



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 11:44 AM UTC

PDB ID : 3J91 / pdb_00003j91
EMDB ID : EMD-6200
Title : Cryo-electron microscopy of Enterovirus 71 (EV71) procapsid in complex with Fab fragments of neutralizing antibody 22A12
Authors : Shingler, K.L.; Cifuentes, J.O.; Ashley, R.E.; Makhov, A.M.; Conway, J.F.; Hafenstein, S.
Deposited on : 2014-11-24
Resolution : 8.80 Å (reported)
Based on initial model : 4GMP

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:

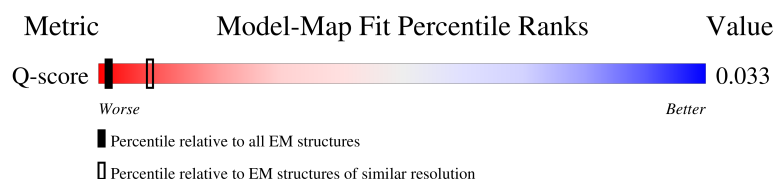
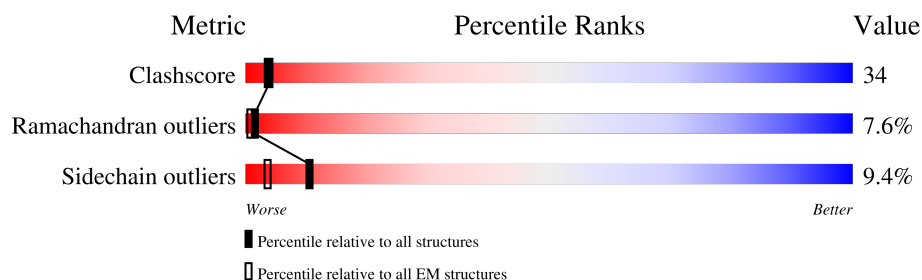
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMD archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (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 8.80 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	265 (8.30 - 9.30)

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	0	323	
2	1	297	
3	3	242	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5395 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 VP0.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	237	Total	C	N	O	S	0	0
			1833	1179	301	345	8		

- Molecule 2 is a protein called VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	218	Total	C	N	O	S	0	0
			1717	1101	289	316	11		

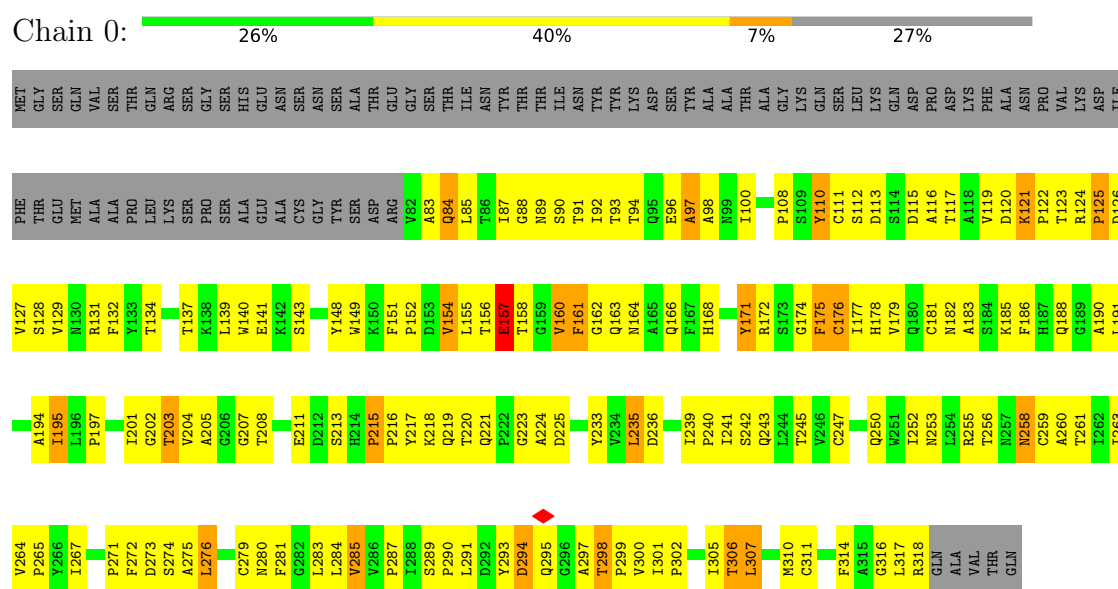
- Molecule 3 is a protein called VP3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	240	Total	C	N	O	S	0	0
			1845	1188	304	342	11		

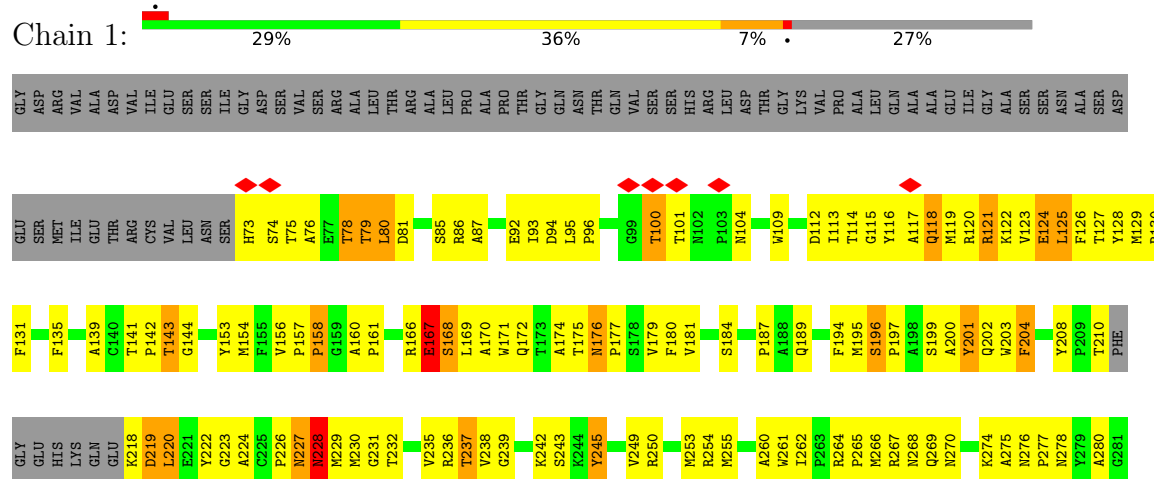
3 Residue-property plots

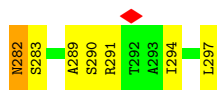
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: VP0

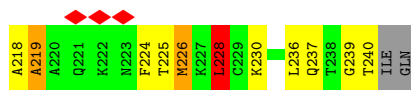
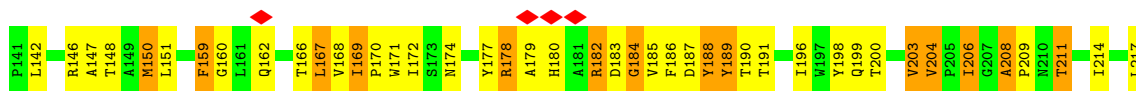
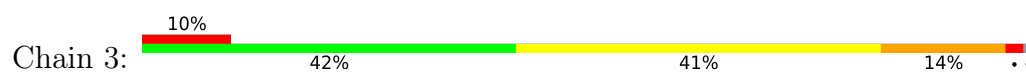


• Molecule 2: VP1





• Molecule 3: VP3



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	15226	Depositor
Resolution determination method	FSC 0.5 CUT-OFF	Depositor
CTF correction method	CTFFind3	Depositor
Microscope	FEI TECNAI F20	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	50000	Depositor
Image detector	KODAK SO-163 FILM	Depositor
Maximum map value	8.483	Depositor
Minimum map value	-4.293	Depositor
Average map value	0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	1	Depositor
Map size ($\text{\AA}$)	585.47, 585.47, 585.47	wwPDB
Map dimensions	461, 461, 461	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.27, 1.27, 1.27	Depositor

5 Model quality

5.1 Standard geometry

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	0	0.57	0/1888	1.05	11/2591 (0.4%)
2	1	0.61	0/1769	1.03	7/2411 (0.3%)
3	3	0.59	0/1897	1.11	23/2596 (0.9%)
All	All	0.59	0/5554	1.06	41/7598 (0.5%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	0	161	PHE	N-CA-C	-8.59	102.24	112.89
3	3	45	GLU	N-CA-C	-8.27	102.22	111.07
3	3	18	ASP	N-CA-C	7.63	120.64	110.88
1	0	294	ASP	N-CA-C	-7.24	99.74	110.23
3	3	74	SER	N-CA-C	7.13	118.00	108.38
2	1	201	TYR	N-CA-C	-7.09	101.69	110.41
2	1	168	SER	N-CA-C	-6.95	100.15	110.23
3	3	2	PHE	CA-C-N	6.69	128.21	119.84
3	3	2	PHE	C-N-CA	6.69	128.21	119.84
3	3	169	ILE	CA-C-N	6.64	128.15	119.84
3	3	169	ILE	C-N-CA	6.64	128.15	119.84
2	1	125	LEU	N-CA-C	-6.39	104.97	112.89
3	3	31	THR	CA-C-N	6.30	126.31	119.89
3	3	31	THR	C-N-CA	6.30	126.31	119.89
3	3	20	VAL	N-CA-C	6.26	116.16	108.53
1	0	84	GLN	N-CA-C	6.21	118.30	108.67
1	0	126	ASP	N-CA-C	-6.19	105.67	113.72
1	0	221	GLN	CA-C-N	6.16	127.55	119.84
1	0	221	GLN	C-N-CA	6.16	127.55	119.84
1	0	175	PHE	N-CA-C	6.07	116.02	108.19
3	3	111	TRP	N-CA-C	6.04	116.39	107.88
1	0	160	VAL	N-CA-C	-6.02	104.89	113.07
2	1	176	ASN	CA-C-N	5.99	127.33	119.84
2	1	176	ASN	C-N-CA	5.99	127.33	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	3	160	GLY	N-CA-C	-5.94	105.60	112.73
1	0	217	TYR	N-CA-C	-5.90	106.13	113.15
2	1	115	GLY	N-CA-C	-5.67	106.73	113.99
3	3	9	GLY	N-CA-C	-5.63	107.24	115.63
1	0	116	ALA	N-CA-C	-5.47	106.64	113.15
3	3	203	VAL	N-CA-C	5.47	115.73	107.75
3	3	52	ILE	N-CA-C	5.45	116.03	108.84
1	0	194	ALA	N-CA-C	5.33	117.17	109.07
2	1	144	GLY	N-CA-C	-5.24	107.82	115.63
3	3	125	PHE	N-CA-C	-5.20	105.23	112.45
3	3	208	ALA	CA-C-N	5.06	125.85	119.98
3	3	208	ALA	C-N-CA	5.06	125.85	119.98
3	3	100	LEU	N-CA-C	-5.04	105.86	111.36
3	3	29	TYR	CA-C-N	5.04	126.14	119.84
3	3	29	TYR	C-N-CA	5.04	126.14	119.84
3	3	19	GLY	CA-C-N	-5.01	116.59	123.46
3	3	19	GLY	C-N-CA	-5.01	116.59	123.46

There are no chirality outliers.

There are no planarity outliers.

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	0	1833	0	1774	132	0
2	1	1717	0	1674	135	0
3	3	1845	0	1822	137	0
All	All	5395	0	5270	359	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 34.

All (359) 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)
3:3:20:VAL:HG22	3:3:21:SER:H	1.17	1.04
1:0:233:TYR:HA	3:3:66:MET:HE3	1.43	1.01
3:3:206:ILE:HD12	3:3:206:ILE:H	1.24	1.00
2:1:229:MET:HE2	2:1:231:GLY:H	1.27	0.98
3:3:58:VAL:HG23	3:3:59:PRO:HD3	1.50	0.93
3:3:109:THR:HB	3:3:228:LEU:HB3	1.52	0.91
2:1:76:ALA:O	2:1:79:THR:HG23	1.73	0.89
3:3:188:TYR:O	3:3:189:TYR:HB3	1.75	0.86
3:3:172:ILE:HD12	3:3:172:ILE:O	1.75	0.86
3:3:167:LEU:HD12	3:3:168:VAL:N	1.91	0.85
2:1:112:ASP:OD2	2:1:114:THR:HG22	1.75	0.85
2:1:218:LYS:HD3	2:1:219:ASP:H	1.43	0.84
2:1:141:THR:OG1	2:1:143:THR:HG23	1.78	0.82
3:3:20:VAL:HG22	3:3:21:SER:N	1.95	0.82
2:1:121:ARG:HG2	2:1:121:ARG:HH11	1.43	0.81
1:0:241:ILE:HG21	3:3:66:MET:HE1	1.60	0.81
1:0:117:THR:HB	1:0:120:ASP:HB3	1.63	0.81
2:1:218:LYS:HD3	2:1:219:ASP:N	1.95	0.81
3:3:58:VAL:HG23	3:3:59:PRO:CD	2.10	0.81
3:3:178:ARG:NH1	3:3:187:ASP:HB3	1.97	0.80
3:3:85:VAL:HG21	3:3:142:LEU:HD12	1.62	0.79
2:1:139:ALA:HB2	2:1:249:VAL:HG22	1.65	0.79
3:3:179:ALA:HA	3:3:184:GLY:HA2	1.66	0.78
2:1:112:ASP:CG	2:1:114:THR:HG22	2.09	0.77
2:1:197:PRO:HD2	2:1:227:ASN:HB2	1.65	0.77
2:1:276:ASN:HB2	2:1:277:PRO:HD2	1.65	0.76
3:3:20:VAL:CG2	3:3:21:SER:H	1.98	0.76
1:0:233:TYR:CA	3:3:66:MET:HE3	2.15	0.76
1:0:188:GLN:NE2	3:3:209:PRO:HB2	2.01	0.75
3:3:84:ALA:HB3	3:3:196:ILE:HD11	1.66	0.75
2:1:100:THR:HG23	2:1:101:THR:H	1.52	0.74
2:1:121:ARG:HD2	2:1:267:ARG:HB3	1.70	0.74
3:3:73:VAL:HA	3:3:198:TYR:OH	1.87	0.73
1:0:139:LEU:HD22	1:0:300:VAL:HB	1.69	0.73
3:3:42:ASN:OD1	3:3:44:LEU:HB2	1.89	0.73
1:0:125:PRO:HB3	1:0:129:VAL:HG21	1.71	0.73
3:3:131:MET:HG2	3:3:159:PHE:HE1	1.53	0.73
2:1:177:PRO:HB2	3:3:24:ILE:HD11	1.70	0.72
2:1:229:MET:HE3	2:1:230:MET:H	1.55	0.72
3:3:178:ARG:HH11	3:3:178:ARG:HB2	1.54	0.72
1:0:128:SER:OG	1:0:160:VAL:HG11	1.91	0.71
1:0:201:ILE:HG22	1:0:202:GLY:N	2.05	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:141:THR:HB	2:1:142:PRO:HD2	1.73	0.70
2:1:229:MET:HE2	2:1:231:GLY:N	2.04	0.70
3:3:239:GLY:O	3:3:240:THR:C	2.33	0.70
3:3:83:CYS:HB3	3:3:196:ILE:HG13	1.73	0.70
2:1:168:SER:C	2:1:170:ALA:H	1.99	0.70
2:1:158:PRO:HD3	2:1:231:GLY:HA2	1.76	0.67
2:1:218:LYS:O	2:1:219:ASP:HB2	1.93	0.67
2:1:229:MET:HE3	2:1:230:MET:N	2.10	0.67
2:1:194:PHE:CZ	2:1:200:ALA:HA	2.30	0.66
2:1:121:ARG:HG2	2:1:121:ARG:NH1	2.08	0.66
2:1:153:TYR:CD1	2:1:235:VAL:HG13	2.31	0.66
3:3:6:LEU:HD12	3:3:10:THR:HG21	1.77	0.65
3:3:204:VAL:HG13	3:3:208:ALA:HB3	1.79	0.65
1:0:179:VAL:HG22	1:0:305:ILE:HG12	1.79	0.64
2:1:78:THR:HG22	3:3:43:LEU:HB2	1.79	0.64
1:0:203:THR:H	1:0:215:PRO:HG3	1.62	0.64
2:1:85:SER:O	2:1:86:ARG:HB2	1.97	0.64
3:3:50:GLU:HA	3:3:219:ALA:HB2	1.80	0.64
2:1:153:TYR:HD1	2:1:235:VAL:HG13	1.64	0.63
2:1:168:SER:O	2:1:170:ALA:N	2.31	0.63
2:1:156:VAL:HB	2:1:232:THR:HB	1.81	0.63
2:1:104:ASN:HA	2:1:242:LYS:NZ	2.14	0.62
1:0:171:TYR:CG	1:0:275:ALA:HB2	2.34	0.62
2:1:112:ASP:OD1	2:1:114:THR:HG22	1.99	0.62
1:0:88:GLY:C	1:0:90:SER:H	2.08	0.62
3:3:178:ARG:HH11	3:3:187:ASP:HB3	1.64	0.61
2:1:189:GLN:HG3	3:3:21:SER:HB3	1.82	0.61
3:3:225:THR:HG22	3:3:226:MET:N	2.15	0.61
2:1:128:TYR:HB2	2:1:261:TRP:HB2	1.82	0.61
2:1:130:ARG:NE	3:3:33:CYS:HB3	2.16	0.61
1:0:129:VAL:HG13	1:0:161:PHE:HD1	1.65	0.60
3:3:225:THR:O	3:3:226:MET:HG2	2.00	0.60
3:3:115:LEU:HD11	3:3:172:ILE:HG12	1.82	0.60
2:1:189:GLN:HG3	3:3:21:SER:CB	2.32	0.60
3:3:2:PHE:CD1	3:3:3:PRO:HD2	2.36	0.60
1:0:120:ASP:O	1:0:122:PRO:HD3	2.03	0.59
1:0:168:HIS:CD2	1:0:314:PHE:HB3	2.38	0.59
3:3:10:THR:HG22	3:3:11:ASN:ND2	2.16	0.59
1:0:177:ILE:HG21	1:0:283:LEU:CD1	2.33	0.59
2:1:139:ALA:CB	2:1:249:VAL:HG22	2.33	0.58
1:0:171:TYR:CB	1:0:275:ALA:HB2	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:148:TYR:HB3	1:0:284:LEU:HD23	1.85	0.58
2:1:158:PRO:HD3	2:1:231:GLY:CA	2.33	0.58
2:1:168:SER:C	2:1:170:ALA:N	2.62	0.58
3:3:178:ARG:HG2	3:3:179:ALA:H	1.68	0.58
1:0:178:HIS:HE1	1:0:259:CYS:HB2	1.69	0.58
1:0:195:ILE:HG13	1:0:281:PHE:CE2	2.39	0.57
3:3:14:LEU:HB3	3:3:17:ASP:HB2	1.86	0.57
2:1:254:ARG:HB3	2:1:254:ARG:HH11	1.70	0.57
3:3:131:MET:HG2	3:3:159:PHE:CE1	2.37	0.57
1:0:92:ILE:HG13	1:0:92:ILE:O	2.04	0.57
2:1:93:ILE:HD11	2:1:109:TRP:HB2	1.86	0.57
1:0:171:TYR:HB3	1:0:275:ALA:HB2	1.85	0.57
2:1:135:PHE:CD1	2:1:253:MET:HB2	2.40	0.57
1:0:201:ILE:CG2	1:0:202:GLY:N	2.68	0.57
2:1:280:ALA:CB	2:1:283:SER:HB3	2.36	0.56
3:3:150:MET:HE2	3:3:150:MET:O	2.05	0.56
2:1:154:MET:HE1	2:1:171:TRP:HA	1.87	0.56
2:1:254:ARG:HB3	2:1:254:ARG:NH1	2.20	0.56
3:3:52:ILE:HA	3:3:217:LEU:HD23	1.87	0.56
3:3:65:LEU:O	3:3:68:ARG:HG3	2.06	0.56
1:0:119:VAL:HG12	1:0:119:VAL:O	2.06	0.55
1:0:90:SER:OG	1:0:132:PHE:HB2	2.06	0.55
3:3:206:ILE:H	3:3:206:ILE:CD1	1.98	0.55
1:0:115:ASP:C	1:0:117:THR:H	2.14	0.55
1:0:188:GLN:HB3	3:3:124:SER:HA	1.89	0.55
2:1:174:ALA:C	2:1:176:ASN:H	2.14	0.55
1:0:155:LEU:C	1:0:157:GLU:N	2.62	0.55
1:0:235:LEU:HD12	1:0:239:ILE:CD1	2.37	0.55
1:0:162:GLY:O	1:0:166:GLN:HG3	2.07	0.55
1:0:250:GLN:HG2	1:0:260:ALA:HB1	1.89	0.55
2:1:120:ARG:HH11	3:3:237:GLN:HE22	1.53	0.55
3:3:6:LEU:CD1	3:3:10:THR:HG21	2.37	0.55
3:3:30:PRO:O	3:3:31:THR:C	2.49	0.55
3:3:50:GLU:HA	3:3:218:ALA:O	2.06	0.55
1:0:137:THR:HG21	1:0:302:PRO:HB2	1.88	0.54
2:1:92:GLU:HG3	2:1:250:ARG:HG2	1.89	0.54
2:1:276:ASN:HB2	2:1:277:PRO:CD	2.34	0.54
3:3:9:GLY:HA2	3:3:12:GLN:OE1	2.08	0.54
1:0:84:GLN:HB3	1:0:93:THR:HA	1.90	0.54
3:3:174:ASN:OD1	3:3:174:ASN:O	2.25	0.54
1:0:100:ILE:HG22	1:0:261:THR:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:125:PRO:HG2	1:0:314:PHE:CD2	2.42	0.54
2:1:93:ILE:HG13	2:1:235:VAL:HG21	1.88	0.54
2:1:94:ASP:O	2:1:95:LEU:HD23	2.07	0.54
1:0:125:PRO:HB3	1:0:129:VAL:CG2	2.37	0.54
1:0:134:THR:HG23	1:0:306:THR:HG23	1.89	0.54
1:0:155:LEU:C	1:0:157:GLU:H	2.14	0.54
1:0:240:PRO:HG3	1:0:243:GLN:NE2	2.22	0.54
2:1:195:MET:O	2:1:196:SER:O	2.25	0.54
2:1:261:TRP:NE1	3:3:39:GLU:HB2	2.23	0.54
1:0:239:ILE:O	1:0:239:ILE:HG13	2.06	0.53
1:0:177:ILE:HG21	1:0:283:LEU:HD11	1.89	0.53
2:1:123:VAL:HG23	2:1:124:GLU:N	2.22	0.53
3:3:91:GLY:HA3	3:3:111:TRP:CZ2	2.43	0.53
3:3:61:ASN:ND2	3:3:64:SER:OG	2.42	0.53
1:0:271:PRO:HG2	1:0:272:PHE:H	1.73	0.53
3:3:18:ASP:O	3:3:19:GLY:O	2.27	0.53
1:0:111:CYS:SG	1:0:115:ASP:OD1	2.67	0.53
2:1:158:PRO:HB3	2:1:229:MET:HG3	1.91	0.53
1:0:197:PRO:HG2	2:1:262:ILE:HG21	1.91	0.53
1:0:208:THR:HA	2:1:282:ASN:OD1	2.08	0.52
1:0:241:ILE:CG2	3:3:66:MET:HE1	2.35	0.52
2:1:208:TYR:HE1	2:1:222:TYR:HB2	1.74	0.52
3:3:56:ASN:O	3:3:68:ARG:HA	2.09	0.52
3:3:225:THR:HG22	3:3:226:MET:H	1.74	0.52
1:0:140:TRP:CE2	1:0:291:LEU:HB2	2.44	0.52
3:3:91:GLY:HA3	3:3:111:TRP:HZ2	1.75	0.52
3:3:79:LYS:HB2	3:3:79:LYS:NZ	2.25	0.52
3:3:74:SER:HB2	3:3:211:THR:OG1	2.10	0.52
3:3:54:GLU:O	3:3:95:PRO:HB3	2.10	0.52
3:3:167:LEU:HD12	3:3:168:VAL:H	1.73	0.52
1:0:183:ALA:HB2	1:0:301:ILE:HG21	1.92	0.52
1:0:201:ILE:HG22	1:0:202:GLY:H	1.75	0.52
2:1:80:LEU:HD21	2:1:260:ALA:HB3	1.92	0.52
2:1:166:ARG:HG2	2:1:166:ARG:HH11	1.74	0.51
2:1:180:PHE:CD1	2:1:180:PHE:N	2.78	0.51
2:1:171:TRP:CZ3	2:1:236:ARG:HG2	2.44	0.51
2:1:194:PHE:HZ	2:1:200:ALA:HA	1.73	0.51
3:3:44:LEU:HD21	3:3:224:PHE:HB3	1.92	0.51
3:3:54:GLU:HG3	3:3:98:SER:CB	2.40	0.51
1:0:123:THR:O	1:0:125:PRO:HD3	2.11	0.51
1:0:188:GLN:HG2	3:3:209:PRO:HG2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:3:167:LEU:HD12	3:3:167:LEU:C	2.36	0.51
1:0:154:VAL:HG23	1:0:224:ALA:HA	1.92	0.51
2:1:135:PHE:CE1	2:1:253:MET:HB2	2.46	0.51
1:0:255:ARG:HG3	1:0:256:THR:HG23	1.93	0.50
2:1:87:ALA:HA	2:1:254:ARG:HB2	1.93	0.50
2:1:204:PHE:CD1	2:1:204:PHE:N	2.77	0.50
2:1:166:ARG:NH2	2:1:237:THR:OG1	2.45	0.50
1:0:182:ASN:O	1:0:301:ILE:HG23	2.11	0.50
2:1:156:VAL:HG22	2:1:176:ASN:ND2	2.26	0.50
1:0:307:LEU:N	1:0:307:LEU:HD23	2.27	0.50
2:1:203:TRP:HB2	2:1:204:PHE:CD1	2.47	0.50
3:3:218:ALA:O	3:3:219:ALA:HB2	2.11	0.50
3:3:128:THR:HG23	3:3:203:VAL:HB	1.93	0.50
1:0:204:VAL:O	1:0:205:ALA:HB3	2.12	0.49
3:3:142:LEU:HD23	3:3:142:LEU:O	2.12	0.49
2:1:73:HIS:HA	3:3:225:THR:HG21	1.94	0.49
2:1:260:ALA:O	3:3:39:GLU:HG3	2.11	0.49
2:1:123:VAL:C	2:1:125:LEU:H	2.20	0.49
1:0:92:ILE:HG12	1:0:132:PHE:CE1	2.47	0.49
2:1:104:ASN:HA	2:1:242:LYS:HZ2	1.77	0.49
2:1:220:LEU:N	2:1:220:LEU:HD23	2.28	0.49
1:0:243:GLN:HA	3:3:51:THR:HG22	1.95	0.49
1:0:265:PRO:O	1:0:267:ILE:HG13	2.11	0.49
2:1:280:ALA:HB1	2:1:283:SER:HB3	1.95	0.49
1:0:293:TYR:CE1	1:0:299:PRO:HA	2.48	0.49
2:1:127:THR:OG1	2:1:264:ARG:NH2	2.46	0.49
1:0:129:VAL:HG22	1:0:164:ASN:ND2	2.29	0.48
1:0:252:ILE:HG22	1:0:252:ILE:O	2.12	0.48
2:1:203:TRP:C	2:1:204:PHE:CD1	2.91	0.48
3:3:130:LYS:HB2	3:3:200:THR:HG23	1.95	0.48
1:0:216:PRO:HD2	1:0:219:GLN:OE1	2.14	0.48
3:3:92:ARG:HD3	3:3:188:TYR:HD2	1.78	0.48
2:1:113:ILE:HG22	2:1:253:MET:SD	2.53	0.48
1:0:276:LEU:HD12	1:0:276:LEU:N	2.29	0.48
2:1:266:MET:O	2:1:267:ARG:C	2.57	0.48
1:0:92:ILE:HG21	1:0:306:THR:HG21	1.96	0.48
2:1:167:GLU:N	2:1:167:GLU:CD	2.71	0.48
1:0:108:PRO:CB	1:0:174:GLY:HA3	2.44	0.47
1:0:125:PRO:CB	1:0:129:VAL:HG21	2.43	0.47
2:1:129:MET:HE2	2:1:131:PHE:CD2	2.48	0.47
1:0:125:PRO:HG2	1:0:314:PHE:CE2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:185:LYS:HD3	3:3:125:PHE:CD1	2.48	0.47
1:0:201:ILE:CG2	1:0:202:GLY:H	2.27	0.47
2:1:229:MET:HE3	2:1:229:MET:HA	1.97	0.47
1:0:195:ILE:HG12	1:0:264:VAL:CG2	2.44	0.47
2:1:122:LYS:HG3	3:3:107:TYR:CE1	2.49	0.47
3:3:109:THR:OG1	3:3:228:LEU:HD12	2.15	0.47
3:3:148:THR:HA	3:3:151:LEU:HD12	1.97	0.47
3:3:170:PRO:HG2	3:3:171:TRP:H	1.79	0.47
3:3:182:ARG:NH2	3:3:182:ARG:HG3	2.29	0.47
2:1:94:ASP:C	2:1:95:LEU:HD23	2.39	0.47
2:1:126:PHE:CG	2:1:260:ALA:HB1	2.50	0.47
2:1:210:THR:O	2:1:210:THR:HG22	2.14	0.47
1:0:96:GLU:O	1:0:97:ALA:HB2	2.15	0.47
2:1:268:ASN:OD1	2:1:269:GLN:HG2	2.14	0.47
1:0:100:ILE:HD12	1:0:100:ILE:O	2.15	0.46
1:0:176:CYS:SG	1:0:263:ILE:HD13	2.55	0.46
2:1:290:SER:OG	3:3:68:ARG:NH2	2.48	0.46
1:0:110:TYR:O	1:0:111:CYS:HB2	2.15	0.46
1:0:258:ASN:CG	1:0:259:CYS:N	2.74	0.46
2:1:238:VAL:HG12	2:1:239:GLY:N	2.29	0.46
3:3:121:PHE:CE2	3:3:123:GLY:HA3	2.50	0.46
1:0:172:ARG:HB3	1:0:272:PHE:CE1	2.50	0.46
2:1:174:ALA:C	2:1:176:ASN:N	2.74	0.46
3:3:71:PHE:CE1	3:3:214:ILE:HB	2.50	0.46
3:3:188:TYR:C	3:3:190:THR:H	2.24	0.46
1:0:83:ALA:HB3	1:0:96:GLU:HA	1.98	0.46
1:0:235:LEU:HD12	1:0:239:ILE:HD12	1.97	0.46
2:1:238:VAL:CG1	2:1:239:GLY:N	2.79	0.46
3:3:22:ALA:HB1	3:3:23:PRO:HD2	1.97	0.46
3:3:24:ILE:HG23	3:3:25:LEU:HD13	1.98	0.46
3:3:115:LEU:HB2	3:3:169:ILE:HB	1.97	0.46
3:3:146:ARG:NH1	3:3:146:ARG:HG2	2.31	0.46
1:0:89:ASN:ND2	1:0:131:ARG:HH21	2.14	0.45
1:0:128:SER:CB	1:0:160:VAL:HG11	2.46	0.45
1:0:129:VAL:CG1	1:0:161:PHE:HD1	2.27	0.45
2:1:154:MET:CE	2:1:171:TRP:HA	2.45	0.45
3:3:58:VAL:H	3:3:59:PRO:HD2	1.81	0.45
2:1:201:TYR:HA	2:1:228:ASN:HD21	1.81	0.45
1:0:233:TYR:HA	3:3:66:MET:CE	2.30	0.45
1:0:156:THR:HA	1:0:162:GLY:HA2	1.99	0.45
3:3:204:VAL:CG1	3:3:208:ALA:HB3	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:0:317:LEU:HD23	1:0:318:ARG:N	2.32	0.45
2:1:74:SER:C	2:1:76:ALA:H	2.24	0.45
3:3:61:ASN:OD1	3:3:62:ALA:N	2.50	0.45
1:0:157:GLU:HB3	1:0:158:THR:H	1.50	0.45
1:0:250:GLN:CG	1:0:260:ALA:HB1	2.46	0.44
1:0:258:ASN:CG	1:0:259:CYS:H	2.24	0.44
2:1:156:VAL:O	2:1:156:VAL:HG12	2.17	0.44
3:3:90:PRO:HD2	3:3:188:TYR:OH	2.17	0.44
1:0:183:ALA:HB2	1:0:301:ILE:CG2	2.47	0.44
1:0:191:LEU:HD23	1:0:287:PRO:HA	1.99	0.44
1:0:283:LEU:C	1:0:283:LEU:HD23	2.42	0.44
2:1:195:MET:O	2:1:196:SER:C	2.60	0.44
2:1:242:LYS:HG3	2:1:243:SER:N	2.32	0.44
3:3:14:LEU:HD22	3:3:16:THR:H	1.82	0.44
2:1:245:TYR:CD1	2:1:245:TYR:N	2.85	0.44
1:0:129:VAL:CG2	1:0:164:ASN:ND2	2.80	0.44
2:1:261:TRP:CD1	3:3:36:ILE:HB	2.52	0.44
3:3:6:LEU:N	3:3:6:LEU:HD23	2.32	0.44
2:1:125:LEU:HD23	2:1:126:PHE:CZ	2.52	0.44
2:1:294:ILE:CD1	3:3:71:PHE:HB3	2.47	0.44
3:3:118:THR:HG23	3:3:166:THR:OG1	2.17	0.44
2:1:116:TYR:HE2	2:1:118:GLN:HE21	1.66	0.44
2:1:229:MET:CE	2:1:230:MET:N	2.79	0.44
1:0:100:ILE:HD12	1:0:100:ILE:C	2.43	0.44
1:0:310:MET:O	1:0:311:CYS:HB2	2.17	0.44
2:1:119:MET:CE	2:1:255:MET:HE3	2.47	0.44
2:1:261:TRP:CD1	3:3:39:GLU:HB2	2.53	0.44
3:3:79:LYS:HB2	3:3:79:LYS:HZ2	1.83	0.44
2:1:104:ASN:O	2:1:166:ARG:HD2	2.18	0.44
3:3:54:GLU:HG2	3:3:69:LEU:HD23	1.99	0.44
3:3:183:ASP:O	3:3:185:VAL:N	2.50	0.44
3:3:46:LEU:O	3:3:49:VAL:HG23	2.18	0.43
1:0:190:ALA:HB3	1:0:289:SER:HB3	1.99	0.43
2:1:123:VAL:CG2	2:1:124:GLU:N	2.81	0.43
2:1:197:PRO:CD	2:1:227:ASN:HB2	2.43	0.43
3:3:182:ARG:HG3	3:3:182:ARG:HH21	1.83	0.43
1:0:152:PRO:HD2	1:0:279:CYS:HA	2.00	0.43
2:1:157:PRO:O	2:1:158:PRO:C	2.60	0.43
2:1:294:ILE:HD12	3:3:56:ASN:HA	1.99	0.43
1:0:235:LEU:O	1:0:236:ASP:HB2	2.19	0.43
1:0:239:ILE:HB	2:1:265:PRO:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:222:TYR:C	2:1:224:ALA:H	2.27	0.43
2:1:297:LEU:HD23	3:3:85:VAL:CG2	2.48	0.43
3:3:132:LEU:C	3:3:132:LEU:HD23	2.44	0.43
3:3:60:THR:O	3:3:61:ASN:C	2.60	0.43
1:0:195:ILE:HG12	1:0:264:VAL:HG21	2.00	0.43
1:0:253:ASN:HD21	3:3:122:THR:HA	1.84	0.43
1:0:273:ASP:OD1	1:0:274:SER:N	2.51	0.43
1:0:129:VAL:CG2	1:0:164:ASN:HD21	2.32	0.43
1:0:141:GLU:C	1:0:143:SER:H	2.27	0.43
1:0:197:PRO:HG2	2:1:262:ILE:CG2	2.48	0.43
1:0:84:GLN:NE2	1:0:91:THR:HG21	2.34	0.43
2:1:120:ARG:HH11	3:3:237:GLN:NE2	2.16	0.43
1:0:88:GLY:C	1:0:90:SER:N	2.74	0.42
1:0:160:VAL:O	1:0:163:GLN:HB3	2.19	0.42
2:1:78:THR:HB	3:3:42:ASN:HD21	1.83	0.42
2:1:235:VAL:HG11	2:1:249:VAL:HG21	2.01	0.42
1:0:216:PRO:C	1:0:218:LYS:H	2.28	0.42
1:0:233:TYR:CD2	3:3:65:LEU:HB3	2.53	0.42
2:1:116:TYR:HE2	2:1:118:GLN:NE2	2.17	0.42
1:0:297:ALA:O	1:0:298:THR:C	2.62	0.42
1:0:175:PHE:HB3	1:0:176:CYS:H	1.67	0.42
1:0:121:LYS:HD2	1:0:121:LYS:H	1.84	0.42
3:3:135:TYR:CD1	3:3:169:ILE:HD12	2.54	0.42
1:0:84:GLN:CB	1:0:93:THR:HA	2.49	0.42
1:0:220:THR:HG23	2:1:222:TYR:CE1	2.55	0.42
3:3:29:TYR:HA	3:3:30:PRO:HD2	1.71	0.42
1:0:110:TYR:CE1	1:0:124:ARG:HB3	2.55	0.42
2:1:81:ASP:O	2:1:85:SER:HB3	2.19	0.42
2:1:120:ARG:HH12	2:1:274:LYS:HA	1.85	0.42
3:3:225:THR:CG2	3:3:226:MET:N	2.83	0.42
2:1:199:SER:OG	3:3:34:ILE:HG12	2.20	0.41
1:0:90:SER:CB	1:0:132:PHE:HB2	2.50	0.41
3:3:89:ASP:HA	3:3:188:TYR:CZ	2.55	0.41
3:3:90:PRO:O	3:3:102:GLY:HA2	2.20	0.41
3:3:178:ARG:HG2	3:3:179:ALA:N	2.34	0.41
2:1:80:LEU:HD12	2:1:80:LEU:HA	1.79	0.41
2:1:126:PHE:CD2	2:1:260:ALA:HB1	2.55	0.41
1:0:121:LYS:HE3	1:0:121:LYS:HB3	1.79	0.41
1:0:186:PHE:CE1	3:3:126:MET:HG3	2.56	0.41
1:0:85:LEU:HD12	1:0:94:THR:HG21	2.01	0.41
2:1:73:HIS:O	2:1:74:SER:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:1:177:PRO:CB	3:3:24:ILE:HD11	2.47	0.41
1:0:141:GLU:C	1:0:143:SER:N	2.75	0.41
2:1:117:ALA:HB1	3:3:236:LEU:HD22	2.03	0.41
2:1:180:PHE:HD1	2:1:180:PHE:H	1.69	0.41
2:1:291:ARG:O	3:3:58:VAL:HG12	2.21	0.41
2:1:297:LEU:HD23	3:3:85:VAL:HG22	2.03	0.41
3:3:74:SER:CB	3:3:211:THR:OG1	2.69	0.41
3:3:133:ILE:HG12	3:3:196:ILE:HG22	2.02	0.41
3:3:187:ASP:O	3:3:188:TYR:C	2.63	0.41
1:0:245:THR:C	1:0:247:CYS:H	2.27	0.41
3:3:25:LEU:HB3	3:3:28:PHE:HB2	2.02	0.41
1:0:211:GLU:C	1:0:213:SER:H	2.29	0.41
2:1:266:MET:HE1	3:3:107:TYR:CE2	2.56	0.41
1:0:151:PHE:N	1:0:151:PHE:CD1	2.88	0.41
1:0:185:LYS:HB3	3:3:125:PHE:HD1	1.85	0.41
1:0:195:ILE:HA	1:0:195:ILE:HD12	1.85	0.41
1:0:289:SER:HA	1:0:290:PRO:HD2	1.82	0.41
2:1:160:ALA:HA	2:1:161:PRO:HD3	1.86	0.41
3:3:162:GLN:O	3:3:162:GLN:HG3	2.21	0.41
1:0:87:ILE:O	1:0:131:ARG:HA	2.21	0.41
3:3:170:PRO:O	3:3:171:TRP:HB2	2.20	0.41
1:0:92:ILE:HG12	1:0:132:PHE:HE1	1.85	0.40
2:1:177:PRO:HB2	3:3:24:ILE:CD1	2.45	0.40
3:3:90:PRO:HG2	3:3:115:LEU:HD11	2.02	0.40
1:0:205:ALA:HB2	2:1:278:ASN:HD22	1.86	0.40
1:0:215:PRO:HA	1:0:216:PRO:HD3	1.85	0.40
2:1:125:LEU:HD13	2:1:266:MET:SD	2.61	0.40
3:3:130:LYS:O	3:3:199:GLN:HB3	2.21	0.40
2:1:94:ASP:C	2:1:96:PRO:HD3	2.47	0.40
1:0:149:TRP:HZ3	1:0:285:VAL:CG1	2.33	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	0	235/323 (73%)	175 (74%)	41 (17%)	19 (8%)	1	9
2	1	214/297 (72%)	171 (80%)	27 (13%)	16 (8%)	1	10
3	3	238/242 (98%)	187 (79%)	34 (14%)	17 (7%)	1	11
All	All	687/862 (80%)	533 (78%)	102 (15%)	52 (8%)	1	10

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	0	97	ALA
1	0	112	SER
1	0	127	VAL
2	1	169	LEU
2	1	196	SER
2	1	226	PRO
2	1	227	ASN
2	1	275	ALA
3	3	92	ARG
3	3	180	HIS
3	3	186	PHE
1	0	98	ALA
1	0	316	GLY
2	1	282	ASN
2	1	289	ALA
3	3	19	GLY
3	3	140	GLY
3	3	177	TYR
1	0	110	TYR
1	0	113	ASP
1	0	176	CYS
1	0	258	ASN
2	1	75	THR
3	3	147	ALA
3	3	182	ARG
3	3	188	TYR
3	3	219	ALA
3	3	228	LEU
1	0	154	VAL
1	0	157	GLU
1	0	223	GLY

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Mol	Chain	Res	Type
1	0	280	ASN
2	1	124	GLU
2	1	167	GLU
2	1	219	ASP
2	1	228	ASN
3	3	43	LEU
1	0	203	THR
1	0	215	PRO
1	0	235	LEU
2	1	79	THR
3	3	189	TYR
1	0	207	GLY
3	3	58	VAL
1	0	125	PRO
2	1	158	PRO
3	3	72	PRO
3	3	184	GLY
3	3	90	PRO
2	1	187	PRO
2	1	223	GLY
1	0	285	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	0	201/272 (74%)	188 (94%)	13 (6%)	15	37
2	1	186/250 (74%)	167 (90%)	19 (10%)	7	23
3	3	200/202 (99%)	177 (88%)	23 (12%)	5	18
All	All	587/724 (81%)	532 (91%)	55 (9%)	10	26

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	0	121	LYS

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Mol	Chain	Res	Type
1	0	157	GLU
1	0	171	TYR
1	0	181	CYS
1	0	195	ILE
1	0	225	ASP
1	0	242	SER
1	0	276	LEU
1	0	294	ASP
1	0	295	GLN
1	0	298	THR
1	0	306	THR
1	0	307	LEU
2	1	78	THR
2	1	80	LEU
2	1	100	THR
2	1	118	GLN
2	1	121	ARG
2	1	143	THR
2	1	167	GLU
2	1	172	GLN
2	1	175	THR
2	1	179	VAL
2	1	181	VAL
2	1	184	SER
2	1	202	GLN
2	1	204	PHE
2	1	220	LEU
2	1	228	ASN
2	1	237	THR
2	1	245	TYR
2	1	270	ASN
3	3	6	LEU
3	3	14	LEU
3	3	44	LEU
3	3	51	THR
3	3	53	LEU
3	3	58	VAL
3	3	61	ASN
3	3	63	THR
3	3	72	PRO
3	3	79	LYS
3	3	87	ARG

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Mol	Chain	Res	Type
3	3	90	PRO
3	3	150	MET
3	3	159	PHE
3	3	167	LEU
3	3	178	ARG
3	3	191	THR
3	3	204	VAL
3	3	206	ILE
3	3	211	THR
3	3	226	MET
3	3	228	LEU
3	3	230	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	0	130	ASN
1	0	164	ASN
1	0	168	HIS
1	0	188	GLN
1	0	214	HIS
1	0	257	ASN
2	1	118	GLN
2	1	172	GLN
2	1	176	ASN
2	1	202	GLN
2	1	278	ASN
3	3	48	GLN
3	3	110	GLN
3	3	176	HIS

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 [i](#)

There are no oligosaccharides 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.

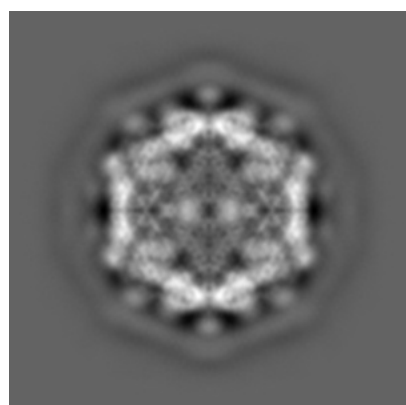
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-6200. These allow visual inspection of the internal detail of the map and identification of artifacts.

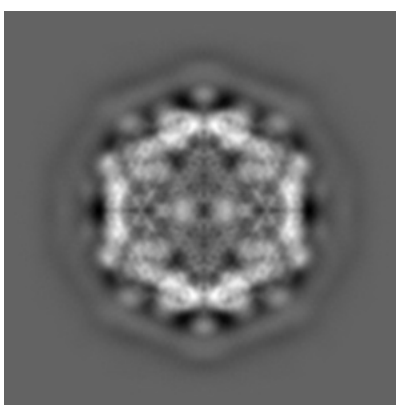
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

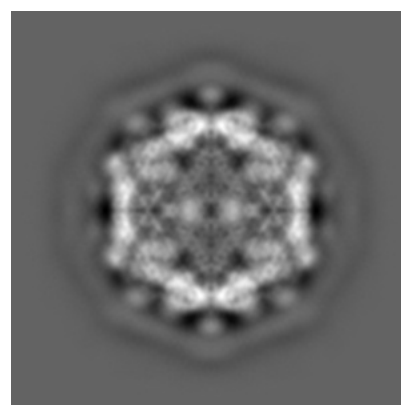
6.1.1 Primary map



X



Y

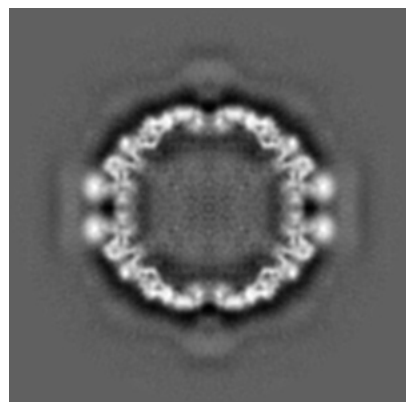


Z

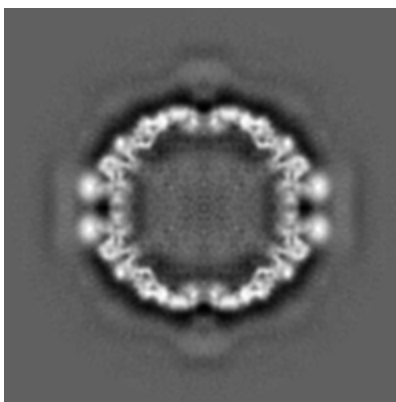
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

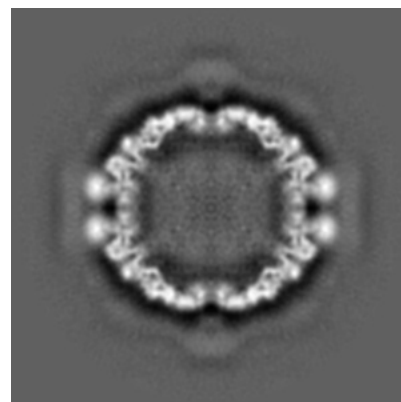
6.2.1 Primary map



X Index: 230



Y Index: 230

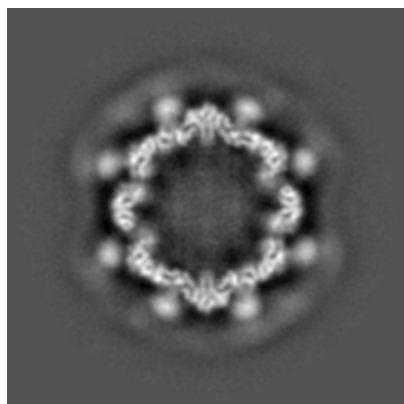


Z Index: 230

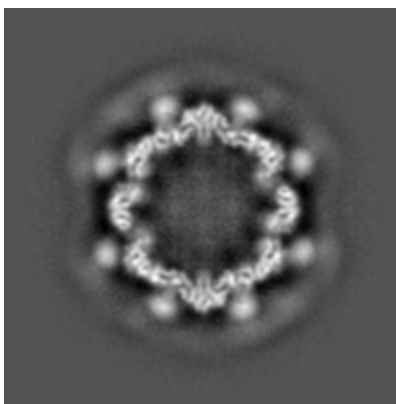
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

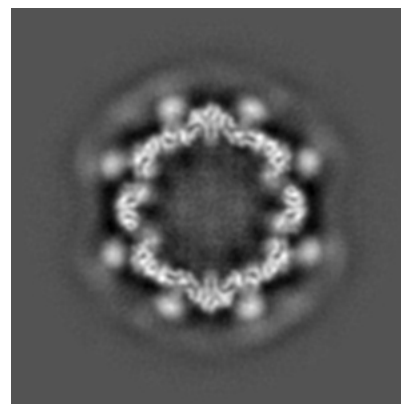
6.3.1 Primary map



X Index: 178



Y Index: 178

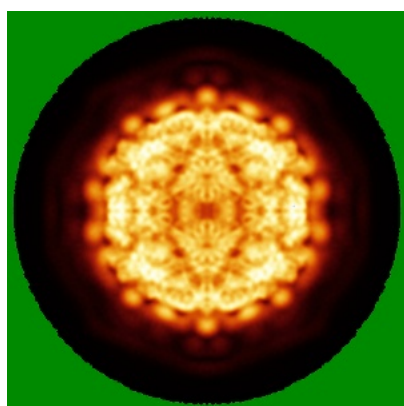


Z Index: 178

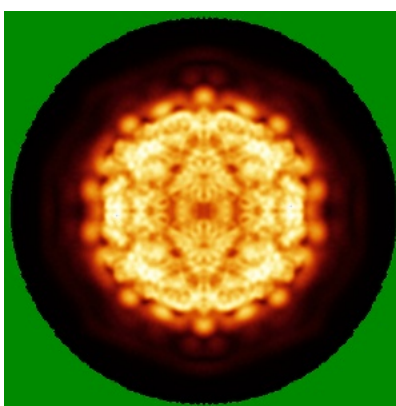
The images above show the largest variance slices of the map in three orthogonal directions.

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

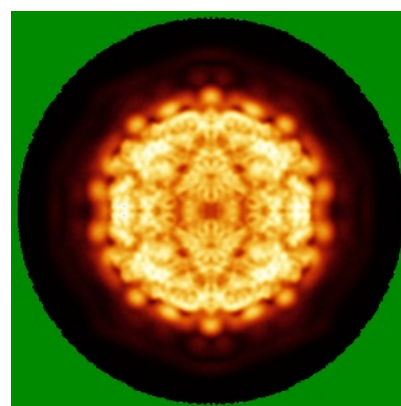
6.4.1 Primary map



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.

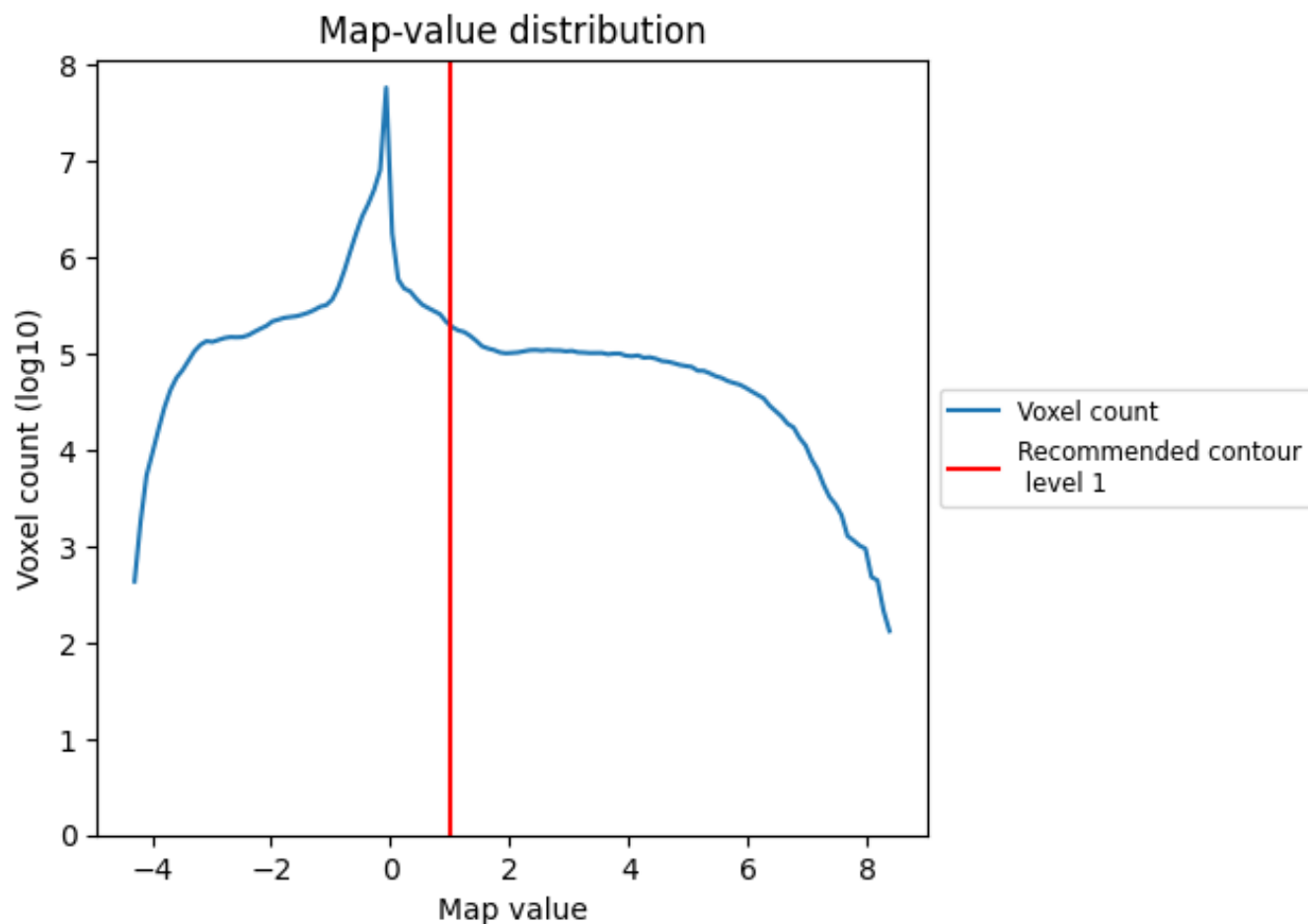
6.6 Mask visualisation

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

7 Map analysis [i](#)

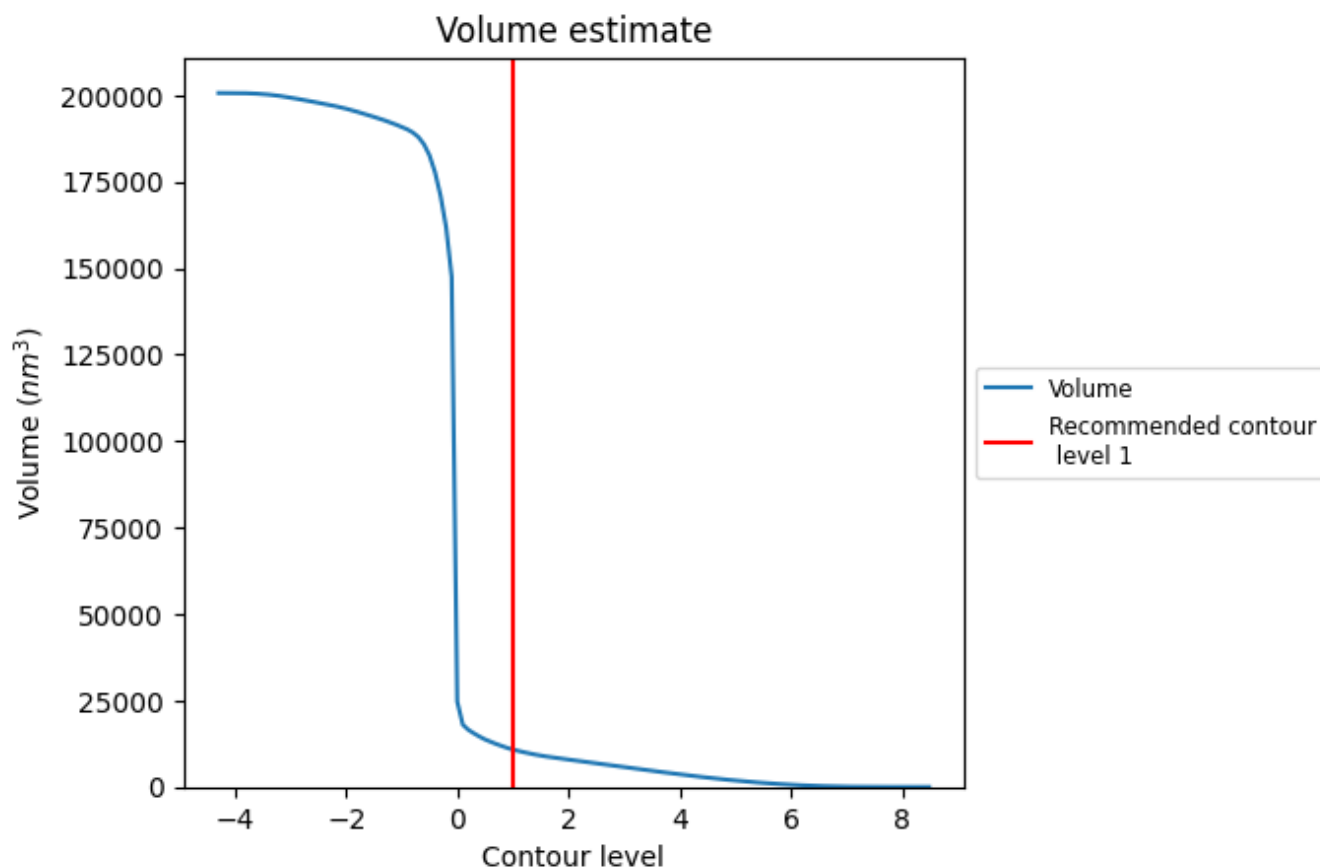
This section contains the results of statistical analysis of the map.

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.

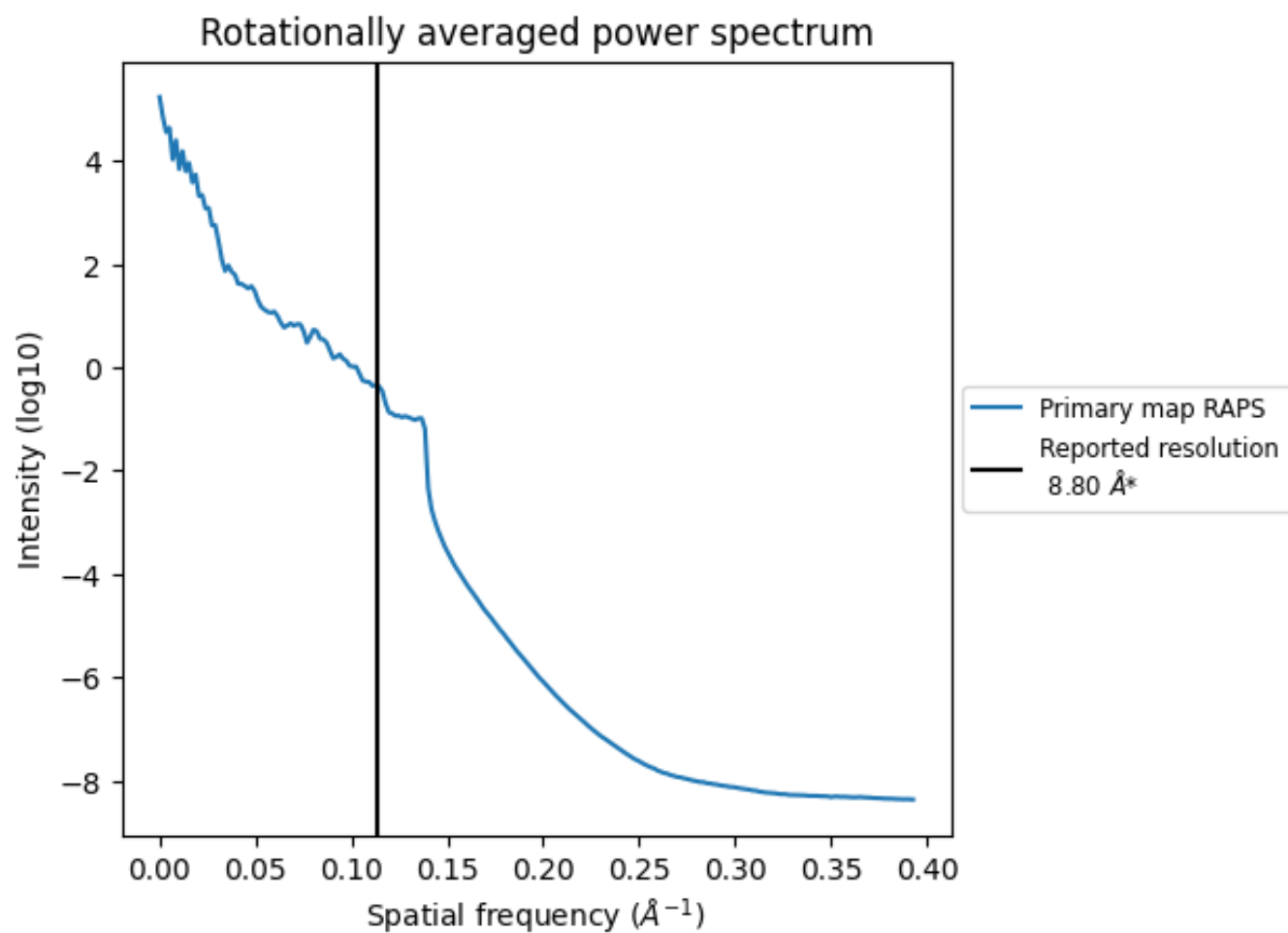
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 10808 nm³; this corresponds to an approximate mass of 9763 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 0.114 Å⁻¹

8 Fourier-Shell correlation ⓘ

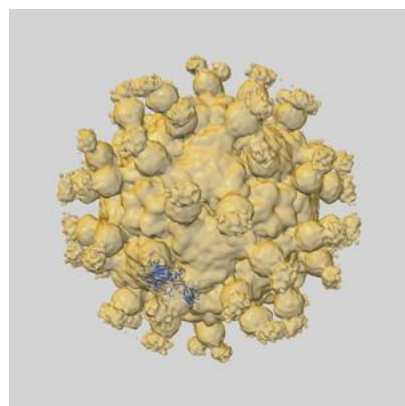
This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

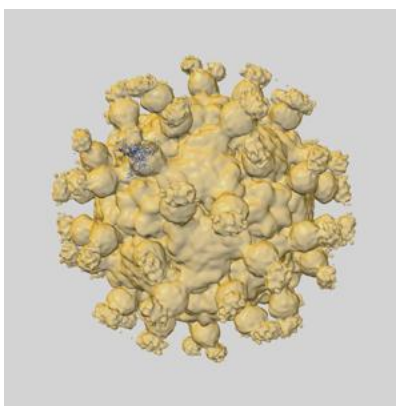
This section contains information regarding the fit between EMDB map EMD-6200 and PDB model 3J91. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlays

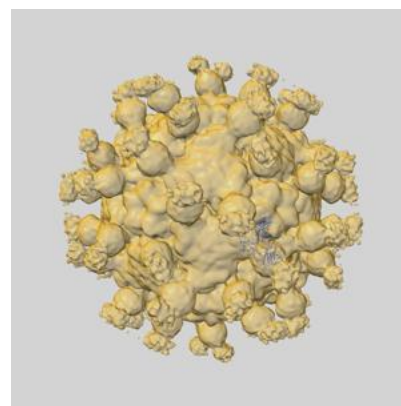
9.1.1 Map-model overlay [i](#)



X

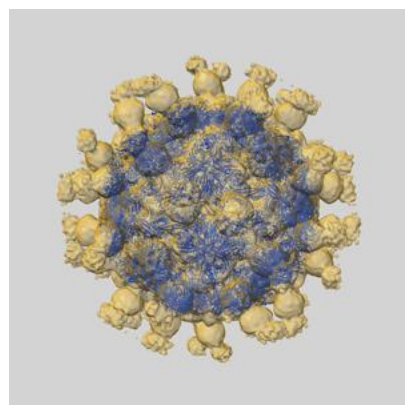


Y

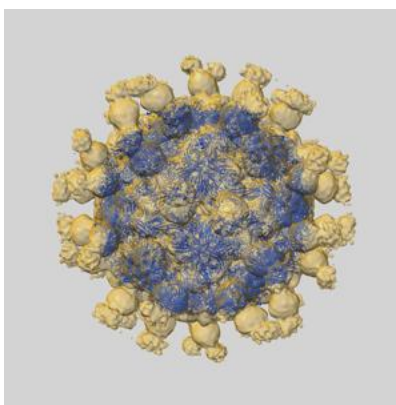


Z

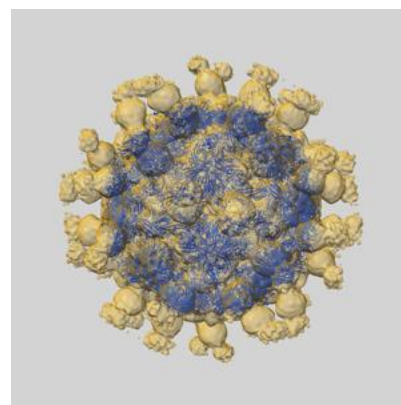
9.1.2 Map-model assembly overlay [i](#)



X



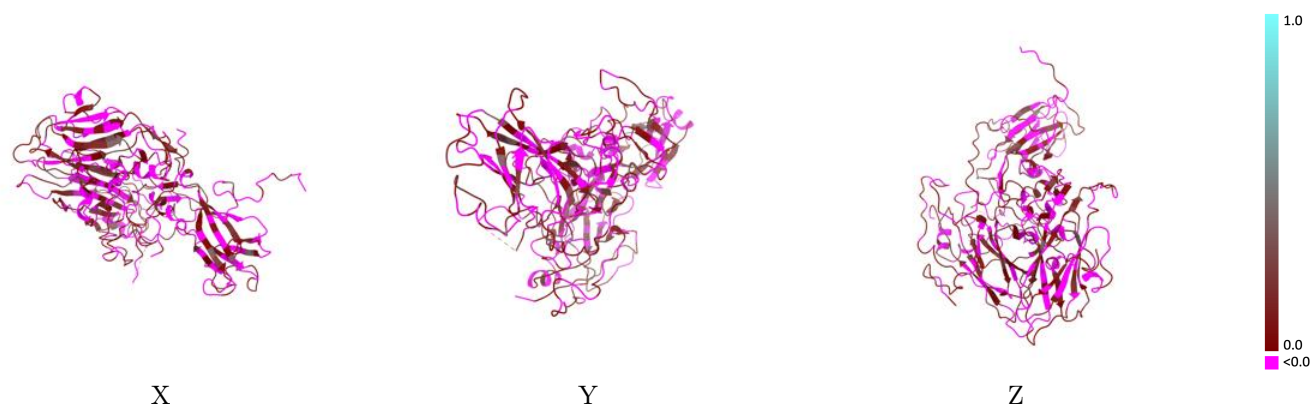
Y



Z

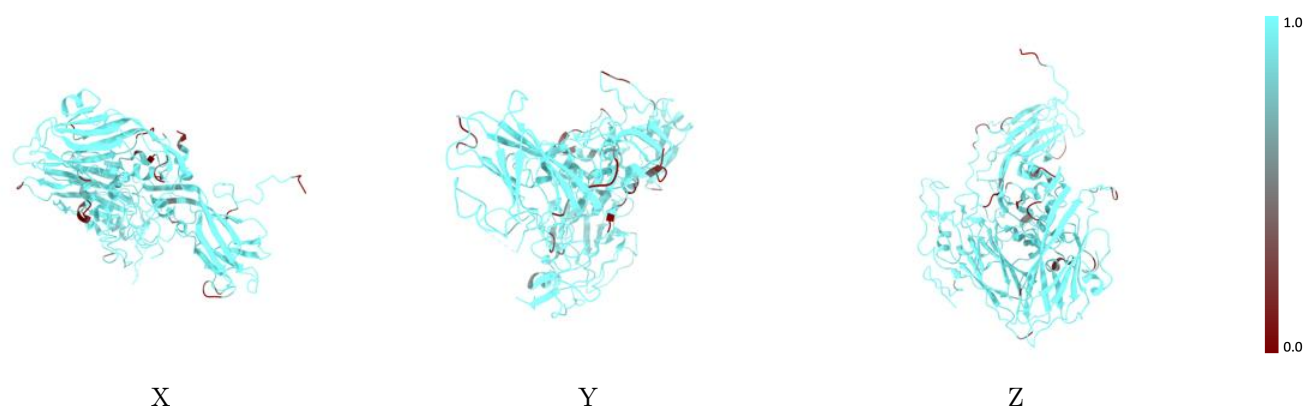
The images above show the 3D surface view of the map at the recommended contour level 1.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



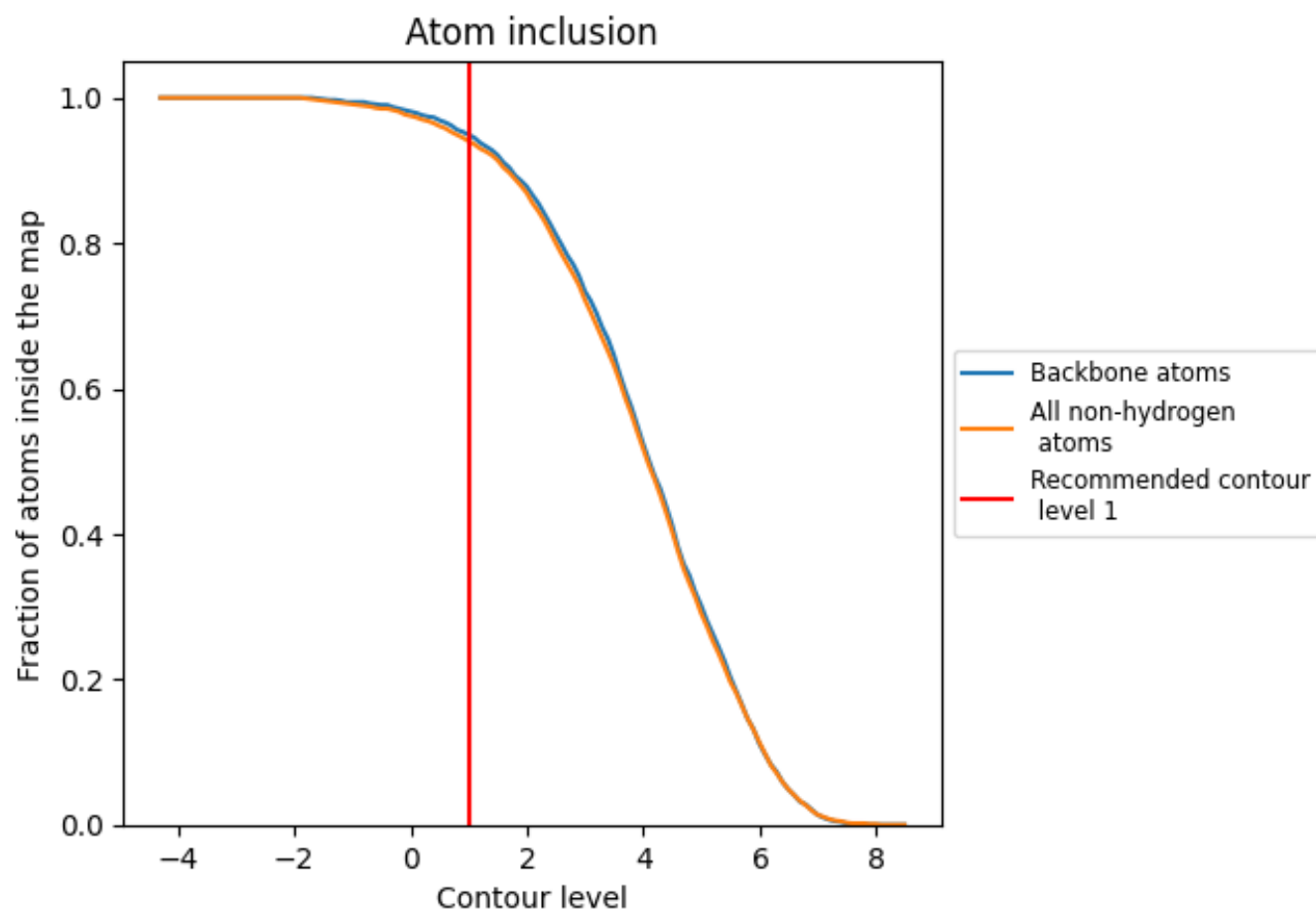
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 94% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

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

Chain	Atom inclusion	Q-score
All	0.9400	0.0330
0	0.9730	0.0420
1	0.9540	0.0360
3	0.8950	0.0210

