



## Full wwPDB EM Validation Report ⓘ

Nov 19, 2022 – 04:07 PM EST

PDB ID : 3J9E  
EMDB ID : EMD-6240  
Title : Atomic structure of a non-enveloped virus reveals pH sensors for a coordinated process of cell entry  
Authors : Zhang, X.; Patel, A.; Celma, C.; Roy, P.; Zhou, Z.H.  
Deposited on : 2015-01-10  
Resolution : 3.50 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](mailto: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>.

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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) : 2.31.3

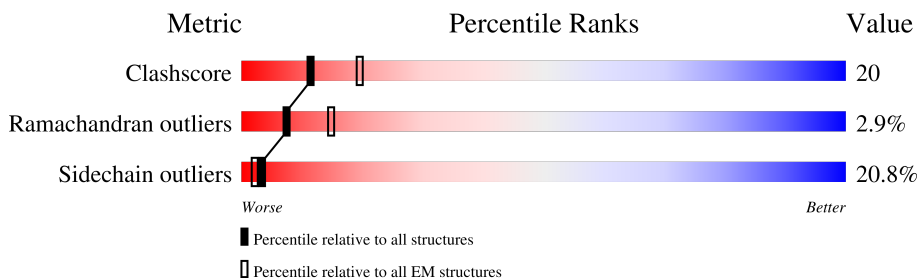
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.50 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  $\geq 3$ , 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	D	526	<div> <div>15%</div> <div>57%</div> <div>30%</div> <div>11%</div> <div>..</div> </div>

## 2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 4098 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.

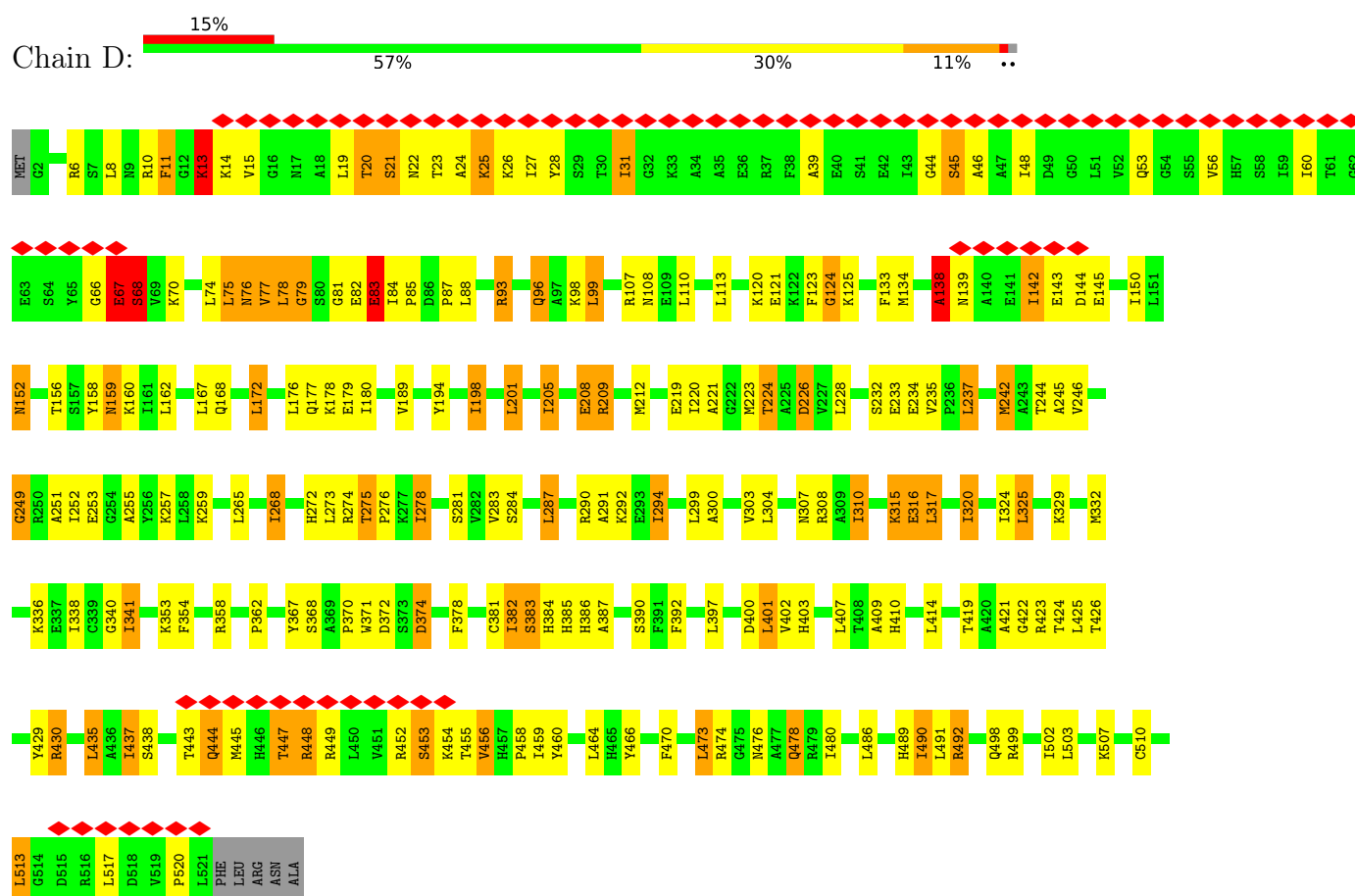
- Molecule 1 is a protein called VP5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	D	520	4098	2584	722	777	15	0	0

### 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: VP5



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, I	Depositor
Number of particles used	5008	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	each particle	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	20	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	3300	Depositor
Magnification	24140	Depositor
Image detector	GATAN K2 (4k x 4k)	Depositor
Maximum map value	5.958	Depositor
Minimum map value	-3.561	Depositor
Average map value	0.008	Depositor
Map value standard deviation	0.212	Depositor
Recommended contour level	0.5	Depositor
Map size ( $\text{\AA}$ )	116.15, 139.38, 150.49	wwPDB
Map dimensions	138, 115, 149	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.01, 1.01, 1.01	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	D	1.00	6/4164 (0.1%)	1.45	23/5605 (0.4%)

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	D	0	10

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	67	GLU	C-N	-25.77	0.74	1.34
1	D	13	LYS	C-N	-21.50	0.84	1.34
1	D	513	LEU	C-N	10.84	1.52	1.33
1	D	83	GLU	CB-CG	10.16	1.71	1.52
1	D	83	GLU	CG-CD	9.13	1.65	1.51
1	D	83	GLU	CD-OE2	6.16	1.32	1.25

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	138	ALA	O-C-N	-52.62	38.51	122.70
1	D	513	LEU	O-C-N	-47.36	42.70	123.20
1	D	13	LYS	O-C-N	-32.50	70.70	122.70
1	D	67	GLU	O-C-N	-19.20	91.99	122.70
1	D	13	LYS	C-N-CA	-16.14	81.35	121.70
1	D	13	LYS	CA-C-N	-14.44	85.42	117.20
1	D	138	ALA	CA-C-N	11.64	142.80	117.20
1	D	99	LEU	CA-CB-CG	-8.28	96.27	115.30
1	D	124	GLY	N-CA-C	7.87	132.78	113.10
1	D	513	LEU	CA-C-N	-7.16	101.88	116.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	67	GLU	CA-C-N	6.83	132.23	117.20
1	D	401	LEU	CA-CB-CG	6.53	130.32	115.30
1	D	45	SER	C-N-CA	5.54	135.56	121.70
1	D	414	LEU	CA-CB-CG	5.53	128.03	115.30
1	D	79	GLY	N-CA-C	-5.49	99.37	113.10
1	D	138	ALA	C-N-CA	-5.35	108.32	121.70
1	D	93	ARG	NE-CZ-NH1	-5.25	117.68	120.30
1	D	226	ASP	CB-CG-OD2	5.23	123.01	118.30
1	D	492	ARG	N-CA-C	5.15	124.92	111.00
1	D	45	SER	N-CA-C	5.14	124.88	111.00
1	D	287	LEU	CA-CB-CG	-5.06	103.67	115.30
1	D	473	LEU	CA-CB-CG	5.02	126.85	115.30
1	D	268	ILE	CB-CA-C	-5.00	101.59	111.60

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	123	PHE	Peptide
1	D	13	LYS	Mainchain
1	D	138	ALA	Peptide,Mainchain
1	D	249	GLY	Peptide
1	D	387	ALA	Peptide
1	D	513	LEU	Mainchain
1	D	67	GLU	Peptide,Mainchain
1	D	77	VAL	Peptide

## 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	D	4098	0	4119	164	0
All	All	4098	0	4119	164	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 20.

All (164) 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:D:453:SER:C	1:D:455:THR:CG2	1.79	1.48
1:D:437:ILE:HD11	1:D:447:THR:CG2	1.48	1.42
1:D:143:GLU:O	1:D:145:GLU:N	1.63	1.31
1:D:453:SER:O	1:D:455:THR:CG2	1.77	1.28
1:D:453:SER:C	1:D:455:THR:HG21	1.09	1.18
1:D:453:SER:O	1:D:455:THR:HG21	1.35	1.18
1:D:453:SER:CA	1:D:455:THR:HG21	1.74	1.16
1:D:144:ASP:CA	1:D:145:GLU:HB2	1.80	1.10
1:D:437:ILE:CD1	1:D:447:THR:HG23	1.81	1.10
1:D:437:ILE:HD11	1:D:447:THR:HG23	1.16	1.07
1:D:437:ILE:HD11	1:D:447:THR:HG21	1.30	1.07
1:D:138:ALA:HB1	1:D:139:ASN:OD1	1.54	1.07
1:D:138:ALA:CB	1:D:139:ASN:OD1	2.04	1.06
1:D:139:ASN:O	1:D:142:ILE:HD12	1.55	1.04
1:D:453:SER:O	1:D:455:THR:HG23	1.54	1.04
1:D:144:ASP:HA	1:D:145:GLU:CB	1.91	1.01
1:D:144:ASP:HA	1:D:145:GLU:HB2	1.02	1.01
1:D:437:ILE:CD1	1:D:447:THR:CG2	2.38	0.93
1:D:142:ILE:HG22	1:D:143:GLU:N	1.84	0.90
1:D:142:ILE:HG22	1:D:143:GLU:H	1.37	0.88
1:D:224:THR:OG1	1:D:226:ASP:OD1	1.93	0.84
1:D:142:ILE:CG2	1:D:143:GLU:H	1.91	0.83
1:D:139:ASN:HA	1:D:142:ILE:HD11	1.62	0.82
1:D:45:SER:N	1:D:46:ALA:HB3	2.00	0.77
1:D:138:ALA:HB3	1:D:139:ASN:OD1	1.85	0.75
1:D:142:ILE:CG2	1:D:143:GLU:N	2.49	0.75
1:D:374:ASP:N	1:D:374:ASP:OD1	2.19	0.74
1:D:158:TYR:HE1	1:D:253:GLU:HG2	1.54	0.73
1:D:144:ASP:O	1:D:145:GLU:OE1	2.07	0.72
1:D:139:ASN:CA	1:D:142:ILE:HD11	2.21	0.71
1:D:249:GLY:O	1:D:253:GLU:HG3	1.90	0.71
1:D:139:ASN:O	1:D:142:ILE:CD1	2.38	0.70
1:D:383:SER:HB2	1:D:470:PHE:CG	2.28	0.69
1:D:139:ASN:C	1:D:142:ILE:HD12	2.14	0.68
1:D:24:ALA:HB2	1:D:60:ILE:HD11	1.76	0.68
1:D:144:ASP:CA	1:D:145:GLU:CB	2.62	0.67
1:D:143:GLU:C	1:D:145:GLU:N	2.48	0.65
1:D:430:ARG:HB2	1:D:430:ARG:CZ	2.24	0.65
1:D:139:ASN:C	1:D:142:ILE:CD1	2.65	0.65
1:D:144:ASP:C	1:D:145:GLU:OE1	2.34	0.65
1:D:340:GLY:HA2	1:D:341:ILE:HB	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:22:ASN:HB3	1:D:26:LYS:HE3	1.79	0.65
1:D:429:TYR:CE1	1:D:458:PRO:HD3	2.33	0.64
1:D:490:ILE:HD11	1:D:502:ILE:HA	1.79	0.64
1:D:437:ILE:CD1	1:D:447:THR:HG21	2.18	0.63
1:D:194:TYR:CE1	1:D:234:GLU:HG3	2.34	0.63
1:D:13:LYS:HB3	1:D:14:LYS:N	2.10	0.62
1:D:453:SER:HA	1:D:455:THR:HG21	1.72	0.62
1:D:144:ASP:C	1:D:145:GLU:HB2	2.21	0.61
1:D:10:ARG:O	1:D:11:PHE:HB2	2.01	0.61
1:D:13:LYS:CB	1:D:14:LYS:N	2.32	0.61
1:D:139:ASN:CA	1:D:142:ILE:CD1	2.79	0.60
1:D:168:GLN:HB3	1:D:235:VAL:HG11	1.83	0.60
1:D:159:ASN:HB2	1:D:317:LEU:HD12	1.82	0.59
1:D:307:ASN:HA	1:D:310:ILE:HG22	1.85	0.59
1:D:120:LYS:HA	1:D:124:GLY:HA2	1.84	0.58
1:D:67:GLU:C	1:D:68:SER:HB3	2.23	0.58
1:D:316:GLU:HG2	1:D:386:HIS:HD2	1.68	0.58
1:D:74:LEU:HD23	1:D:510:CYS:HB2	1.86	0.58
1:D:437:ILE:CG1	1:D:447:THR:HG23	2.32	0.58
1:D:370:PRO:HA	1:D:459:ILE:HD12	1.84	0.58
1:D:381:CYS:O	1:D:390:SER:HB2	2.05	0.57
1:D:8:LEU:HD21	1:D:402:VAL:HB	1.87	0.57
1:D:77:VAL:HG22	1:D:221:ALA:HB3	1.87	0.57
1:D:87:PRO:HB3	1:D:503:LEU:HD22	1.87	0.55
1:D:120:LYS:O	1:D:124:GLY:HA3	2.06	0.55
1:D:82:GLU:HG2	1:D:83:GLU:O	2.06	0.54
1:D:219:GLU:OE1	1:D:219:GLU:N	2.38	0.54
1:D:244:THR:HG22	1:D:245:ALA:H	1.72	0.54
1:D:208:GLU:HG2	1:D:221:ALA:HA	1.90	0.54
1:D:409:ALA:O	1:D:410:HIS:HB2	2.08	0.54
1:D:316:GLU:OE1	1:D:385:HIS:N	2.41	0.54
1:D:144:ASP:C	1:D:145:GLU:CB	2.77	0.54
1:D:66:GLY:O	1:D:68:SER:HB3	2.08	0.53
1:D:67:GLU:C	1:D:68:SER:CB	2.49	0.53
1:D:421:ALA:HB3	1:D:422:GLY:HA3	1.91	0.53
1:D:316:GLU:HG2	1:D:384:HIS:HB3	1.91	0.53
1:D:294:ILE:HD11	1:D:299:LEU:HD21	1.90	0.52
1:D:237:LEU:HD22	1:D:249:GLY:HA3	1.91	0.52
1:D:275:THR:HG22	1:D:276:PRO:HD2	1.93	0.51
1:D:74:LEU:O	1:D:76:ASN:N	2.43	0.51
1:D:172:LEU:HD23	1:D:303:VAL:HG11	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:172:LEU:O	1:D:172:LEU:HG	2.11	0.50
1:D:84:ILE:HG22	1:D:85:PRO:HD2	1.92	0.50
1:D:290:ARG:HG2	1:D:290:ARG:HH11	1.76	0.50
1:D:272:HIS:CG	1:D:362:PRO:HG2	2.47	0.49
1:D:426:THR:HA	1:D:429:TYR:CE2	2.47	0.49
1:D:249:GLY:HA2	1:D:252:ILE:HG13	1.93	0.49
1:D:152:ASN:O	1:D:156:THR:HG23	2.11	0.49
1:D:110:LEU:HA	1:D:113:LEU:HD12	1.95	0.48
1:D:367:TYR:CE2	1:D:460:TYR:HD1	2.31	0.48
1:D:371:TRP:HD1	1:D:372:ASP:H	1.56	0.48
1:D:476:ASN:O	1:D:480:ILE:HG12	2.13	0.48
1:D:315:LYS:HE3	1:D:315:LYS:HB2	1.60	0.48
1:D:299:LEU:HA	1:D:299:LEU:HD23	1.63	0.48
1:D:320:ILE:HD12	1:D:324:ILE:HD12	1.95	0.48
1:D:168:GLN:HG2	1:D:235:VAL:HG13	1.96	0.47
1:D:287:LEU:HD23	1:D:287:LEU:HA	1.44	0.47
1:D:304:LEU:HD23	1:D:304:LEU:HA	1.69	0.47
1:D:11:PHE:HZ	1:D:489:HIS:CD2	2.33	0.47
1:D:27:ILE:O	1:D:31:ILE:HB	2.15	0.47
1:D:489:HIS:O	1:D:492:ARG:HD3	2.14	0.47
1:D:176:LEU:HD23	1:D:176:LEU:HA	1.70	0.46
1:D:201:LEU:HD22	1:D:205:ILE:HD12	1.97	0.46
1:D:371:TRP:CD1	1:D:372:ASP:N	2.82	0.46
1:D:316:GLU:HA	1:D:386:HIS:HB3	1.97	0.46
1:D:329:LYS:HA	1:D:332:MET:HE2	1.97	0.46
1:D:139:ASN:HB3	1:D:142:ILE:HD12	1.98	0.45
1:D:362:PRO:HD3	1:D:385:HIS:NE2	2.32	0.45
1:D:198:ILE:HD12	1:D:198:ILE:HA	1.58	0.45
1:D:21:SER:C	1:D:23:THR:H	2.20	0.45
1:D:82:GLU:O	1:D:96:GLN:NE2	2.49	0.45
1:D:162:LEU:HD23	1:D:162:LEU:HA	1.61	0.45
1:D:362:PRO:HB3	1:D:383:SER:O	2.17	0.45
1:D:44:GLY:HA2	1:D:48:ILE:HB	1.98	0.45
1:D:44:GLY:CA	1:D:48:ILE:HB	2.47	0.44
1:D:142:ILE:C	1:D:144:ASP:H	2.21	0.44
1:D:237:LEU:CD2	1:D:249:GLY:HA3	2.48	0.44
1:D:88:LEU:HD22	1:D:499:ARG:HD3	2.00	0.44
1:D:194:TYR:HE1	1:D:234:GLU:HG3	1.79	0.44
1:D:99:LEU:HA	1:D:99:LEU:HD12	1.73	0.44
1:D:320:ILE:HD12	1:D:320:ILE:HA	1.84	0.44
1:D:492:ARG:N	1:D:498:GLN:HG2	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:25:LYS:HB2	1:D:25:LYS:NZ	2.32	0.43
1:D:362:PRO:HA	1:D:382:ILE:HB	2.00	0.43
1:D:78:LEU:HB3	1:D:79:GLY:HA2	1.99	0.43
1:D:278:ILE:HD12	1:D:278:ILE:H	1.82	0.43
1:D:28:TYR:CE1	1:D:56:VAL:HG21	2.52	0.43
1:D:158:TYR:CE1	1:D:253:GLU:HG2	2.44	0.43
1:D:139:ASN:HB3	1:D:142:ILE:CD1	2.48	0.43
1:D:10:ARG:O	1:D:11:PHE:CB	2.67	0.43
1:D:316:GLU:HG3	1:D:384:HIS:CG	2.53	0.43
1:D:142:ILE:O	1:D:144:ASP:N	2.44	0.43
1:D:325:LEU:HD13	1:D:325:LEU:HA	1.73	0.43
1:D:251:ALA:O	1:D:255:ALA:N	2.52	0.43
1:D:444:GLN:HG2	1:D:445:MET:HG3	2.01	0.43
1:D:478:GLN:HE21	1:D:478:GLN:HA	1.84	0.43
1:D:108:ASN:HB3	1:D:255:ALA:HA	2.01	0.42
1:D:265:LEU:HD23	1:D:265:LEU:HA	1.70	0.42
1:D:133:PHE:CD1	1:D:150:ILE:HD13	2.54	0.42
1:D:307:ASN:HA	1:D:310:ILE:CG2	2.47	0.42
1:D:490:ILE:CD1	1:D:502:ILE:HA	2.48	0.42
1:D:209:ARG:HA	1:D:212:MET:HE3	2.02	0.41
1:D:383:SER:HA	1:D:466:TYR:OH	2.20	0.41
1:D:378:PHE:CD1	1:D:392:PHE:CE2	3.08	0.41
1:D:316:GLU:HG2	1:D:386:HIS:CD2	2.53	0.41
1:D:294:ILE:HD11	1:D:299:LEU:CD2	2.50	0.41
1:D:392:PHE:HB2	1:D:407:LEU:HD11	2.02	0.41
1:D:226:ASP:OD1	1:D:226:ASP:N	2.53	0.41
1:D:449:ARG:HH11	1:D:456:VAL:HG12	1.85	0.41
1:D:474:ARG:HH21	1:D:474:ARG:HG2	1.85	0.41
1:D:503:LEU:HD12	1:D:503:LEU:HA	1.88	0.41
1:D:242:MET:O	1:D:242:MET:HE3	2.21	0.41
1:D:435:LEU:HD22	1:D:435:LEU:HA	1.69	0.41
1:D:453:SER:HA	1:D:455:THR:CG2	2.46	0.41
1:D:445:MET:O	1:D:448:ARG:HB3	2.21	0.41
1:D:453:SER:CA	1:D:455:THR:CG2	2.65	0.41
1:D:491:LEU:HA	1:D:491:LEU:HD23	1.91	0.41
1:D:220:ILE:HG21	1:D:283:VAL:HG11	2.03	0.41
1:D:252:ILE:HG13	1:D:252:ILE:H	1.64	0.41
1:D:176:LEU:HB3	1:D:300:ALA:HB2	2.02	0.40
1:D:19:LEU:HB2	1:D:20:THR:HA	2.04	0.40
1:D:78:LEU:HB3	1:D:79:GLY:CA	2.52	0.40
1:D:354:PHE:CD1	1:D:354:PHE:N	2.89	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	D	512/526 (97%)	447 (87%)	50 (10%)	15 (3%)	4	31

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	39	ALA
1	D	68	SER
1	D	142	ILE
1	D	180	ILE
1	D	400	ASP
1	D	11	PHE
1	D	81	GLY
1	D	21	SER
1	D	75	LEU
1	D	291	ALA
1	D	453	SER
1	D	15	VAL
1	D	83	GLU
1	D	341	ILE
1	D	520	PRO

### 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	D	437/442 (99%)	346 (79%)	91 (21%)	<b>1</b> <b>6</b>

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	6	ARG
1	D	20	THR
1	D	25	LYS
1	D	31	ILE
1	D	53	GLN
1	D	67	GLU
1	D	68	SER
1	D	70	LYS
1	D	75	LEU
1	D	76	ASN
1	D	78	LEU
1	D	93	ARG
1	D	96	GLN
1	D	98	LYS
1	D	107	ARG
1	D	121	GLU
1	D	125	LYS
1	D	134	MET
1	D	152	ASN
1	D	159	ASN
1	D	160	LYS
1	D	167	LEU
1	D	172	LEU
1	D	177	GLN
1	D	178	LYS
1	D	179	GLU
1	D	189	VAL
1	D	198	ILE
1	D	201	LEU
1	D	205	ILE
1	D	208	GLU
1	D	209	ARG
1	D	223	MET
1	D	224	THR
1	D	228	LEU
1	D	232	SER
1	D	233	GLU
1	D	237	LEU

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Mol	Chain	Res	Type
1	D	242	MET
1	D	246	VAL
1	D	257	LYS
1	D	259	LYS
1	D	268	ILE
1	D	273	LEU
1	D	274	ARG
1	D	275	THR
1	D	278	ILE
1	D	281	SER
1	D	284	SER
1	D	292	LYS
1	D	294	ILE
1	D	308	ARG
1	D	310	ILE
1	D	315	LYS
1	D	316	GLU
1	D	317	LEU
1	D	320	ILE
1	D	325	LEU
1	D	336	LYS
1	D	338	ILE
1	D	353	LYS
1	D	358	ARG
1	D	368	SER
1	D	374	ASP
1	D	382	ILE
1	D	383	SER
1	D	397	LEU
1	D	401	LEU
1	D	403	HIS
1	D	419	THR
1	D	423	ARG
1	D	424	THR
1	D	425	LEU
1	D	430	ARG
1	D	435	LEU
1	D	437	ILE
1	D	438	SER
1	D	443	THR
1	D	444	GLN
1	D	447	THR

*Continued on next page...*

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Mol	Chain	Res	Type
1	D	448	ARG
1	D	452	ARG
1	D	454	LYS
1	D	456	VAL
1	D	464	LEU
1	D	473	LEU
1	D	478	GLN
1	D	486	LEU
1	D	490	ILE
1	D	507	LYS
1	D	517	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	489	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 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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	5

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	144:ASP	C	145:GLU	N	4.38
1	D	454:LYS	C	455:THR	N	2.55
1	D	442:GLY	C	443:THR	N	2.16
1	D	13:LYS	C	14:LYS	N	0.84
1	D	67:GLU	C	68:SER	N	0.74



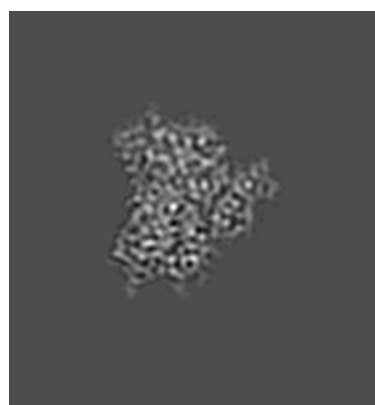
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-6240. 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: 57



Y Index: 69



Z Index: 74

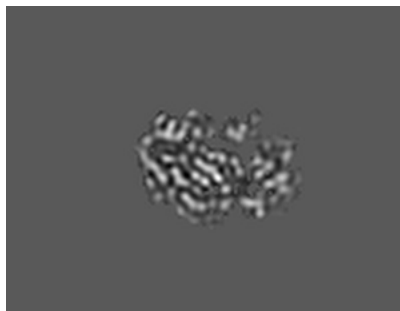
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

### 6.3.1 Primary map



X Index: 56



Y Index: 57

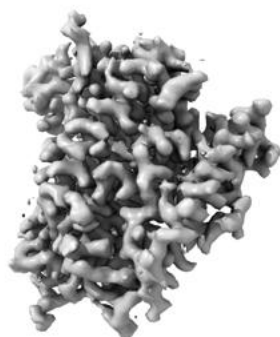


Z Index: 67

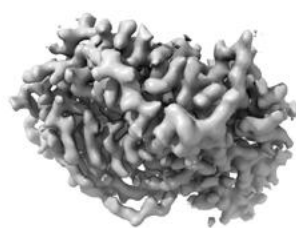
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

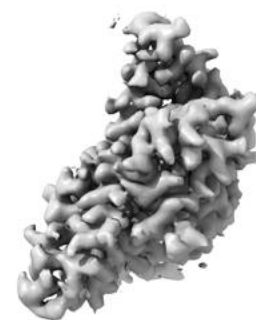
### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

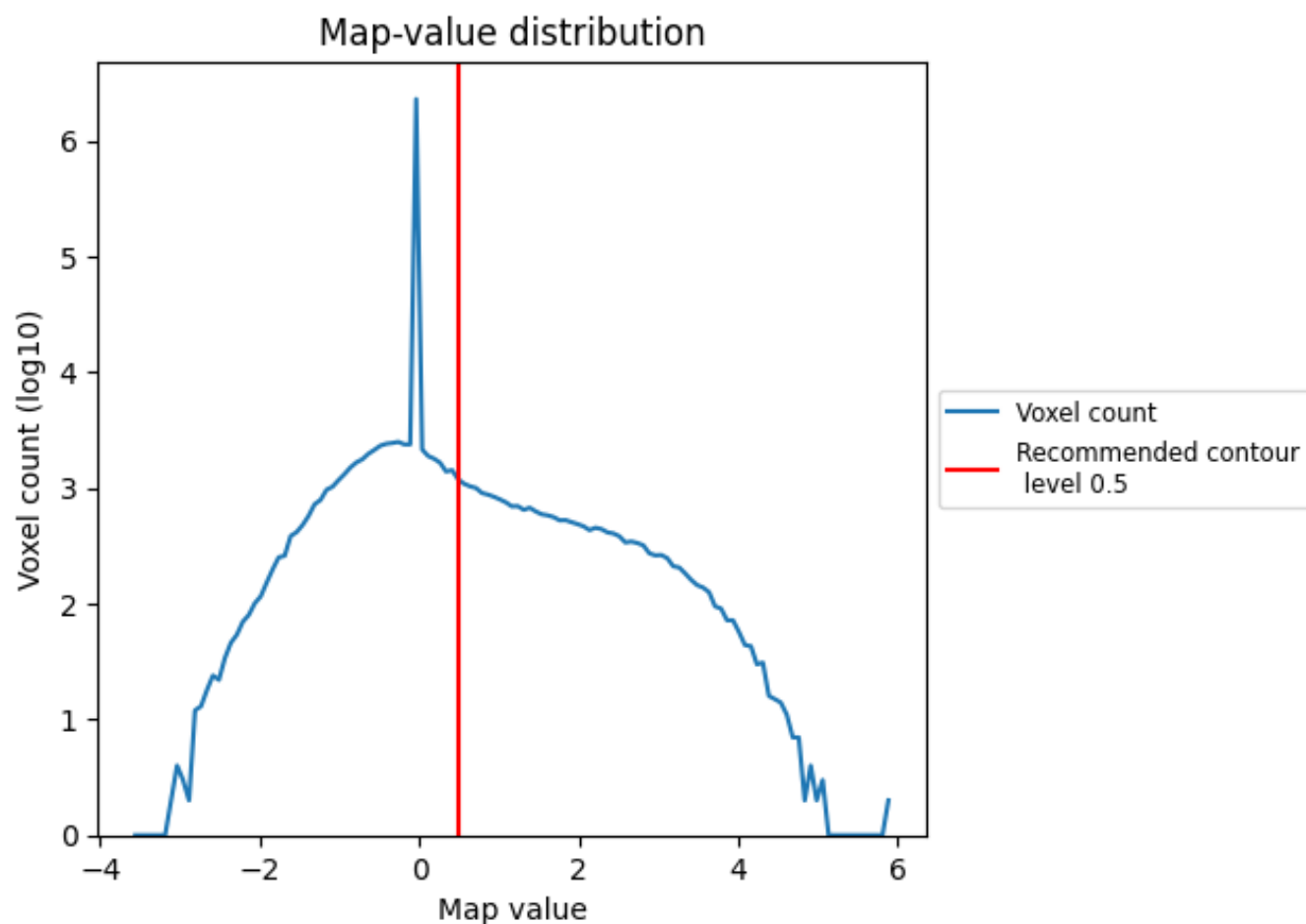
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

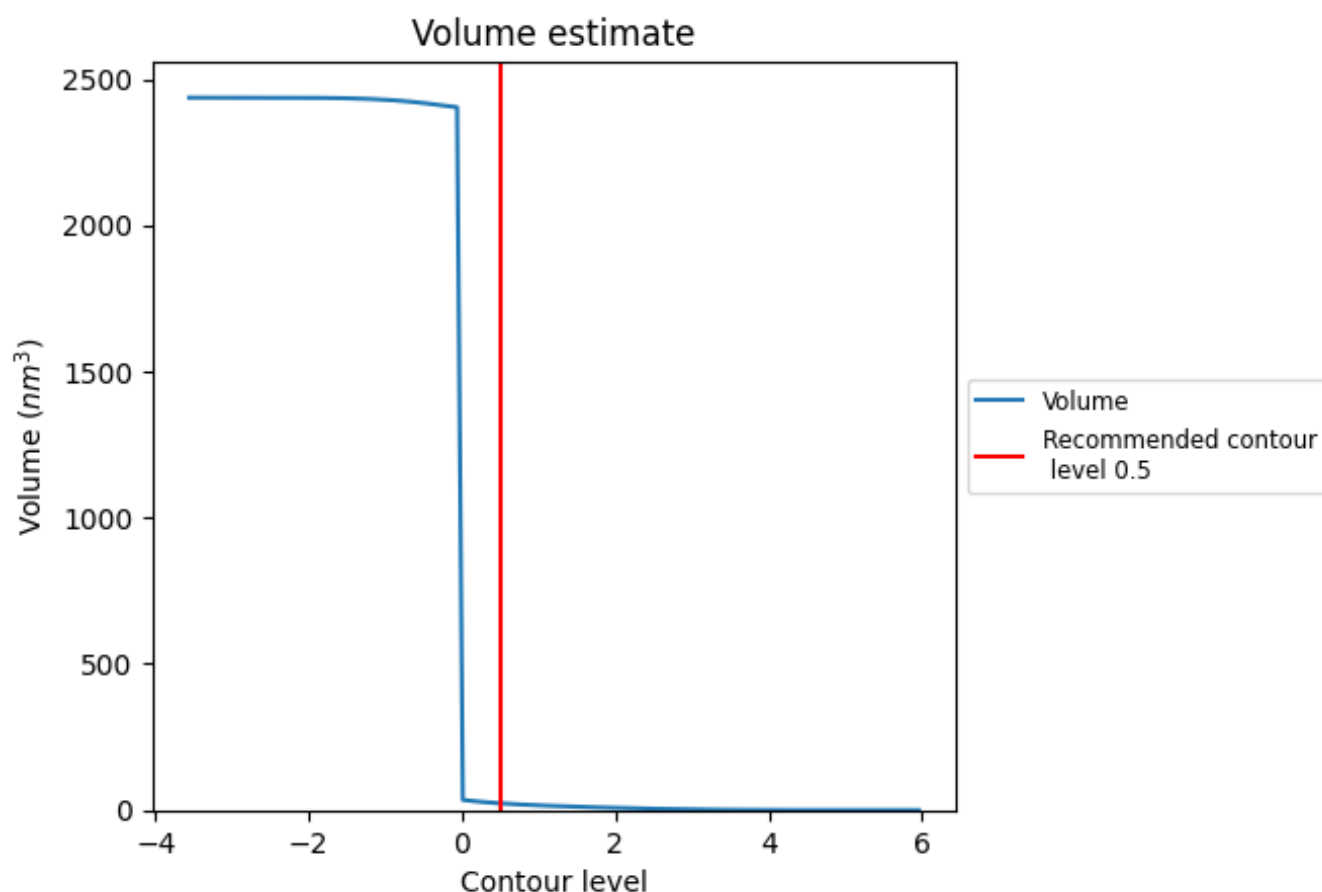
This section contains the results of statistical analysis of the map.

### 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 23 nm<sup>3</sup>; this corresponds to an approximate mass of 21 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](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

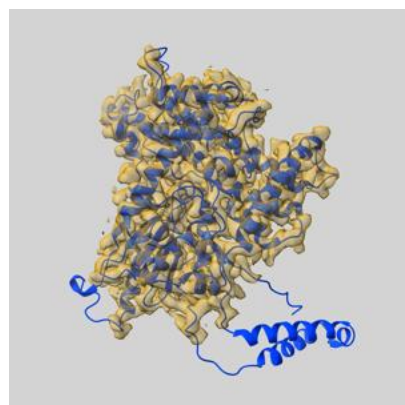
## 8 Fourier-Shell correlation

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

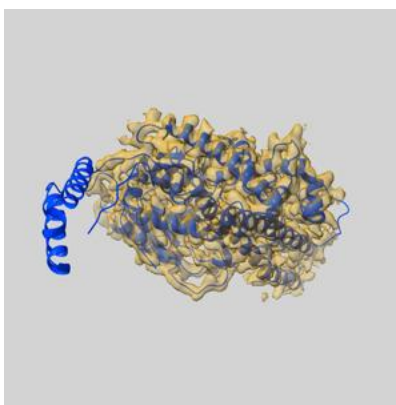
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-6240 and PDB model 3J9E. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

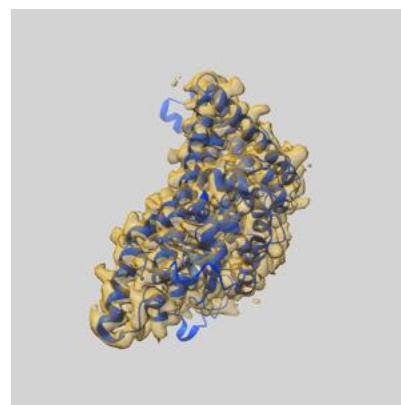
### 9.1 Map-model overlay [i](#)



X



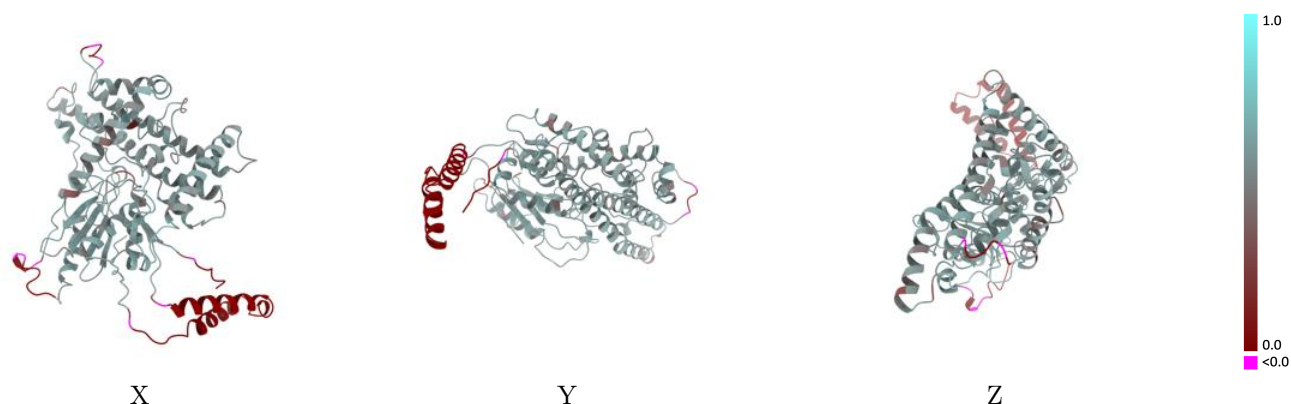
Y



Z

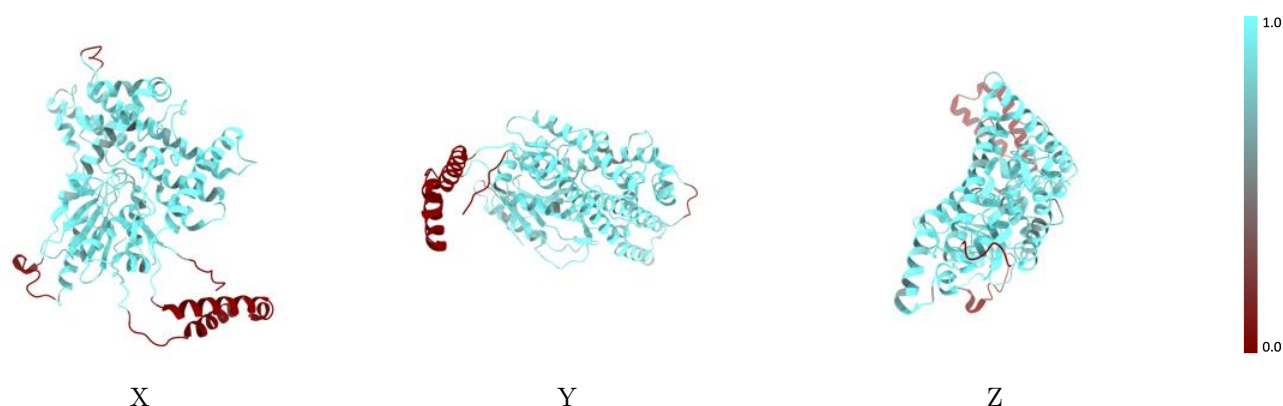
The images above show the 3D surface view of the map at the recommended contour level 0.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

## 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.

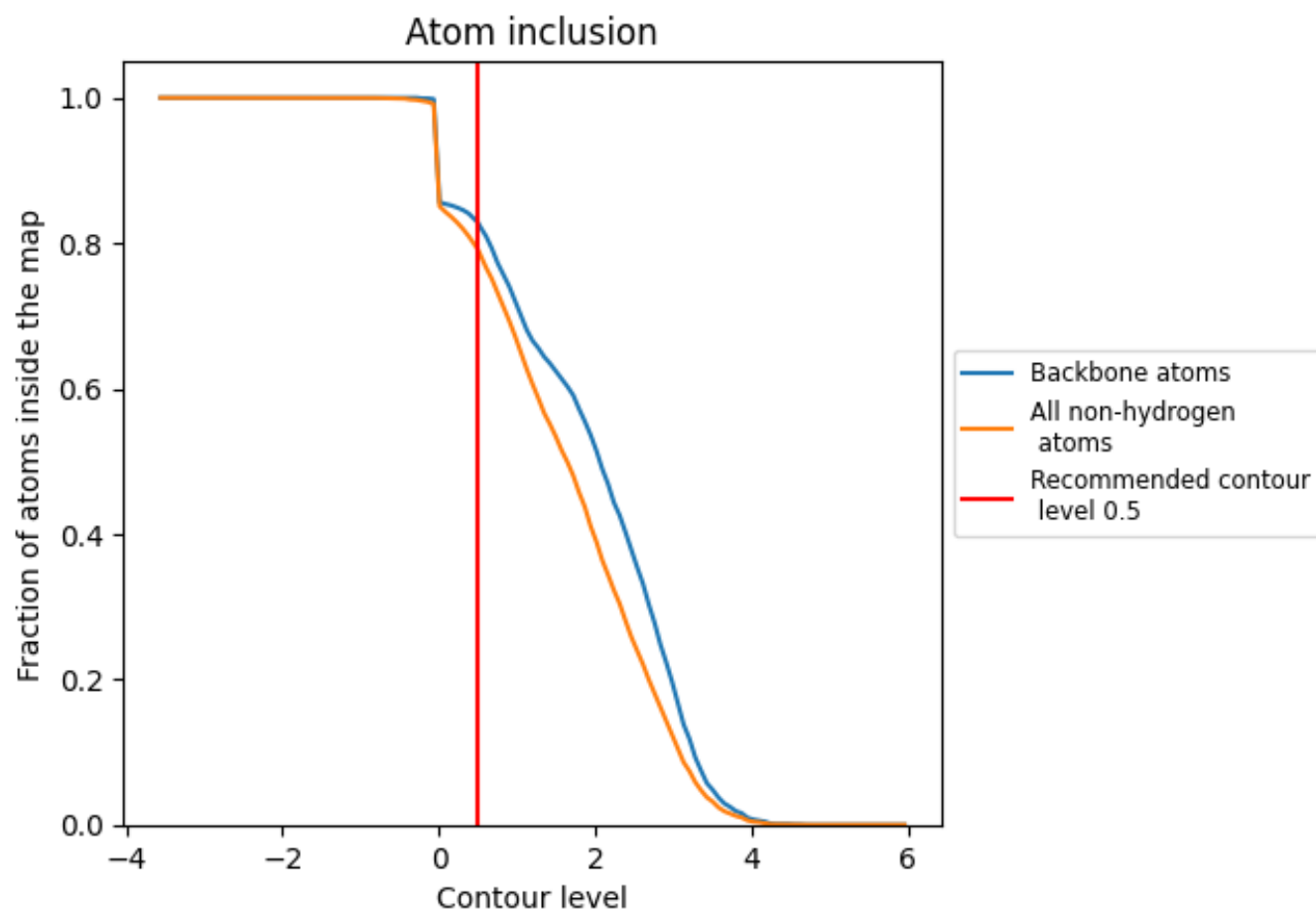
## 9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.5).



## 9.4 Atom inclusion [i](#)



At the recommended contour level, 83% of all backbone atoms, 79% 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 (0.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.7914	<div><div></div></div> 0.4740
D	<div><div></div></div> 0.7914	<div><div></div></div> 0.4740

