



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 05:46 AM UTC

PDB ID : 9O58 / pdb_00009o58
EMDB ID : EMD-70127
Title : Zymogen ADAM17-iRhom2 complex bound by the MEDI3622 Fab
Authors : Slone, C.E.; Maciag, J.J.
Deposited on : 2025-04-09
Resolution : 3.53 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

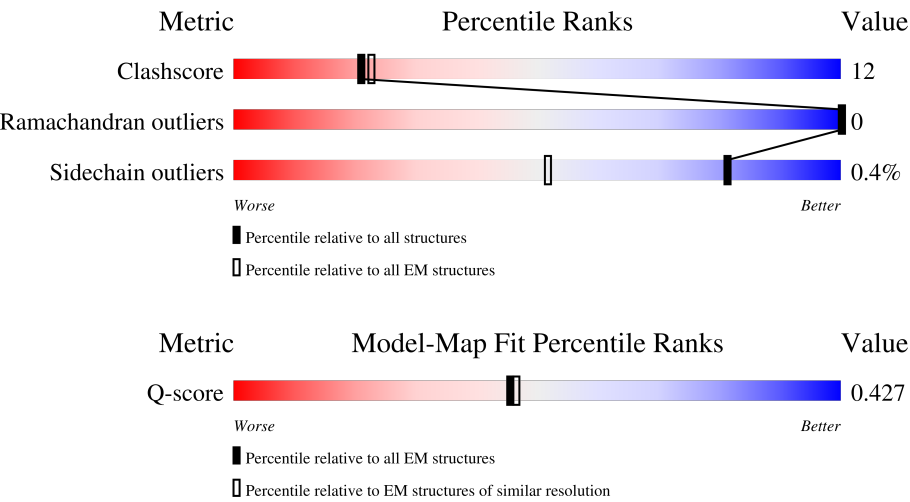
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.53 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	12947 (3.03 - 4.02)

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 <=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	807	5% 58% 21% 20%
2	B	741	50% 16% 34%
3	C	214	36% 74% 26%
4	D	221	29% 73% 27%

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Mol	Chain	Length	Quality of chain
5	E	2	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 12422 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 Disintegrin and metalloproteinase domain-containing protein 17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	642	5082	3172	873	993	44	0	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	406	ALA	GLU	conflict	UNP P78536
A	825	GLU	-	expression tag	UNP P78536
A	826	GLN	-	expression tag	UNP P78536
A	827	LYS	-	expression tag	UNP P78536
A	828	LEU	-	expression tag	UNP P78536
A	829	ILE	-	expression tag	UNP P78536
A	830	SER	-	expression tag	UNP P78536
A	831	GLU	-	expression tag	UNP P78536
A	832	GLU	-	expression tag	UNP P78536
A	833	ASP	-	expression tag	UNP P78536
A	834	LEU	-	expression tag	UNP P78536

- Molecule 2 is a protein called GFP-iRhom2 fusion protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	490	3941	2564	666	685	26	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	116	MET	-	initiating methionine	UNP P42212
B	117	VAL	-	expression tag	UNP P42212
B	162	LEU	PHE	conflict	UNP P42212
B	180	LEU	PHE	conflict	UNP P42212
B	181	GLY	SER	conflict	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
B	184	LEU	VAL	conflict	UNP P42212
B	188	ALA	SER	conflict	UNP P42212
B	269	THR	MET	conflict	UNP P42212
B	279	ALA	VAL	conflict	UNP P42212
B	291	GLY	SER	conflict	UNP P42212
B	319	TYR	THR	conflict	UNP P42212
B	322	LYS	ALA	conflict	UNP P42212
B	347	LEU	HIS	conflict	UNP P42212
B	355	ASP	-	linker	UNP P42212
B	356	TYR	-	linker	UNP P42212
B	357	LYS	-	linker	UNP P42212
B	358	ASP	-	linker	UNP P42212
B	359	ASP	-	linker	UNP P42212
B	360	ASP	-	linker	UNP P42212
B	361	LYS	-	linker	UNP P42212
B	362	SER	-	linker	UNP P42212
B	363	GLY	-	linker	UNP P42212

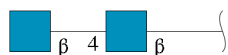
- Molecule 3 is a protein called MEDI3622 Fab Light Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	214	Total	C	N	O	S	0	0
			1639	1023	275	335	6		

- Molecule 4 is a protein called MEDI3622 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	221	Total	C	N	O	S	0	0
			1646	1035	278	325	8		

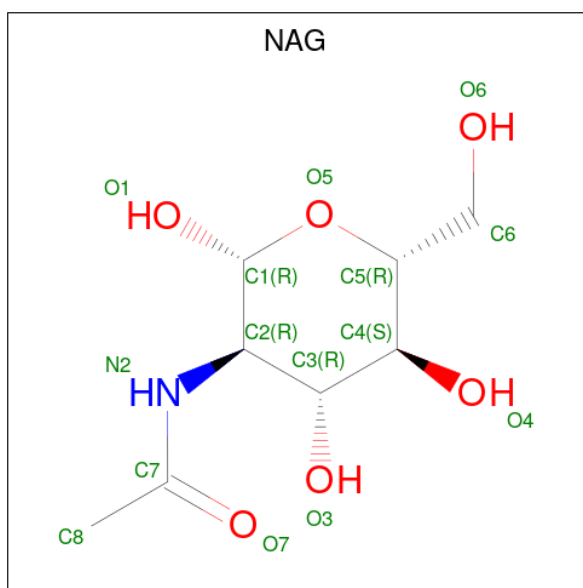
- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				AltConf	Trace
5	E	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula:

C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	A	1	Total	C	N	O	0
			14	8	1	5	
6	A	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
7	A	1	Total	Ca	0
			1	1	

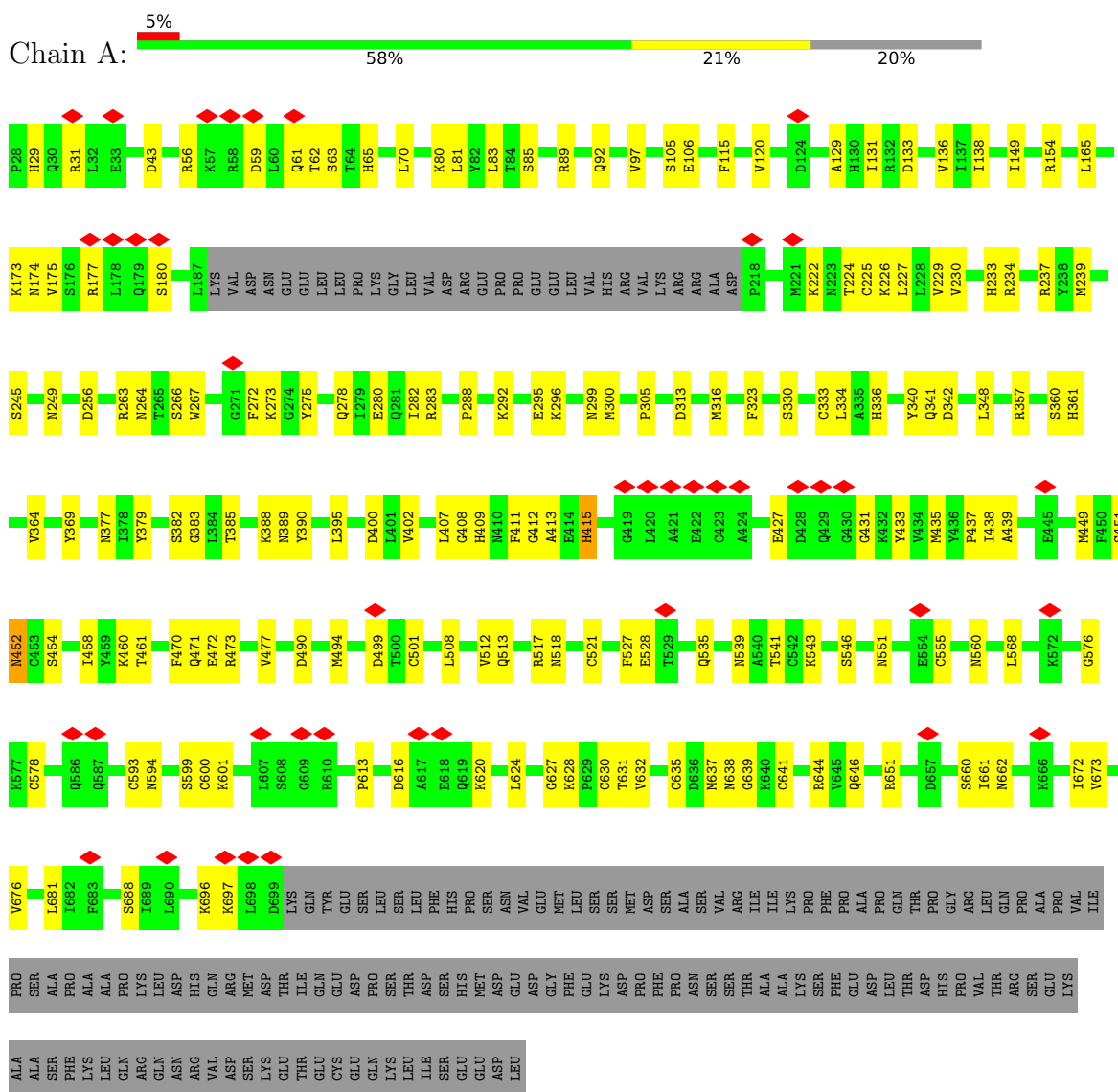
- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
8	A	1	Total	Zn	0
			1	1	

3 Residue-property plots

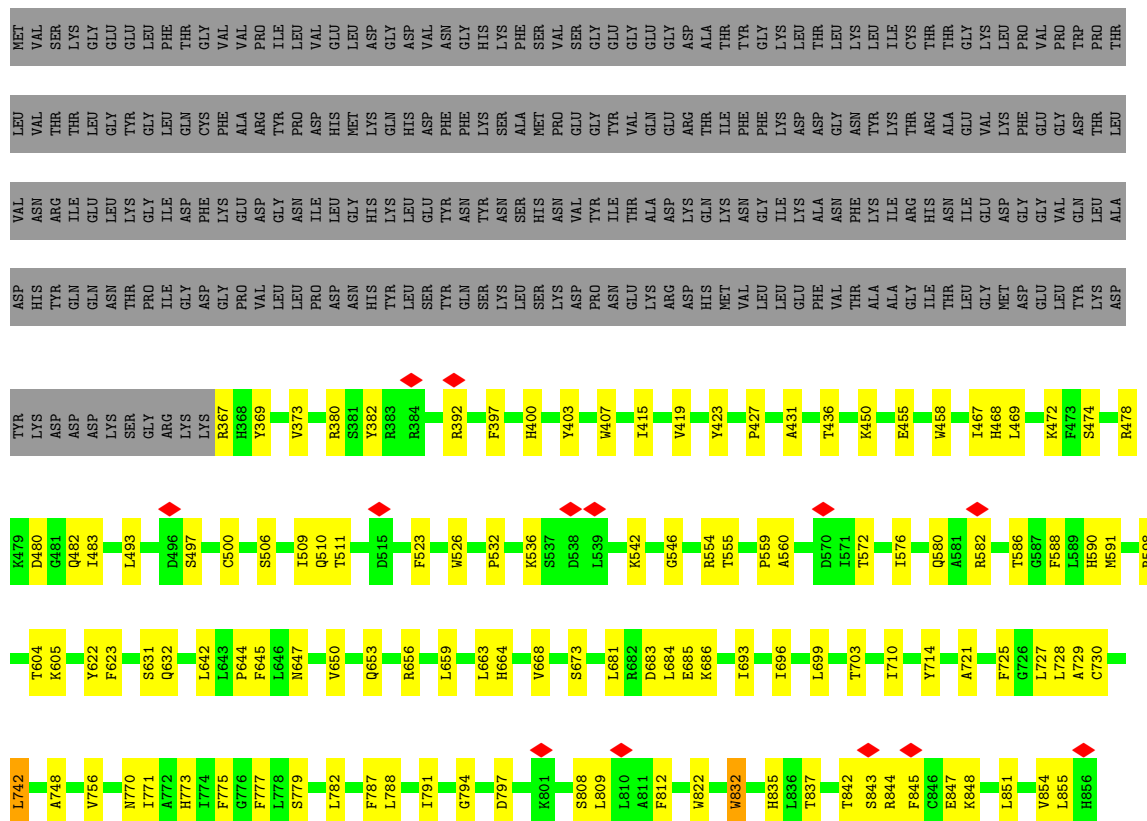
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Disintegrin and metalloproteinase domain-containing protein 17

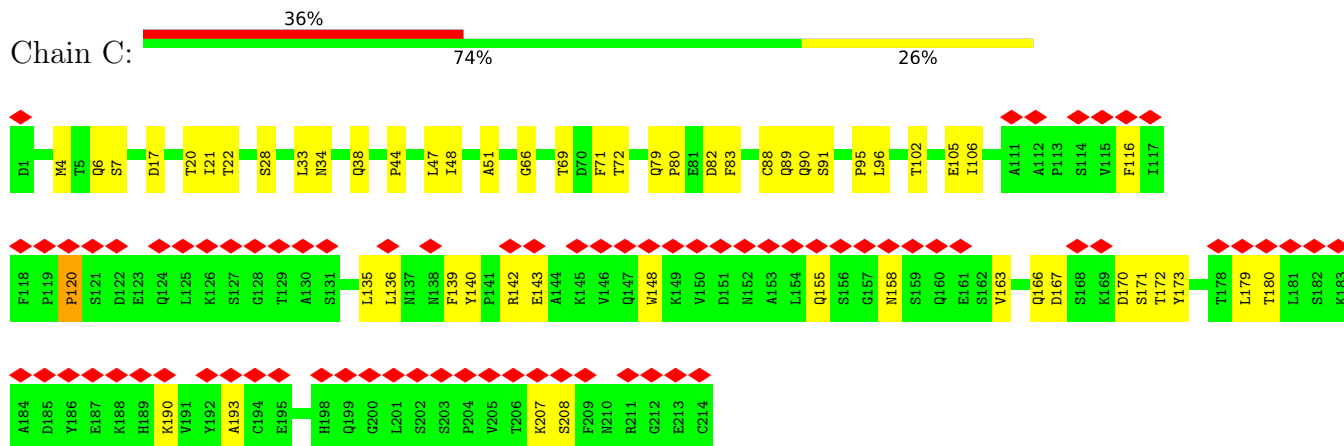


- Molecule 2: GFP-iRhom2 fusion protein

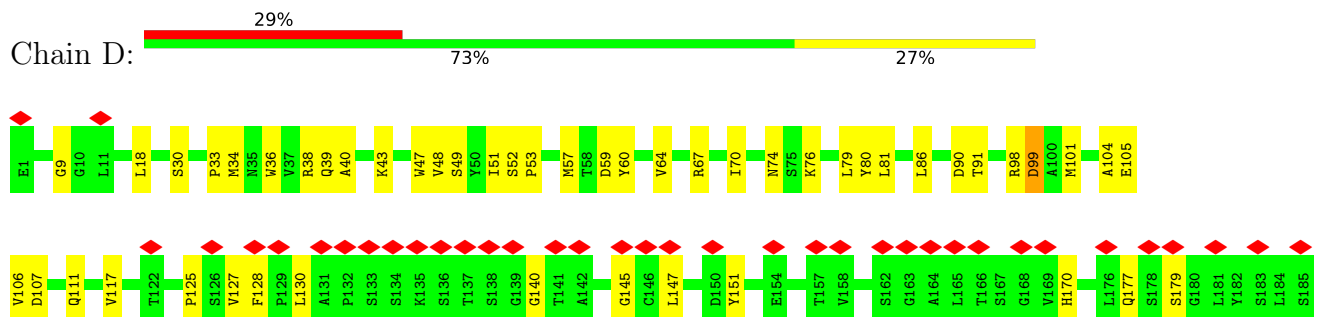




• Molecule 3: MEDI3622 Fab Light Chain



• Molecule 4: MEDI3622 Fab Heavy Chain





● Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	46753	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	58.3	Depositor
Minimum defocus (nm)	400	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.598	Depositor
Minimum map value	-0.305	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.014	Depositor
Recommended contour level	0.102	Depositor
Map size (Å)	361.68002, 361.68002, 361.68002	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.822, 0.822, 0.822	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: NAG, CA, ZN

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	0.16	0/5183	0.39	1/6995 (0.0%)
2	B	0.12	0/4058	0.30	0/5521
3	C	0.24	1/1674 (0.1%)	0.53	4/2271 (0.2%)
4	D	0.16	0/1685	0.40	0/2294
All	All	0.16	1/12600 (0.0%)	0.39	5/17081 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	120	PRO	CG-CD	-7.52	1.25	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	120	PRO	N-CD-CG	-13.05	83.63	103.20
3	C	120	PRO	CA-CB-CG	-10.34	84.86	104.50
1	A	452	ASN	CB-CA-C	-8.45	100.75	111.22
3	C	120	PRO	N-CA-CB	-6.66	96.26	103.25
3	C	120	PRO	CA-N-CD	-6.11	103.45	112.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	594	ASN	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5082	0	4889	133	0
2	B	3941	0	3876	81	0
3	C	1639	0	1599	40	0
4	D	1646	0	1612	53	0
5	E	28	0	25	2	0
6	A	84	0	78	1	0
7	A	1	0	0	0	0
8	A	1	0	0	0	0
All	All	12422	0	12079	292	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (292) 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:278:GLN:OE1	1:A:473:ARG:NH1	1.80	1.13
1:A:263:ARG:HH22	1:A:473:ARG:HD3	1.22	1.04
1:A:409:HIS:CE1	1:A:415:HIS:HD1	1.91	0.80
1:A:278:GLN:CD	1:A:473:ARG:NH1	2.44	0.76
1:A:412:GLY:HA3	1:A:461:THR:HG21	1.69	0.75
1:A:263:ARG:NH2	1:A:473:ARG:HD3	2.01	0.74
4:D:105:GLU:N	4:D:105:GLU:OE1	2.20	0.74
1:A:472:GLU:OE1	1:A:472:GLU:N	2.17	0.72
1:A:313:ASP:HB2	1:A:316:MET:HB2	1.72	0.72
3:C:33:LEU:HB3	3:C:51:ALA:HB2	1.71	0.71
1:A:413:ALA:HB1	1:A:454:SER:HB3	1.71	0.71
3:C:95:PRO:HA	4:D:47:TRP:HZ3	1.57	0.70
1:A:89:ARG:NH1	1:A:256:ASP:OD2	2.24	0.69
3:C:116:PHE:HB3	3:C:135:LEU:HB3	1.75	0.68
4:D:99:ASP:HA	4:D:106:VAL:HA	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:127:VAL:HG11	4:D:215:LYS:HG2	1.73	0.68
3:C:143:GLU:OE1	3:C:143:GLU:N	2.21	0.67
1:A:83:LEU:HB3	1:A:115:PHE:HB3	1.76	0.67
2:B:845:PHE:HA	2:B:848:LYS:HE3	1.75	0.67
1:A:578:CYS:SG	2:B:536:LYS:NZ	2.68	0.66
4:D:101:MET:HE2	4:D:101:MET:HA	1.77	0.65
2:B:458:TRP:HE3	2:B:668:VAL:HG13	1.61	0.65
4:D:30:SER:HA	4:D:74:ASN:HD21	1.61	0.65
1:A:70:LEU:HB3	1:A:81:LEU:HB2	1.78	0.65
2:B:756:VAL:HG21	2:B:777:PHE:HE2	1.63	0.64
4:D:47:TRP:CD1	4:D:48:VAL:N	2.66	0.63
4:D:91:THR:HG22	4:D:117:VAL:H	1.63	0.63
2:B:851:LEU:HD13	2:B:854:VAL:HB	1.80	0.62
1:A:92:GLN:HE21	1:A:477:VAL:CG1	2.12	0.62
2:B:605:LYS:HE3	2:B:714:TYR:HB3	1.81	0.61
4:D:59:ASP:OD1	4:D:60:TYR:N	2.33	0.61
2:B:367:ARG:HD3	2:B:686:LYS:HG2	1.82	0.61
3:C:6:GLN:NE2	3:C:88:CYS:SG	2.74	0.61
2:B:576:ILE:HG12	2:B:632:GLN:HE21	1.65	0.61
1:A:348:LEU:HD11	1:A:389:ASN:HB3	1.84	0.60
4:D:216:ARG:HH12	4:D:219:PRO:HD3	1.67	0.60
3:C:158:ASN:HB3	3:C:180:THR:HB	1.84	0.60
1:A:305:PRO:HG3	1:A:316:MET:HG3	1.84	0.59
1:A:225:CYS:HB3	1:A:470:PHE:CD1	2.38	0.59
1:A:233:HIS:HD2	1:A:237:ARG:HH21	1.51	0.59
2:B:400:HIS:O	2:B:686:LYS:NZ	2.36	0.58
2:B:647:ASN:HB3	2:B:650:VAL:HG22	1.84	0.58
2:B:681:LEU:HD13	2:B:727:LEU:HD21	1.85	0.58
4:D:33:PRO:HG3	4:D:53:PRO:HD3	1.84	0.58
3:C:91:SER:HB3	4:D:104:ALA:HB3	1.86	0.58
1:A:651:ARG:NH2	2:B:644:PRO:O	2.37	0.58
2:B:572:THR:HG23	2:B:631:SER:HB3	1.85	0.58
1:A:382:SER:OG	1:A:383:GLY:N	2.37	0.57
4:D:86:LEU:HD22	4:D:117:VAL:HG11	1.85	0.57
1:A:599:SER:OG	1:A:641:CYS:SG	2.57	0.57
2:B:369:TYR:O	2:B:382:TYR:OH	2.18	0.57
2:B:380:ARG:HD2	2:B:742:LEU:HD11	1.86	0.57
3:C:105:GLU:N	3:C:105:GLU:OE2	2.37	0.57
1:A:568:LEU:HD12	2:B:560:ALA:HB1	1.85	0.57
2:B:526:TRP:H	2:B:546:GLY:HA2	1.69	0.57
1:A:593:CYS:HB2	1:A:599:SER:HA	1.84	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:340:TYR:CE2	1:A:388:LYS:HD3	2.40	0.57
1:A:154:ARG:NH1	1:A:239:MET:O	2.37	0.56
4:D:128:PHE:HB3	4:D:147:LEU:HB2	1.87	0.56
3:C:167:ASP:OD2	3:C:170:ASP:N	2.39	0.56
1:A:427:GLU:HA	1:A:431:GLY:HA2	1.88	0.56
1:A:230:VAL:HG22	1:A:283:ARG:HB2	1.88	0.56
2:B:847:GLU:HG3	2:B:851:LEU:HD21	1.88	0.56
3:C:80:PRO:HA	3:C:83:PHE:HE1	1.71	0.55
1:A:535:GLN:HE22	2:B:586:THR:C	2.13	0.55
4:D:47:TRP:HE1	4:D:49:SER:C	2.13	0.55
4:D:51:ILE:HD11	4:D:70:ILE:HG23	1.89	0.55
1:A:59:ASP:N	1:A:59:ASP:OD1	2.39	0.55
1:A:92:GLN:HE21	1:A:477:VAL:HG12	1.71	0.55
1:A:499:ASP:OD1	1:A:517:ARG:NH1	2.39	0.55
1:A:234:ARG:NH1	1:A:341:GLN:OE1	2.40	0.55
4:D:177:GLN:NE2	4:D:179:SER:H	2.05	0.55
1:A:129:ALA:HB2	1:A:138:ILE:HG23	1.89	0.54
1:A:245:SER:O	1:A:249:ASN:ND2	2.40	0.54
1:A:131:ILE:HA	1:A:136:VAL:HG12	1.89	0.54
1:A:62:THR:OG1	1:A:63:SER:N	2.39	0.54
2:B:656:ARG:HA	2:B:659:LEU:HB3	1.89	0.53
2:B:506:SER:O	2:B:506:SER:OG	2.25	0.53
1:A:229:VAL:HB	1:A:282:ILE:HG12	1.89	0.53
3:C:89:GLN:NE2	3:C:96:LEU:HD13	2.24	0.53
2:B:532:PRO:HB3	2:B:554:ARG:HH21	1.73	0.53
1:A:278:GLN:CD	1:A:473:ARG:HH11	2.14	0.53
1:A:413:ALA:CB	1:A:454:SER:HB3	2.37	0.52
1:A:630:CYS:SG	1:A:631:THR:N	2.82	0.52
3:C:47:LEU:HD23	3:C:48:ILE:HG12	1.92	0.52
2:B:832:TRP:HA	2:B:835:HIS:CE1	2.44	0.52
3:C:79:GLN:N	3:C:82:ASP:OD2	2.39	0.52
3:C:142:ARG:HD3	3:C:163:VAL:HG11	1.90	0.52
1:A:334:LEU:HD11	1:A:364:VAL:HG12	1.91	0.52
1:A:348:LEU:HD23	1:A:402:VAL:HG21	1.90	0.52
2:B:419:VAL:HG22	2:B:469:LEU:HD21	1.91	0.52
1:A:342:ASP:OD2	1:A:388:LYS:NZ	2.43	0.52
3:C:148:TRP:HB2	3:C:155:GLN:HB2	1.91	0.52
1:A:174:ASN:O	1:A:177:ARG:HG2	2.10	0.51
1:A:501:CYS:HB2	1:A:512:VAL:HG11	1.92	0.51
1:A:637:MET:HE3	1:A:637:MET:H	1.76	0.51
4:D:98:ARG:NH1	4:D:107:ASP:OD2	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:HIS:NE2	1:A:43:ASP:OD1	2.41	0.51
2:B:782:LEU:O	2:B:808:SER:OG	2.29	0.50
1:A:578:CYS:O	2:B:536:LYS:NZ	2.33	0.50
1:A:628:LYS:NZ	2:B:560:ALA:O	2.43	0.50
2:B:373:VAL:N	2:B:683:ASP:OD2	2.42	0.50
1:A:296:LYS:HD2	1:A:300:MET:HE1	1.93	0.50
2:B:392:ARG:NH1	2:B:392:ARG:HA	2.26	0.50
2:B:725:PHE:HE2	2:B:773:HIS:ND1	2.08	0.50
1:A:61:GLN:OE1	1:A:494:MET:HE1	2.11	0.50
1:A:454:SER:O	1:A:458:ILE:HG13	2.11	0.50
1:A:323:PHE:CD2	4:D:57:MET:HE1	2.47	0.50
1:A:234:ARG:HD2	1:A:341:GLN:HG2	1.95	0.49
1:A:267:TRP:HE1	1:A:275:TYR:HB2	1.77	0.49
2:B:730:CYS:SG	2:B:788:LEU:HD11	2.53	0.49
3:C:38:GLN:OE1	4:D:39:GLN:NE2	2.45	0.49
4:D:47:TRP:HD1	4:D:48:VAL:H	1.61	0.49
1:A:407:LEU:O	1:A:411:PHE:N	2.43	0.49
3:C:47:LEU:O	3:C:48:ILE:HD13	2.13	0.49
1:A:472:GLU:H	1:A:472:GLU:CD	2.15	0.49
3:C:106:ILE:HD11	3:C:171:SER:HB3	1.94	0.49
1:A:535:GLN:HB2	1:A:546:SER:HB3	1.95	0.49
4:D:38:ARG:NH1	4:D:90:ASP:OD1	2.46	0.49
1:A:551:ASN:OD1	6:A:902:NAG:N2	2.46	0.49
3:C:17:ASP:N	3:C:17:ASP:OD1	2.43	0.49
4:D:140:GLY:HA2	4:D:192:SER:HB3	1.94	0.48
1:A:173:LYS:HB3	1:A:175:VAL:HG22	1.95	0.48
1:A:513:GLN:OE1	1:A:527:PHE:N	2.42	0.48
3:C:20:THR:OG1	3:C:72:THR:OG1	2.31	0.48
4:D:34:MET:HG3	4:D:79:LEU:HD22	1.93	0.48
4:D:60:TYR:CE1	4:D:70:ILE:HG22	2.48	0.48
1:A:673:VAL:HG13	2:B:415:ILE:HD13	1.95	0.48
1:A:70:LEU:O	1:A:81:LEU:N	2.41	0.48
3:C:193:ALA:HA	3:C:208:SER:HB2	1.95	0.48
3:C:38:GLN:HB2	3:C:44:PRO:HG3	1.96	0.48
4:D:86:LEU:HD21	4:D:117:VAL:HG21	1.95	0.48
1:A:263:ARG:HH22	1:A:473:ARG:CD	2.10	0.47
2:B:480:ASP:HB3	2:B:483:ILE:HB	1.96	0.47
1:A:226:LYS:HB3	1:A:280:GLU:HB2	1.96	0.47
1:A:616:ASP:OD1	1:A:620:LYS:N	2.44	0.47
1:A:336:HIS:HE2	1:A:385:THR:HG1	1.40	0.47
2:B:493:LEU:HD23	2:B:493:LEU:HA	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:33:LEU:HD12	3:C:34:ASN:H	1.80	0.47
1:A:227:LEU:N	1:A:278:GLN:O	2.48	0.47
1:A:541:THR:OG1	1:A:543:LYS:HG2	2.14	0.47
2:B:403:TYR:H	2:B:685:GLU:CD	2.22	0.47
3:C:155:GLN:HB3	3:C:158:ASN:ND2	2.30	0.47
1:A:80:LYS:HD2	1:A:120:VAL:HG21	1.96	0.47
1:A:539:ASN:OD1	5:E:1:NAG:N2	2.47	0.47
2:B:794:GLY:H	2:B:797:ASP:HB3	1.79	0.47
1:A:177:ARG:HH21	1:A:438:ILE:HG21	1.79	0.47
1:A:224:THR:O	1:A:471:GLN:O	2.33	0.47
2:B:710:ILE:HD11	2:B:822:TRP:C	2.40	0.47
4:D:51:ILE:HG22	4:D:52:SER:O	2.15	0.47
1:A:632:VAL:HG12	1:A:641:CYS:HB3	1.96	0.46
3:C:7:SER:OG	3:C:22:THR:OG1	2.33	0.46
4:D:206:HIS:CE1	4:D:208:PRO:HD2	2.51	0.46
1:A:233:HIS:H	1:A:299:ASN:HD21	1.63	0.46
1:A:435:MET:HA	1:A:435:MET:HE3	1.97	0.46
1:A:646:GLN:HE21	1:A:651:ARG:HE	1.62	0.46
1:A:672:ILE:O	1:A:676:VAL:HG23	2.16	0.46
2:B:843:SER:OG	2:B:844:ARG:NH1	2.47	0.46
1:A:635:CYS:HB3	1:A:641:CYS:HA	1.97	0.46
2:B:699:LEU:O	2:B:703:THR:HG23	2.16	0.46
4:D:36:TRP:CG	4:D:81:LEU:HD12	2.51	0.46
2:B:693:ILE:HA	2:B:696:ILE:HG22	1.98	0.46
4:D:128:PHE:O	4:D:147:LEU:N	2.46	0.46
2:B:645:PHE:HE1	2:B:653:GLN:HE21	1.64	0.46
1:A:624:LEU:HB3	1:A:628:LYS:HG3	1.97	0.45
4:D:34:MET:HB3	4:D:79:LEU:HD22	1.98	0.45
1:A:560:ASN:HD21	1:A:576:GLY:N	2.13	0.45
1:A:233:HIS:CD2	1:A:237:ARG:HH21	2.31	0.45
2:B:380:ARG:HD3	2:B:748:ALA:HB1	1.98	0.45
1:A:681:LEU:HG	2:B:407:TRP:CH2	2.51	0.45
4:D:64:VAL:HG12	4:D:67:ARG:HH21	1.81	0.45
5:E:1:NAG:H62	5:E:2:NAG:H82	1.98	0.45
2:B:472:LYS:HB3	2:B:472:LYS:HE3	1.83	0.45
2:B:555:THR:HG21	2:B:591:MET:SD	2.56	0.45
4:D:151:TYR:HB2	4:D:206:HIS:NE2	2.31	0.45
1:A:508:LEU:HD23	1:A:512:VAL:HG12	1.99	0.45
3:C:106:ILE:O	3:C:140:TYR:OH	2.35	0.45
4:D:220:LYS:HD2	4:D:220:LYS:HA	1.71	0.45
1:A:264:ASN:HA	1:A:273:LYS:HE3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:725:PHE:HA	2:B:728:LEU:HB3	1.98	0.45
3:C:179:LEU:HD12	3:C:180:THR:N	2.32	0.45
1:A:97:VAL:HG13	1:A:283:ARG:HG2	1.98	0.45
1:A:357:ARG:HB2	1:A:360:SER:HB3	1.98	0.45
2:B:725:PHE:O	2:B:729:ALA:N	2.43	0.44
3:C:136:LEU:HD11	3:C:139:PHE:CD1	2.52	0.44
2:B:431:ALA:O	2:B:455:GLU:HG2	2.16	0.44
2:B:469:LEU:HD23	2:B:663:LEU:HD21	1.99	0.44
1:A:180:SER:HB3	1:A:390:TYR:CE2	2.52	0.44
1:A:333:CYS:SG	1:A:334:LEU:HD12	2.57	0.44
2:B:588:PHE:O	2:B:590:HIS:N	2.48	0.44
3:C:136:LEU:HD11	3:C:139:PHE:HD1	1.82	0.44
1:A:490:ASP:H	1:A:518:ASN:ND2	2.16	0.44
1:A:521:CYS:O	1:A:528:GLU:HB3	2.17	0.44
2:B:497:SER:OG	2:B:510:GLN:OE1	2.30	0.44
2:B:721:ALA:HB1	2:B:773:HIS:NE2	2.33	0.44
1:A:361:HIS:CE1	1:A:369:TYR:HB3	2.53	0.44
3:C:4:MET:HE1	3:C:33:LEU:HD13	2.00	0.44
4:D:207:LYS:HA	4:D:207:LYS:HD3	1.63	0.44
2:B:771:ILE:HG12	2:B:837:THR:OG1	2.17	0.44
2:B:436:THR:HG22	2:B:450:LYS:HB3	1.99	0.44
4:D:130:LEU:HD23	4:D:145:GLY:C	2.43	0.44
1:A:105:SER:OG	1:A:106:GLU:N	2.51	0.43
1:A:408:GLY:O	1:A:413:ALA:HB3	2.18	0.43
4:D:52:SER:OG	4:D:53:PRO:HD2	2.17	0.43
1:A:56:ARG:HE	1:A:61:GLN:HA	1.82	0.43
2:B:392:ARG:HA	2:B:392:ARG:CZ	2.49	0.43
4:D:125:PRO:HB3	4:D:151:TYR:HB3	2.00	0.43
1:A:616:ASP:OD1	1:A:616:ASP:N	2.39	0.43
2:B:542:LYS:HD2	2:B:542:LYS:HA	1.83	0.43
4:D:151:TYR:HB2	4:D:206:HIS:HE2	1.82	0.43
1:A:292:LYS:HB2	1:A:295:GLU:HG2	2.01	0.43
2:B:500:CYS:HB3	2:B:523:PHE:HD1	1.82	0.43
1:A:460:LYS:HE2	1:A:460:LYS:HB2	1.80	0.43
1:A:688:SER:O	2:B:403:TYR:OH	2.29	0.43
2:B:844:ARG:N	2:B:844:ARG:HD2	2.34	0.43
3:C:21:ILE:HG21	3:C:102:THR:HG21	2.01	0.43
2:B:419:VAL:O	2:B:423:TYR:HB2	2.19	0.43
2:B:482:GLN:HB3	2:B:855:LEU:HD12	2.01	0.43
3:C:120:PRO:O	3:C:120:PRO:HG2	2.14	0.43
4:D:111:GLN:HA	4:D:111:GLN:OE1	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:ARG:H	1:A:357:ARG:HG3	1.58	0.43
3:C:28:SER:HA	3:C:69:THR:HG22	2.00	0.43
1:A:660:SER:OG	1:A:662:ASN:OD1	2.22	0.43
2:B:468:HIS:HE1	2:B:622:TYR:OH	2.02	0.43
1:A:222:LYS:HA	1:A:472:GLU:HA	2.01	0.42
2:B:427:PRO:O	2:B:455:GLU:HB2	2.19	0.42
3:C:90:GLN:NE2	3:C:96:LEU:HA	2.34	0.42
4:D:79:LEU:HD12	4:D:80:TYR:N	2.35	0.42
1:A:330:SER:HB2	1:A:379:TYR:CE1	2.54	0.42
2:B:604:THR:HG23	2:B:605:LYS:HD2	2.01	0.42
1:A:31:ARG:HA	1:A:31:ARG:HD3	1.81	0.42
2:B:632:GLN:H	2:B:632:GLN:HG2	1.68	0.42
1:A:521:CYS:SG	1:A:555:CYS:HB3	2.60	0.42
1:A:644:ARG:HA	1:A:644:ARG:HD2	1.81	0.42
3:C:66:GLY:HA3	3:C:71:PHE:HA	2.02	0.42
2:B:851:LEU:HD22	2:B:854:VAL:HG21	2.02	0.42
4:D:107:ASP:OD1	4:D:107:ASP:C	2.61	0.42
1:A:92:GLN:HE21	1:A:477:VAL:HG11	1.82	0.42
2:B:642:LEU:HD23	2:B:642:LEU:HA	1.82	0.42
1:A:400:ASP:OD1	1:A:400:ASP:C	2.63	0.42
1:A:600:CYS:HB3	1:A:639:GLY:HA2	2.01	0.42
2:B:787:PHE:HE1	2:B:809:LEU:HD21	1.85	0.42
3:C:208:SER:O	3:C:208:SER:OG	2.32	0.42
4:D:170:HIS:HB3	4:D:187:VAL:O	2.19	0.42
4:D:33:PRO:HB3	4:D:51:ILE:O	2.20	0.42
4:D:212:LYS:HA	4:D:212:LYS:HD3	1.81	0.42
2:B:598:ARG:HB2	2:B:623:PHE:CZ	2.55	0.41
4:D:40:ALA:HB3	4:D:43:LYS:HB2	2.02	0.41
1:A:237:ARG:NH2	1:A:288:PRO:HD3	2.34	0.41
1:A:646:GLN:NE2	1:A:651:ARG:HE	2.18	0.41
2:B:580:GLN:HE21	2:B:582:ARG:HB2	1.84	0.41
2:B:699:LEU:HD12	2:B:812:PHE:CZ	2.55	0.41
2:B:770:ASN:HA	2:B:773:HIS:HB2	2.02	0.41
3:C:172:THR:OG1	3:C:173:TYR:N	2.53	0.41
1:A:415:HIS:NE2	1:A:437:PRO:HA	2.35	0.41
1:A:661:ILE:HD12	1:A:661:ILE:HA	1.97	0.41
4:D:206:HIS:ND1	4:D:208:PRO:HD2	2.35	0.41
1:A:449:MET:HE3	1:A:449:MET:HB3	1.89	0.41
1:A:601:LYS:HG3	1:A:613:PRO:HB3	2.02	0.41
1:A:696:LYS:HG3	1:A:697:LYS:HD3	2.02	0.41
2:B:509:ILE:O	2:B:511:THR:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:9:GLY:H	4:D:18:LEU:HD21	1.85	0.41
2:B:397:PHE:CE2	2:B:791:ILE:HD11	2.55	0.41
1:A:133:ASP:OD1	1:A:133:ASP:N	2.51	0.41
2:B:467:ILE:H	2:B:467:ILE:HG12	1.66	0.41
2:B:842:THR:O	2:B:845:PHE:HD1	2.04	0.41
1:A:180:SER:O	1:A:439:ALA:HB2	2.21	0.41
1:A:377:ASN:OD1	1:A:377:ASN:N	2.53	0.41
1:A:688:SER:HB2	2:B:403:TYR:CE2	2.56	0.41
3:C:166:GLN:OE1	3:C:167:ASP:N	2.54	0.41
1:A:433:TYR:CD2	1:A:451:SER:HB3	2.57	0.41
1:A:637:MET:SD	1:A:638:ASN:N	2.94	0.41
1:A:662:ASN:OD1	1:A:662:ASN:N	2.54	0.41
2:B:664:HIS:HE1	2:B:673:SER:HB3	1.85	0.41
2:B:775:PHE:O	2:B:779:SER:OG	2.38	0.41
1:A:149:ILE:HA	1:A:165:LEU:O	2.21	0.40
1:A:336:HIS:NE2	1:A:385:THR:OG1	2.34	0.40
4:D:48:VAL:HG11	4:D:81:LEU:HD11	2.03	0.40
4:D:91:THR:HG22	4:D:117:VAL:N	2.34	0.40
1:A:266:SER:HA	1:A:272:PHE:O	2.21	0.40
1:A:323:PHE:CG	4:D:57:MET:HE1	2.56	0.40
1:A:535:GLN:N	1:A:546:SER:OG	2.38	0.40
1:A:627:GLY:HA3	2:B:559:PRO:HG3	2.03	0.40
3:C:207:LYS:HB2	3:C:207:LYS:HE3	1.83	0.40
1:A:65:HIS:HB2	1:A:85:SER:O	2.21	0.40
1:A:237:ARG:HH22	1:A:288:PRO:HD3	1.86	0.40
1:A:628:LYS:HA	1:A:628:LYS:HD3	1.86	0.40
3:C:190:LYS:HD3	3:C:190:LYS:HA	1.92	0.40
1:A:395:LEU:H	1:A:395:LEU:HD12	1.86	0.40
2:B:474:SER:HB2	2:B:478:ARG:HE	1.86	0.40
4:D:76:LYS:HE2	4:D:76:LYS:HB3	1.85	0.40
4:D:48:VAL:HG12	4:D:49:SER:OG	2.21	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	A	638/807 (79%)	614 (96%)	24 (4%)	0	100	100
2	B	488/741 (66%)	464 (95%)	24 (5%)	0	100	100
3	C	212/214 (99%)	199 (94%)	13 (6%)	0	100	100
4	D	219/221 (99%)	208 (95%)	11 (5%)	0	100	100
All	All	1557/1983 (78%)	1485 (95%)	72 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	575/726 (79%)	573 (100%)	2 (0%)	86	83
2	B	430/647 (66%)	427 (99%)	3 (1%)	76	78
3	C	189/189 (100%)	189 (100%)	0	100	100
4	D	184/184 (100%)	183 (100%)	1 (0%)	81	81
All	All	1378/1746 (79%)	1372 (100%)	6 (0%)	81	81

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	415	HIS
1	A	452	ASN
2	B	684	LEU
2	B	742	LEU
2	B	832	TRP
4	D	99	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	92	GLN
1	A	233	HIS
1	A	297	HIS
1	A	299	ASN
1	A	320	GLN
1	A	415	HIS
1	A	426	ASN
1	A	446	ASN
1	A	447	ASN
1	A	518	ASN
1	A	535	GLN
1	A	560	ASN
1	A	598	ASN
1	A	619	GLN
1	A	621	ASN
1	A	646	GLN
2	B	376	ASN
2	B	391	GLN
2	B	393	GLN
2	B	400	HIS
2	B	412	HIS
2	B	468	HIS
2	B	485	GLN
2	B	565	HIS
2	B	632	GLN
2	B	669	HIS
3	C	155	GLN
3	C	158	ASN
3	C	198	HIS
4	D	39	GLN
4	D	161	ASN
4	D	170	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

2 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	E	1	1,5	14,14,15	0.28	0	17,19,21	0.62	1 (5%)
5	NAG	E	2	5	14,14,15	0.27	0	17,19,21	0.46	0

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
5	NAG	E	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	E	2	5	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	1	NAG	C1-O5-C5	2.02	114.89	112.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

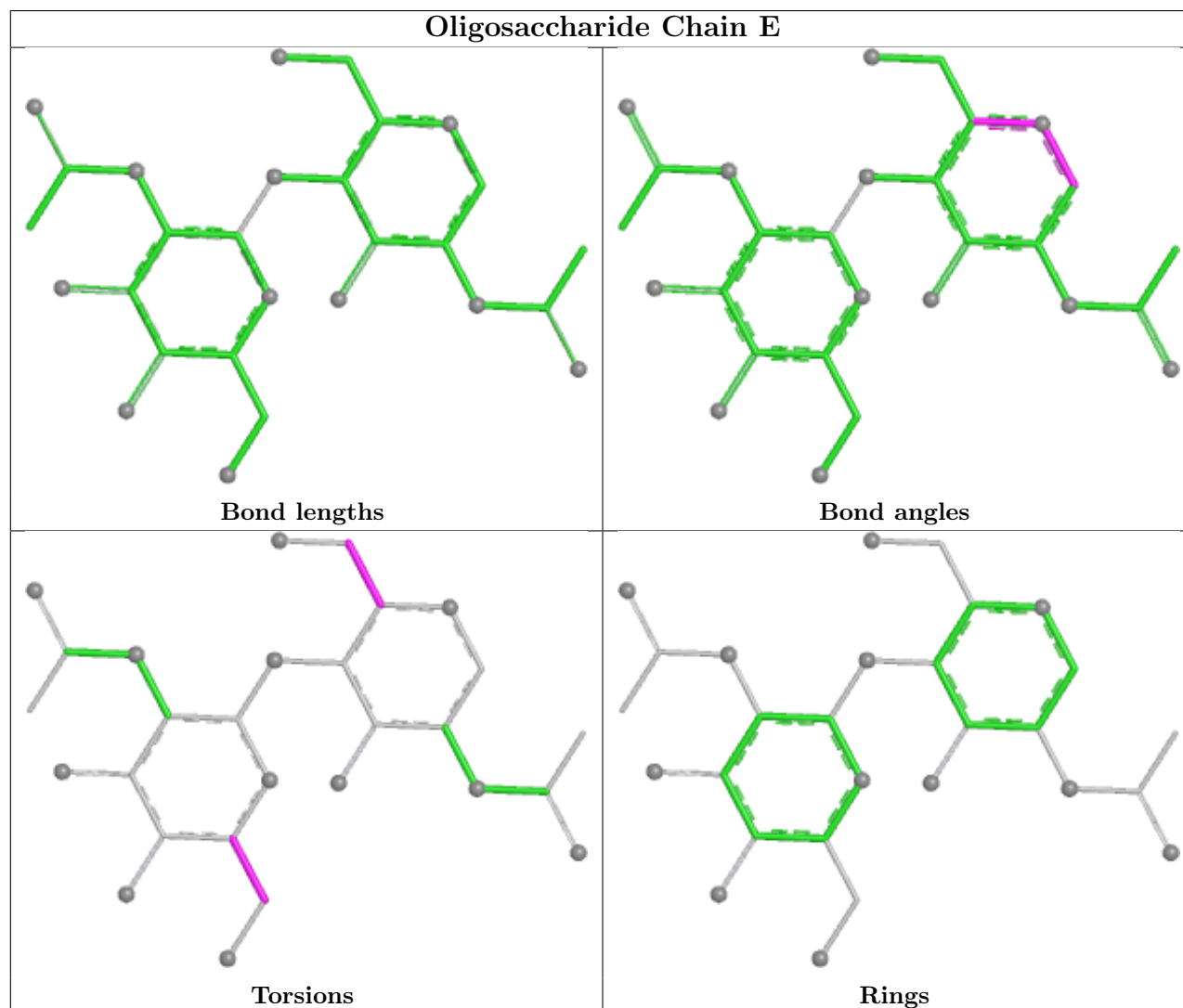
Mol	Chain	Res	Type	Atoms
5	E	1	NAG	O5-C5-C6-O6
5	E	1	NAG	C4-C5-C6-O6
5	E	2	NAG	C4-C5-C6-O6
5	E	2	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	1	NAG	2	0
5	E	2	NAG	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 oligosaccharide.



5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	$\# Z > 2$
6	NAG	A	904	1	14,14,15	0.22	0	17,19,21	0.45	0
6	NAG	A	905	1	14,14,15	0.69	1 (7%)	17,19,21	0.64	0
6	NAG	A	906	1	14,14,15	0.57	0	17,19,21	0.70	1 (5%)
6	NAG	A	902	1	14,14,15	0.35	0	17,19,21	0.69	1 (5%)
6	NAG	A	903	1	14,14,15	0.26	0	17,19,21	0.41	0
6	NAG	A	901	1	14,14,15	0.93	1 (7%)	17,19,21	0.61	0

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
6	NAG	A	904	1	-	2/6/23/26	0/1/1/1
6	NAG	A	905	1	-	2/6/23/26	0/1/1/1
6	NAG	A	906	1	-	2/6/23/26	0/1/1/1
6	NAG	A	902	1	-	4/6/23/26	0/1/1/1
6	NAG	A	903	1	-	2/6/23/26	0/1/1/1
6	NAG	A	901	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	901	NAG	O5-C1	-2.76	1.39	1.43
6	A	905	NAG	C1-C2	2.03	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	902	NAG	C1-O5-C5	2.45	115.46	112.19
6	A	906	NAG	C1-O5-C5	2.22	115.16	112.19

There are no chirality outliers.

All (12) torsion outliers are listed below:

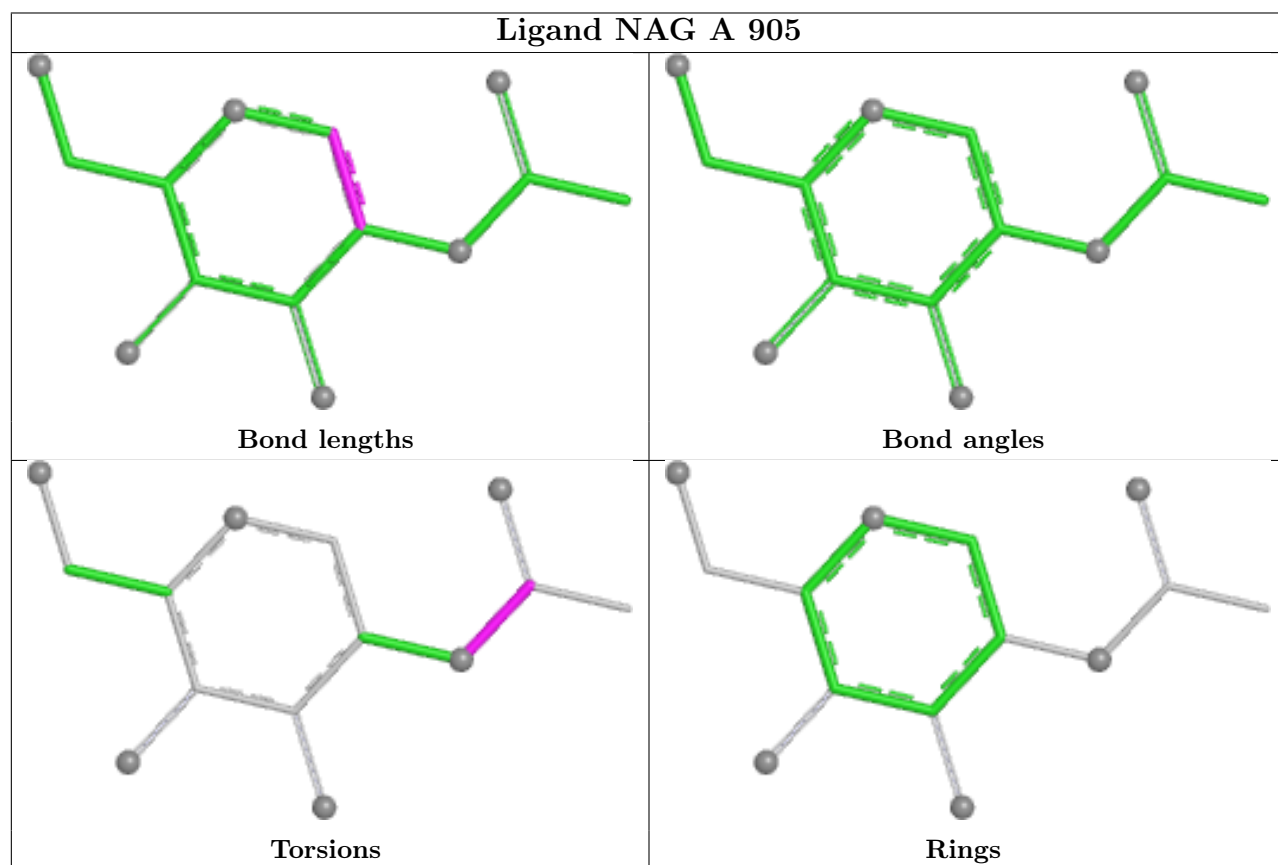
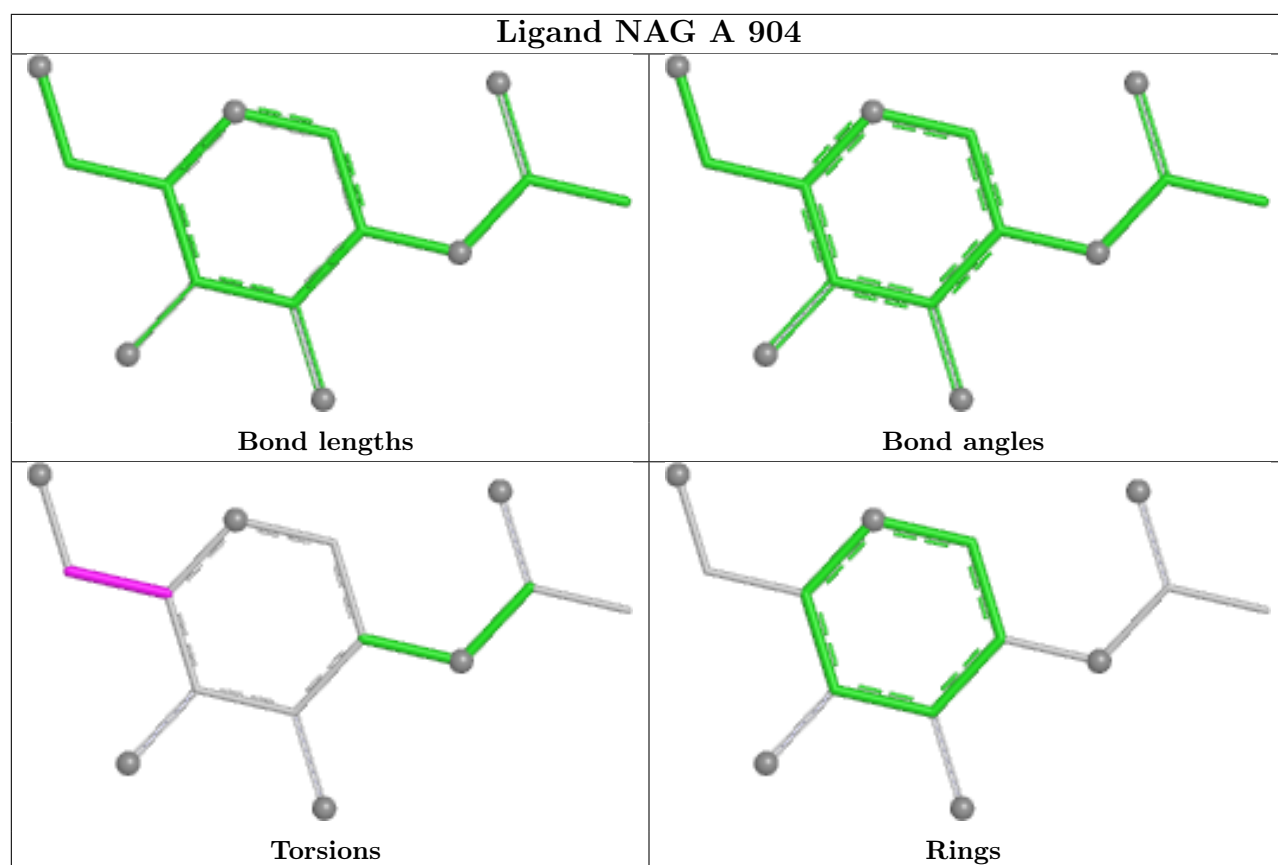
Mol	Chain	Res	Type	Atoms
6	A	904	NAG	C4-C5-C6-O6
6	A	904	NAG	O5-C5-C6-O6
6	A	906	NAG	C4-C5-C6-O6
6	A	906	NAG	O5-C5-C6-O6
6	A	902	NAG	C4-C5-C6-O6
6	A	902	NAG	O5-C5-C6-O6
6	A	902	NAG	C8-C7-N2-C2
6	A	902	NAG	O7-C7-N2-C2
6	A	905	NAG	C8-C7-N2-C2
6	A	905	NAG	O7-C7-N2-C2
6	A	903	NAG	O5-C5-C6-O6
6	A	903	NAG	C1-C2-N2-C7

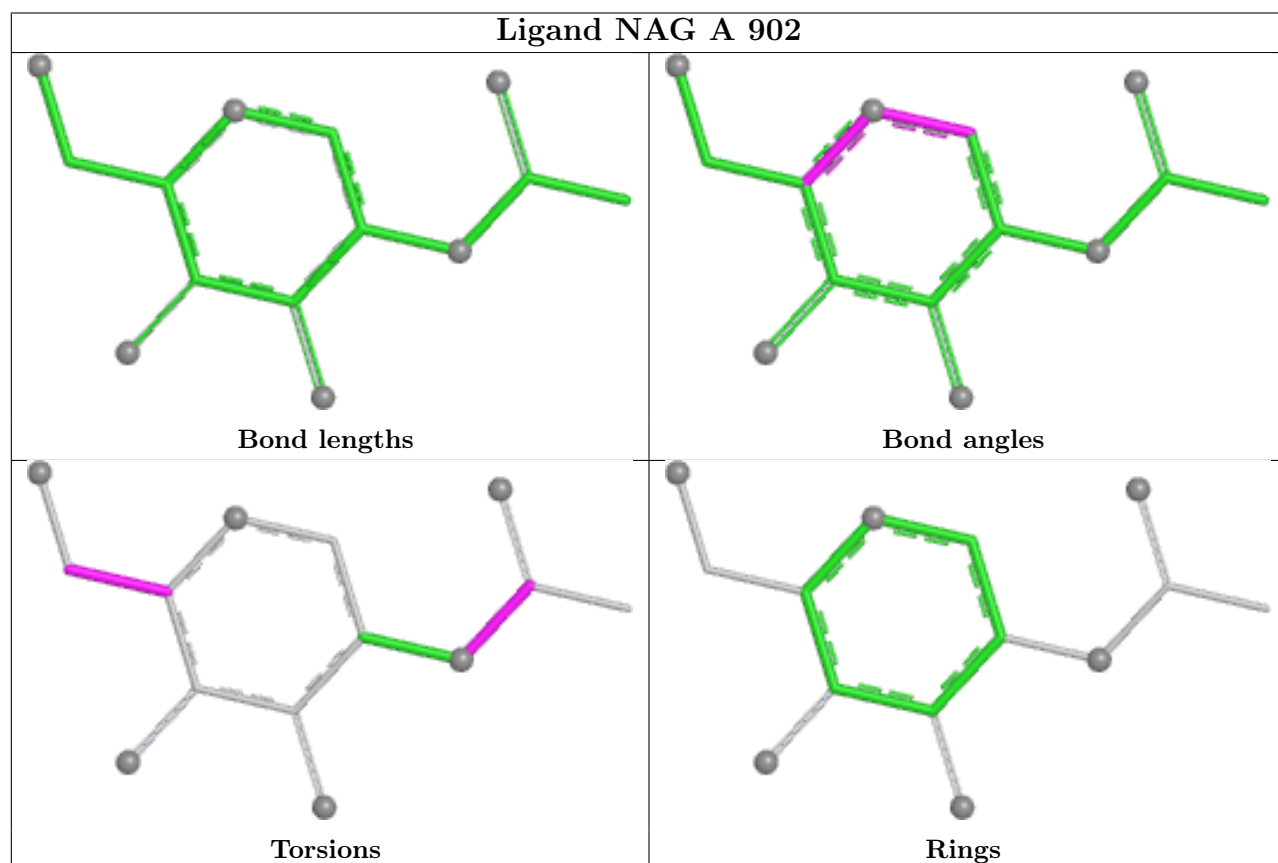
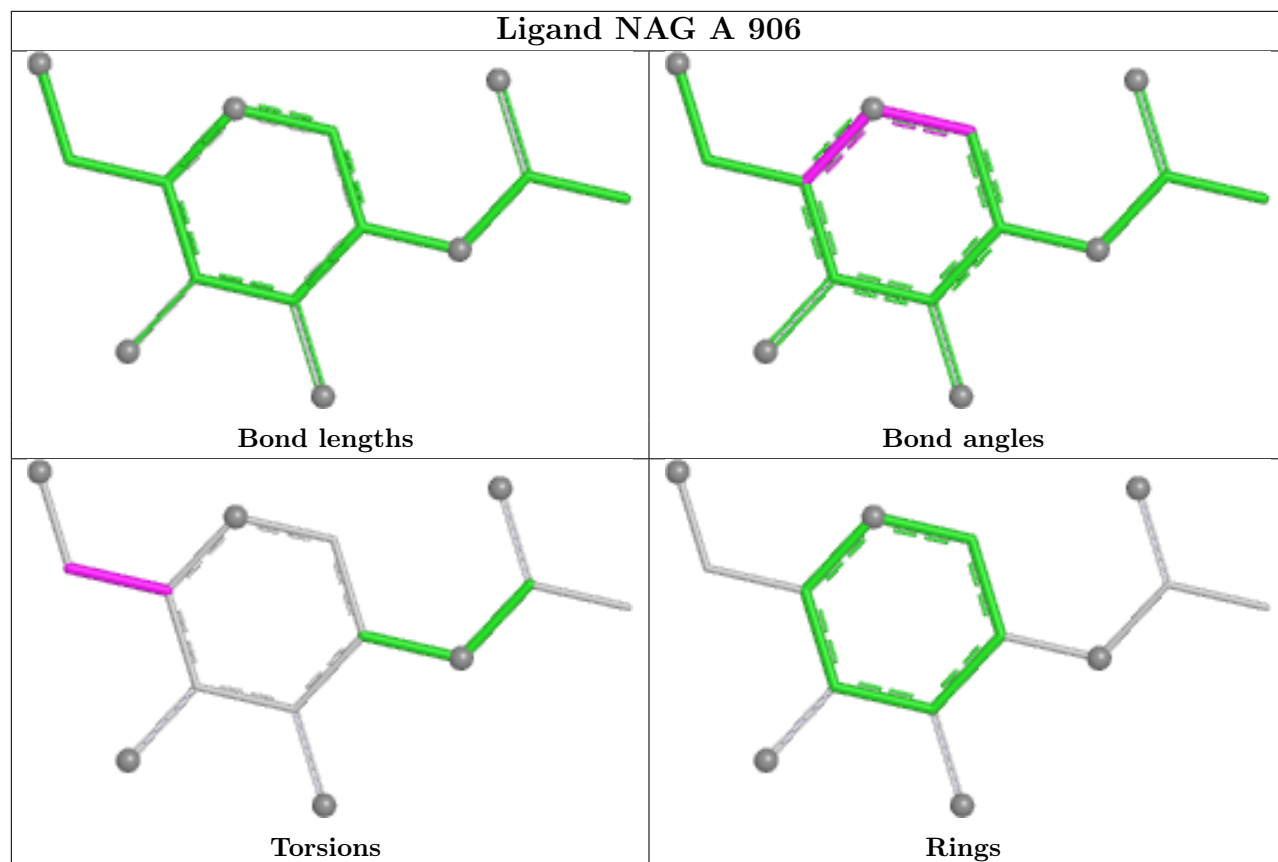
There are no ring outliers.

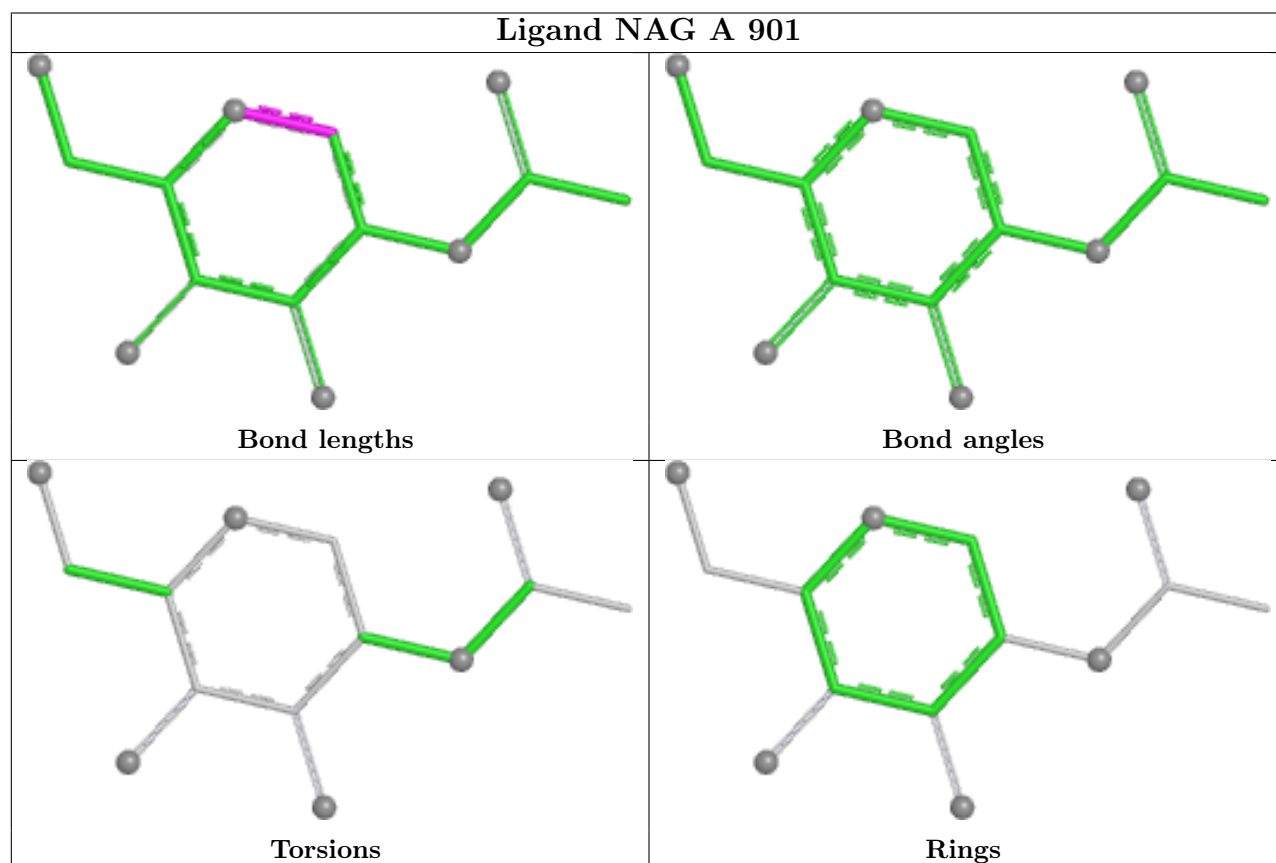
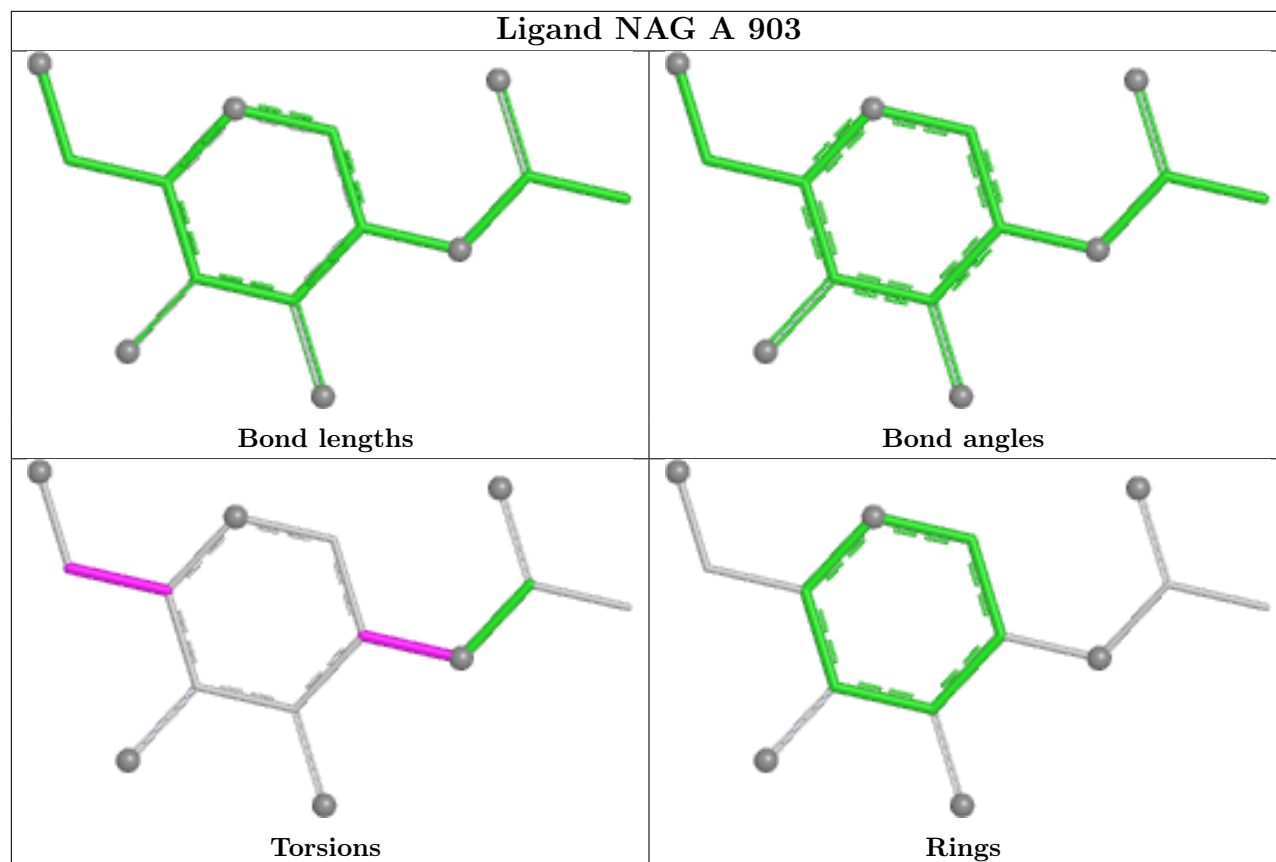
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	902	NAG	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.

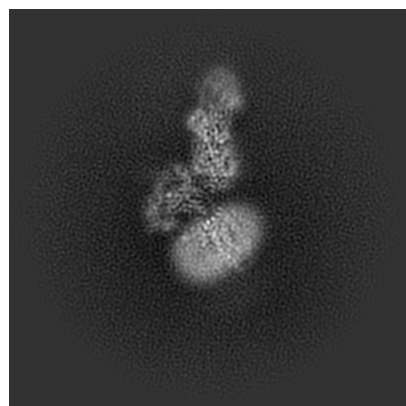
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-70127. These allow visual inspection of the internal detail of the map and identification of artifacts.

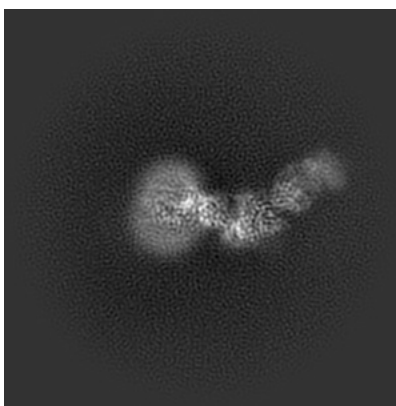
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

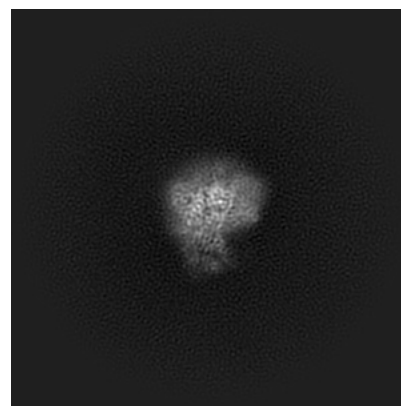
6.1.1 Primary map



X

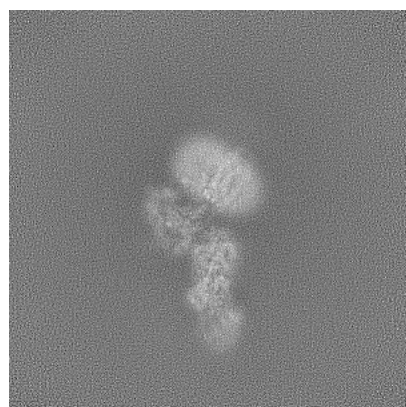


Y

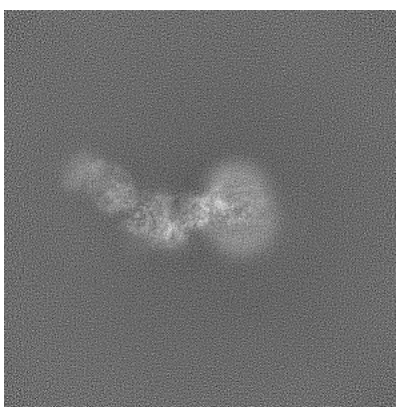


Z

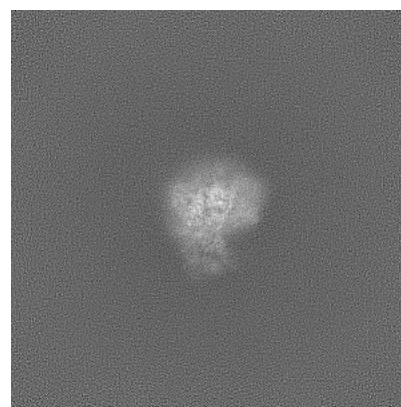
6.1.2 Raw map



X



Y

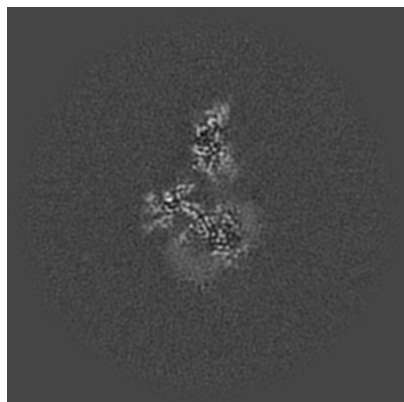


Z

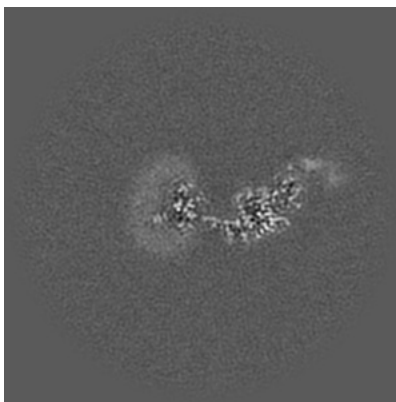
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

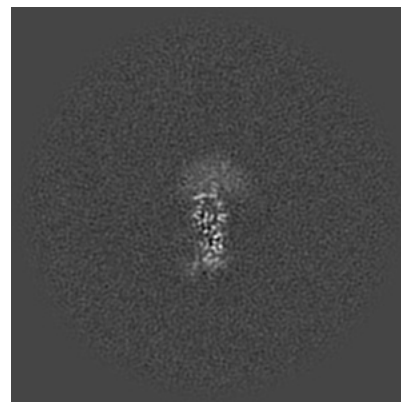
6.2.1 Primary map



X Index: 220

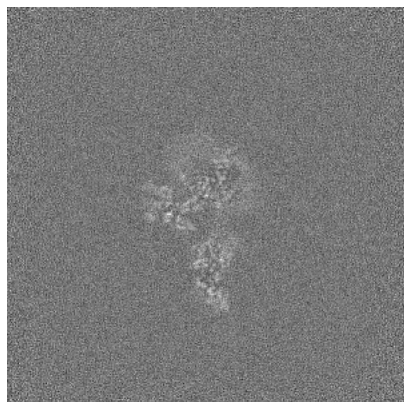


Y Index: 220

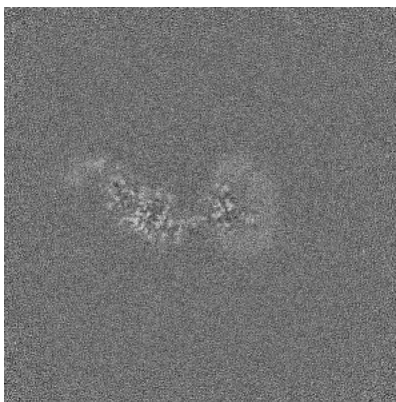


Z Index: 220

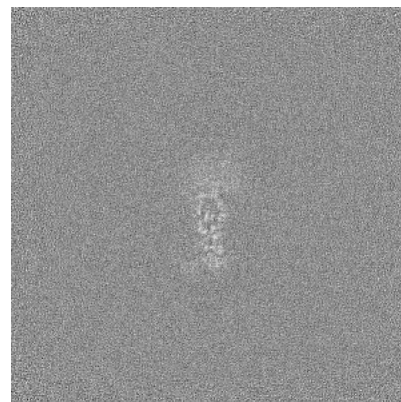
6.2.2 Raw map



X Index: 220



Y Index: 220

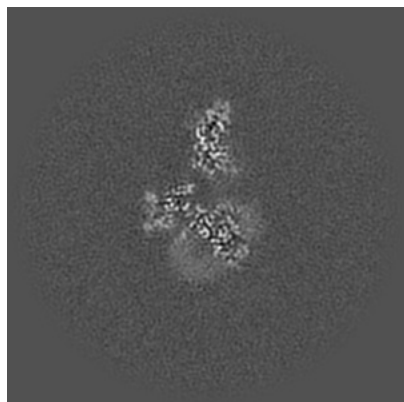


Z Index: 220

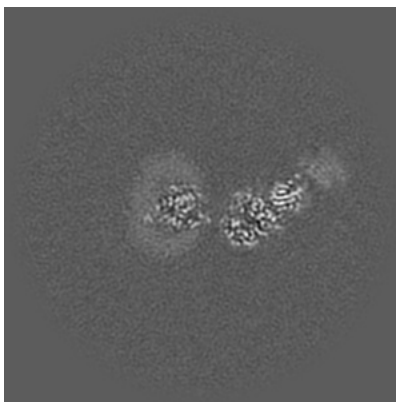
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

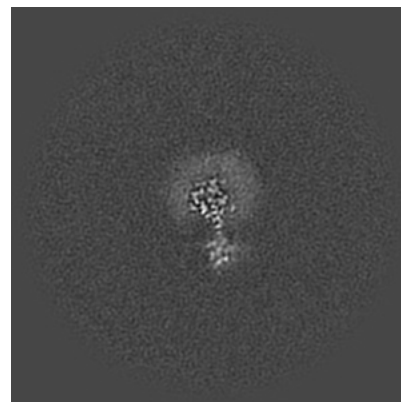
6.3.1 Primary map



X Index: 222

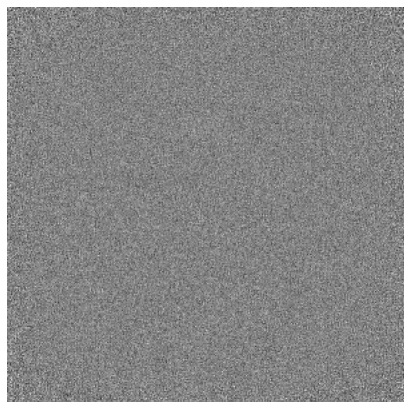


Y Index: 227

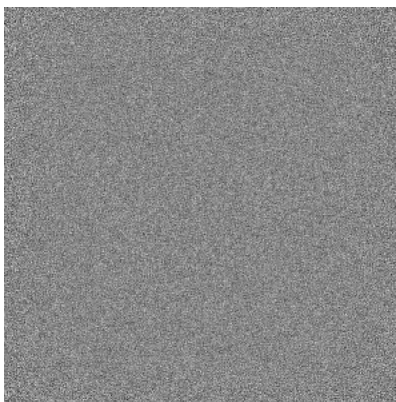


Z Index: 203

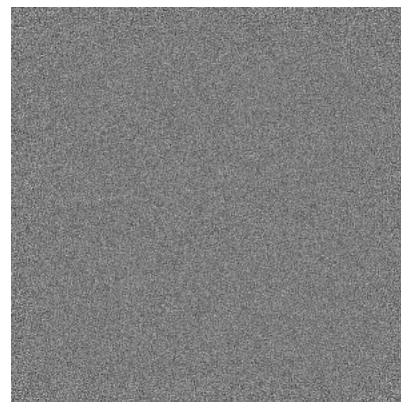
6.3.2 Raw map



X Index: 0



Y Index: 0

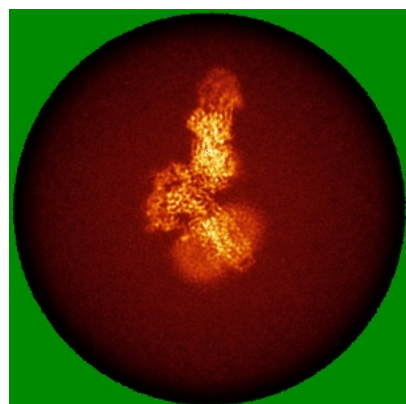


Z Index: 0

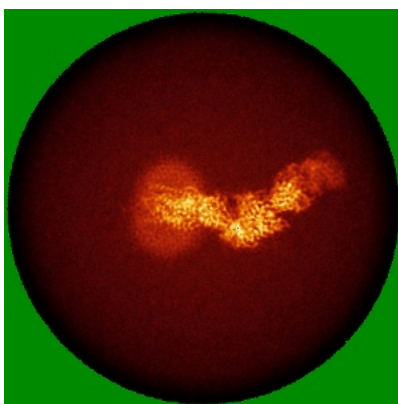
The images above show the largest variance slices of the map in three orthogonal directions.

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

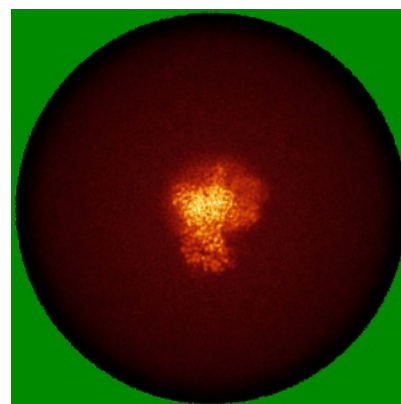
6.4.1 Primary map



X

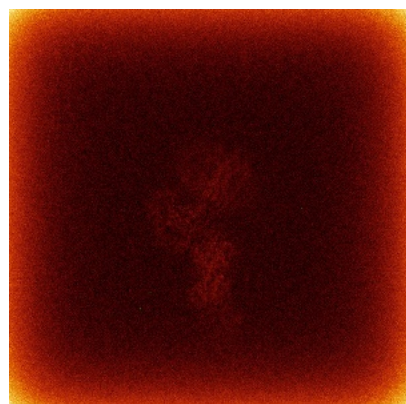


Y

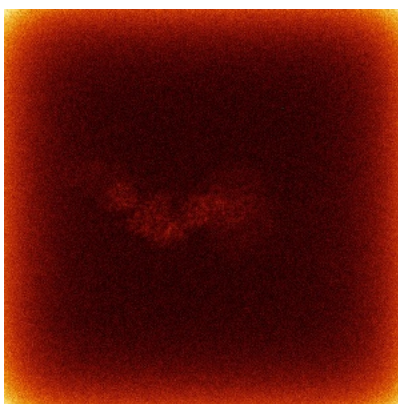


Z

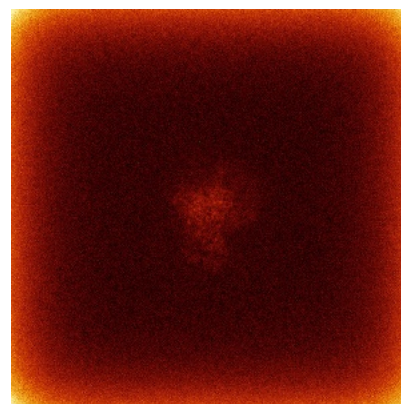
6.4.2 Raw map



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](#)

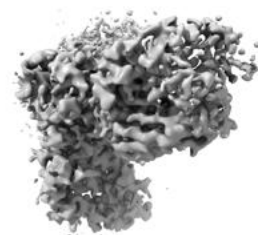
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.102. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

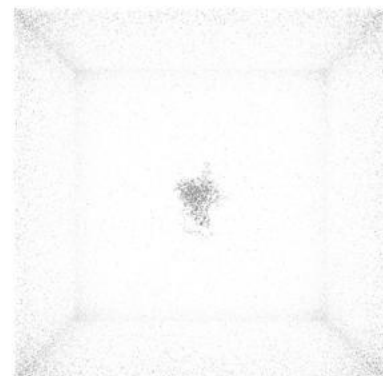
6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

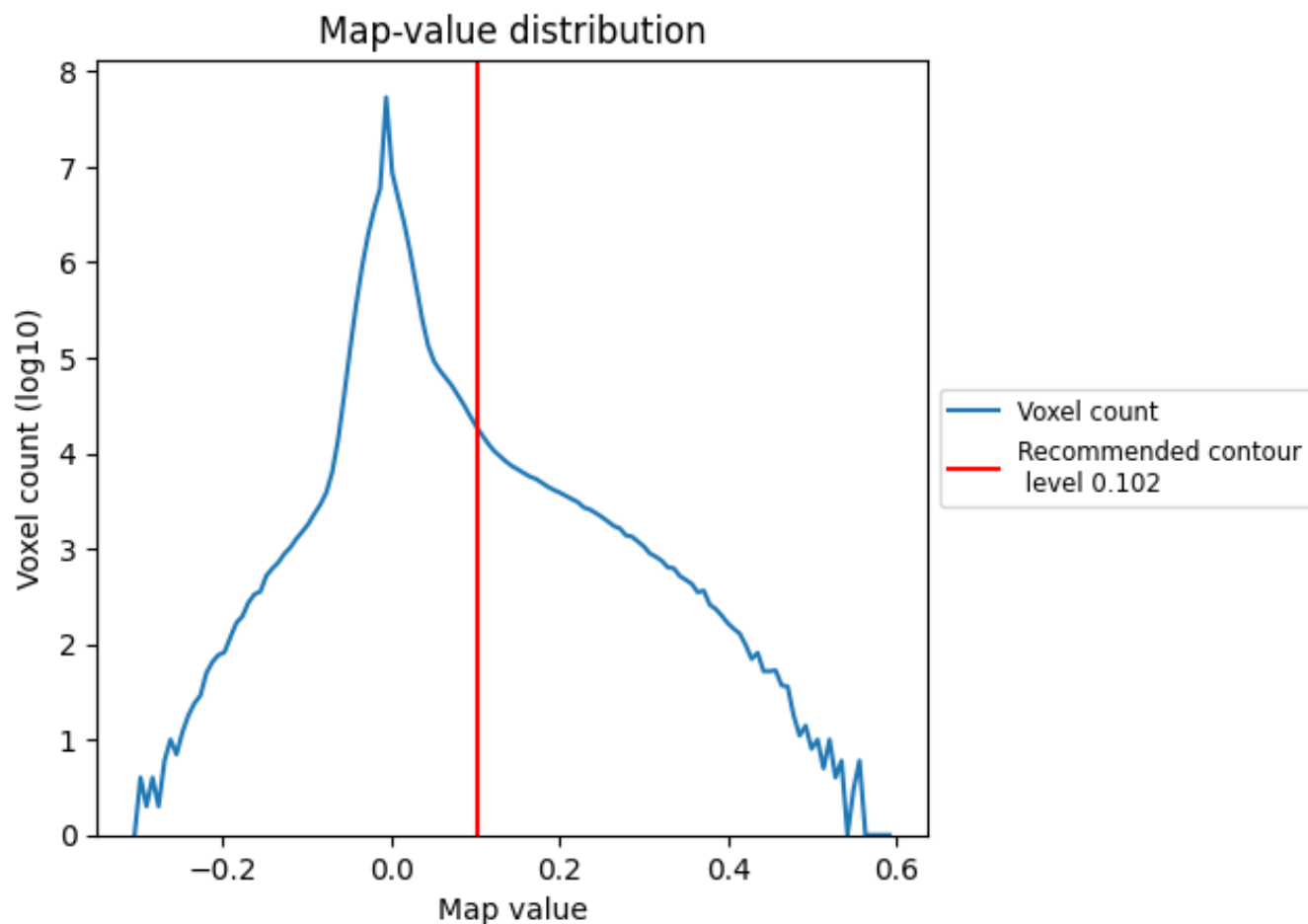
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

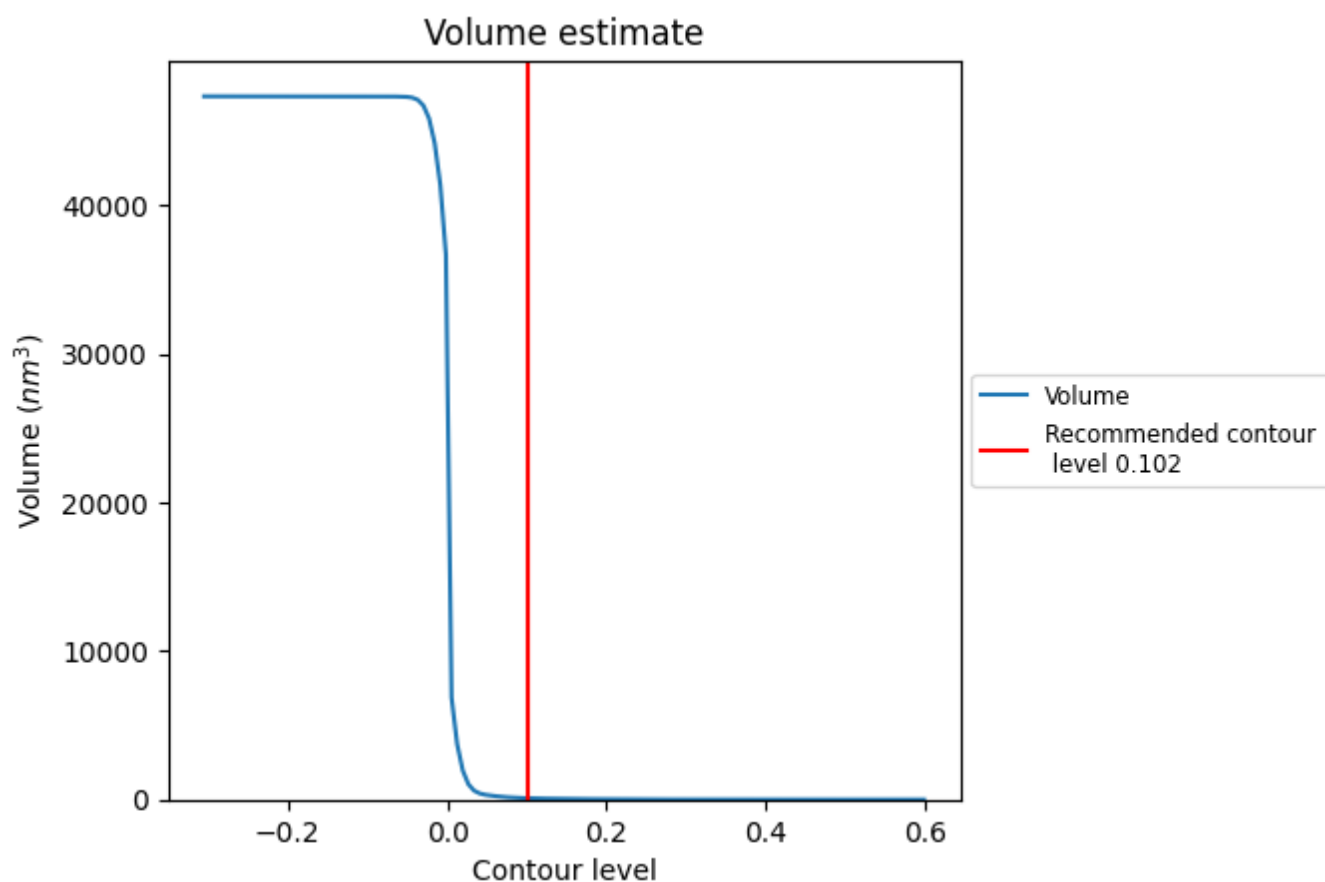
This section contains the results of statistical analysis of the map.

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.

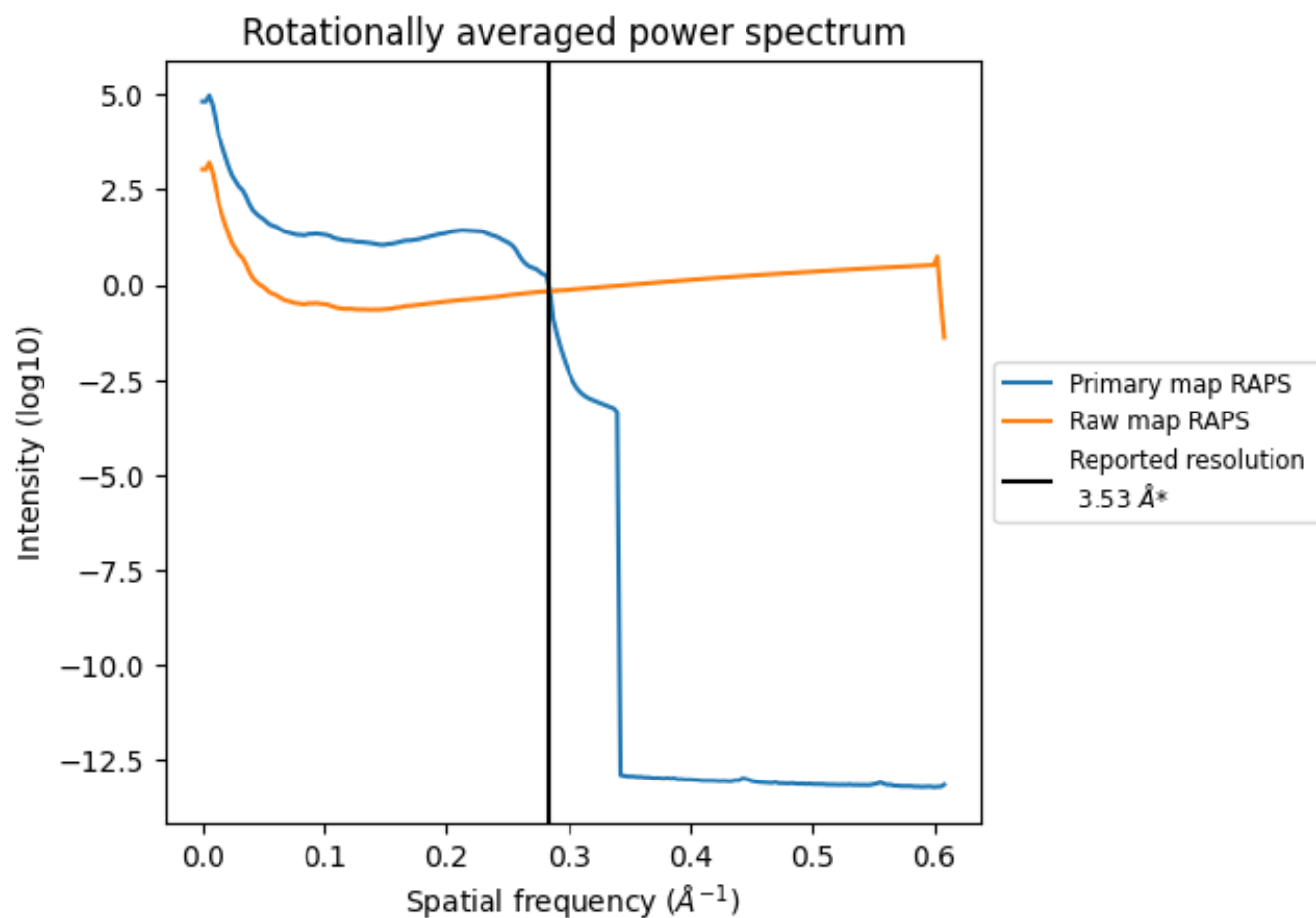
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 87 nm³; this corresponds to an approximate mass of 79 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 ⓘ

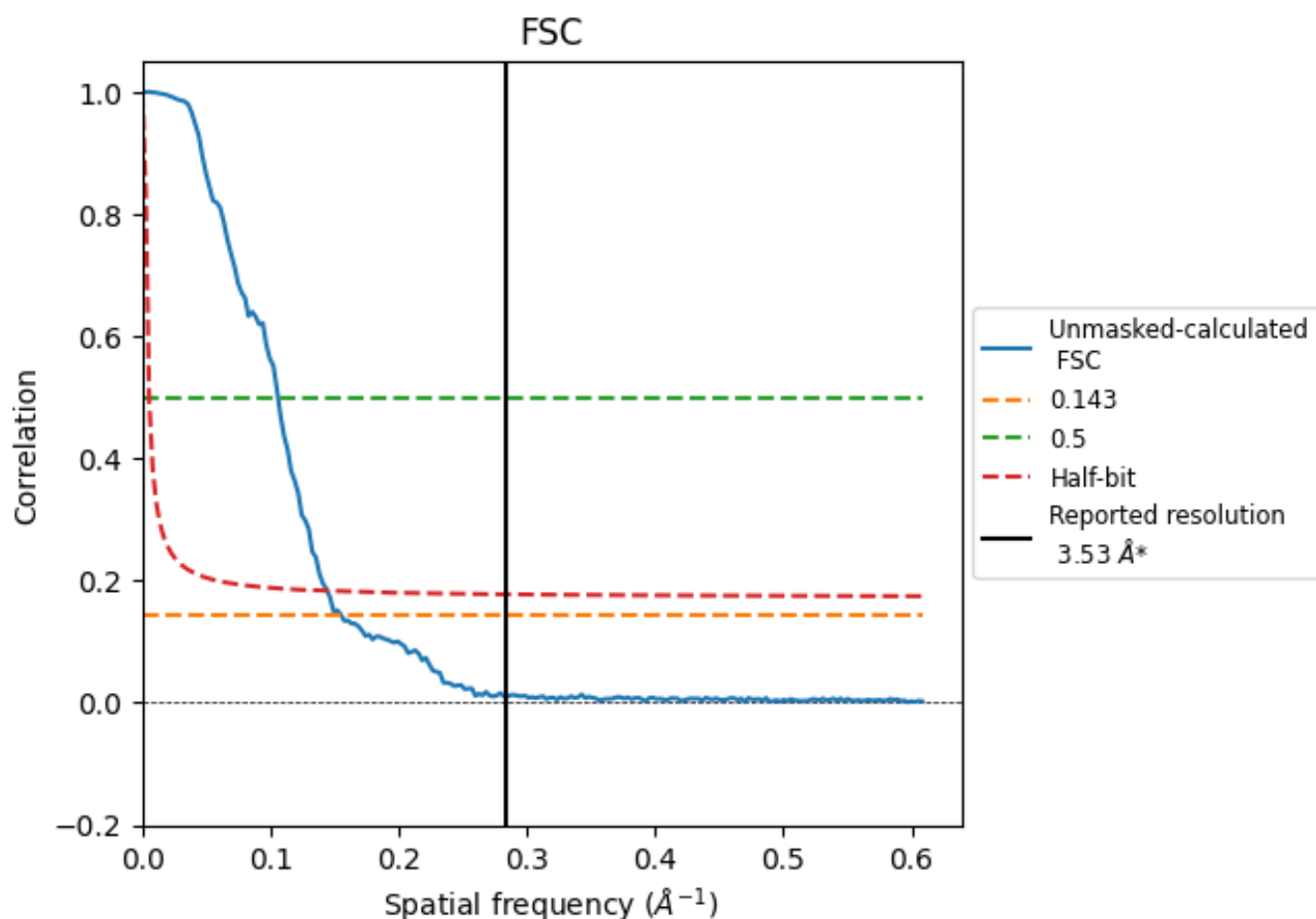


*Reported resolution corresponds to spatial frequency of 0.283 Å⁻¹

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 $0.283 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.53	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	6.44	9.44	6.93

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 6.44 differs from the reported value 3.53 by more than 10 %

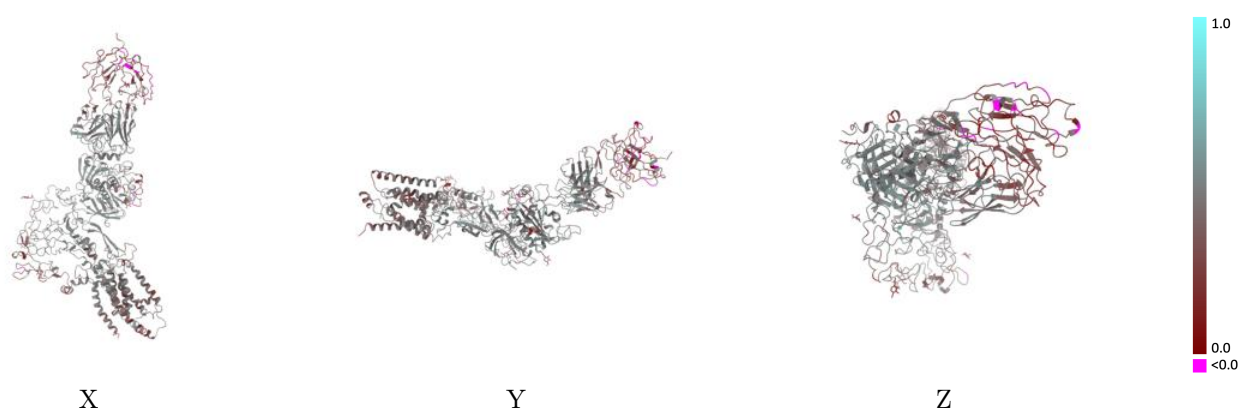
9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-70127 and PDB model 9O58. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)

This section was not generated.

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

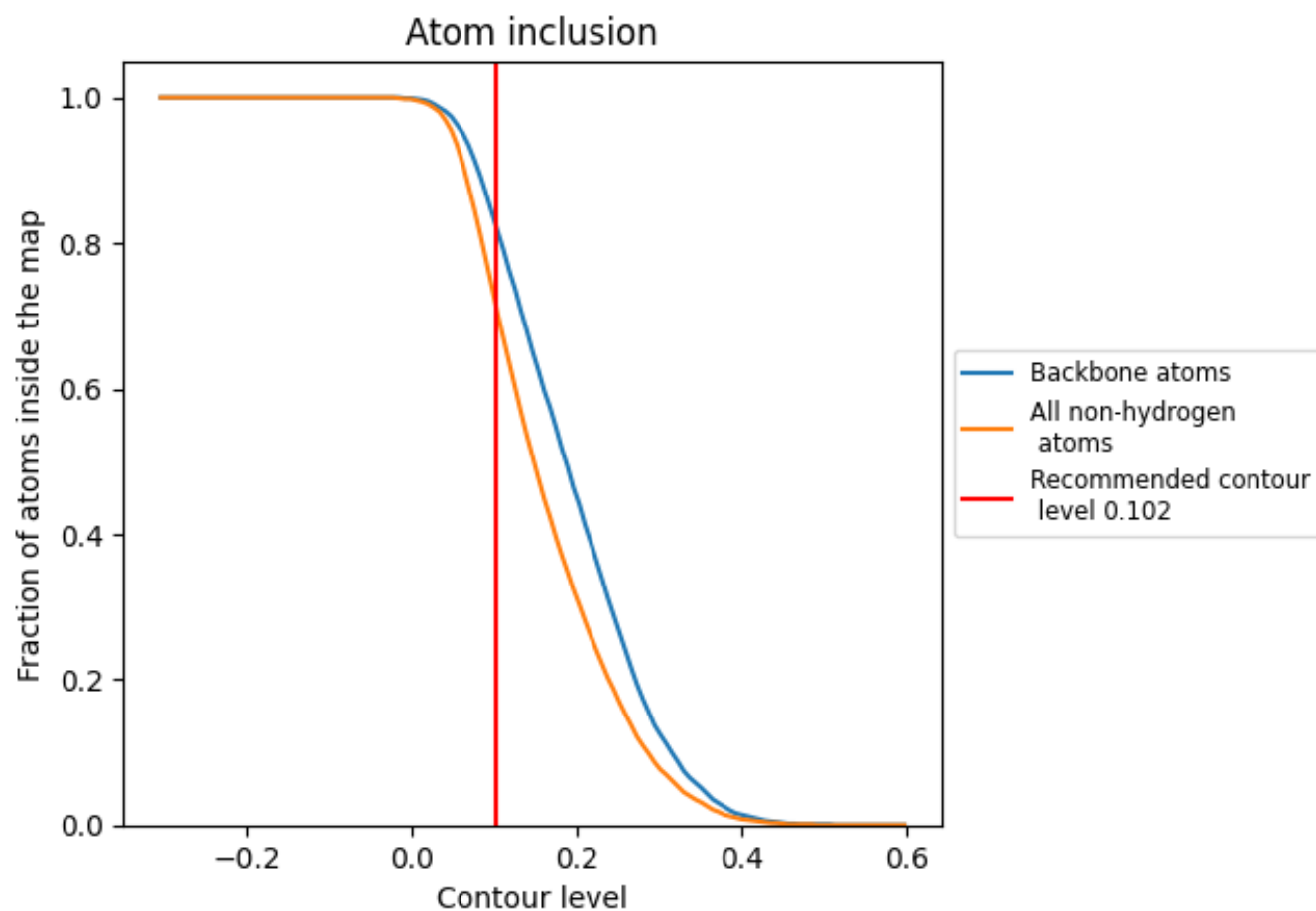


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 72% 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 (0.102) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7160	0.4270
A	0.7570	0.4480
B	0.7780	0.4530
C	0.5410	0.3510
D	0.6130	0.3740
E	0.6430	0.3990

