



Full wwPDB EM Validation Report ⓘ

Mar 26, 2026 – 03:48 PM UTC

PDB ID : 7MI3 / pdb_00007mi3
EMDB ID : EMD-23838
Title : Signal subtracted reconstruction of AAA2, AAA3, and AAA4 domains of dynein in the presence of a pyrazolo-pyrimidinone-based compound, Model 4
Authors : Santarossa, C.C.; Coudray, N.; Urnavicius, L.; Ekiert, D.C.; Bhabha, G.; Kapoor, T.M.
Deposited on : 2021-04-16
Resolution : 3.50 Å (reported)
Based on initial model : 7MI1

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

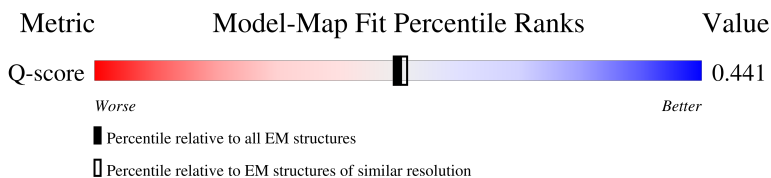
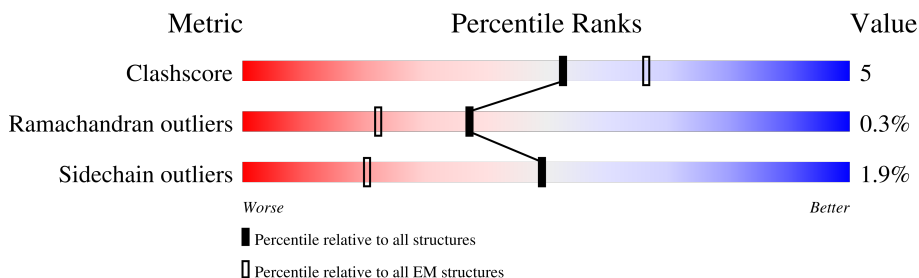
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.50 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	13950 (3.00 - 4.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	2661	 38% 6% 55%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9682 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 Fusion protein of Dynein and Endolysin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1192	Total	C	N	O	S	0	0
			9590	6143	1611	1787	49		

There are 33 discrepancies between the modelled and reference sequences:

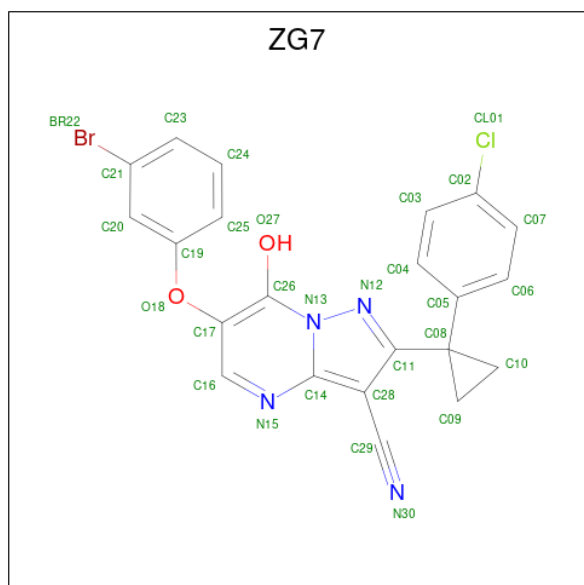
Chain	Residue	Modelled	Actual	Comment	Reference
A	1363	GLY	-	expression tag	UNP P36022
A	1849	GLN	GLU	conflict	UNP P36022
A	3120	GLY	-	linker	UNP P36022
A	3121	SER	-	linker	UNP P36022
A	3122	GLY	-	linker	UNP P36022
A	3123	SER	-	linker	UNP P36022
A	3124	GLY	-	linker	UNP P36022
A	3125	SER	-	linker	UNP P36022
A	3136	GLY	ARG	conflict	UNP P00720
A	3178	THR	CYS	conflict	UNP P00720
A	3221	ALA	CYS	conflict	UNP P00720
A	3261	ARG	ILE	conflict	UNP P00720
A	3286	GLY	-	linker	UNP P00720
A	3287	SER	-	linker	UNP P00720
A	3288	GLY	-	linker	UNP P00720
A	3289	SER	-	linker	UNP P00720
A	3290	GLY	-	linker	UNP P00720
A	3291	SER	-	linker	UNP P00720
A	3742	ASP	ASN	conflict	UNP P36022
A	3895	VAL	PHE	conflict	UNP P36022
A	4072	ASP	ASN	conflict	UNP P36022
A	4093	GLY	-	expression tag	UNP P36022
A	4094	SER	-	expression tag	UNP P36022
A	4095	GLY	-	expression tag	UNP P36022
A	4096	SER	-	expression tag	UNP P36022
A	4097	GLY	-	expression tag	UNP P36022
A	4098	SER	-	expression tag	UNP P36022
A	4099	HIS	-	expression tag	UNP P36022

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Chain	Residue	Modelled	Actual	Comment	Reference
A	4100	HIS	-	expression tag	UNP P36022
A	4101	HIS	-	expression tag	UNP P36022
A	4102	HIS	-	expression tag	UNP P36022
A	4103	HIS	-	expression tag	UNP P36022
A	4104	HIS	-	expression tag	UNP P36022

- Molecule 2 is (8S)-6-(3-bromophenoxy)-2-[1-(4-chlorophenyl)cyclopropyl]-7-hydroxypyrazolo[1,5-a]pyrimidine-3-carbonitrile (CCD ID: ZG7) (formula: C₂₂H₁₄BrClN₄O₂).



Mol	Chain	Residues	Atoms						AltConf
2	A	1	Total	Br	C	Cl	N	O	0
			30	1	22	1	4	2	
2	A	1	Total	Br	C	Cl	N	O	0
			30	1	22	1	4	2	

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C	N	O	P	0
			31	10	5	13	3	

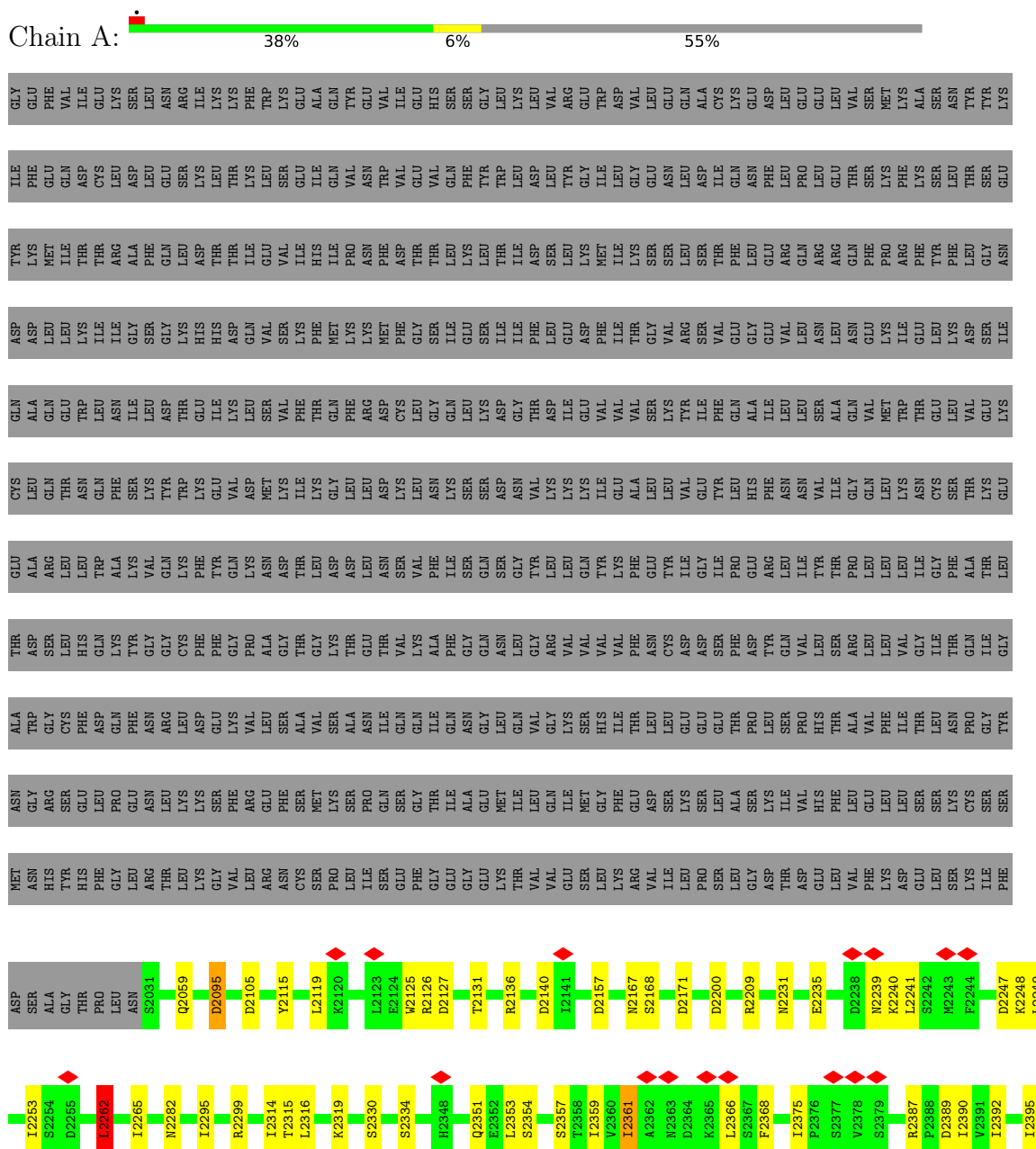
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	Mg	0
			1	1	

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Fusion protein of Dynein and Endolysin





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	131702	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	44	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	3700	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.020	Depositor
Minimum map value	-0.009	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.004	Depositor
Map size (Å)	190.62401, 190.62401, 190.62401	wwPDB
Map dimensions	368, 368, 368	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.518, 0.518, 0.518	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MG, ZG7, ATP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.45	0/9788	0.83	10/13249 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2115	TYR	N-CA-C	-6.89	99.76	110.17
1	A	3456	GLU	CA-C-N	6.41	136.22	121.87
1	A	3456	GLU	C-N-CA	6.41	136.22	121.87
1	A	2684	GLN	N-CA-C	6.07	119.88	112.23
1	A	2262	LEU	CD1-CG-CD2	-5.48	98.75	110.80
1	A	3467	SER	N-CA-C	-5.41	98.47	107.99
1	A	2985	ASN	N-CA-C	5.22	121.36	109.81
1	A	2491	LEU	N-CA-C	5.16	121.21	109.81
1	A	2115	TYR	CA-C-N	-5.14	116.89	122.55
1	A	2115	TYR	C-N-CA	-5.14	116.89	122.55

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9590	0	9637	103	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	60	0	0	3	0
3	A	31	0	12	1	0
4	A	1	0	0	0	0
All	All	9682	0	9649	103	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 5.

All (103) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2476:LYS:HE3	1:A:2528:ARG:HD2	1.59	0.85
1:A:2838:ALA:HB3	1:A:2878:VAL:HG13	1.76	0.65
1:A:2763:ARG:NH1	1:A:3511:SER:O	2.30	0.64
1:A:2517:LYS:O	1:A:2521:ASN:N	2.32	0.60
1:A:2979:LYS:HG3	1:A:3345:MET:HE1	1.84	0.59
1:A:2366:LEU:HD23	1:A:2368:PHE:HE2	1.67	0.59
1:A:2395:ILE:HD12	1:A:2395:ILE:H	1.68	0.58
1:A:2696:PHE:HA	1:A:2706:GLU:HA	1.86	0.57
1:A:2786:ILE:HD11	1:A:2821:ASN:HA	1.86	0.57
1:A:2316:LEU:HD23	1:A:2359:ILE:HD11	1.87	0.57
1:A:3342:ARG:NH1	1:A:3392:GLU:OE1	2.37	0.56
1:A:3504:ASP:OD1	1:A:3504:ASP:N	2.40	0.55
1:A:3409:ASP:HB3	1:A:3518:PHE:HB2	1.89	0.54
1:A:2136:ARG:O	1:A:2140:ASP:HB2	2.09	0.53
1:A:3408:LEU:HD23	1:A:3517:HIS:CE1	2.43	0.52
1:A:2476:LYS:HD3	1:A:2476:LYS:H	1.74	0.52
1:A:2728:LEU:HD13	1:A:2730:VAL:HG23	1.92	0.52
1:A:3385:ASP:OD1	1:A:3385:ASP:N	2.43	0.52
1:A:2642:ARG:O	1:A:2646:ARG:HG2	2.11	0.51
1:A:2984:VAL:O	1:A:2987:ARG:NH1	2.44	0.51
1:A:2984:VAL:HG13	1:A:2985:ASN:H	1.75	0.51
1:A:2832:ASN:OD1	1:A:2905:SER:OG	2.26	0.51
1:A:2241:LEU:HD13	1:A:2299:ARG:HH11	1.76	0.50
1:A:2765:GLY:H	2:A:4201:ZG7:C26	2.24	0.50
1:A:2936:ILE:HD12	1:A:3000:VAL:HA	1.93	0.50
1:A:2157:ASP:N	1:A:2157:ASP:OD1	2.43	0.50
1:A:2401:GLU:OE2	1:A:2430:ASN:ND2	2.32	0.50
1:A:2758:LEU:HD23	1:A:2915:ASN:HB3	1.93	0.50
1:A:2262:LEU:HA	1:A:2265:ILE:HD12	1.93	0.49
1:A:2944:ILE:HG21	1:A:2961:ILE:HG12	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2387:ARG:NH2	1:A:2389:ASP:OD2	2.33	0.49
1:A:2494:LEU:HD22	1:A:2498:GLY:HA2	1.94	0.49
1:A:2552:ARG:NH1	3:A:4203:ATP:O1G	2.42	0.49
1:A:2330:SER:HB3	1:A:2334:SER:HB2	1.93	0.49
1:A:3387:ASN:HB3	1:A:3390:PHE:HD1	1.77	0.49
1:A:2646:ARG:NH1	1:A:2706:GLU:OE1	2.46	0.48
1:A:3414:MET:C	1:A:3416:THR:H	2.20	0.48
1:A:2354:SER:HG	1:A:2357:SER:HB2	1.79	0.48
1:A:2842:ASP:OD1	1:A:2842:ASP:N	2.38	0.48
1:A:2471:LEU:HD12	1:A:2523:TRP:HB3	1.95	0.48
1:A:2646:ARG:HH11	1:A:2646:ARG:HG3	1.79	0.47
1:A:2171:ASP:OD1	1:A:2209:ARG:NH1	2.46	0.47
1:A:2951:GLU:OE1	1:A:2951:GLU:N	2.39	0.47
1:A:2441:VAL:HG23	1:A:2482:LEU:HD11	1.96	0.47
1:A:2929:ALA:HB2	1:A:2992:PHE:HZ	1.80	0.47
1:A:3394:MET:HE3	1:A:3398:MET:HE3	1.97	0.47
1:A:2472:THR:O	1:A:2524:VAL:HA	2.14	0.46
1:A:2933:VAL:O	1:A:2936:ILE:HG22	2.14	0.46
1:A:2941:THR:O	1:A:2944:ILE:HG12	2.14	0.46
1:A:3379:TRP:NE1	1:A:3398:MET:SD	2.88	0.46
1:A:2491:LEU:HD21	1:A:2543:ARG:NH1	2.29	0.46
1:A:2792:LEU:HA	1:A:2827:PHE:HE1	1.79	0.46
1:A:2105:ASP:N	1:A:2105:ASP:OD1	2.48	0.46
1:A:2730:VAL:HG12	1:A:2732:MET:HG3	1.98	0.46
1:A:2319:LYS:NZ	1:A:2351:GLN:O	2.43	0.46
1:A:2728:LEU:HD22	1:A:3469:GLU:HG2	1.98	0.46
1:A:2424:LYS:HG2	1:A:2559:LEU:HD12	1.97	0.46
1:A:2700:LEU:HD22	1:A:2707:VAL:HG11	1.97	0.46
1:A:3468:ARG:HA	1:A:3468:ARG:HD3	1.80	0.46
1:A:3483:ASP:OD1	1:A:3484:HIS:ND1	2.47	0.46
1:A:2445:PHE:HA	1:A:2449:THR:HG21	1.97	0.45
1:A:2856:LEU:HD21	1:A:2877:PHE:HB2	1.97	0.45
1:A:2319:LYS:HB3	1:A:2353:LEU:HD22	1.98	0.45
1:A:2536:ASN:HB2	1:A:2543:ARG:HE	1.81	0.45
1:A:2253:ILE:HD11	1:A:2314:ILE:HD12	1.99	0.45
1:A:2375:ILE:HD11	1:A:2395:ILE:HG13	1.98	0.45
1:A:2247:ASP:OD1	1:A:2248:LYS:N	2.50	0.45
1:A:2249:LEU:O	1:A:2253:ILE:HG12	2.17	0.45
1:A:2241:LEU:HD13	1:A:2299:ARG:HD2	1.99	0.45
1:A:2095:ASP:N	1:A:2095:ASP:OD1	2.51	0.44
1:A:2200:ASP:N	1:A:2200:ASP:OD1	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3370:LEU:HD21	1:A:3399:ASN:HD22	1.83	0.44
1:A:3512:ARG:HH22	2:A:4201:ZG7:C25	2.31	0.44
1:A:2792:LEU:HD11	1:A:2830:ARG:HE	1.84	0.43
1:A:2977:TYR:CE2	1:A:2987:ARG:HB2	2.53	0.43
1:A:3339:GLU:H	1:A:3339:GLU:CD	2.26	0.43
1:A:2282:ASN:HB3	1:A:2552:ARG:HG3	1.99	0.43
1:A:2576:LYS:HA	1:A:2589:THR:HG21	2.00	0.43
1:A:2844:PHE:O	1:A:2849:TYR:HB2	2.18	0.43
1:A:2392:ILE:HG13	1:A:2570:ILE:HG23	2.01	0.43
1:A:2787:HIS:HB3	1:A:3461:ILE:HG23	2.01	0.43
1:A:2858:ASN:OD1	1:A:2861:ARG:NH2	2.53	0.42
1:A:2992:PHE:HD2	1:A:2993:ILE:HD12	1.84	0.42
1:A:2315:THR:O	1:A:2319:LYS:HG3	2.20	0.42
1:A:2389:ASP:OD1	1:A:2389:ASP:N	2.52	0.42
1:A:2239:ASN:OD1	1:A:2240:LYS:N	2.53	0.42
1:A:2684:GLN:N	1:A:2684:GLN:OE1	2.54	0.41
1:A:2689:ILE:HD12	1:A:2689:ILE:HG23	1.76	0.41
1:A:2423:GLY:N	2:A:4202:ZG7:O27	2.48	0.41
1:A:2125:TRP:CZ3	1:A:2127:ASP:HA	2.56	0.41
1:A:3415:ILE:HD11	1:A:3497:HIS:ND1	2.36	0.41
1:A:2956:GLU:HG3	1:A:2957:PRO:HD2	2.02	0.41
1:A:2240:LYS:NZ	1:A:2299:ARG:HH22	2.18	0.41
1:A:2786:ILE:HG13	1:A:2821:ASN:HD22	1.85	0.41
1:A:3374:ASP:OD1	1:A:3375:GLU:N	2.52	0.41
1:A:2361:ILE:HG22	1:A:2368:PHE:CE1	2.56	0.40
1:A:2749:LEU:HD23	1:A:2749:LEU:HA	1.91	0.40
1:A:3408:LEU:HD23	1:A:3517:HIS:ND1	2.36	0.40
1:A:2119:LEU:HD21	1:A:2126:ARG:NE	2.37	0.40
1:A:2167:ASN:OD1	1:A:2168:SER:N	2.55	0.40
1:A:2867:LEU:HD22	1:A:2872:GLU:HB3	2.03	0.40
1:A:2231:ASN:O	1:A:2235:GLU:HG3	2.22	0.40
1:A:3429:LEU:HD21	1:A:3439:ARG:HD2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM

entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	1188/2661 (45%)	1103 (93%)	81 (7%)	4 (0%)	36 67

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3413	HIS
1	A	2984	VAL
1	A	2959	GLN
1	A	2985	ASN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	1084/2406 (45%)	1063 (98%)	21 (2%)	50 68

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2059	GLN
1	A	2095	ASP
1	A	2131	THR
1	A	2262	LEU
1	A	2295	ILE
1	A	2361	ILE
1	A	2390	ILE
1	A	2476	LYS
1	A	2702	LEU
1	A	2709	LYS
1	A	2728	LEU
1	A	2842	ASP
1	A	2942	ASP

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Mol	Chain	Res	Type
1	A	2945	VAL
1	A	2958	ILE
1	A	2959	GLN
1	A	2960	THR
1	A	3385	ASP
1	A	3402	ASP
1	A	3409	ASP
1	A	3513	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2363	ASN
1	A	2481	ASN
1	A	2500	GLN
1	A	2508	GLN
1	A	2536	ASN
1	A	2634	ASN
1	A	2714	ASN
1	A	2741	HIS
1	A	2783	GLN
1	A	2854	ASN
1	A	2959	GLN
1	A	3442	ASN
1	A	3472	HIS
1	A	3517	HIS
1	A	3521	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ZG7	A	4201	-	30,34,34	2.44	6 (20%)	35,51,51	1.78	7 (20%)
3	ATP	A	4203	4	32,33,33	1.30	6 (18%)	48,52,52	1.93	11 (22%)
2	ZG7	A	4202	-	30,34,34	2.61	6 (20%)	35,51,51	1.86	6 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ZG7	A	4201	-	-	3/10/22/22	0/5/5/5
3	ATP	A	4203	4	-	7/22/38/38	0/3/3/3
2	ZG7	A	4202	-	-	1/10/22/22	0/5/5/5

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	4202	ZG7	C14-N15	8.82	1.53	1.35
2	A	4201	ZG7	C14-N15	8.26	1.52	1.35
2	A	4202	ZG7	C16-N15	7.50	1.45	1.32
2	A	4201	ZG7	C16-N15	6.93	1.44	1.32
2	A	4201	ZG7	C29-C28	4.87	1.53	1.43
2	A	4202	ZG7	C29-C28	4.76	1.53	1.43
2	A	4202	ZG7	O27-C26	-4.58	1.17	1.31
2	A	4201	ZG7	O27-C26	-3.90	1.19	1.31
3	A	4203	ATP	C5-C4	3.65	1.45	1.39
2	A	4201	ZG7	C11-N12	2.89	1.37	1.32
2	A	4202	ZG7	C11-N12	2.78	1.37	1.32
3	A	4203	ATP	C5-C6	2.63	1.48	1.41
3	A	4203	ATP	C5-N7	-2.56	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	4203	ATP	C4-N9	-2.45	1.32	1.37
2	A	4202	ZG7	BR22-C21	2.35	1.94	1.90
3	A	4203	ATP	C8-N9	-2.31	1.33	1.37
2	A	4201	ZG7	BR22-C21	2.18	1.94	1.90
3	A	4203	ATP	C8-N7	2.07	1.35	1.31

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	4203	ATP	C5-C4-N3	-6.03	118.41	126.72
2	A	4202	ZG7	C16-N15-C14	5.87	121.81	115.71
2	A	4201	ZG7	C16-N15-C14	4.64	120.53	115.71
3	A	4203	ATP	N3-C4-N9	4.41	134.67	127.17
2	A	4202	ZG7	N13-C14-N15	-4.24	118.04	122.20
3	A	4203	ATP	C4-C5-N7	-4.23	105.74	110.58
2	A	4201	ZG7	N13-C14-N15	-4.12	118.16	122.20
2	A	4201	ZG7	C10-C08-C05	3.94	123.18	118.58
3	A	4203	ATP	C2-N3-C4	3.92	121.41	111.83
2	A	4202	ZG7	C09-C08-C10	3.81	60.15	58.45
2	A	4202	ZG7	C28-C14-N15	3.46	137.04	131.48
2	A	4201	ZG7	C09-C08-C10	3.31	59.93	58.45
3	A	4203	ATP	N3-C2-N1	-3.22	123.70	128.58
3	A	4203	ATP	C4-N9-C8	3.16	109.05	105.74
3	A	4203	ATP	C5-N7-C8	3.08	108.28	103.45
3	A	4203	ATP	N9-C8-N7	-2.77	110.01	113.94
3	A	4203	ATP	C3'-C2'-C1'	2.73	106.62	101.46
3	A	4203	ATP	C6-C5-N7	2.66	137.23	132.09
3	A	4203	ATP	O5'-C5'-C4'	2.61	117.89	108.99
2	A	4201	ZG7	C09-C08-C05	2.59	121.61	118.58
2	A	4201	ZG7	C10-C09-C08	-2.52	59.63	60.78
2	A	4201	ZG7	C28-C14-N15	2.50	135.49	131.48
2	A	4202	ZG7	C19-C20-C21	2.34	120.71	118.07
2	A	4202	ZG7	C06-C05-C04	2.28	121.43	118.03

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	4202	ZG7	C16-C17-O18-C19
3	A	4203	ATP	PB-O3B-PG-O3G
2	A	4201	ZG7	C06-C05-C08-C10
3	A	4203	ATP	PB-O3B-PG-O1G

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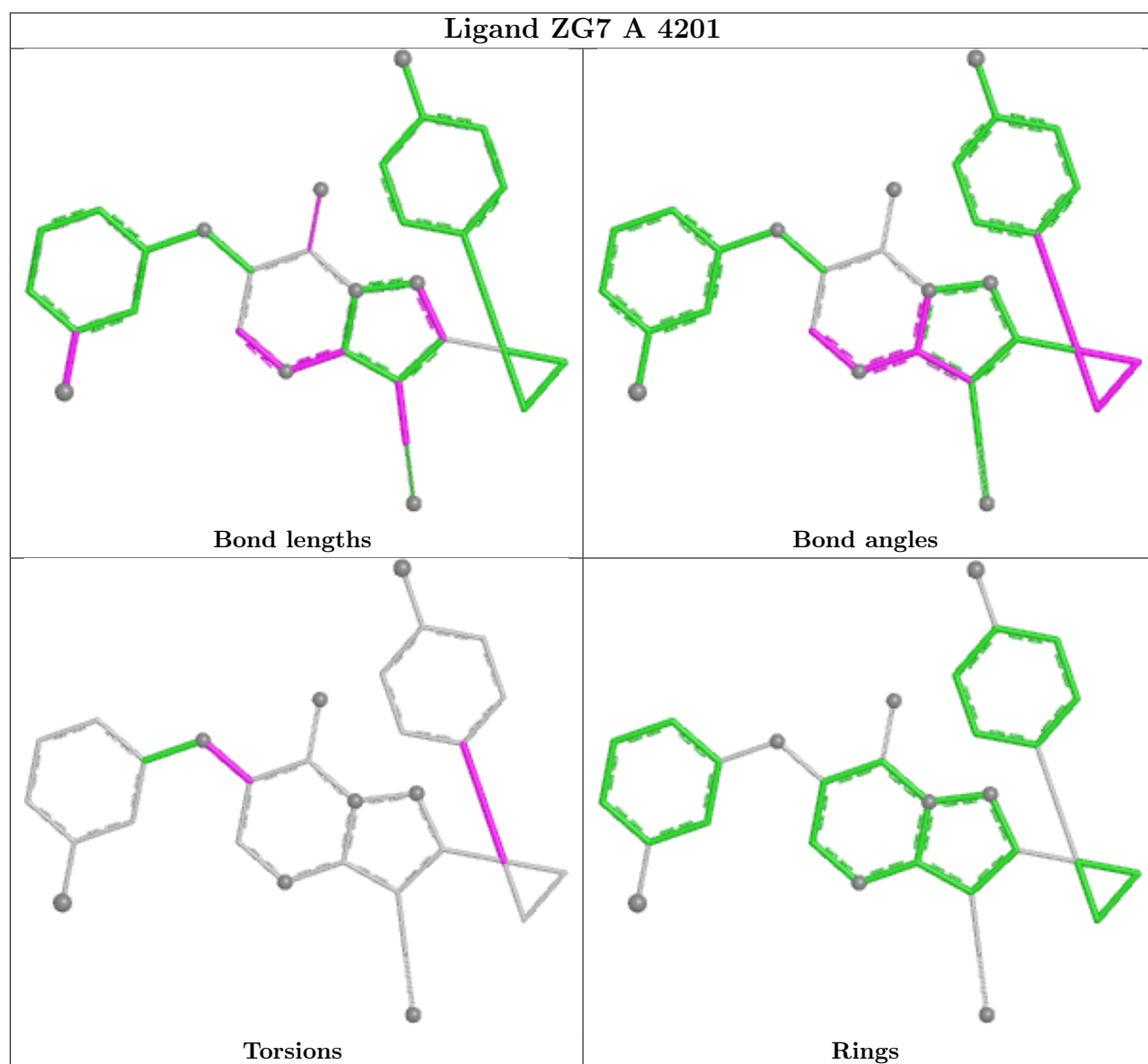
Mol	Chain	Res	Type	Atoms
3	A	4203	ATP	PG-O3B-PB-O1B
3	A	4203	ATP	PB-O3A-PA-O1A
2	A	4201	ZG7	C26-C17-O18-C19
3	A	4203	ATP	PB-O3A-PA-O2A
3	A	4203	ATP	C4'-C5'-O5'-PA
2	A	4201	ZG7	C04-C05-C08-C10
3	A	4203	ATP	PG-O3B-PB-O2B

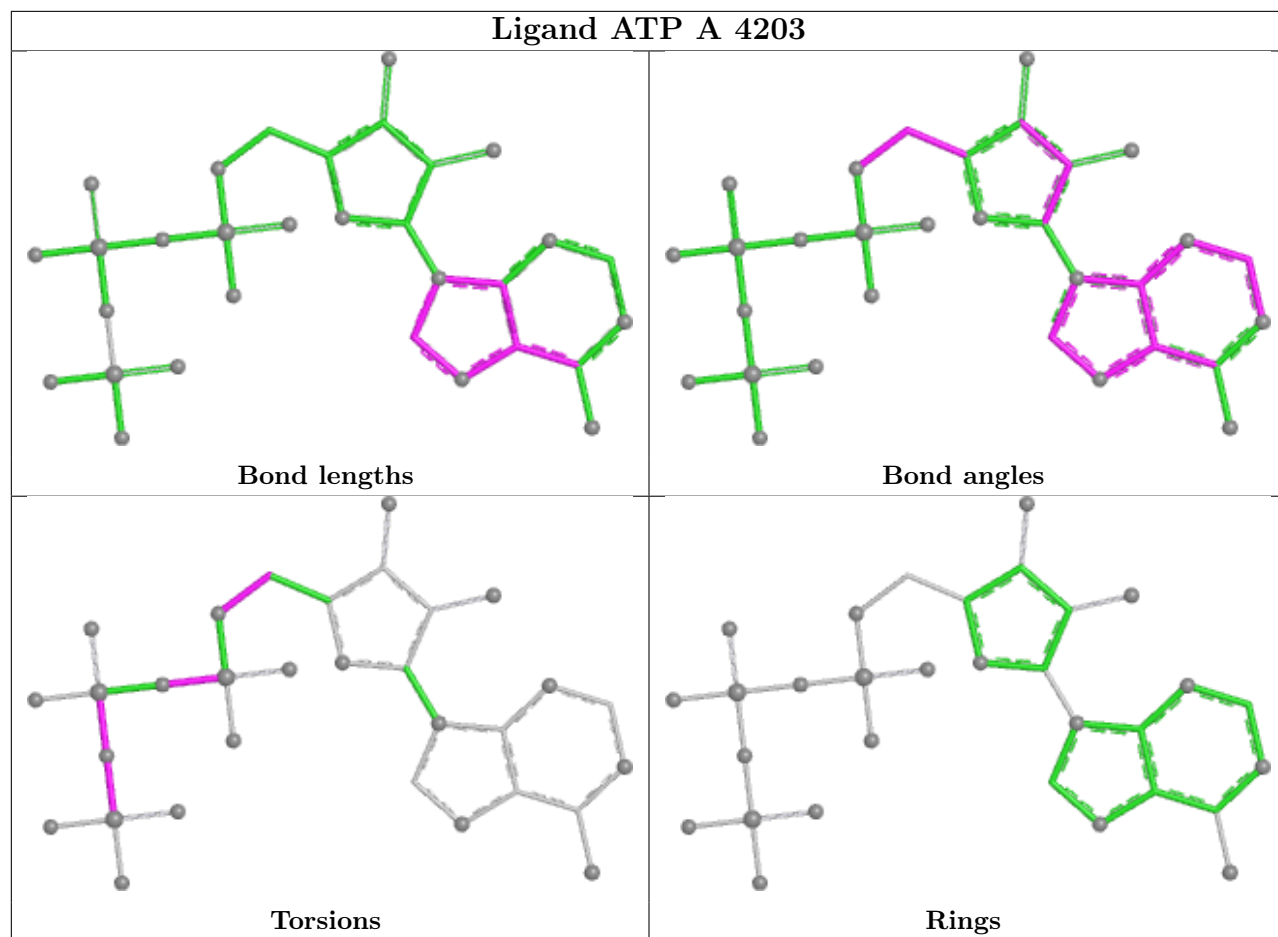
There are no ring outliers.

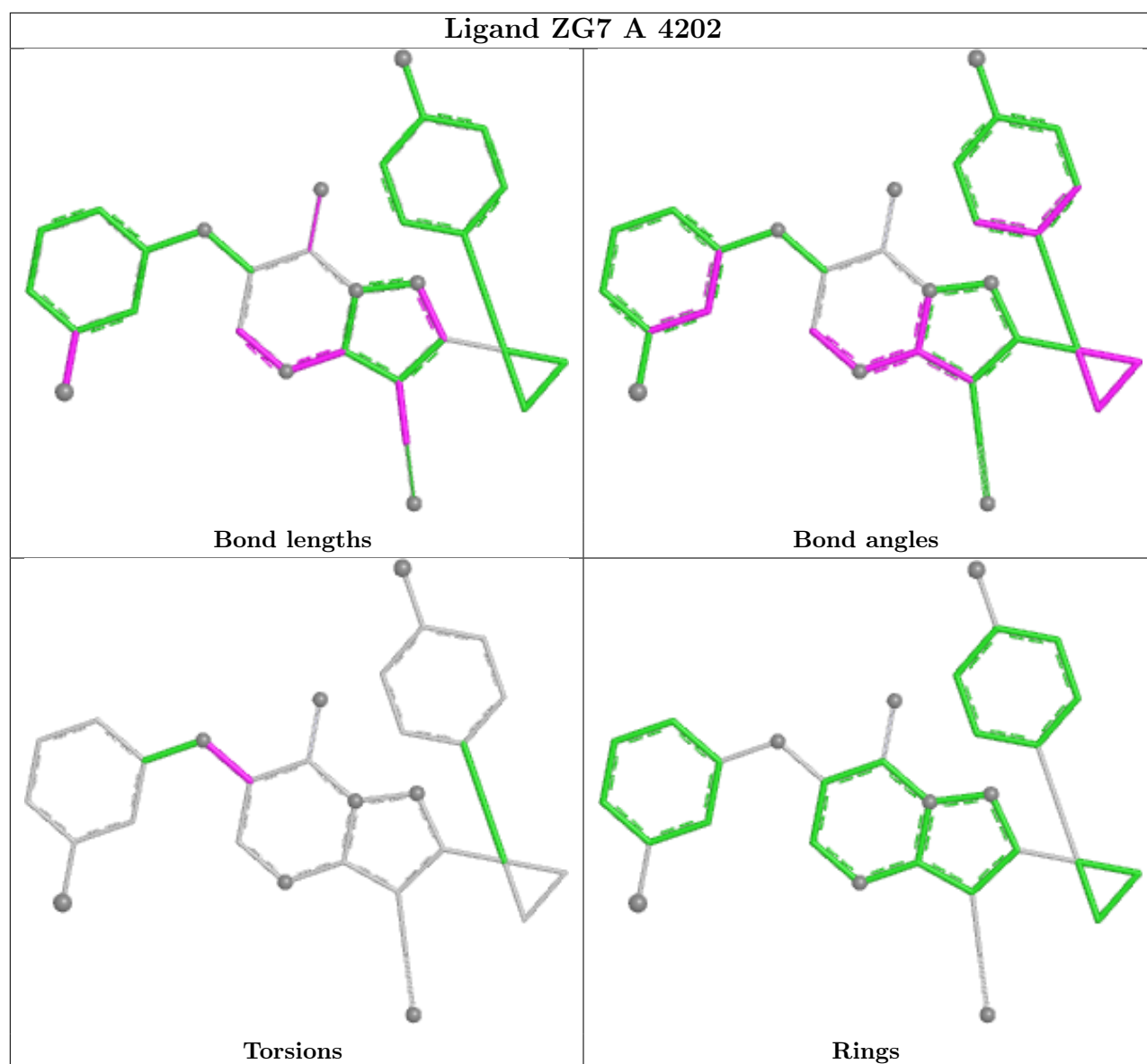
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	4201	ZG7	2	0
3	A	4203	ATP	1	0
2	A	4202	ZG7	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







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.

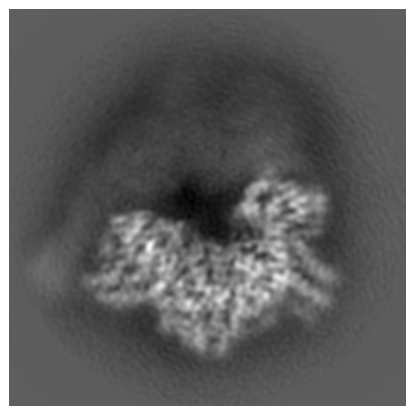
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23838. These allow visual inspection of the internal detail of the map and identification of artifacts.

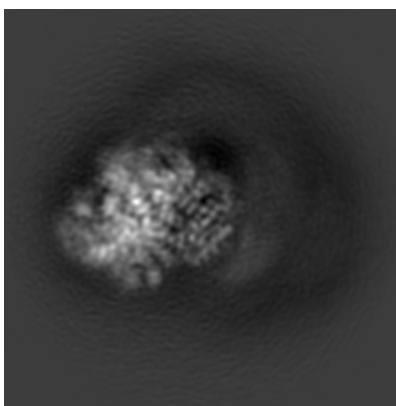
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

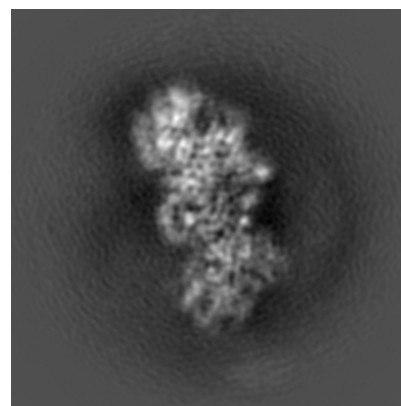
6.1.1 Primary map



X

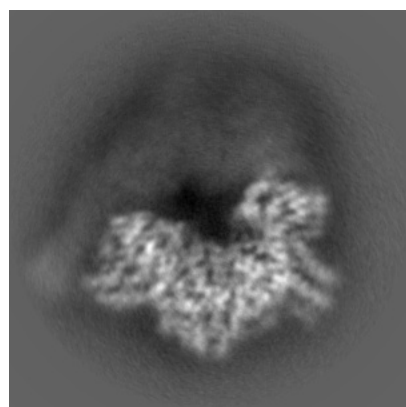


Y

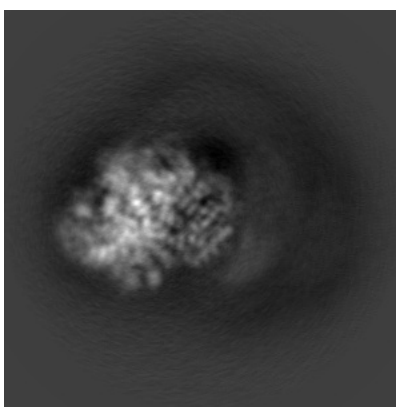


Z

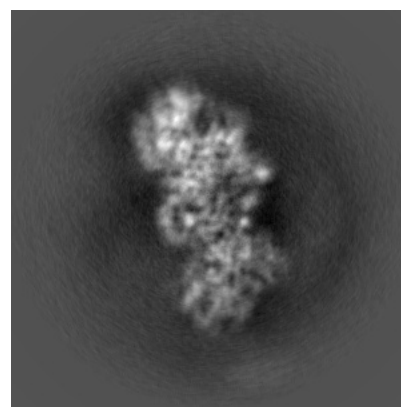
6.1.2 Raw map



X



Y

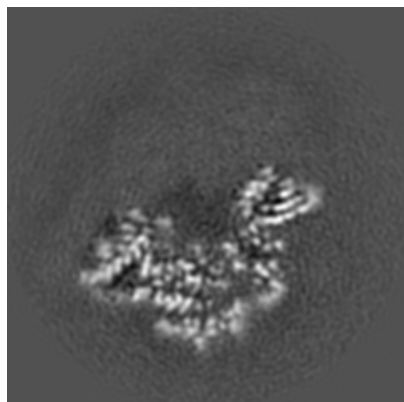


Z

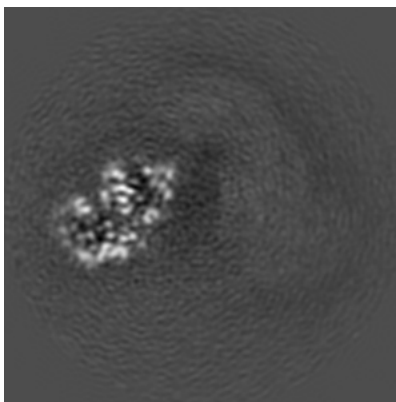
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

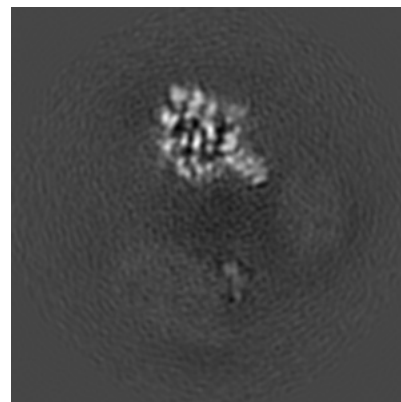
6.2.1 Primary map



X Index: 184

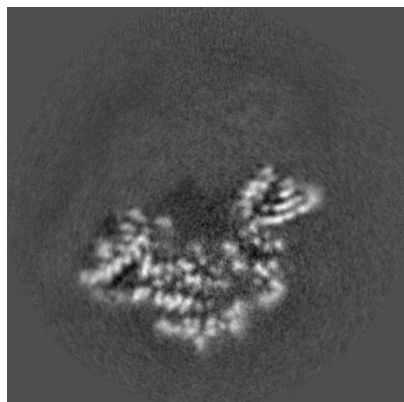


Y Index: 184

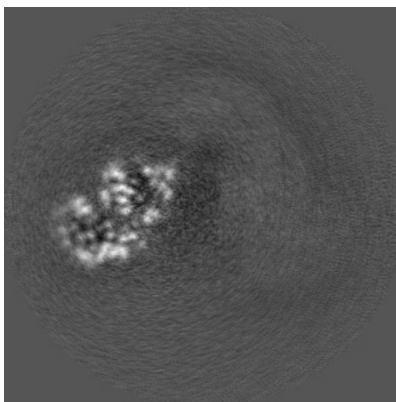


Z Index: 184

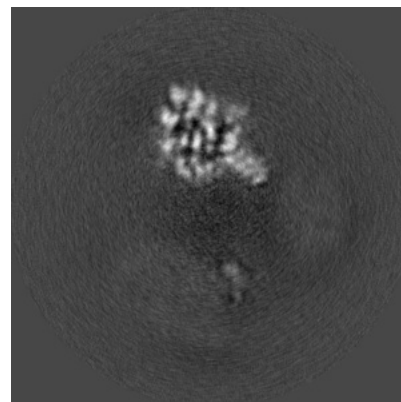
6.2.2 Raw map



X Index: 184



Y Index: 184

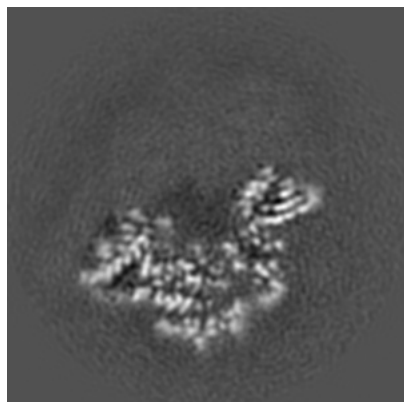


Z Index: 184

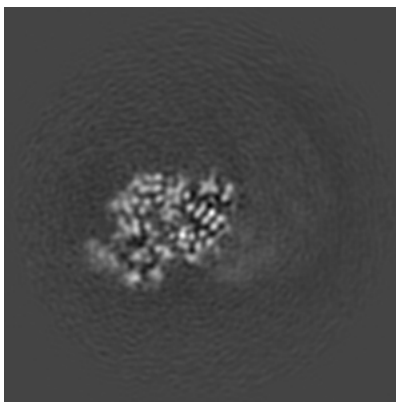
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

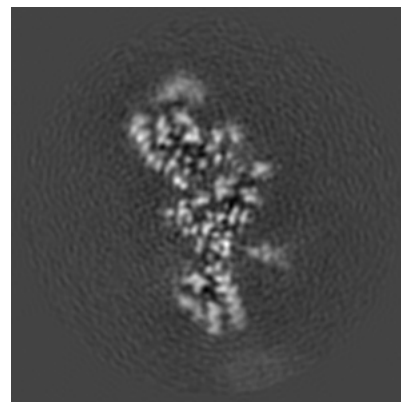
6.3.1 Primary map



X Index: 184

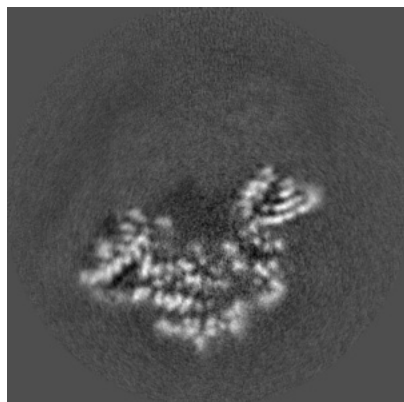


Y Index: 247

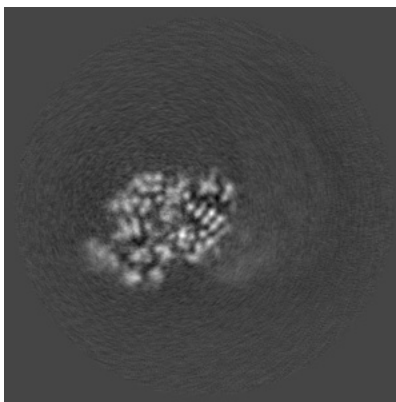


Z Index: 120

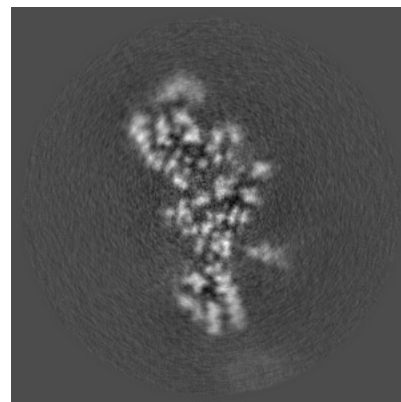
6.3.2 Raw map



X Index: 184



Y Index: 246

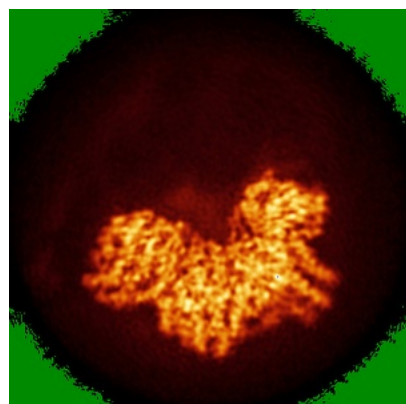


Z Index: 120

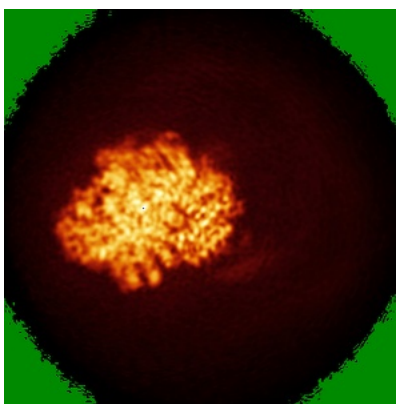
The images above show the largest variance slices of the map in three orthogonal directions.

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

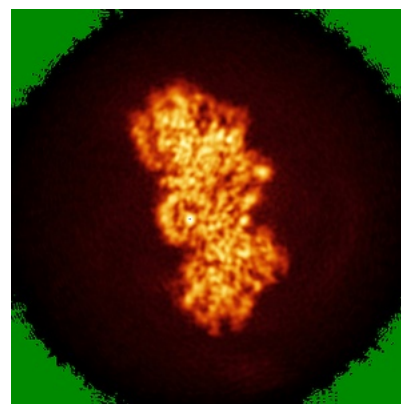
6.4.1 Primary map



X

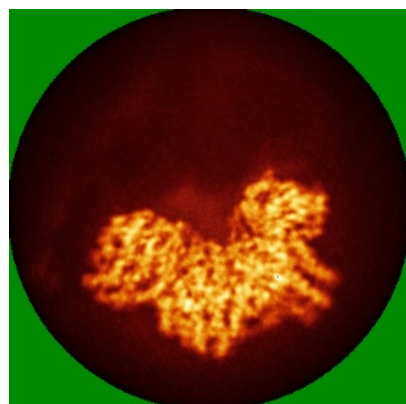


Y

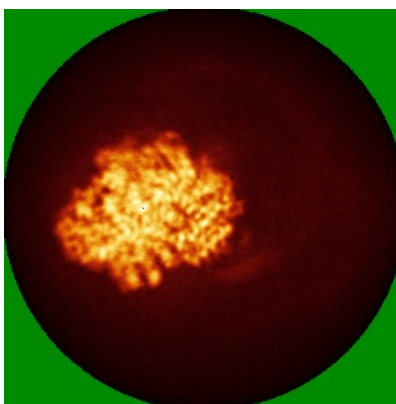


Z

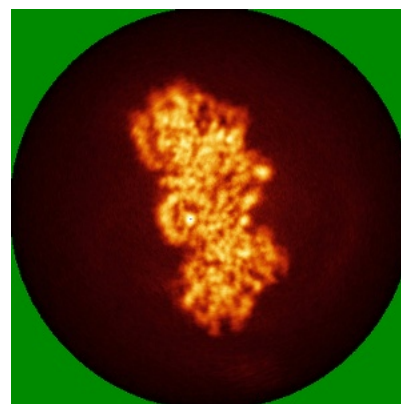
6.4.2 Raw map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.004. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

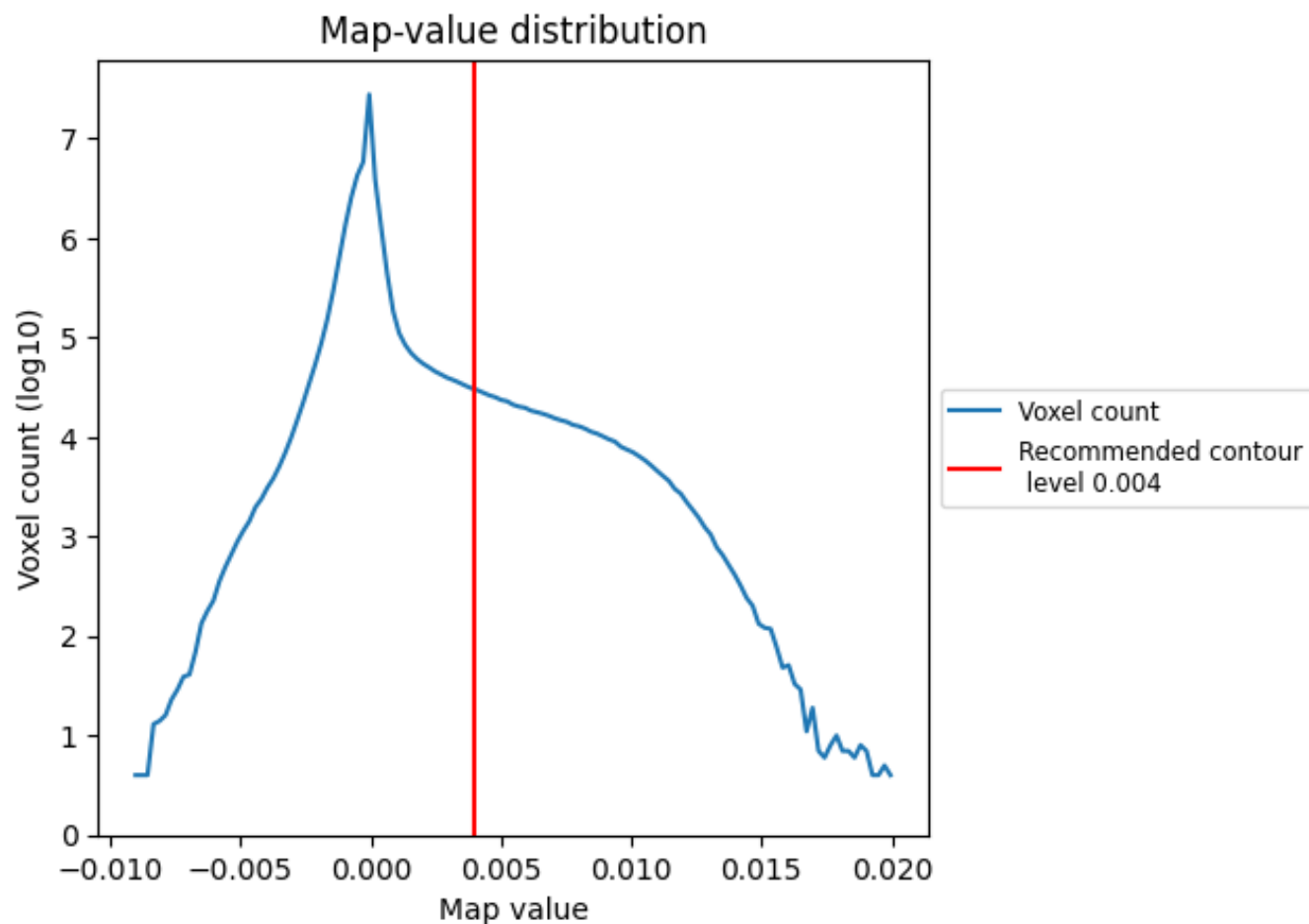
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

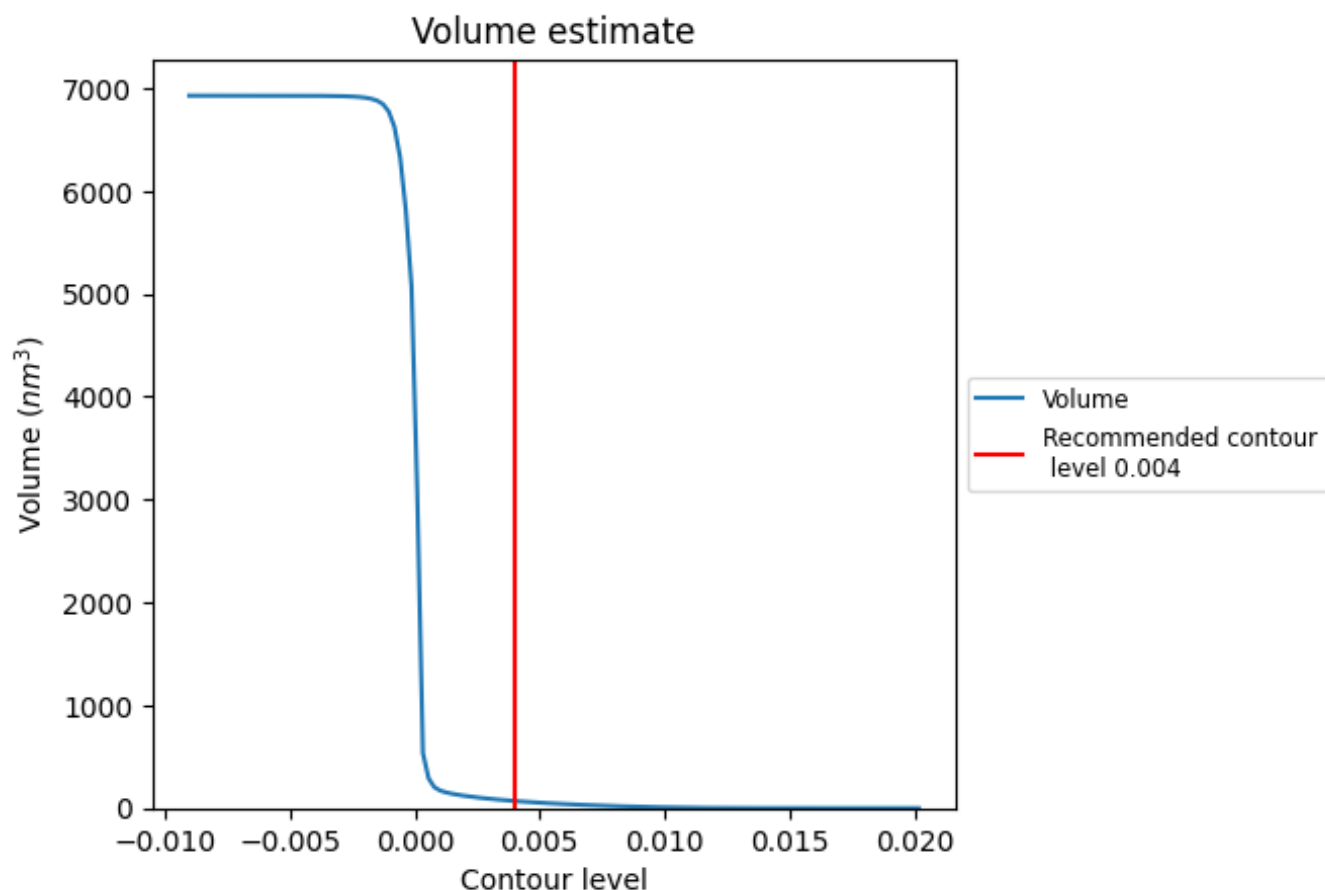
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

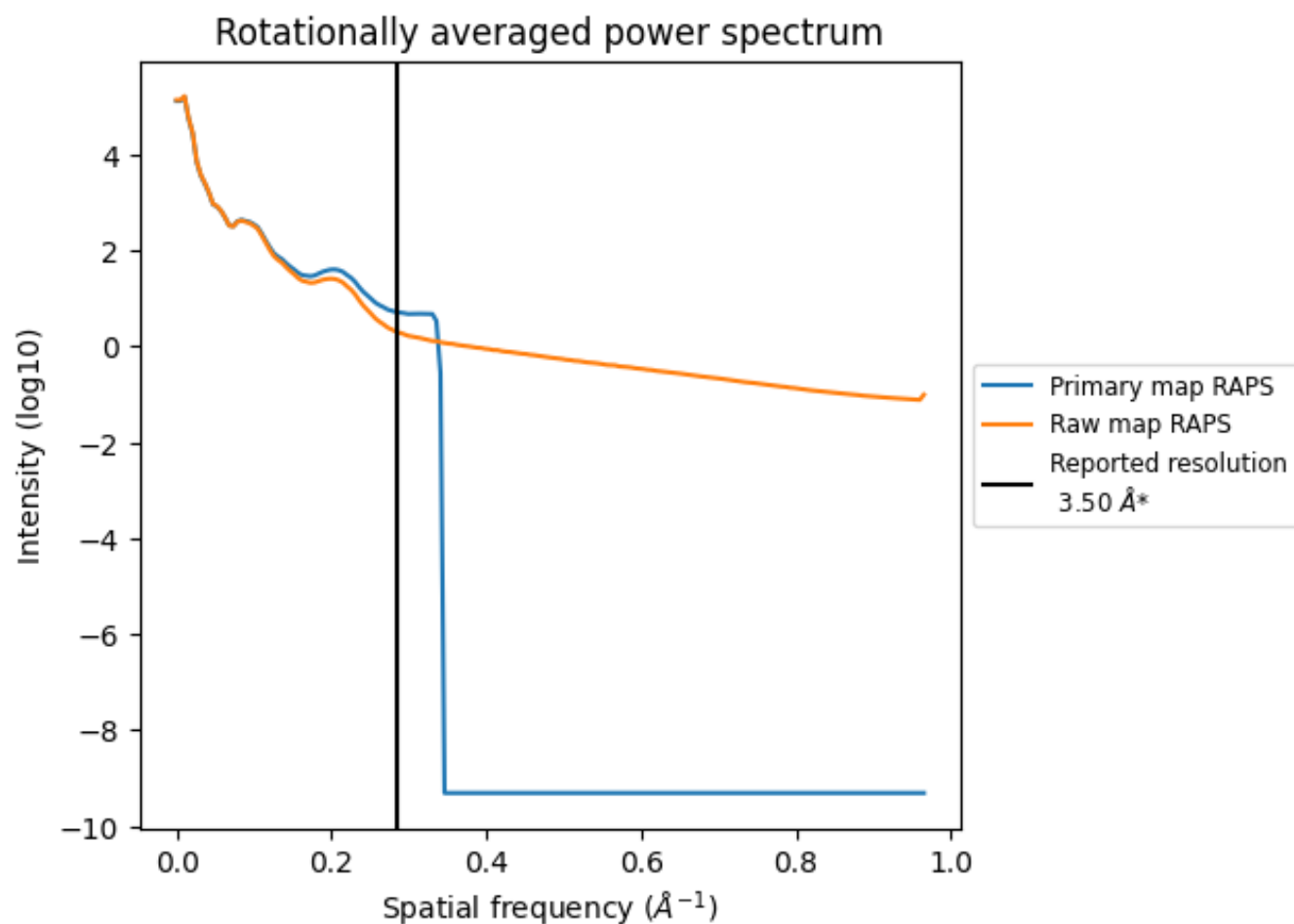
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 69 nm^3 ; this corresponds to an approximate mass of 63 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 ⓘ

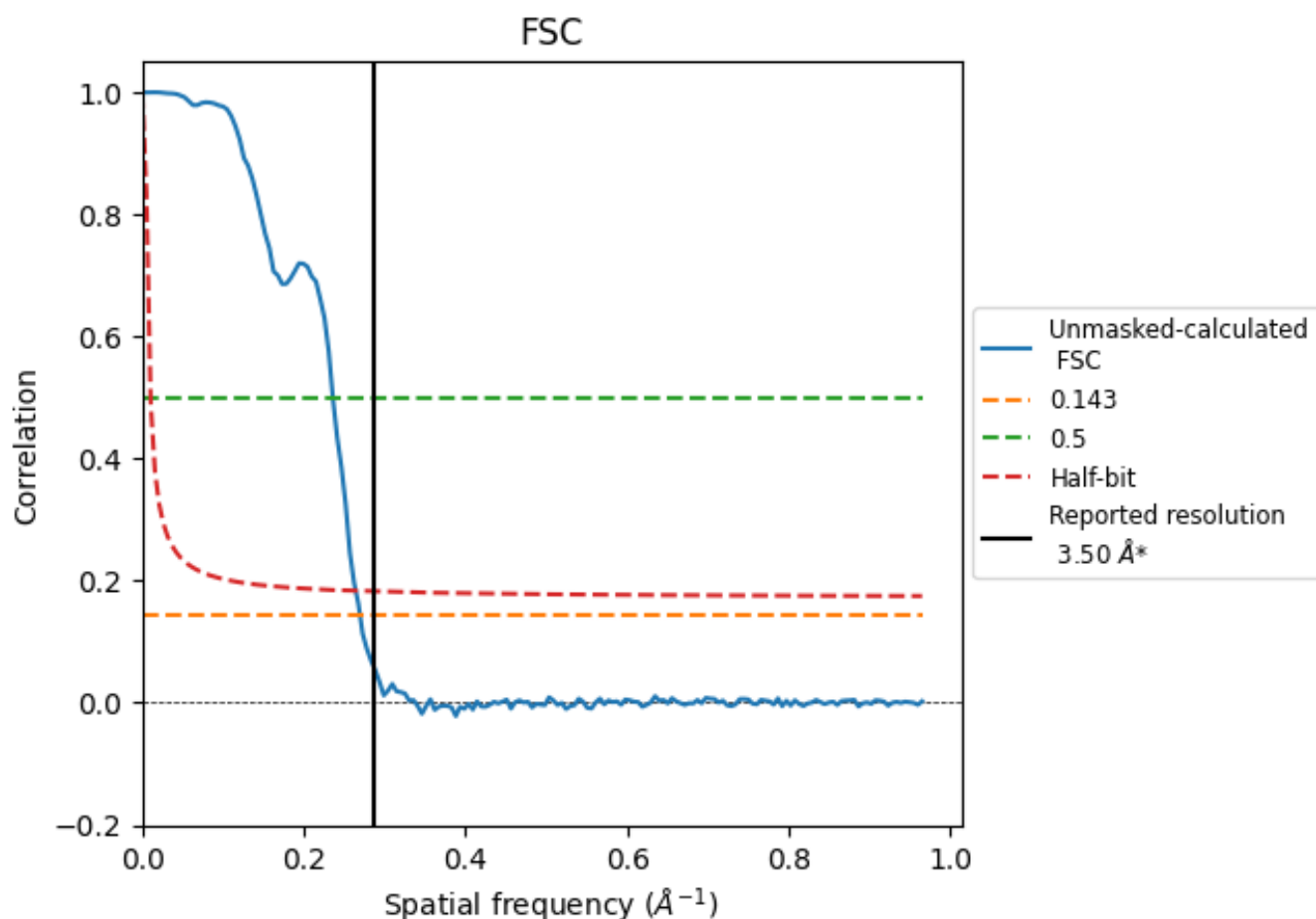


*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.286 Å⁻¹

8.2 Resolution estimates [i](#)

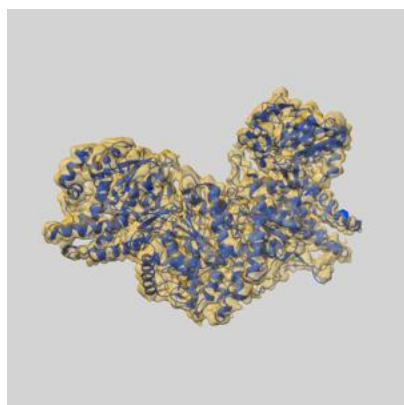
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.50	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.71	4.24	3.78

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

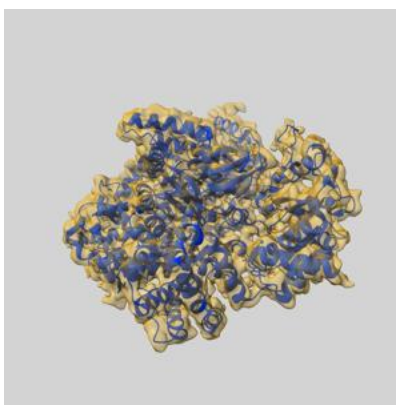
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-23838 and PDB model 7MI3. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

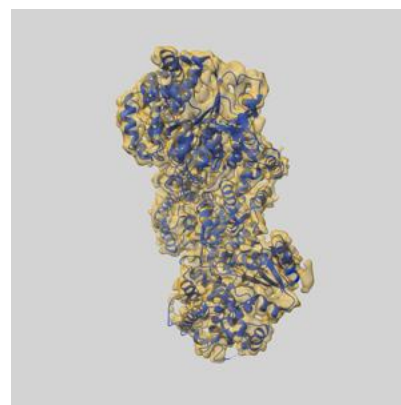
9.1 Map-model overlay [i](#)



X



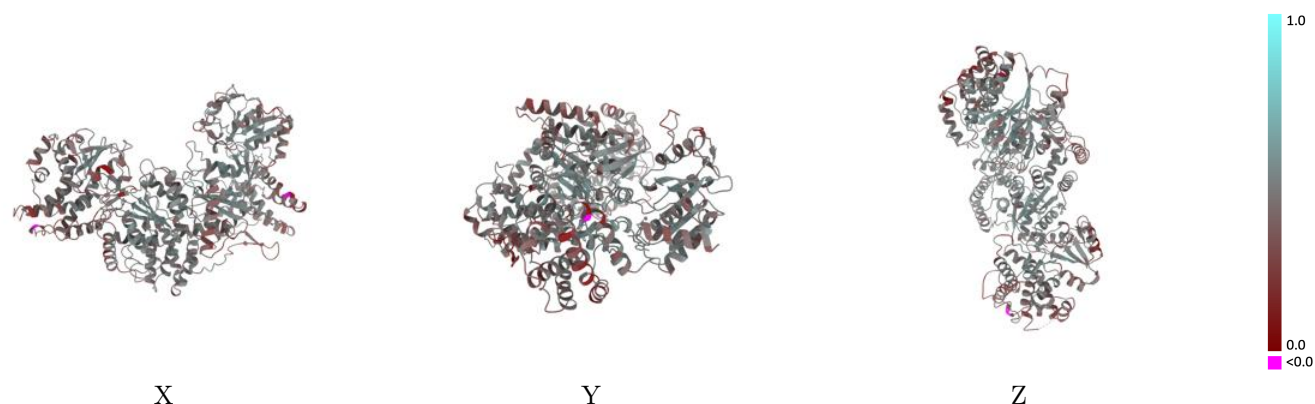
Y



Z

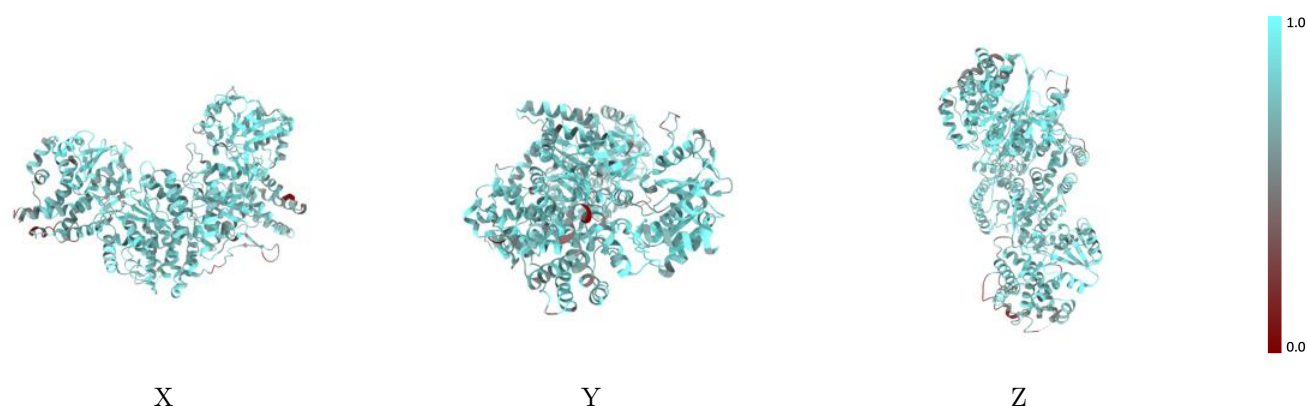
The images above show the 3D surface view of the map at the recommended contour level 0.004 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



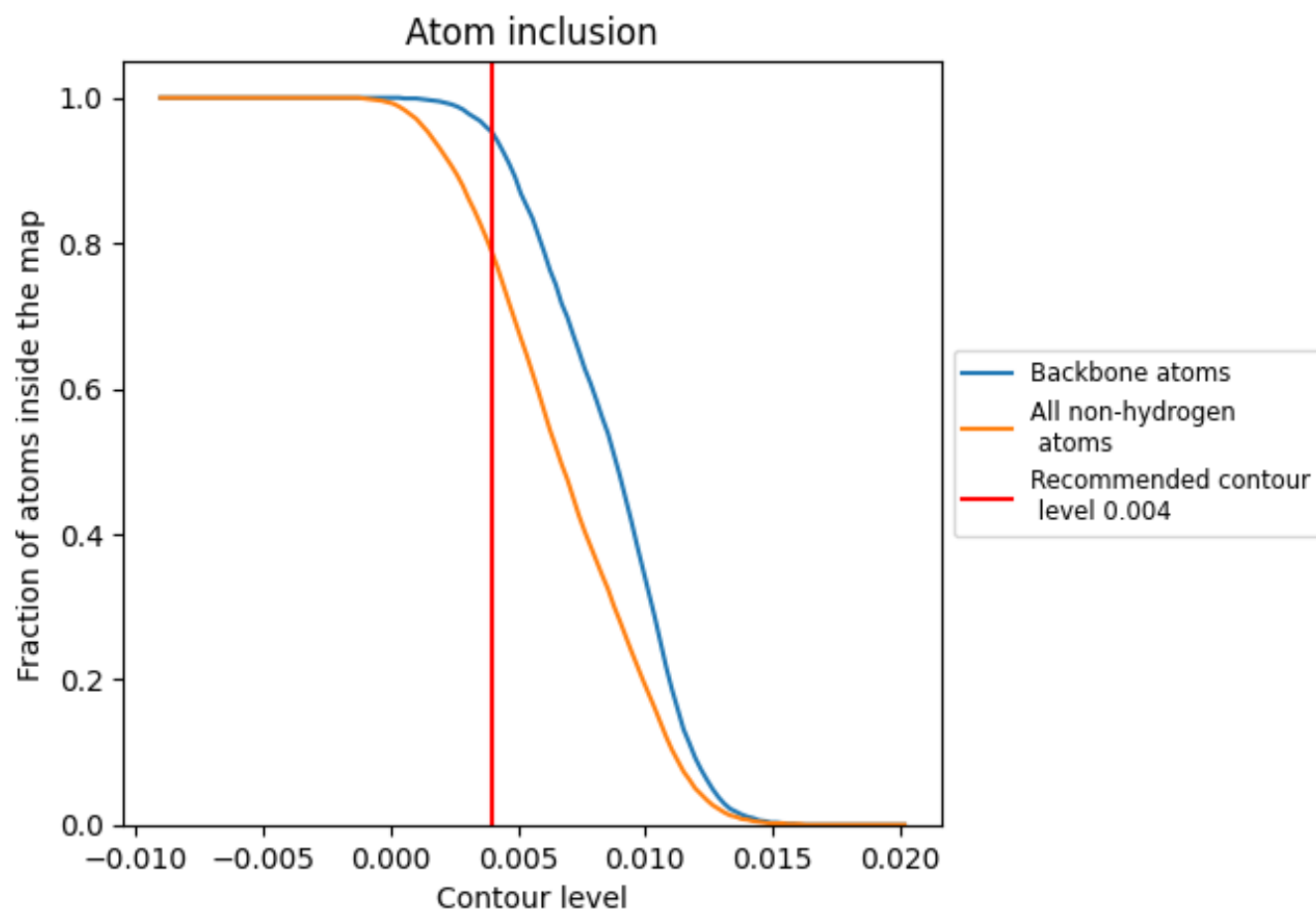
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.004).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% 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 (0.004) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7860	0.4410
A	0.7860	0.4410

