN	GLU	conflict	UNP Q04IN8

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: Pneumolysin



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C42	Depositor
Number of particles used	6461	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	1.02	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.080	Depositor
Minimum map value	-0.023	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.005	Depositor
Recommended contour level	0.0282	Depositor
Map size (Å)	504.0, 504.0, 504.0	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.4, 1.4, 1.4	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	B	0.29	1/3798 (0.0%)	0.47	7/5164 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	57	THR	C-N	11.34	1.60	1.34

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	39	LEU	C-N-CD	-10.37	97.79	120.60
1	B	117	ASP	C-N-CD	-9.52	99.67	120.60
1	B	273	ALA	C-N-CD	-8.36	102.20	120.60
1	B	71	TYR	C-N-CD	-7.96	103.09	120.60
1	B	57	THR	O-C-N	-7.08	111.38	122.70
1	B	295	ASP	C-N-CD	-5.37	108.79	120.60
1	B	59	ASP	CB-CG-OD2	5.22	123.00	118.30

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	B	3725	0	3673	691	0
All	All	3725	0	3673	691	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 93.

All (691) 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:B:302:VAL:HG21	1:B:313:LEU:CD1	1.12	1.56
1:B:302:VAL:CG2	1:B:313:LEU:CD1	1.85	1.55
1:B:302:VAL:CG2	1:B:313:LEU:HD13	1.37	1.54
1:B:244:ARG:HB3	1:B:293:GLY:CA	1.43	1.46
1:B:235:VAL:HG22	1:B:335:PHE:CD1	1.46	1.46
1:B:39:LEU:HB3	1:B:40:PRO:CD	1.40	1.40
1:B:74:ALA:CB	1:B:237:ILE:HG13	1.49	1.39
1:B:455:ILE:CG2	1:B:464:VAL:HG13	1.52	1.38
1:B:235:VAL:HG22	1:B:335:PHE:CE1	1.60	1.37
1:B:72:PRO:HB2	1:B:99:TYR:CE1	1.59	1.35
1:B:235:VAL:CG2	1:B:335:PHE:CE1	2.10	1.35
1:B:233:PRO:HB3	1:B:337:ARG:CZ	1.60	1.32
1:B:135:HIS:CE1	1:B:322:ALA:HB2	1.66	1.29
1:B:244:ARG:CB	1:B:293:GLY:CA	2.11	1.28
1:B:135:HIS:CE1	1:B:322:ALA:CB	2.16	1.26
1:B:371:TYR:CE1	1:B:427:GLU:HB2	1.71	1.26
1:B:371:TYR:OH	1:B:455:ILE:HD12	1.27	1.24
1:B:379:TRP:HB3	1:B:421:LEU:CD1	1.70	1.20
1:B:233:PRO:CB	1:B:337:ARG:CZ	2.18	1.20
1:B:244:ARG:CB	1:B:293:GLY:HA3	1.71	1.19
1:B:39:LEU:CB	1:B:40:PRO:HD2	1.72	1.19
1:B:104:PRO:CG	1:B:144:VAL:HA	1.71	1.19
1:B:74:ALA:O	1:B:236:TYR:HB2	1.41	1.18
1:B:455:ILE:HG23	1:B:464:VAL:HG13	1.22	1.18
1:B:233:PRO:HB3	1:B:337:ARG:NH1	1.57	1.18
1:B:51:ARG:HH21	1:B:351:VAL:CG2	1.57	1.17
1:B:244:ARG:CB	1:B:293:GLY:HA2	1.73	1.17
1:B:244:ARG:HB3	1:B:293:GLY:HA3	1.22	1.16
1:B:74:ALA:HB2	1:B:237:ILE:HG13	1.27	1.15
1:B:103:LEU:HG	1:B:104:PRO:CD	1.76	1.15
1:B:228:ILE:HD12	1:B:234:LEU:HD22	1.29	1.14
1:B:81:THR:HB	1:B:86:ASN:O	1.43	1.13
1:B:295:ASP:OD2	1:B:324:HIS:ND1	1.81	1.13
1:B:70:LEU:O	1:B:237:ILE:HD12	1.50	1.11
1:B:195:PHE:HB3	1:B:246:VAL:HB	1.26	1.11
1:B:158:MET:HB2	1:B:187:GLU:CB	1.80	1.11
1:B:161:LEU:HB3	1:B:184:HIS:HB2	1.20	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:TYR:HB3	1:B:242:TYR:HB2	1.15	1.11
1:B:233:PRO:CB	1:B:337:ARG:NH2	2.12	1.11
1:B:51:ARG:HH21	1:B:351:VAL:HG23	1.08	1.10
1:B:235:VAL:HG21	1:B:335:PHE:CE1	1.87	1.10
1:B:244:ARG:HB3	1:B:293:GLY:HA2	1.17	1.09
1:B:71:TYR:HB2	1:B:72:PRO:HD3	1.27	1.09
1:B:455:ILE:HG22	1:B:464:VAL:HG13	1.32	1.09
1:B:154:THR:HG22	1:B:191:GLN:HB2	1.11	1.08
1:B:158:MET:CB	1:B:187:GLU:HB3	1.82	1.08
1:B:71:TYR:HB2	1:B:72:PRO:CD	1.83	1.08
1:B:379:TRP:CB	1:B:421:LEU:HD12	1.83	1.07
1:B:57:THR:HG21	1:B:60:ILE:HB	1.35	1.07
1:B:456:TRP:O	1:B:463:GLN:HB2	1.55	1.07
1:B:76:LEU:N	1:B:235:VAL:O	1.88	1.07
1:B:85:ASN:HD21	1:B:345:GLN:HB2	1.10	1.07
1:B:371:TYR:HE1	1:B:373:ALA:HB2	1.18	1.06
1:B:448:VAL:HG21	1:B:468:VAL:CG2	1.84	1.06
1:B:302:VAL:CG2	1:B:313:LEU:HD12	1.83	1.05
1:B:17:LYS:NZ	1:B:21:LEU:HB2	1.71	1.05
1:B:74:ALA:CB	1:B:237:ILE:CG1	2.35	1.05
1:B:103:LEU:CG	1:B:104:PRO:HD3	1.87	1.04
1:B:161:LEU:HB3	1:B:184:HIS:CB	1.88	1.04
1:B:51:ARG:NH2	1:B:351:VAL:HG23	1.72	1.03
1:B:74:ALA:HB3	1:B:237:ILE:HG13	1.36	1.03
1:B:57:THR:CG2	1:B:60:ILE:HB	1.89	1.03
1:B:39:LEU:HB3	1:B:40:PRO:HD3	1.38	1.02
1:B:150:TYR:HB2	1:B:195:PHE:HA	1.40	1.02
1:B:371:TYR:OH	1:B:455:ILE:CD1	2.07	1.02
1:B:72:PRO:HB2	1:B:99:TYR:CZ	1.95	1.02
1:B:375:TYR:HE2	1:B:408:PHE:CE2	1.77	1.02
1:B:104:PRO:HG3	1:B:144:VAL:HA	1.03	1.01
1:B:185:SER:OG	1:B:256:SER:HB2	1.60	1.01
1:B:104:PRO:HG3	1:B:145:PRO:CD	1.91	1.01
1:B:179:ASP:HB3	1:B:262:ALA:HB3	1.38	1.01
1:B:74:ALA:HB3	1:B:237:ILE:CG1	1.89	1.00
1:B:455:ILE:CG2	1:B:464:VAL:CG1	2.38	1.00
1:B:72:PRO:CB	1:B:99:TYR:CE1	2.45	1.00
1:B:74:ALA:O	1:B:236:TYR:CB	2.09	1.00
1:B:103:LEU:HG	1:B:104:PRO:HD3	1.03	1.00
1:B:133:LYS:HE3	1:B:137:ASP:HB2	1.44	1.00
1:B:3:ASN:O	1:B:6:VAL:HG13	1.59	1.00

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:235:VAL:CG2	1:B:335:PHE:CZ	2.44	1.00
1:B:455:ILE:HG22	1:B:464:VAL:CG1	1.92	1.00
1:B:95:ALA:HB1	1:B:96:PRO:HD2	1.40	1.00
1:B:15:TYR:CZ	1:B:337:ARG:CZ	2.44	1.00
1:B:448:VAL:CG1	1:B:451:ARG:HB2	1.92	0.99
1:B:103:LEU:CG	1:B:104:PRO:CD	2.39	0.99
1:B:235:VAL:CG2	1:B:335:PHE:CD1	2.39	0.99
1:B:418:VAL:HG12	1:B:419:ARG:N	1.77	0.99
1:B:448:VAL:HG21	1:B:468:VAL:HG23	1.04	0.99
1:B:154:THR:CG2	1:B:191:GLN:HB2	1.92	0.99
1:B:104:PRO:HG3	1:B:145:PRO:HD2	1.45	0.98
1:B:74:ALA:HB3	1:B:237:ILE:H	1.28	0.98
1:B:355:VAL:HG12	1:B:356:THR:H	1.26	0.98
1:B:71:TYR:CB	1:B:72:PRO:HD3	1.90	0.98
1:B:244:ARG:CA	1:B:293:GLY:HA2	1.94	0.97
1:B:302:VAL:HG23	1:B:313:LEU:CD1	1.94	0.96
1:B:209:ASN:HB3	1:B:210:PRO:CD	1.96	0.96
1:B:362:ASP:CB	1:B:413:PRO:HA	1.95	0.96
1:B:154:THR:HG22	1:B:191:GLN:CB	1.95	0.96
1:B:135:HIS:CE1	1:B:322:ALA:HB3	1.98	0.95
1:B:375:TYR:HE2	1:B:408:PHE:HE2	1.11	0.95
1:B:65:THR:O	1:B:69:ARG:HG2	1.67	0.95
1:B:235:VAL:HG21	1:B:335:PHE:CZ	2.02	0.95
1:B:202:VAL:HB	1:B:240:VAL:HG12	1.47	0.95
1:B:209:ASN:HB3	1:B:210:PRO:HD2	1.49	0.95
1:B:39:LEU:CB	1:B:40:PRO:CD	2.19	0.95
1:B:81:THR:OG1	1:B:87:PRO:HA	1.67	0.94
1:B:81:THR:OG1	1:B:87:PRO:HB3	1.65	0.94
1:B:233:PRO:HB2	1:B:337:ARG:NH2	1.82	0.94
1:B:118:PRO:HG3	1:B:122:SER:HB2	1.48	0.94
1:B:375:TYR:CE2	1:B:408:PHE:CE2	2.56	0.94
1:B:76:LEU:HD23	1:B:87:PRO:CG	1.98	0.93
1:B:81:THR:CB	1:B:87:PRO:HA	1.99	0.93
1:B:440:TYR:HE1	1:B:442:LYS:HE2	1.31	0.93
1:B:302:VAL:HG23	1:B:313:LEU:HD12	1.47	0.93
1:B:376:TYR:HD1	1:B:424:LYS:O	1.50	0.92
1:B:414:LEU:HG	1:B:447:LEU:HD21	1.50	0.92
1:B:228:ILE:HD12	1:B:234:LEU:CD2	1.98	0.92
1:B:244:ARG:HA	1:B:293:GLY:HA2	1.51	0.92
1:B:81:THR:OG1	1:B:87:PRO:CB	2.18	0.92
1:B:104:PRO:HG3	1:B:144:VAL:CA	1.97	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:418:VAL:HG12	1:B:419:ARG:H	1.34	0.91
1:B:81:THR:OG1	1:B:87:PRO:CA	2.19	0.91
1:B:233:PRO:HB2	1:B:337:ARG:CZ	2.00	0.91
1:B:236:TYR:O	1:B:334:SER:N	2.02	0.91
1:B:355:VAL:HG12	1:B:356:THR:N	1.82	0.91
1:B:39:LEU:HB3	1:B:40:PRO:HD2	0.93	0.91
1:B:85:ASN:ND2	1:B:345:GLN:HB2	1.85	0.90
1:B:156:HIS:HD2	1:B:189:GLN:HG3	1.36	0.90
1:B:371:TYR:CE1	1:B:373:ALA:HB2	2.07	0.90
1:B:418:VAL:CG1	1:B:419:ARG:H	1.84	0.90
1:B:99:TYR:HD2	1:B:115:VAL:HG21	1.37	0.90
1:B:200:TYR:CB	1:B:242:TYR:HB2	2.01	0.90
1:B:379:TRP:HB3	1:B:421:LEU:HD12	0.92	0.90
1:B:85:ASN:ND2	1:B:345:GLN:HE21	1.67	0.90
1:B:103:LEU:CB	1:B:104:PRO:CD	2.48	0.90
1:B:376:TYR:HB2	1:B:424:LYS:HB3	1.53	0.89
1:B:158:MET:HB2	1:B:187:GLU:HB3	0.92	0.89
1:B:76:LEU:HD23	1:B:87:PRO:HG3	1.55	0.89
1:B:370:ALA:O	1:B:459:THR:CG2	2.21	0.88
1:B:455:ILE:HG22	1:B:464:VAL:CB	2.02	0.88
1:B:448:VAL:CG2	1:B:468:VAL:HG23	1.98	0.88
1:B:135:HIS:ND1	1:B:322:ALA:HB2	1.87	0.88
1:B:302:VAL:HG21	1:B:313:LEU:HD11	1.52	0.88
1:B:51:ARG:NH2	1:B:351:VAL:CG2	2.31	0.88
1:B:17:LYS:HZ2	1:B:21:LEU:CB	1.88	0.87
1:B:244:ARG:HB2	1:B:293:GLY:HA3	1.54	0.87
1:B:44:VAL:HG12	1:B:356:THR:HG23	1.57	0.86
1:B:15:TYR:CZ	1:B:337:ARG:NH2	2.43	0.86
1:B:235:VAL:HG22	1:B:335:PHE:CG	2.11	0.86
1:B:236:TYR:N	1:B:334:SER:O	2.08	0.86
1:B:236:TYR:HE1	1:B:334:SER:HB3	1.41	0.85
1:B:99:TYR:CD2	1:B:115:VAL:HG21	2.10	0.85
1:B:455:ILE:HG22	1:B:464:VAL:HG22	1.56	0.85
1:B:418:VAL:CG1	1:B:419:ARG:N	2.41	0.84
1:B:74:ALA:O	1:B:236:TYR:HA	1.76	0.84
1:B:135:HIS:HE1	1:B:322:ALA:CB	1.90	0.84
1:B:355:VAL:CG1	1:B:356:THR:H	1.91	0.83
1:B:17:LYS:HZ3	1:B:21:LEU:HB2	1.43	0.83
1:B:448:VAL:HG11	1:B:451:ARG:HB2	1.59	0.83
1:B:309:MET:SD	1:B:310:VAL:N	2.51	0.83
1:B:362:ASP:HB3	1:B:413:PRO:HA	1.58	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:193:VAL:HG11	1:B:309:MET:HE1	1.61	0.83
1:B:208:LYS:HA	1:B:208:LYS:HZ3	1.42	0.83
1:B:455:ILE:HG22	1:B:464:VAL:CG2	2.09	0.83
1:B:78:VAL:HA	1:B:82:LEU:HD12	1.61	0.82
1:B:421:LEU:HD23	1:B:445:LEU:CD1	2.09	0.82
1:B:448:VAL:HG12	1:B:451:ARG:HB2	1.59	0.82
1:B:44:VAL:HG12	1:B:356:THR:CG2	2.09	0.82
1:B:156:HIS:CD2	1:B:189:GLN:HG3	2.14	0.82
1:B:329:ILE:O	1:B:347:SER:HB3	1.78	0.82
1:B:374:GLN:HG2	1:B:402:GLN:N	1.93	0.82
1:B:23:HIS:CD2	1:B:79:ASP:O	2.32	0.82
1:B:21:LEU:HD13	1:B:21:LEU:O	1.79	0.82
1:B:72:PRO:HB2	1:B:99:TYR:CD1	2.15	0.82
1:B:199:TYR:OH	1:B:244:ARG:NE	2.12	0.82
1:B:377:ILE:HD13	1:B:412:ILE:HD11	1.60	0.82
1:B:247:TYR:O	1:B:289:ALA:HB1	1.78	0.81
1:B:302:VAL:HG22	1:B:309:MET:HG2	1.61	0.81
1:B:236:TYR:CE1	1:B:334:SER:CB	2.63	0.81
1:B:74:ALA:O	1:B:236:TYR:CA	2.28	0.81
1:B:3:ASN:O	1:B:6:VAL:CG1	2.28	0.80
1:B:6:VAL:HG21	1:B:214:PHE:CD2	2.16	0.80
1:B:57:THR:HG21	1:B:60:ILE:CB	2.10	0.80
1:B:214:PHE:HE1	1:B:223:LEU:HD13	1.46	0.80
1:B:193:VAL:HB	1:B:248:LEU:HB3	1.61	0.80
1:B:228:ILE:CD1	1:B:234:LEU:HD22	2.12	0.80
1:B:103:LEU:CG	1:B:104:PRO:HD2	2.12	0.80
1:B:371:TYR:HH	1:B:455:ILE:HD12	1.45	0.80
1:B:15:TYR:CE2	1:B:337:ARG:NH2	2.50	0.80
1:B:17:LYS:HZ2	1:B:21:LEU:HB2	1.44	0.80
1:B:20:LEU:O	1:B:20:LEU:HD22	1.81	0.80
1:B:236:TYR:CE1	1:B:334:SER:HB3	2.17	0.80
1:B:379:TRP:CB	1:B:421:LEU:CD1	2.54	0.80
1:B:70:LEU:HB2	1:B:237:ILE:CD1	2.11	0.79
1:B:421:LEU:HD23	1:B:445:LEU:HD11	1.65	0.79
1:B:179:ASP:CB	1:B:262:ALA:HB3	2.13	0.79
1:B:376:TYR:CD1	1:B:424:LYS:O	2.36	0.79
1:B:440:TYR:CE1	1:B:442:LYS:HE2	2.16	0.79
1:B:219:THR:HB	1:B:222:ASP:HB2	1.65	0.78
1:B:103:LEU:HB3	1:B:104:PRO:HD2	1.63	0.78
1:B:15:TYR:CE1	1:B:337:ARG:NH1	2.52	0.78
1:B:214:PHE:CE1	1:B:223:LEU:HD13	2.18	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:LYS:NZ	1:B:21:LEU:CB	2.43	0.77
1:B:202:VAL:HB	1:B:240:VAL:CG1	2.12	0.77
1:B:295:ASP:OD2	1:B:324:HIS:CE1	2.37	0.77
1:B:70:LEU:HB3	1:B:74:ALA:HB2	1.67	0.77
1:B:309:MET:SD	1:B:310:VAL:HG13	2.24	0.77
1:B:61:SER:O	1:B:65:THR:HG23	1.85	0.76
1:B:377:ILE:HD13	1:B:397:TRP:CE3	2.21	0.76
1:B:103:LEU:CB	1:B:104:PRO:HD2	2.15	0.76
1:B:158:MET:HA	1:B:187:GLU:HA	1.66	0.76
1:B:213:VAL:HG22	1:B:336:LEU:HG	1.68	0.76
1:B:371:TYR:CD1	1:B:427:GLU:HB2	2.21	0.76
1:B:59:ASP:O	1:B:63:THR:HG23	1.85	0.75
1:B:75:LEU:HA	1:B:235:VAL:O	1.86	0.75
1:B:89:LEU:N	1:B:89:LEU:HD13	2.01	0.75
1:B:70:LEU:HB2	1:B:237:ILE:HD11	1.69	0.75
1:B:103:LEU:HB3	1:B:104:PRO:CD	2.17	0.75
1:B:372:VAL:HG23	1:B:428:CYS:HB3	1.70	0.74
1:B:214:PHE:HB3	1:B:218:VAL:HG21	1.69	0.74
1:B:236:TYR:HE1	1:B:334:SER:CB	2.00	0.74
1:B:455:ILE:CG2	1:B:464:VAL:HG22	2.17	0.74
1:B:51:ARG:HH21	1:B:351:VAL:HG21	1.50	0.74
1:B:104:PRO:CG	1:B:145:PRO:HD2	2.15	0.74
1:B:72:PRO:CB	1:B:99:TYR:CD1	2.69	0.74
1:B:239:SER:OG	1:B:332:THR:HG21	1.88	0.74
1:B:375:TYR:CE2	1:B:408:PHE:HE2	1.99	0.74
1:B:34:LYS:HG3	1:B:34:LYS:O	1.87	0.73
1:B:374:GLN:HG2	1:B:402:GLN:H	1.53	0.73
1:B:107:ALA:O	1:B:111:SER:HB2	1.89	0.73
1:B:421:LEU:HB3	1:B:445:LEU:CD1	2.19	0.73
1:B:133:LYS:HE3	1:B:137:ASP:CB	2.17	0.73
1:B:327:LEU:HD22	1:B:328:PRO:HD2	1.70	0.73
1:B:16:ASP:OD1	1:B:16:ASP:N	2.21	0.73
1:B:15:TYR:CZ	1:B:337:ARG:NH1	2.56	0.73
1:B:57:THR:OG1	1:B:60:ILE:HD12	1.88	0.73
1:B:208:LYS:HA	1:B:208:LYS:NZ	2.04	0.72
1:B:233:PRO:CG	1:B:337:ARG:NH2	2.51	0.72
1:B:298:SER:O	1:B:313:LEU:CD1	2.37	0.72
1:B:70:LEU:O	1:B:237:ILE:HG21	1.90	0.72
1:B:23:HIS:HD2	1:B:79:ASP:O	1.72	0.72
1:B:9:PHE:CZ	1:B:13:MET:HE2	2.24	0.72
1:B:76:LEU:HD23	1:B:87:PRO:HG2	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:ASN:HA	1:B:6:VAL:CG1	2.20	0.72
1:B:161:LEU:N	1:B:184:HIS:O	2.22	0.72
1:B:233:PRO:HG3	1:B:337:ARG:HH22	1.54	0.72
1:B:248:LEU:HA	1:B:289:ALA:HB2	1.72	0.72
1:B:168:ASP:HB3	1:B:177:ASP:OD1	1.90	0.71
1:B:236:TYR:CE1	1:B:334:SER:HB2	2.24	0.71
1:B:81:THR:HB	1:B:87:PRO:HA	1.73	0.71
1:B:148:MET:HG3	1:B:197:GLN:HG3	1.71	0.71
1:B:205:ASP:O	1:B:206:ALA:O	2.07	0.71
1:B:171:LYS:CG	1:B:172:ALA:N	2.53	0.70
1:B:381:GLU:O	1:B:381:GLU:HG3	1.89	0.70
1:B:47:GLU:OE1	1:B:470:ASN:HB3	1.91	0.70
1:B:76:LEU:O	1:B:235:VAL:HB	1.90	0.70
1:B:250:LEU:HB3	1:B:287:VAL:HG23	1.74	0.70
1:B:327:LEU:HD22	1:B:328:PRO:CD	2.22	0.70
1:B:377:ILE:HG21	1:B:412:ILE:HD13	1.74	0.69
1:B:2:ALA:O	1:B:6:VAL:HG12	1.92	0.69
1:B:437:ARG:HH21	1:B:437:ARG:HG3	1.55	0.69
1:B:355:VAL:CG1	1:B:449:ARG:NH2	2.55	0.69
1:B:219:THR:CB	1:B:222:ASP:HB2	2.22	0.69
1:B:62:VAL:O	1:B:65:THR:OG1	2.11	0.69
1:B:71:TYR:HB2	1:B:72:PRO:HD2	1.72	0.69
1:B:248:LEU:HD23	1:B:309:MET:CE	2.23	0.69
1:B:259:VAL:HB	1:B:278:TRP:HB2	1.73	0.69
1:B:362:ASP:HB3	1:B:413:PRO:CA	2.23	0.69
1:B:440:TYR:HE1	1:B:442:LYS:CE	2.04	0.69
1:B:60:ILE:O	1:B:63:THR:OG1	2.11	0.68
1:B:207:VAL:HG13	1:B:212:ASP:OD2	1.93	0.68
1:B:371:TYR:HB3	1:B:459:THR:HG22	1.74	0.68
1:B:9:PHE:CZ	1:B:13:MET:CE	2.76	0.68
1:B:147:ARG:H	1:B:147:ARG:HD3	1.57	0.68
1:B:51:ARG:HD2	1:B:348:THR:HG23	1.75	0.68
1:B:15:TYR:HD1	1:B:15:TYR:H	1.41	0.68
1:B:30:ASN:CB	1:B:45:VAL:HG12	2.23	0.68
1:B:70:LEU:O	1:B:237:ILE:CD1	2.36	0.67
1:B:29:GLU:HB3	1:B:46:ILE:HG13	1.74	0.67
1:B:188:LYS:HG2	1:B:253:THR:HG22	1.77	0.67
1:B:44:VAL:CG1	1:B:356:THR:HG23	2.24	0.67
1:B:85:ASN:HD21	1:B:345:GLN:CB	1.99	0.67
1:B:237:ILE:HA	1:B:333:THR:HA	1.77	0.67
1:B:446:PRO:O	1:B:451:ARG:NH2	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:VAL:O	1:B:232:ARG:HB3	1.94	0.67
1:B:213:VAL:HG21	1:B:336:LEU:CD1	2.24	0.66
1:B:298:SER:O	1:B:313:LEU:HD12	1.96	0.66
1:B:452:THR:N	1:B:467:LYS:O	2.27	0.66
1:B:89:LEU:HD13	1:B:89:LEU:H	1.56	0.66
1:B:95:ALA:HB1	1:B:96:PRO:CD	2.22	0.66
1:B:378:THR:O	1:B:422:SER:N	2.29	0.66
1:B:429:THR:HG22	1:B:431:LEU:HD23	1.76	0.66
1:B:179:ASP:HB3	1:B:262:ALA:CB	2.21	0.66
1:B:79:ASP:OD2	1:B:232:ARG:NH2	2.29	0.66
1:B:193:VAL:HG11	1:B:309:MET:CE	2.26	0.66
1:B:238:SER:H	1:B:333:THR:HA	1.61	0.66
1:B:415:LYS:NZ	1:B:415:LYS:HB2	2.11	0.66
1:B:421:LEU:HB3	1:B:445:LEU:HD11	1.77	0.65
1:B:149:GLN:HG3	1:B:149:GLN:O	1.96	0.65
1:B:171:LYS:HG2	1:B:172:ALA:N	2.11	0.65
1:B:437:ARG:HG3	1:B:437:ARG:NH2	2.09	0.65
1:B:209:ASN:CB	1:B:210:PRO:CD	2.70	0.65
1:B:246:VAL:CG2	1:B:313:LEU:HD22	2.27	0.65
1:B:158:MET:CB	1:B:187:GLU:CB	2.59	0.65
1:B:214:PHE:HB3	1:B:218:VAL:CG2	2.26	0.65
1:B:310:VAL:O	1:B:314:ILE:HG22	1.97	0.65
1:B:79:ASP:OD2	1:B:232:ARG:CZ	2.45	0.65
1:B:46:ILE:HG13	1:B:46:ILE:O	1.97	0.65
1:B:70:LEU:C	1:B:237:ILE:HD12	2.18	0.65
1:B:133:LYS:CE	1:B:137:ASP:HB2	2.24	0.64
1:B:177:ASP:HB3	1:B:264:GLU:HB2	1.80	0.64
1:B:329:ILE:O	1:B:347:SER:CB	2.44	0.64
1:B:361:GLY:O	1:B:414:LEU:HD23	1.98	0.64
1:B:85:ASN:ND2	1:B:345:GLN:NE2	2.45	0.64
1:B:195:PHE:HB3	1:B:246:VAL:CB	2.16	0.64
1:B:370:ALA:O	1:B:459:THR:HG21	1.97	0.64
1:B:84:GLU:HG2	1:B:350:TYR:HD2	1.62	0.64
1:B:355:VAL:HG12	1:B:449:ARG:NH2	2.13	0.64
1:B:361:GLY:O	1:B:414:LEU:HB2	1.97	0.64
1:B:195:PHE:CB	1:B:246:VAL:HB	2.15	0.64
1:B:207:VAL:C	1:B:208:LYS:HE2	2.17	0.64
1:B:248:LEU:CD2	1:B:309:MET:HE2	2.27	0.64
1:B:355:VAL:CG1	1:B:449:ARG:HH21	2.10	0.64
1:B:104:PRO:CG	1:B:145:PRO:CD	2.70	0.64
1:B:185:SER:HG	1:B:256:SER:HB2	1.63	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:74:ALA:HB3	1:B:237:ILE:N	2.07	0.63
1:B:207:VAL:HG21	1:B:336:LEU:HD23	1.79	0.63
1:B:236:TYR:HD1	1:B:334:SER:O	1.79	0.63
1:B:247:TYR:HB2	1:B:290:VAL:O	1.98	0.63
1:B:94:ARG:HD3	1:B:117:ASP:HB2	1.81	0.63
1:B:259:VAL:O	1:B:278:TRP:HB2	1.98	0.63
1:B:118:PRO:CG	1:B:122:SER:HB2	2.26	0.63
1:B:159:GLU:N	1:B:186:GLY:O	2.31	0.63
1:B:17:LYS:HD2	1:B:17:LYS:O	1.98	0.63
1:B:376:TYR:CB	1:B:424:LYS:HB3	2.29	0.63
1:B:214:PHE:CB	1:B:218:VAL:HG21	2.29	0.62
1:B:78:VAL:O	1:B:78:VAL:HG13	2.00	0.62
1:B:74:ALA:C	1:B:236:TYR:HA	2.20	0.62
1:B:155:ALA:N	1:B:190:ILE:O	2.26	0.62
1:B:199:TYR:OH	1:B:244:ARG:CZ	2.47	0.61
1:B:248:LEU:HD23	1:B:309:MET:HE2	1.81	0.61
1:B:373:ALA:CB	1:B:427:GLU:HA	2.30	0.61
1:B:76:LEU:CD2	1:B:87:PRO:HG2	2.29	0.61
1:B:355:VAL:HG12	1:B:449:ARG:HH21	1.65	0.61
1:B:15:TYR:OH	1:B:337:ARG:NH2	2.34	0.61
1:B:135:HIS:HE1	1:B:322:ALA:HB3	1.56	0.61
1:B:233:PRO:CG	1:B:337:ARG:HH22	2.12	0.61
1:B:419:ARG:O	1:B:420:ASN:HB2	2.00	0.61
1:B:10:ILE:HG23	1:B:337:ARG:HG2	1.82	0.61
1:B:207:VAL:HG11	1:B:336:LEU:HD21	1.83	0.61
1:B:250:LEU:CB	1:B:287:VAL:HG23	2.30	0.61
1:B:196:LYS:HD3	1:B:245:GLN:HE21	1.64	0.61
1:B:228:ILE:HG22	1:B:228:ILE:O	2.01	0.61
1:B:17:LYS:HD3	1:B:341:VAL:HG13	1.83	0.61
1:B:104:PRO:HG3	1:B:145:PRO:HD3	1.80	0.61
1:B:21:LEU:HD13	1:B:21:LEU:C	2.21	0.61
1:B:161:LEU:O	1:B:184:HIS:N	2.18	0.60
1:B:98:THR:HG22	1:B:113:LEU:O	2.02	0.60
1:B:426:ARG:O	1:B:426:ARG:HG3	2.01	0.60
1:B:75:LEU:CA	1:B:235:VAL:O	2.49	0.60
1:B:65:THR:O	1:B:69:ARG:CG	2.47	0.60
1:B:72:PRO:HB3	1:B:99:TYR:CD1	2.36	0.60
1:B:355:VAL:HG11	1:B:449:ARG:NH2	2.16	0.60
1:B:39:LEU:CB	1:B:40:PRO:HD3	2.12	0.59
1:B:253:THR:OG1	1:B:284:ASN:HB2	2.02	0.59
1:B:117:ASP:N	1:B:118:PRO:HD3	2.18	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:362:ASP:CA	1:B:413:PRO:HA	2.33	0.59
1:B:379:TRP:CG	1:B:421:LEU:CD1	2.86	0.58
1:B:150:TYR:CB	1:B:195:PHE:HA	2.25	0.58
1:B:213:VAL:HG21	1:B:336:LEU:HD12	1.86	0.58
1:B:94:ARG:HD3	1:B:117:ASP:CB	2.34	0.58
1:B:207:VAL:CB	1:B:336:LEU:HD21	2.33	0.58
1:B:247:TYR:HD2	1:B:290:VAL:HG23	1.68	0.58
1:B:150:TYR:HD1	1:B:194:ASN:O	1.87	0.58
1:B:66:ASN:HA	1:B:69:ARG:HG3	1.85	0.57
1:B:15:TYR:OH	1:B:341:VAL:HG21	2.05	0.57
1:B:291:ILE:O	1:B:291:ILE:HG23	2.03	0.57
1:B:421:LEU:HB3	1:B:445:LEU:HD12	1.86	0.57
1:B:455:ILE:HG22	1:B:464:VAL:CA	2.34	0.57
1:B:15:TYR:OH	1:B:337:ARG:CZ	2.52	0.57
1:B:85:ASN:HD22	1:B:345:GLN:HE21	1.49	0.57
1:B:101:ILE:HG21	1:B:200:TYR:CE2	2.40	0.57
1:B:259:VAL:HB	1:B:278:TRP:CB	2.35	0.57
1:B:298:SER:O	1:B:313:LEU:HD11	2.04	0.57
1:B:71:TYR:HD2	1:B:240:VAL:HG21	1.68	0.57
1:B:455:ILE:HA	1:B:464:VAL:HA	1.87	0.57
1:B:17:LYS:NZ	1:B:21:LEU:CG	2.67	0.56
1:B:30:ASN:O	1:B:30:ASN:ND2	2.34	0.56
1:B:259:VAL:O	1:B:278:TRP:N	2.38	0.56
1:B:364:LEU:HD12	1:B:452:THR:OG1	2.05	0.56
1:B:81:THR:CB	1:B:87:PRO:CA	2.78	0.56
1:B:78:VAL:CA	1:B:82:LEU:HD12	2.34	0.56
1:B:154:THR:HA	1:B:191:GLN:HA	1.87	0.56
1:B:104:PRO:HG2	1:B:144:VAL:HA	1.76	0.56
1:B:147:ARG:H	1:B:147:ARG:CD	2.19	0.56
1:B:246:VAL:HG22	1:B:291:ILE:HD12	1.88	0.56
1:B:370:ALA:O	1:B:459:THR:HG22	2.05	0.56
1:B:205:ASP:O	1:B:206:ALA:C	2.43	0.56
1:B:170:GLU:O	1:B:170:GLU:HG3	2.05	0.56
1:B:373:ALA:HB2	1:B:427:GLU:HA	1.86	0.56
1:B:15:TYR:CE1	1:B:337:ARG:CZ	2.89	0.56
1:B:202:VAL:N	1:B:240:VAL:O	2.34	0.56
1:B:228:ILE:CD1	1:B:234:LEU:CD2	2.76	0.56
1:B:81:THR:HB	1:B:86:ASN:C	2.22	0.56
1:B:385:ASP:OD2	1:B:391:VAL:HG13	2.06	0.56
1:B:377:ILE:HD13	1:B:397:TRP:CZ3	2.41	0.55
1:B:372:VAL:O	1:B:428:CYS:CB	2.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:ARG:CA	1:B:293:GLY:CA	2.70	0.55
1:B:76:LEU:N	1:B:76:LEU:HD12	2.22	0.55
1:B:51:ARG:HD2	1:B:348:THR:CG2	2.37	0.55
1:B:144:VAL:HG11	1:B:320:PHE:CB	2.36	0.55
1:B:207:VAL:HB	1:B:336:LEU:CD2	2.37	0.55
1:B:17:LYS:HZ2	1:B:21:LEU:CG	2.20	0.55
1:B:55:THR:HG23	1:B:55:THR:O	2.07	0.55
1:B:231:GLU:OE1	1:B:231:GLU:HA	2.07	0.55
1:B:246:VAL:HG21	1:B:313:LEU:HD22	1.89	0.55
1:B:196:LYS:HG3	1:B:245:GLN:HG3	1.89	0.55
1:B:336:LEU:C	1:B:336:LEU:HD13	2.28	0.55
1:B:57:THR:HG21	1:B:60:ILE:CG2	2.36	0.54
1:B:193:VAL:CG1	1:B:309:MET:HE1	2.34	0.54
1:B:363:LEU:O	1:B:411:SER:HA	2.07	0.54
1:B:208:LYS:HE2	1:B:208:LYS:N	2.22	0.54
1:B:239:SER:O	1:B:332:THR:CG2	2.55	0.54
1:B:44:VAL:CB	1:B:356:THR:HG23	2.38	0.54
1:B:58:SER:O	1:B:61:SER:OG	2.18	0.54
1:B:213:VAL:CG2	1:B:336:LEU:HG	2.35	0.54
1:B:72:PRO:HD2	1:B:123:VAL:CG2	2.36	0.54
1:B:248:LEU:HD23	1:B:309:MET:HE3	1.89	0.54
1:B:448:VAL:HG11	1:B:451:ARG:CB	2.33	0.54
1:B:311:GLU:HA	1:B:314:ILE:HG22	1.90	0.54
1:B:207:VAL:CG1	1:B:336:LEU:HD21	2.38	0.54
1:B:244:ARG:HB2	1:B:293:GLY:CA	2.20	0.54
1:B:9:PHE:CE2	1:B:13:MET:CE	2.90	0.54
1:B:377:ILE:HG22	1:B:379:TRP:CZ3	2.43	0.54
1:B:58:SER:O	1:B:62:VAL:HG23	2.08	0.54
1:B:82:LEU:HD11	1:B:235:VAL:HB	1.88	0.54
1:B:252:THR:HA	1:B:284:ASN:O	2.07	0.54
1:B:302:VAL:CG2	1:B:313:LEU:HD11	2.19	0.54
1:B:379:TRP:N	1:B:379:TRP:CE3	2.76	0.54
1:B:6:VAL:HG21	1:B:214:PHE:HD2	1.70	0.53
1:B:362:ASP:CB	1:B:413:PRO:CA	2.77	0.53
1:B:144:VAL:HG11	1:B:320:PHE:HB3	1.89	0.53
1:B:415:LYS:HB2	1:B:415:LYS:HZ3	1.73	0.53
1:B:449:ARG:HB3	1:B:470:ASN:HA	1.90	0.53
1:B:161:LEU:HB3	1:B:184:HIS:HB3	1.86	0.53
1:B:171:LYS:CG	1:B:172:ALA:H	2.22	0.53
1:B:362:ASP:HB3	1:B:413:PRO:CB	2.38	0.53
1:B:426:ARG:HB2	1:B:437:ARG:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:PHE:CE1	1:B:13:MET:HE2	2.43	0.53
1:B:235:VAL:HA	1:B:335:PHE:HA	1.91	0.53
1:B:239:SER:H	1:B:332:THR:HG23	1.73	0.53
1:B:314:ILE:O	1:B:314:ILE:HG23	2.07	0.53
1:B:402:GLN:HG3	1:B:403:ASP:N	2.23	0.53
1:B:371:TYR:HE1	1:B:373:ALA:CB	2.07	0.53
1:B:207:VAL:HB	1:B:336:LEU:HD21	1.90	0.53
1:B:151:GLU:HG2	1:B:151:GLU:O	2.08	0.53
1:B:171:LYS:HG3	1:B:172:ALA:H	1.73	0.52
1:B:207:VAL:HG21	1:B:336:LEU:CD2	2.39	0.52
1:B:314:ILE:C	1:B:314:ILE:HD12	2.30	0.52
1:B:199:TYR:CD1	1:B:199:TYR:N	2.77	0.52
1:B:363:LEU:O	1:B:412:ILE:N	2.41	0.52
1:B:362:ASP:HA	1:B:413:PRO:HA	1.91	0.52
1:B:426:ARG:CZ	1:B:436:TRP:CG	2.92	0.52
1:B:66:ASN:O	1:B:69:ARG:HG3	2.09	0.52
1:B:44:VAL:HG12	1:B:356:THR:HG21	1.90	0.52
1:B:3:ASN:HA	1:B:6:VAL:HG12	1.92	0.52
1:B:20:LEU:HD22	1:B:20:LEU:C	2.28	0.52
1:B:29:GLU:HB3	1:B:46:ILE:CG1	2.40	0.52
1:B:85:ASN:HD22	1:B:345:GLN:NE2	2.04	0.52
1:B:146:ALA:HB2	1:B:199:TYR:HD2	1.75	0.52
1:B:155:ALA:O	1:B:190:ILE:HB	2.10	0.52
1:B:103:LEU:CD2	1:B:104:PRO:HD2	2.40	0.52
1:B:9:PHE:CZ	1:B:13:MET:HE1	2.44	0.52
1:B:43:PHE:CZ	1:B:447:LEU:HD22	2.45	0.52
1:B:165:PHE:HD2	1:B:180:PHE:HB2	1.75	0.51
1:B:239:SER:N	1:B:332:THR:HG23	2.25	0.51
1:B:378:THR:HG23	1:B:396:ALA:HB2	1.93	0.51
1:B:426:ARG:HD2	1:B:436:TRP:HB3	1.91	0.51
1:B:4:LYS:O	1:B:8:ASP:OD1	2.28	0.51
1:B:301:ARG:N	1:B:301:ARG:CD	2.73	0.51
1:B:70:LEU:CB	1:B:237:ILE:CD1	2.85	0.51
1:B:376:TYR:O	1:B:423:VAL:HA	2.10	0.51
1:B:414:LEU:CG	1:B:447:LEU:HD21	2.32	0.51
1:B:455:ILE:HG22	1:B:464:VAL:HA	1.93	0.51
1:B:195:PHE:HD1	1:B:196:LYS:N	2.08	0.51
1:B:103:LEU:HD23	1:B:104:PRO:HD2	1.92	0.51
1:B:146:ALA:CB	1:B:199:TYR:CD2	2.93	0.51
1:B:10:ILE:HG23	1:B:337:ARG:CG	2.41	0.51
1:B:177:ASP:N	1:B:264:GLU:O	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:TYR:CD1	1:B:334:SER:O	2.61	0.51
1:B:333:THR:OG1	1:B:343:THR:HG22	2.10	0.51
1:B:221:GLU:HA	1:B:224:LYS:HE2	1.92	0.51
1:B:3:ASN:CA	1:B:6:VAL:CG1	2.88	0.50
1:B:99:TYR:CE1	1:B:204:VAL:HG23	2.46	0.50
1:B:148:MET:HA	1:B:197:GLN:HG2	1.91	0.50
1:B:301:ARG:N	1:B:301:ARG:HD2	2.27	0.50
1:B:439:VAL:CG1	1:B:464:VAL:HG21	2.42	0.50
1:B:154:THR:HG22	1:B:191:GLN:CA	2.42	0.50
1:B:3:ASN:HA	1:B:6:VAL:HG11	1.94	0.50
1:B:456:TRP:CD1	1:B:456:TRP:C	2.86	0.50
1:B:104:PRO:CD	1:B:145:PRO:HD2	2.42	0.50
1:B:411:SER:O	1:B:413:PRO:HD3	2.12	0.50
1:B:66:ASN:HA	1:B:69:ARG:CG	2.41	0.49
1:B:118:PRO:HG3	1:B:122:SER:CB	2.33	0.49
1:B:70:LEU:HB2	1:B:237:ILE:HD12	1.93	0.49
1:B:72:PRO:HD2	1:B:123:VAL:HG23	1.94	0.49
1:B:156:HIS:HB2	1:B:189:GLN:HG2	1.93	0.49
1:B:374:GLN:HG2	1:B:402:GLN:CA	2.42	0.49
1:B:377:ILE:CD1	1:B:397:TRP:CE3	2.94	0.49
1:B:451:ARG:HG3	1:B:468:VAL:HB	1.93	0.49
1:B:44:VAL:HG23	1:B:45:VAL:N	2.25	0.49
1:B:218:VAL:HG12	1:B:219:THR:N	2.27	0.49
1:B:356:THR:O	1:B:449:ARG:NH2	2.45	0.49
1:B:17:LYS:NZ	1:B:21:LEU:HG	2.28	0.49
1:B:69:ARG:NH2	1:B:328:PRO:HG2	2.28	0.49
1:B:105:GLY:O	1:B:134:TRP:HZ3	1.96	0.49
1:B:118:PRO:HG3	1:B:122:SER:H	1.77	0.49
1:B:28:ILE:HG22	1:B:47:GLU:HB2	1.95	0.49
1:B:92:VAL:O	1:B:94:ARG:HG3	2.11	0.49
1:B:327:LEU:O	1:B:329:ILE:HG12	2.13	0.49
1:B:76:LEU:N	1:B:76:LEU:CD1	2.76	0.48
1:B:104:PRO:CB	1:B:145:PRO:HD3	2.43	0.48
1:B:76:LEU:CD2	1:B:87:PRO:CG	2.81	0.48
1:B:213:VAL:CG2	1:B:336:LEU:CD1	2.91	0.48
1:B:425:ILE:HG22	1:B:425:ILE:O	2.13	0.48
1:B:418:VAL:HG13	1:B:419:ARG:H	1.75	0.48
1:B:366:ASP:OD1	1:B:409:THR:HG23	2.13	0.48
1:B:104:PRO:CG	1:B:145:PRO:HD3	2.41	0.48
1:B:364:LEU:HD23	1:B:411:SER:OG	2.13	0.48
1:B:377:ILE:CD1	1:B:412:ILE:HD11	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:ASN:HB2	1:B:45:VAL:HG12	1.96	0.47
1:B:335:PHE:HB2	1:B:341:VAL:O	2.14	0.47
1:B:3:ASN:CA	1:B:6:VAL:HG12	2.44	0.47
1:B:97:MET:O	1:B:115:VAL:N	2.35	0.47
1:B:146:ALA:HB2	1:B:199:TYR:CD2	2.49	0.47
1:B:32:PHE:CZ	1:B:418:VAL:O	2.66	0.47
1:B:81:THR:HB	1:B:87:PRO:CA	2.44	0.47
1:B:90:LEU:HD11	1:B:226:ARG:HD3	1.97	0.47
1:B:367:HIS:HB2	1:B:375:TYR:OH	2.12	0.47
1:B:104:PRO:CG	1:B:144:VAL:CA	2.66	0.47
1:B:372:VAL:O	1:B:428:CYS:HB3	2.13	0.47
1:B:442:LYS:HB2	1:B:445:LEU:HD23	1.97	0.47
1:B:78:VAL:HG12	1:B:233:PRO:O	2.14	0.47
1:B:158:MET:HA	1:B:187:GLU:CA	2.39	0.47
1:B:179:ASP:HB3	1:B:262:ALA:O	2.15	0.47
1:B:238:SER:N	1:B:333:THR:HA	2.28	0.47
1:B:37:ASN:HB2	1:B:384:TYR:CD2	2.49	0.47
1:B:70:LEU:HD23	1:B:70:LEU:N	2.30	0.47
1:B:198:ILE:HG23	1:B:242:TYR:O	2.15	0.47
1:B:452:THR:O	1:B:467:LYS:N	2.24	0.47
1:B:17:LYS:HZ2	1:B:21:LEU:HG	1.80	0.47
1:B:208:LYS:HA	1:B:208:LYS:CE	2.45	0.46
1:B:3:ASN:C	1:B:6:VAL:CG1	2.83	0.46
1:B:148:MET:CE	1:B:314:ILE:HD11	2.45	0.46
1:B:235:VAL:CG2	1:B:335:PHE:CE2	2.96	0.46
1:B:431:LEU:HB3	1:B:432:ALA:H	1.54	0.46
1:B:208:LYS:CE	1:B:208:LYS:CA	2.94	0.46
1:B:381:GLU:HG2	1:B:393:THR:O	2.14	0.46
1:B:191:GLN:O	1:B:191:GLN:HG3	2.15	0.46
1:B:379:TRP:CG	1:B:421:LEU:HD12	2.48	0.46
1:B:57:THR:HG23	1:B:60:ILE:HB	1.87	0.46
1:B:239:SER:O	1:B:332:THR:HG22	2.14	0.46
1:B:309:MET:SD	1:B:310:VAL:CA	3.04	0.46
1:B:426:ARG:CZ	1:B:436:TRP:CD2	2.99	0.46
1:B:30:ASN:HB3	1:B:45:VAL:HG12	1.95	0.46
1:B:86:ASN:N	1:B:87:PRO:HD3	2.31	0.46
1:B:70:LEU:CB	1:B:237:ILE:HD12	2.45	0.46
1:B:421:LEU:CD2	1:B:445:LEU:HD11	2.42	0.46
1:B:246:VAL:HG22	1:B:313:LEU:HD22	1.98	0.46
1:B:302:VAL:HG21	1:B:313:LEU:HD13	0.48	0.46
1:B:81:THR:HG1	1:B:87:PRO:HA	1.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:455:ILE:CG2	1:B:464:VAL:CG2	2.83	0.45
1:B:453:ILE:O	1:B:453:ILE:HG22	2.16	0.45
1:B:74:ALA:HB3	1:B:237:ILE:CB	2.43	0.45
1:B:259:VAL:HG23	1:B:278:TRP:HB3	1.98	0.45
1:B:29:GLU:N	1:B:46:ILE:O	2.49	0.45
1:B:75:LEU:C	1:B:235:VAL:O	2.51	0.45
1:B:199:TYR:CD1	1:B:243:GLY:HA2	2.51	0.45
1:B:335:PHE:HB2	1:B:341:VAL:HB	1.97	0.45
1:B:434:GLU:OE2	1:B:437:ARG:NH2	2.33	0.45
1:B:214:PHE:CG	1:B:218:VAL:HG11	2.52	0.45
1:B:302:VAL:HA	1:B:309:MET:CG	2.47	0.45
1:B:38:GLN:O	1:B:38:GLN:HG3	2.15	0.45
1:B:53:LEU:C	1:B:56:ASN:HD21	2.18	0.45
1:B:84:GLU:HA	1:B:350:TYR:CE2	2.52	0.45
1:B:177:ASP:CB	1:B:264:GLU:HB2	2.46	0.45
1:B:207:VAL:CB	1:B:336:LEU:CD2	2.93	0.45
1:B:214:PHE:HB3	1:B:218:VAL:HG11	1.99	0.45
1:B:336:LEU:CD1	1:B:336:LEU:C	2.86	0.45
1:B:362:ASP:HB2	1:B:413:PRO:HA	1.90	0.45
1:B:440:TYR:HE1	1:B:442:LYS:CD	2.29	0.45
1:B:118:PRO:CG	1:B:122:SER:H	2.30	0.45
1:B:244:ARG:NH1	1:B:317:GLY:O	2.50	0.45
1:B:417:ASN:O	1:B:418:VAL:HG23	2.17	0.45
1:B:76:LEU:CG	1:B:87:PRO:HG2	2.47	0.44
1:B:179:ASP:CB	1:B:262:ALA:O	2.65	0.44
1:B:223:LEU:O	1:B:223:LEU:HG	2.16	0.44
1:B:236:TYR:O	1:B:334:SER:O	2.34	0.44
1:B:15:TYR:CD1	1:B:15:TYR:N	2.73	0.44
1:B:315:GLN:NE2	1:B:315:GLN:HA	2.32	0.44
1:B:379:TRP:HB2	1:B:420:ASN:O	2.16	0.44
1:B:76:LEU:O	1:B:235:VAL:N	2.44	0.44
1:B:150:TYR:CD1	1:B:194:ASN:O	2.69	0.44
1:B:238:SER:HB3	1:B:239:SER:H	1.62	0.44
1:B:51:ARG:HH22	1:B:60:ILE:HG21	1.82	0.44
1:B:401:GLY:O	1:B:402:GLN:HB2	2.18	0.44
1:B:404:LEU:HD22	1:B:408:PHE:CE1	2.52	0.44
1:B:15:TYR:CD1	1:B:337:ARG:NH1	2.85	0.44
1:B:21:LEU:C	1:B:21:LEU:CD1	2.86	0.44
1:B:375:TYR:CD1	1:B:425:ILE:HG23	2.52	0.44
1:B:452:THR:HB	1:B:467:LYS:HB2	1.98	0.44
1:B:94:ARG:HB3	1:B:212:ASP:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:239:SER:O	1:B:332:THR:HG23	2.18	0.44
1:B:248:LEU:HA	1:B:289:ALA:CB	2.42	0.44
1:B:371:TYR:HE1	1:B:427:GLU:HB2	1.60	0.44
1:B:439:VAL:HG11	1:B:464:VAL:CG2	2.48	0.44
1:B:95:ALA:CB	1:B:96:PRO:CD	2.92	0.44
1:B:46:ILE:HB	1:B:354:LYS:HG2	1.98	0.44
1:B:374:GLN:O	1:B:376:TYR:CD1	2.70	0.44
1:B:426:ARG:NE	1:B:436:TRP:CD2	2.86	0.44
1:B:374:GLN:O	1:B:376:TYR:CE1	2.71	0.43
1:B:71:TYR:HB3	1:B:240:VAL:HG11	2.00	0.43
1:B:153:ILE:O	1:B:153:ILE:HG12	2.18	0.43
1:B:215:GLN:O	1:B:218:VAL:HG23	2.18	0.43
1:B:415:LYS:HB2	1:B:415:LYS:HZ2	1.83	0.43
1:B:437:ARG:HH21	1:B:437:ARG:CG	2.23	0.43
1:B:426:ARG:NH1	1:B:436:TRP:CG	2.87	0.43
1:B:84:GLU:HG2	1:B:350:TYR:CD2	2.48	0.43
1:B:74:ALA:HB3	1:B:237:ILE:HG12	1.94	0.43
1:B:379:TRP:N	1:B:379:TRP:HE3	2.17	0.43
1:B:439:VAL:HG11	1:B:464:VAL:HG21	2.01	0.43
1:B:176:LEU:N	1:B:176:LEU:HD12	2.33	0.42
1:B:66:ASN:CA	1:B:69:ARG:HG3	2.47	0.42
1:B:83:LEU:HD13	1:B:83:LEU:HA	1.77	0.42
1:B:252:THR:OG1	1:B:285:THR:HG23	2.18	0.42
1:B:105:GLY:O	1:B:134:TRP:CZ3	2.72	0.42
1:B:214:PHE:HB3	1:B:218:VAL:CG1	2.50	0.42
1:B:217:THR:HG22	1:B:217:THR:O	2.19	0.42
1:B:327:LEU:HD22	1:B:328:PRO:HD3	1.99	0.42
1:B:379:TRP:CG	1:B:421:LEU:HD13	2.53	0.42
1:B:150:TYR:HB2	1:B:195:PHE:CA	2.29	0.42
1:B:398:ASP:OD1	1:B:398:ASP:N	2.52	0.42
1:B:460:LEU:HB3	1:B:461:TYR:CD1	2.55	0.42
1:B:69:ARG:HH21	1:B:328:PRO:HG2	1.84	0.42
1:B:76:LEU:HG	1:B:87:PRO:HG2	2.01	0.42
1:B:78:VAL:O	1:B:232:ARG:HA	2.19	0.42
1:B:95:ALA:CB	1:B:96:PRO:HD2	2.27	0.42
1:B:147:ARG:CD	1:B:147:ARG:N	2.81	0.42
1:B:219:THR:OG1	1:B:222:ASP:HB2	2.19	0.42
1:B:362:ASP:HB3	1:B:413:PRO:HB3	2.00	0.42
1:B:17:LYS:HD2	1:B:17:LYS:HA	1.82	0.42
1:B:57:THR:CB	1:B:60:ILE:HD12	2.49	0.42
1:B:333:THR:OG1	1:B:343:THR:CG2	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:403:ASP:C	1:B:404:LEU:HG	2.39	0.42
1:B:124:ARG:O	1:B:128:ASN:HB2	2.19	0.42
1:B:199:TYR:CE1	1:B:243:GLY:HA2	2.55	0.42
1:B:376:TYR:HB2	1:B:424:LYS:CB	2.36	0.42
1:B:43:PHE:HZ	1:B:447:LEU:HD22	1.84	0.42
1:B:85:ASN:ND2	1:B:345:GLN:CB	2.68	0.42
1:B:70:LEU:N	1:B:70:LEU:CD2	2.82	0.42
1:B:236:TYR:O	1:B:236:TYR:CD1	2.73	0.42
1:B:375:TYR:CE1	1:B:425:ILE:HG23	2.55	0.42
1:B:309:MET:SD	1:B:309:MET:C	2.98	0.41
1:B:148:MET:HE3	1:B:148:MET:HB2	1.89	0.41
1:B:106:LEU:HD12	1:B:106:LEU:H	1.85	0.41
1:B:161:LEU:CB	1:B:184:HIS:CB	2.80	0.41
1:B:195:PHE:CD1	1:B:196:LYS:N	2.87	0.41
1:B:259:VAL:H	1:B:278:TRP:HB3	1.86	0.41
1:B:363:LEU:O	1:B:411:SER:CA	2.68	0.41
1:B:233:PRO:HB3	1:B:337:ARG:HH12	1.70	0.41
1:B:327:LEU:HD22	1:B:327:LEU:HA	1.80	0.41
1:B:7:ASN:HA	1:B:210:PRO:HG2	2.02	0.41
1:B:375:TYR:CD1	1:B:375:TYR:N	2.88	0.41
1:B:78:VAL:HA	1:B:82:LEU:CD1	2.42	0.41
1:B:215:GLN:O	1:B:218:VAL:CG2	2.69	0.41
1:B:221:GLU:H	1:B:221:GLU:HG3	1.73	0.41
1:B:171:LYS:HB2	1:B:171:LYS:HE3	1.91	0.41
1:B:218:VAL:HG12	1:B:219:THR:H	1.84	0.41
1:B:207:VAL:CG2	1:B:336:LEU:CD2	2.98	0.41
1:B:367:HIS:HE1	1:B:371:TYR:CE2	2.39	0.41
1:B:454:SER:HB2	1:B:465:GLU:OE2	2.21	0.41
1:B:244:ARG:HA	1:B:293:GLY:CA	2.37	0.40
1:B:346:ASN:HB3	1:B:347:SER:H	1.68	0.40
1:B:415:LYS:NZ	1:B:415:LYS:CB	2.78	0.40
1:B:18:LYS:H	1:B:18:LYS:HG3	1.49	0.40
1:B:99:TYR:CD2	1:B:115:VAL:CG2	2.96	0.40
1:B:302:VAL:CG2	1:B:309:MET:HG2	2.43	0.40
1:B:379:TRP:HB3	1:B:421:LEU:HA	2.03	0.40
1:B:255:LYS:HG2	1:B:256:SER:N	2.35	0.40
1:B:275:GLN:O	1:B:275:GLN:HG3	2.21	0.40
1:B:421:LEU:HD23	1:B:445:LEU:HD13	1.96	0.40
1:B:140:GLN:HB3	1:B:141:VAL:H	1.59	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	B	469/471 (100%)	348 (74%)	98 (21%)	23 (5%)	2 23

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	39	LEU
1	B	40	PRO
1	B	72	PRO
1	B	103	LEU
1	B	118	PRO
1	B	206	ALA
1	B	209	ASN
1	B	274	PRO
1	B	296	PRO
1	B	141	VAL
1	B	87	PRO
1	B	383	SER
1	B	388	GLY
1	B	414	LEU
1	B	415	LYS
1	B	411	SER
1	B	419	ARG
1	B	447	LEU
1	B	54	SER
1	B	382	LEU
1	B	233	PRO
1	B	341	VAL
1	B	418	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	B	413/413 (100%)	269 (65%)	144 (35%)	0 1

All (144) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	6	VAL
1	B	7	ASN
1	B	11	LEU
1	B	13	MET
1	B	15	TYR
1	B	16	ASP
1	B	17	LYS
1	B	18	LYS
1	B	19	LYS
1	B	20	LEU
1	B	30	ASN
1	B	31	ARG
1	B	34	LYS
1	B	35	GLU
1	B	38	GLN
1	B	42	GLU
1	B	43	PHE
1	B	44	VAL
1	B	45	VAL
1	B	46	ILE
1	B	51	ARG
1	B	52	SER
1	B	53	LEU
1	B	56	ASN
1	B	67	ASP
1	B	69	ARG
1	B	70	LEU
1	B	75	LEU
1	B	83	LEU
1	B	86	ASN

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Mol	Chain	Res	Type
1	B	89	LEU
1	B	90	LEU
1	B	92	VAL
1	B	94	ARG
1	B	97	MET
1	B	102	ASP
1	B	106	LEU
1	B	109	SER
1	B	110	ASP
1	B	114	GLN
1	B	119	SER
1	B	122	SER
1	B	123	VAL
1	B	129	ASP
1	B	130	LEU
1	B	133	LYS
1	B	136	GLN
1	B	140	GLN
1	B	144	VAL
1	B	147	ARG
1	B	149	GLN
1	B	150	TYR
1	B	151	GLU
1	B	153	ILE
1	B	156	HIS
1	B	159	GLU
1	B	160	GLN
1	B	161	LEU
1	B	169	PHE
1	B	175	SER
1	B	177	ASP
1	B	180	PHE
1	B	183	VAL
1	B	184	HIS
1	B	187	GLU
1	B	189	GLN
1	B	191	GLN
1	B	194	ASN
1	B	195	PHE
1	B	197	GLN
1	B	199	TYR
1	B	204	VAL

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Mol	Chain	Res	Type
1	B	205	ASP
1	B	208	LYS
1	B	220	VAL
1	B	221	GLU
1	B	223	LEU
1	B	226	ARG
1	B	232	ARG
1	B	234	LEU
1	B	236	TYR
1	B	237	ILE
1	B	238	SER
1	B	244	ARG
1	B	249	LYS
1	B	250	LEU
1	B	254	SER
1	B	257	ASP
1	B	268	LEU
1	B	270	VAL
1	B	272	VAL
1	B	278	TRP
1	B	281	ILE
1	B	282	LEU
1	B	284	ASN
1	B	286	GLU
1	B	288	LYS
1	B	291	ILE
1	B	297	SER
1	B	301	ARG
1	B	307	VAL
1	B	310	VAL
1	B	313	LEU
1	B	314	ILE
1	B	316	GLU
1	B	318	SER
1	B	319	ARG
1	B	323	ASP
1	B	327	LEU
1	B	330	SER
1	B	334	SER
1	B	337	ARG
1	B	340	VAL
1	B	343	THR

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Mol	Chain	Res	Type
1	B	344	PHE
1	B	345	GLN
1	B	346	ASN
1	B	359	ARG
1	B	364	LEU
1	B	372	VAL
1	B	380	ASP
1	B	383	SER
1	B	386	HIS
1	B	391	VAL
1	B	392	LEU
1	B	395	LYS
1	B	398	ASP
1	B	399	ARG
1	B	403	ASP
1	B	404	LEU
1	B	415	LYS
1	B	419	ARG
1	B	422	SER
1	B	427	GLU
1	B	434	GLU
1	B	437	ARG
1	B	438	THR
1	B	445	LEU
1	B	447	LEU
1	B	456	TRP
1	B	460	LEU
1	B	465	GLU
1	B	468	VAL
1	B	471	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	23	HIS
1	B	37	ASN
1	B	38	GLN
1	B	85	ASN
1	B	135	HIS
1	B	140	GLN
1	B	156	HIS
1	B	245	GLN

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Mol	Chain	Res	Type
1	B	275	GLN
1	B	284	ASN
1	B	367	HIS
1	B	407	HIS
1	B	463	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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.

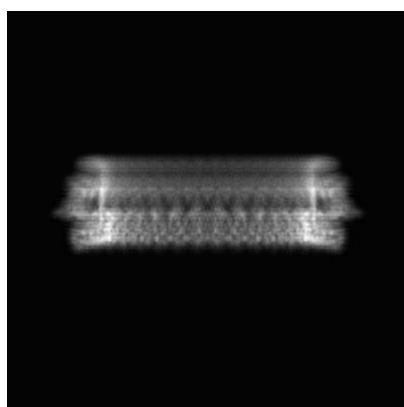
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-4118. These allow visual inspection of the internal detail of the map and identification of artifacts.

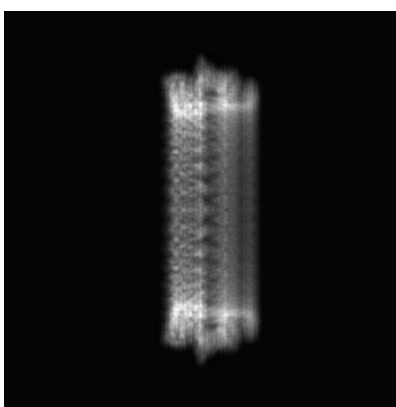
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

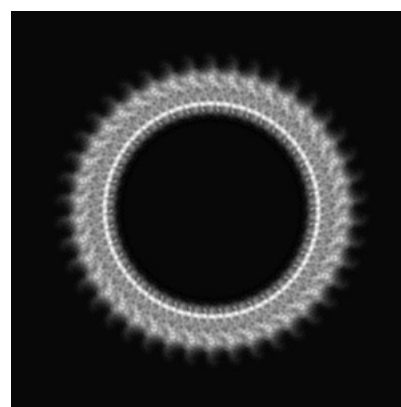
6.1.1 Primary map



X



Y

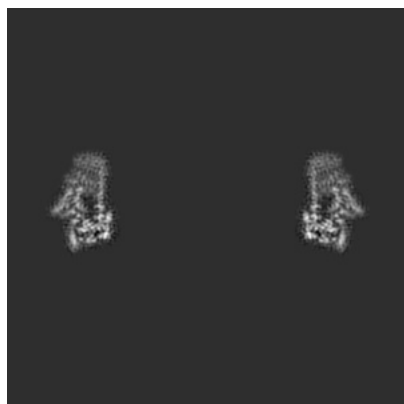


Z

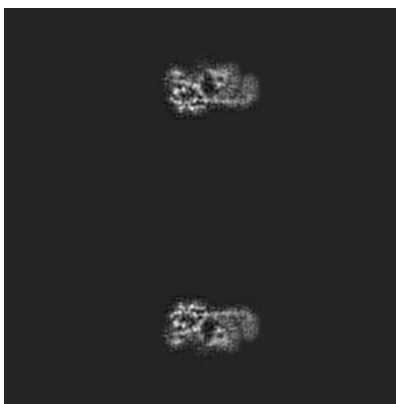
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

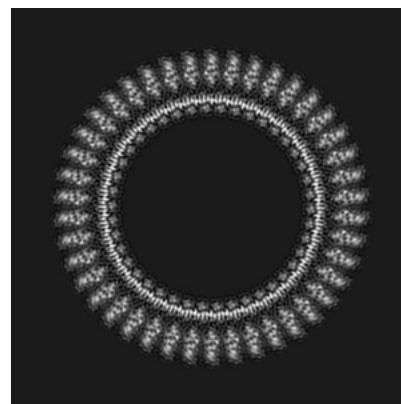
6.2.1 Primary map



X Index: 180



Y Index: 180

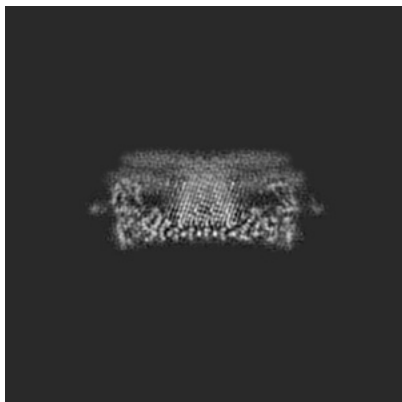


Z Index: 180

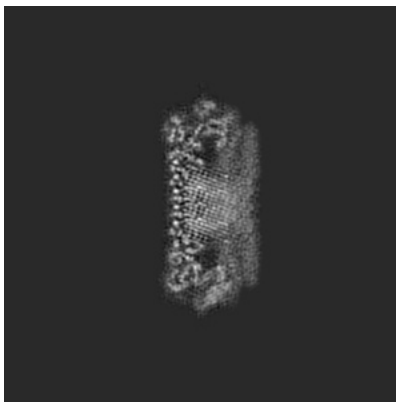
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

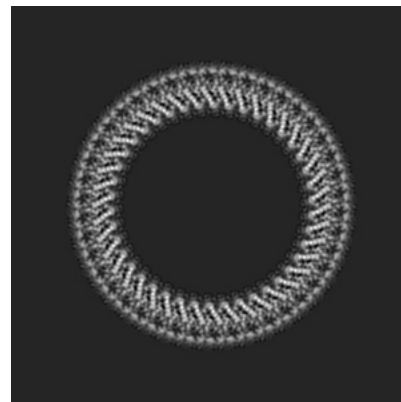
6.3.1 Primary map



X Index: 275



Y Index: 275

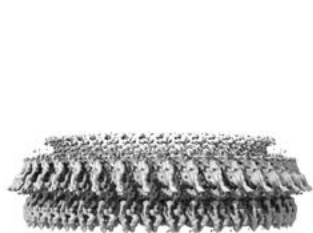


Z Index: 155

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

6.4 Orthogonal surface views [i](#)

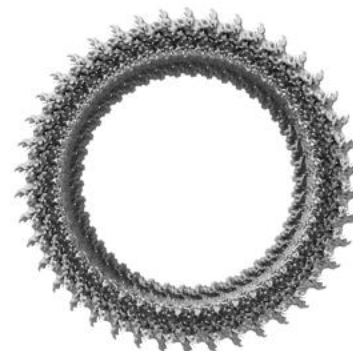
6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0282. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

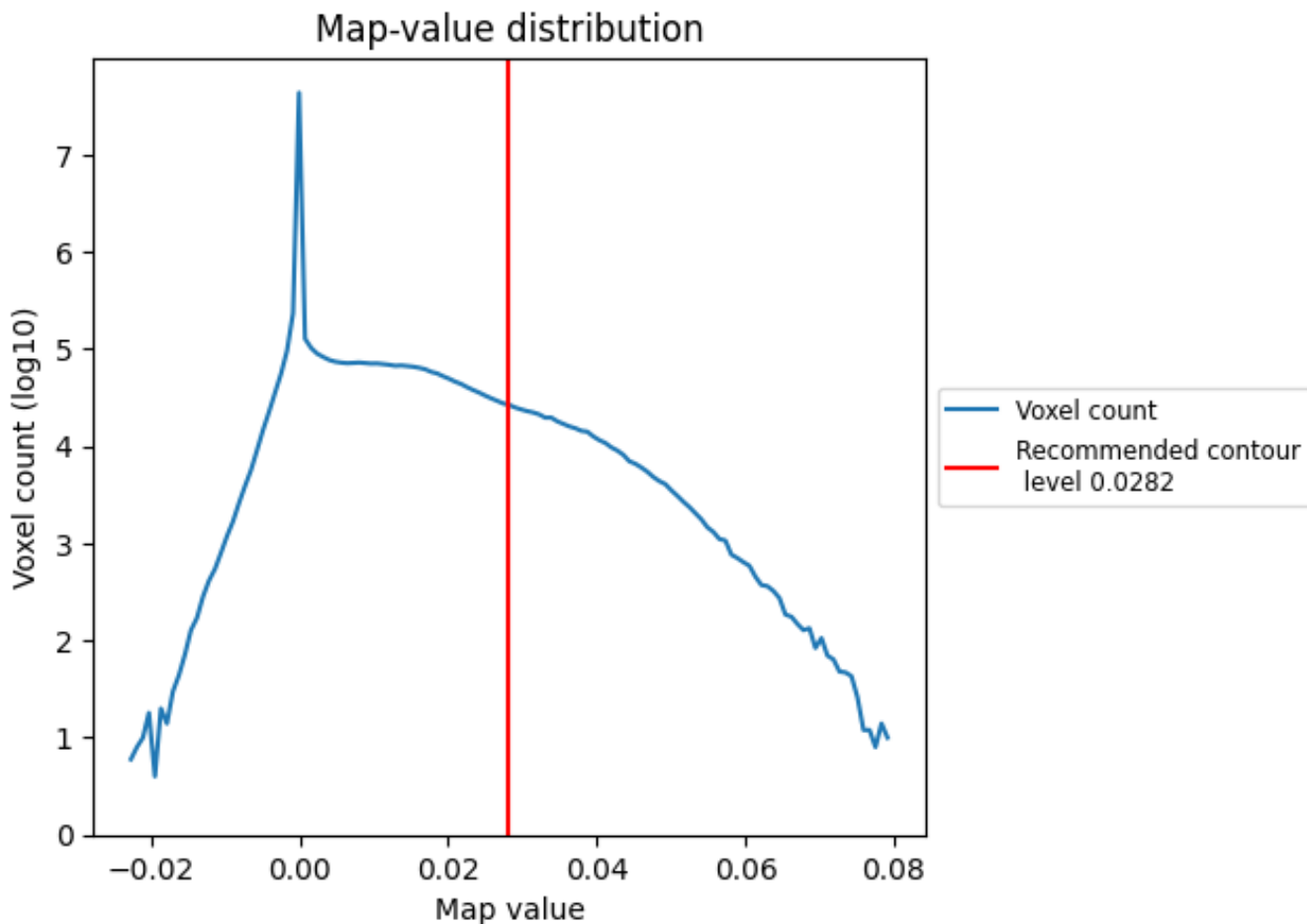
6.5 Mask visualisation

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

7 Map analysis [i](#)

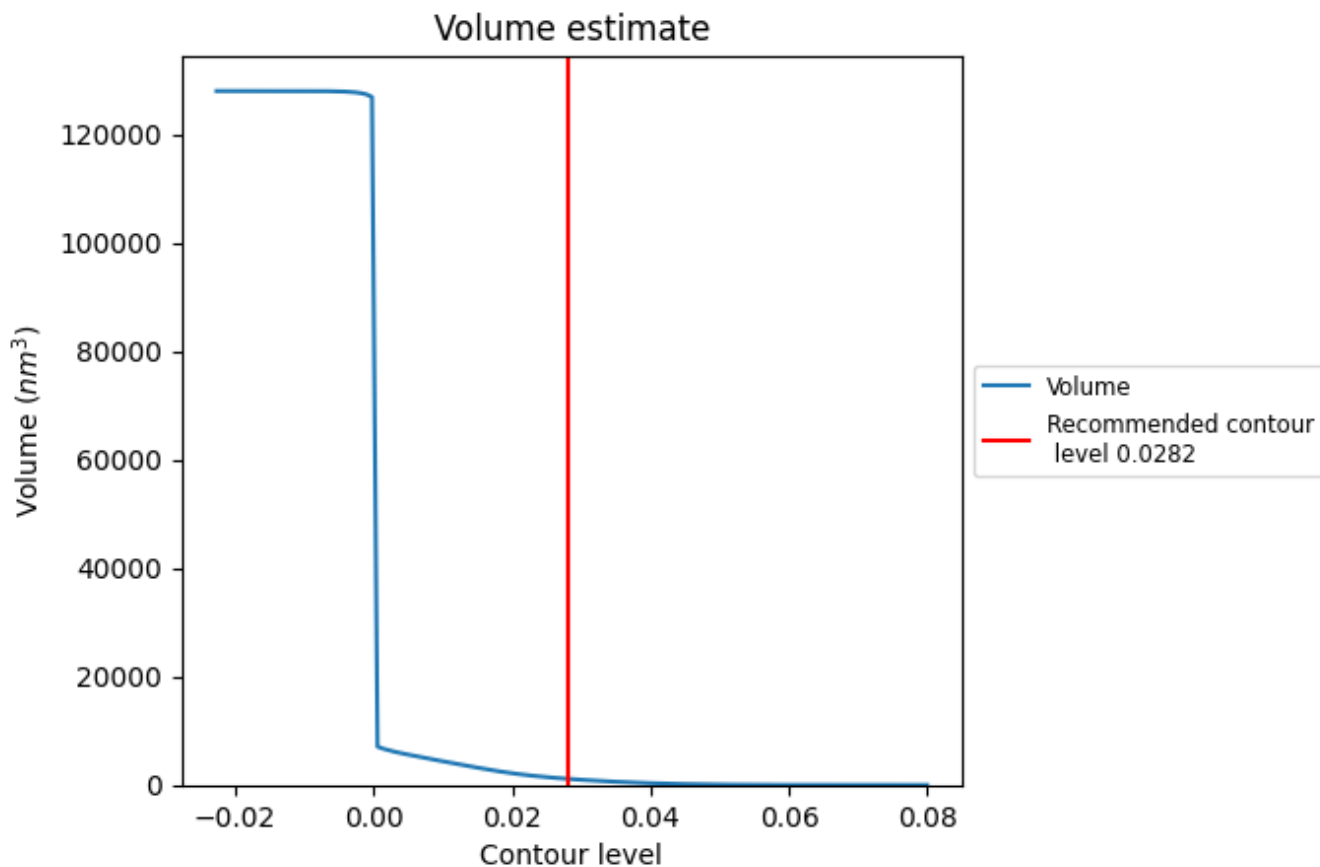
This section contains the results of statistical analysis of the map.

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.

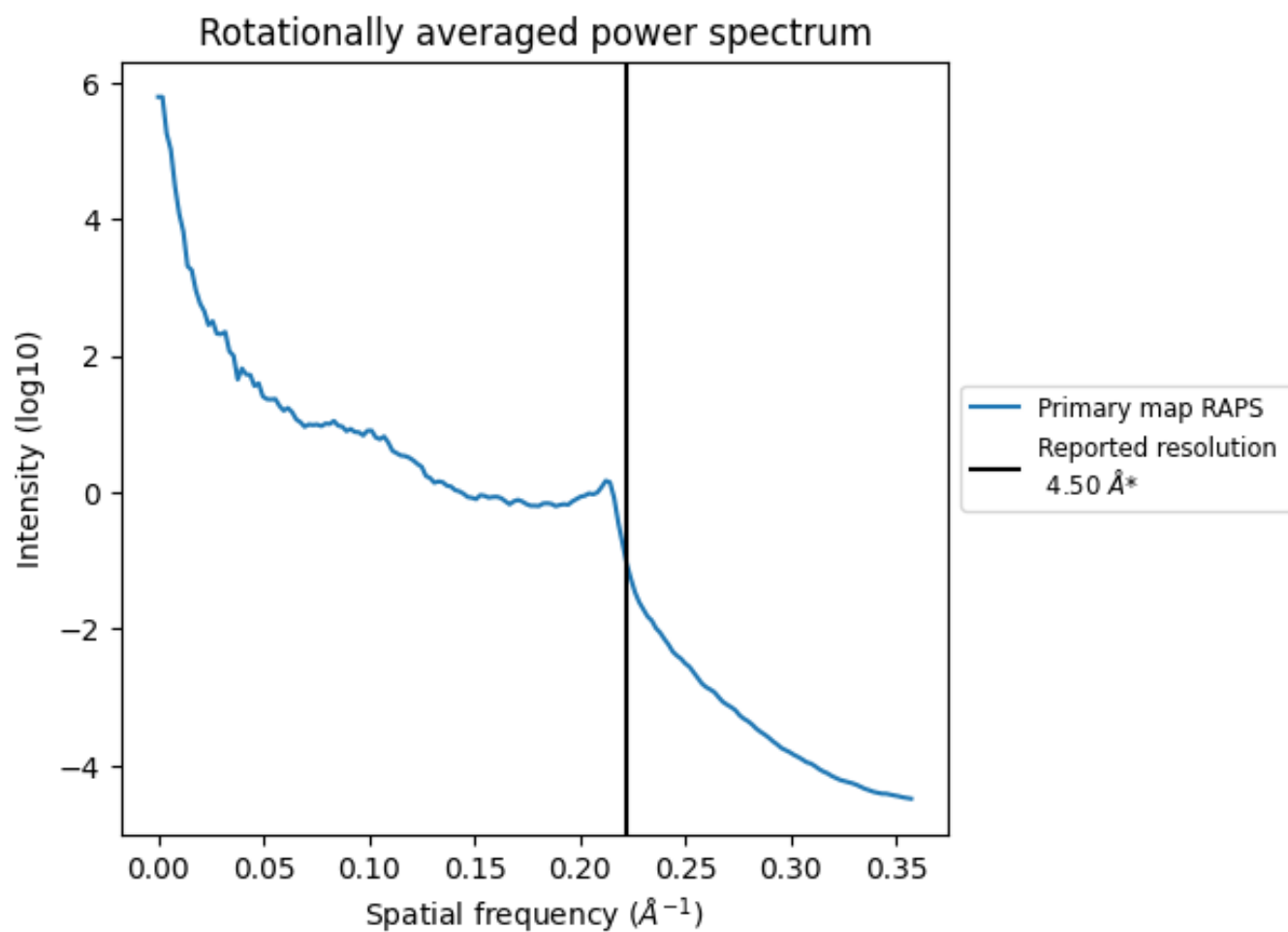
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 1109 nm³; this corresponds to an approximate mass of 1002 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 [i](#)

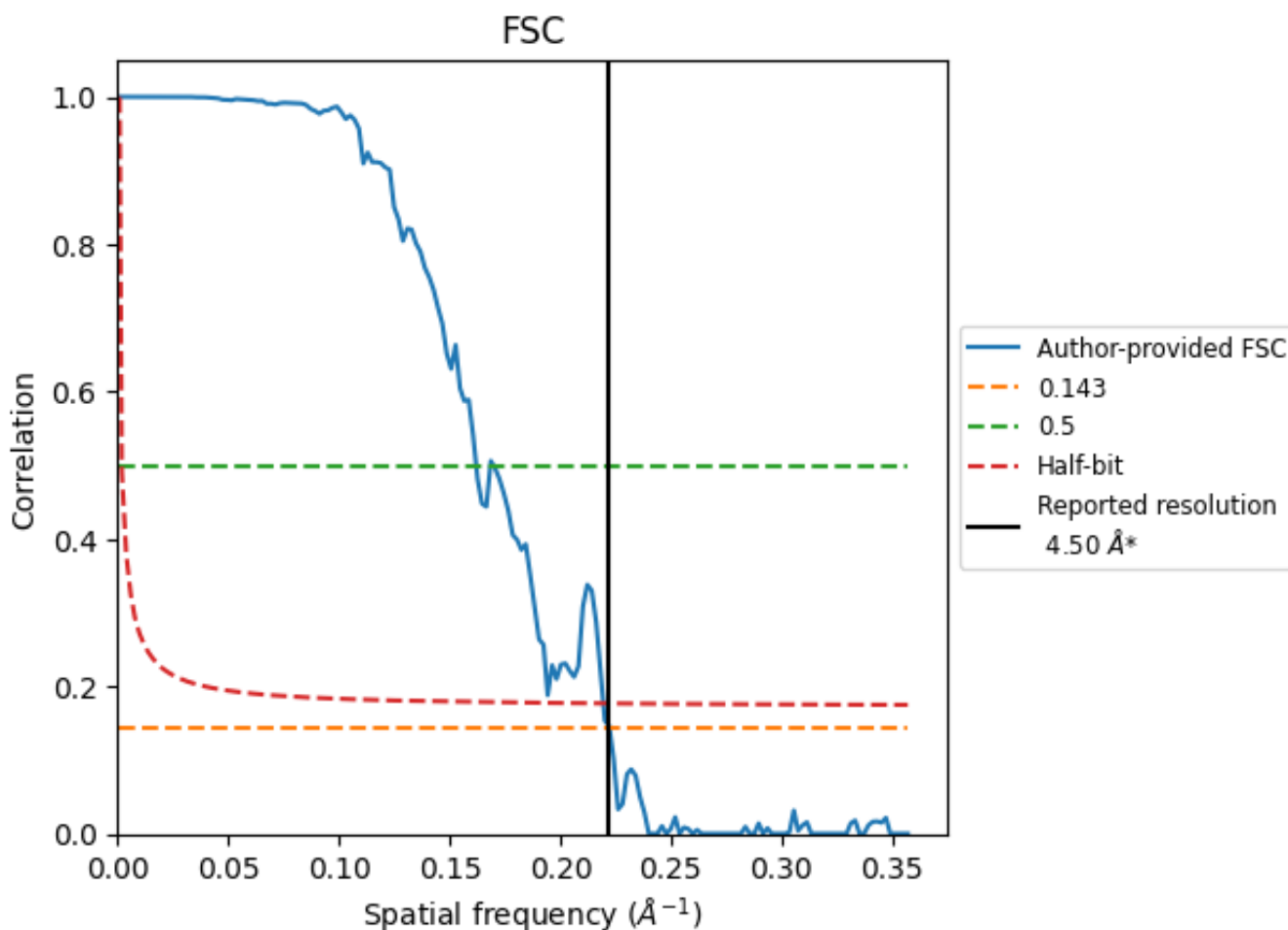


*Reported resolution corresponds to spatial frequency of 0.222\AA^{-1}

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 0.222 Å⁻¹

8.2 Resolution estimates [i](#)

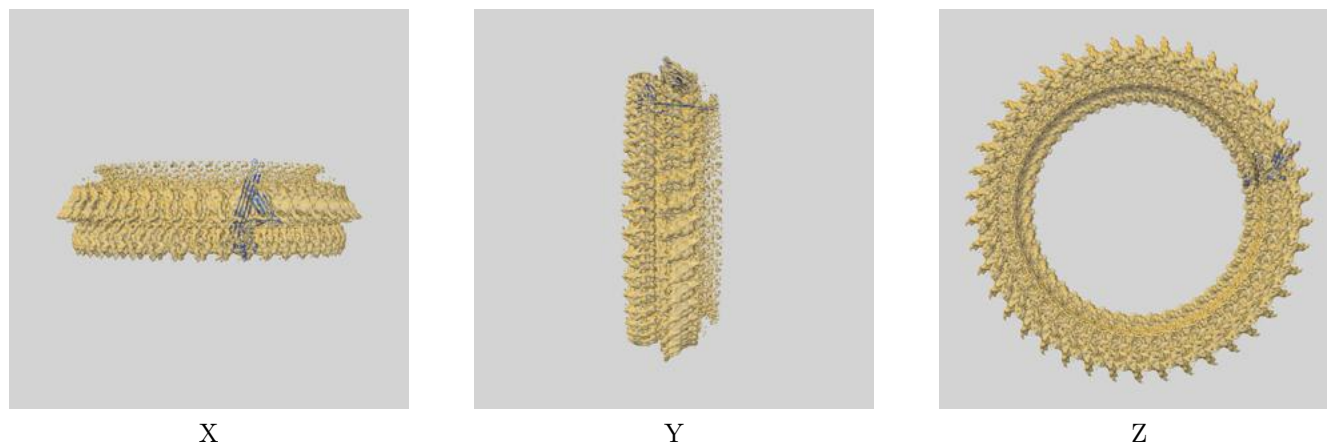
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.50	-	-
Author-provided FSC curve	4.50	6.17	4.56
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

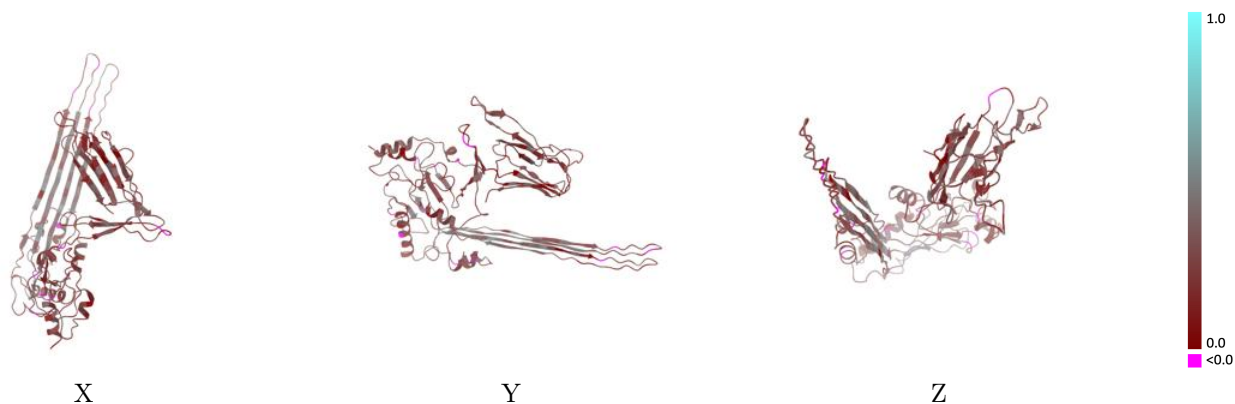
This section contains information regarding the fit between EMDB map EMD-4118 and PDB model 5LY6. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)



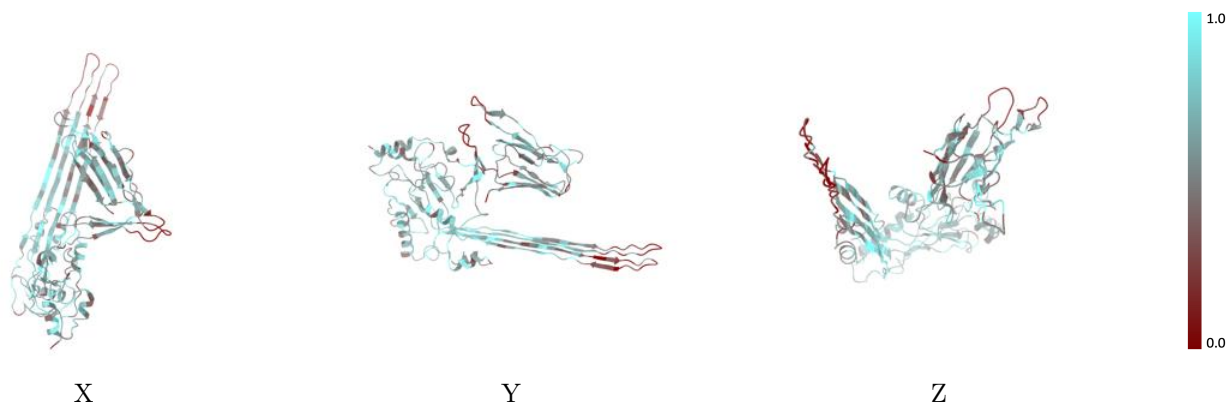
The images above show the 3D surface view of the map at the recommended contour level 0.0282 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



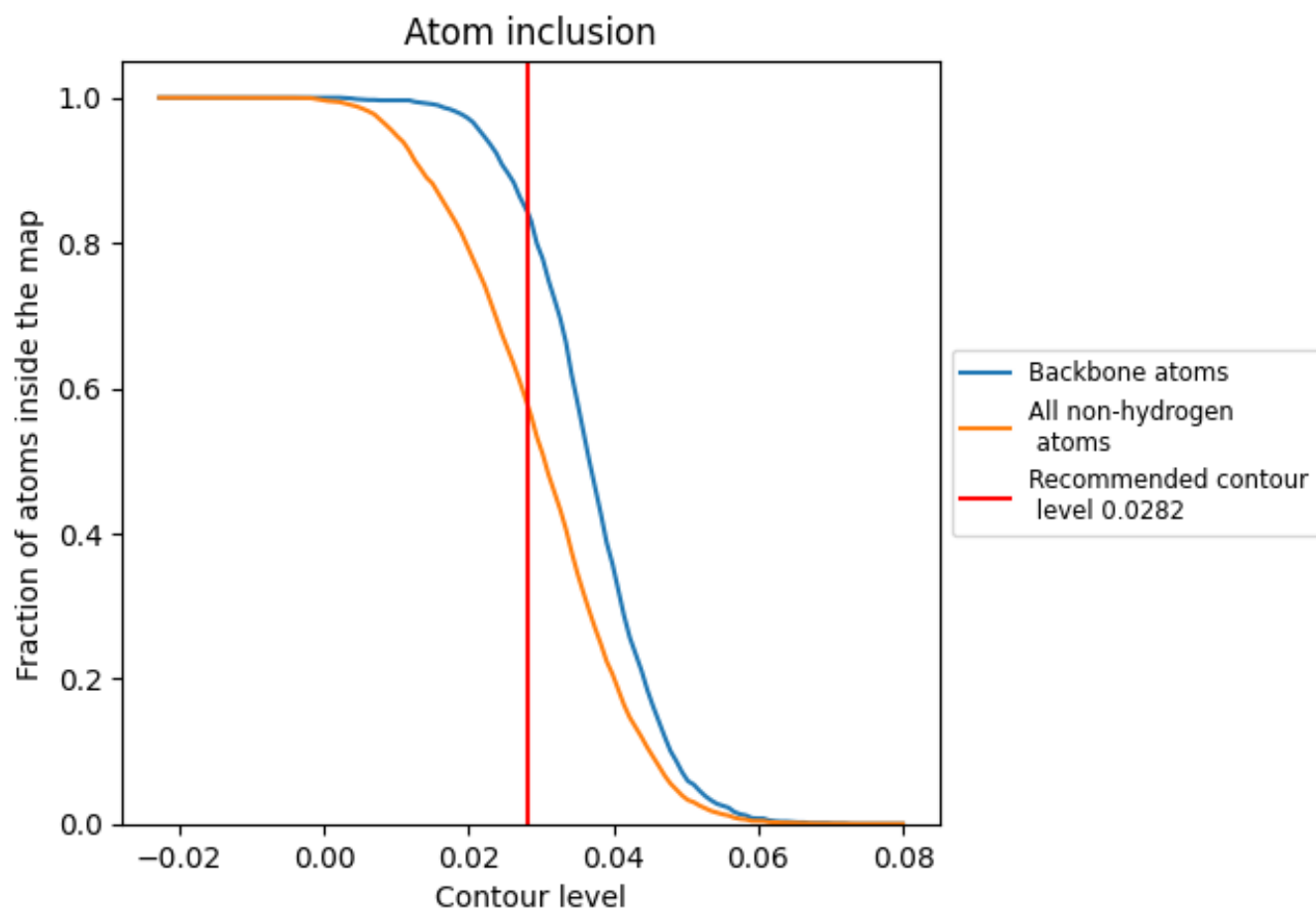
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0282).





9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 58% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.0282) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.5752	 0.2590
B	 0.5752	 0.2590

