



Full wwPDB EM Validation Report ⓘ

Mar 5, 2026 – 08:16 AM UTC

PDB ID : 8DDX / pdb_00008ddx
EMDB ID : EMD-27345
Title : cryo-EM structure of TRPM3 ion channel in complex with Gbg in the presence of PIP2, tethered by ALFA-nanobody
Authors : Zhao, C.; MacKinnon, R.
Deposited on : 2022-06-19
Resolution : 3.80 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

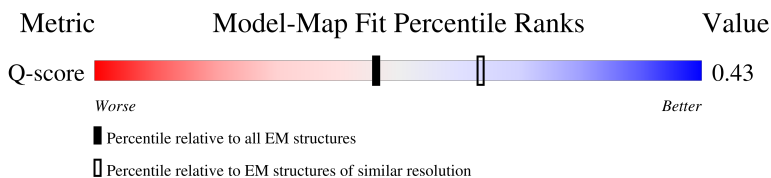
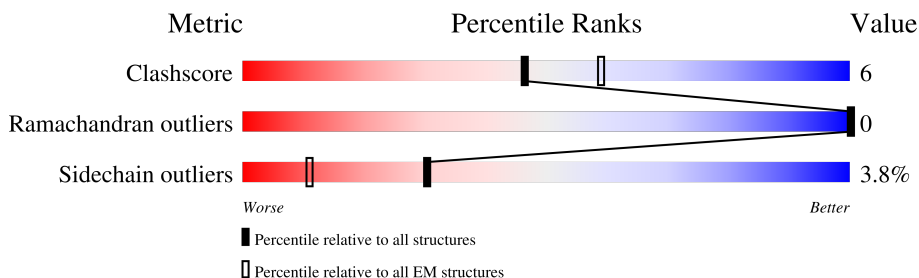
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10198 (3.30 - 4.30)

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

Mol	Chain	Length	Quality of chain
1	A	1370	
1	B	1370	
1	C	1370	
1	D	1370	

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Mol	Chain	Length	Quality of chain
2	E	17	88% 12%
2	F	17	88% 12%
2	G	17	82% 18%
2	H	17	94% 6%
3	I	340	27% 68% 29% ..
4	J	70	51% 66% 23% 11%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 35500 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 Transient receptor potential cation channel, subfamily M, member 3.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	968	Total	C	N	O	S	2	0
			7841	5083	1327	1373	58		
1	B	968	Total	C	N	O	S	2	0
			7841	5083	1327	1373	58		
1	C	968	Total	C	N	O	S	2	0
			7841	5083	1327	1373	58		
1	D	992	Total	C	N	O	S	2	0
			8030	5205	1357	1409	59		

- Molecule 2 is a protein called Unidentified segment at the N-terminus of TRPM3.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	E	17	Total	C	N	O	0	0
			85	51	17	17		
2	F	17	Total	C	N	O	0	0
			85	51	17	17		
2	G	17	Total	C	N	O	0	0
			85	51	17	17		
2	H	17	Total	C	N	O	0	0
			85	51	17	17		

- Molecule 3 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	I	337	Total	C	N	O	S	2	0
			2598	1603	466	506	23		

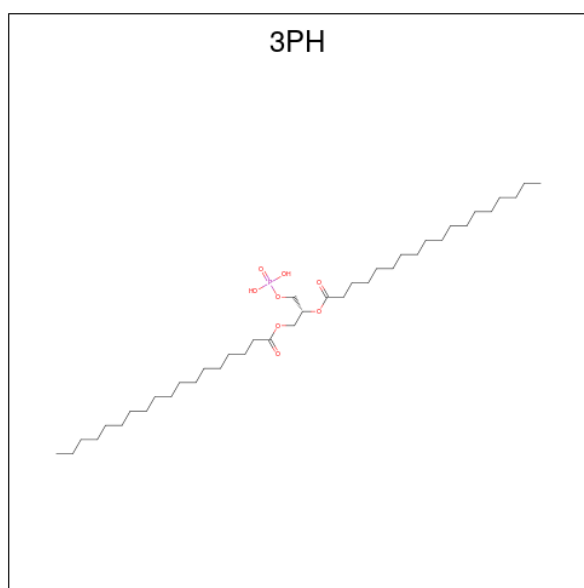
- Molecule 4 is a protein called Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	J	62	Total	C	N	O	S	0	0
			484	307	84	90	3		

There is a discrepancy between the modelled and reference sequences:

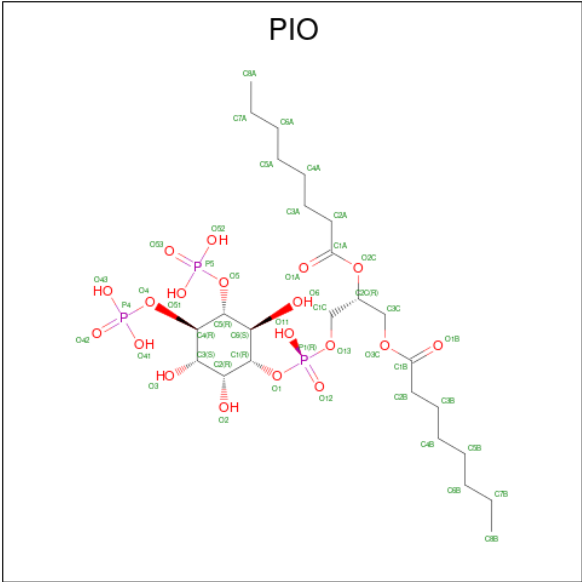
Chain	Residue	Modelled	Actual	Comment	Reference
J	68	SER	CYS	engineered mutation	UNP P59768

- Molecule 5 is 1,2-DIACYL-GLYCEROL-3-SN-PHOSPHATE (CCD ID: 3PH) (formula: $C_{39}H_{77}O_8P$).



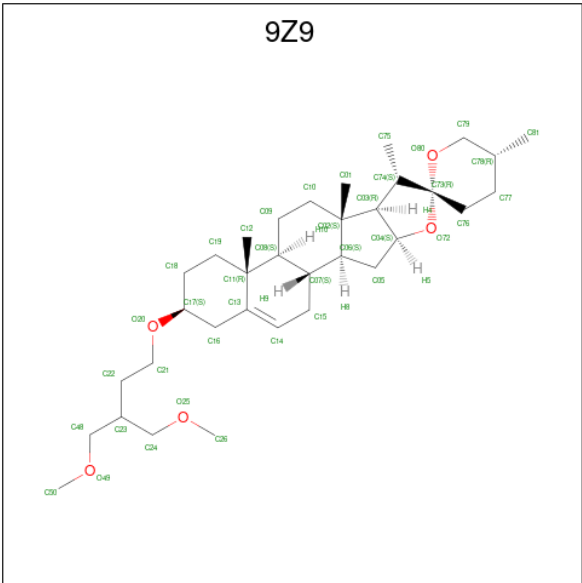
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	O	P	0
			45	36	8	1	
5	B	1	Total	C	O	P	0
			45	36	8	1	
5	C	1	Total	C	O	P	0
			45	36	8	1	
5	D	1	Total	C	O	P	0
			45	36	8	1	

- Molecule 6 is [(2R)-2-octanoyloxy-3-[oxidanyl-[(1R,2R,3S,4R,5R,6S)-2,3,6-tris(oxidanyl)-4,5-diphosphonooxy-cyclohexyl]oxy-phosphoryl]oxy-propyl] octanoate (CCD ID: PIO) (formula: $C_{25}H_{49}O_{19}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
6	A	1	Total	C	O	P	0
			47	25	19	3	
6	B	1	Total	C	O	P	0
			47	25	19	3	
6	C	1	Total	C	O	P	0
			47	25	19	3	
6	D	1	Total	C	O	P	0
			47	25	19	3	

- Molecule 7 is (3beta,14beta,17beta,25R)-3-[4-methoxy-3-(methoxymethyl)butoxy]spirost-5-en (CCD ID: 9Z9) (formula: C₃₄H₅₆O₅).



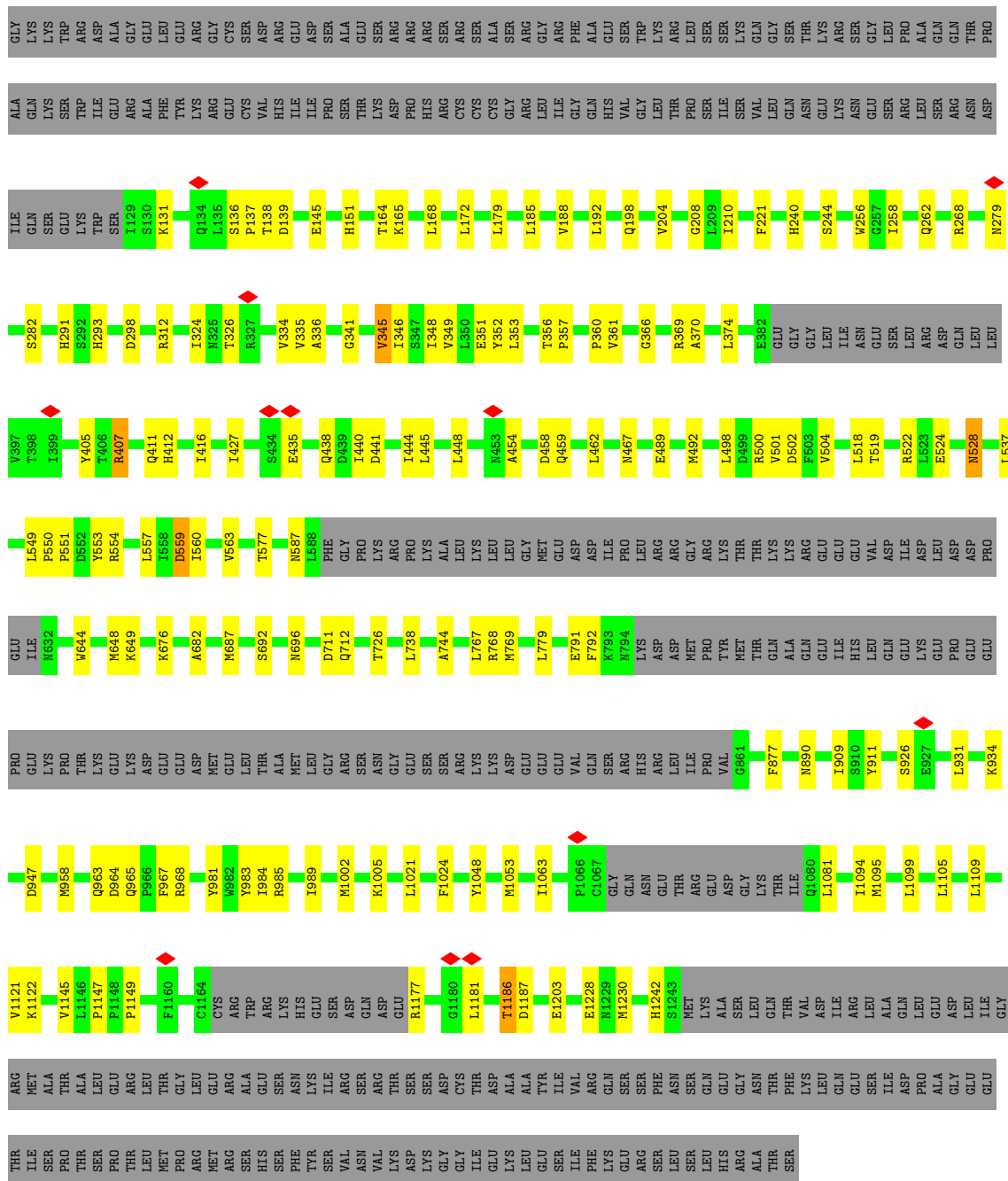
Mol	Chain	Residues	Atoms			AltConf
7	A	1	Total	C	O	0
			39	34	5	
7	B	1	Total	C	O	0
			39	34	5	
7	C	1	Total	C	O	0
			39	34	5	
7	C	1	Total	C	O	0
			39	34	5	

- Molecule 8 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		AltConf
8	C	1	Total	Na	0
			1	1	

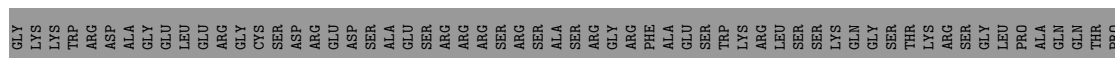
- Molecule 1: Transient receptor potential cation channel, subfamily M, member 3

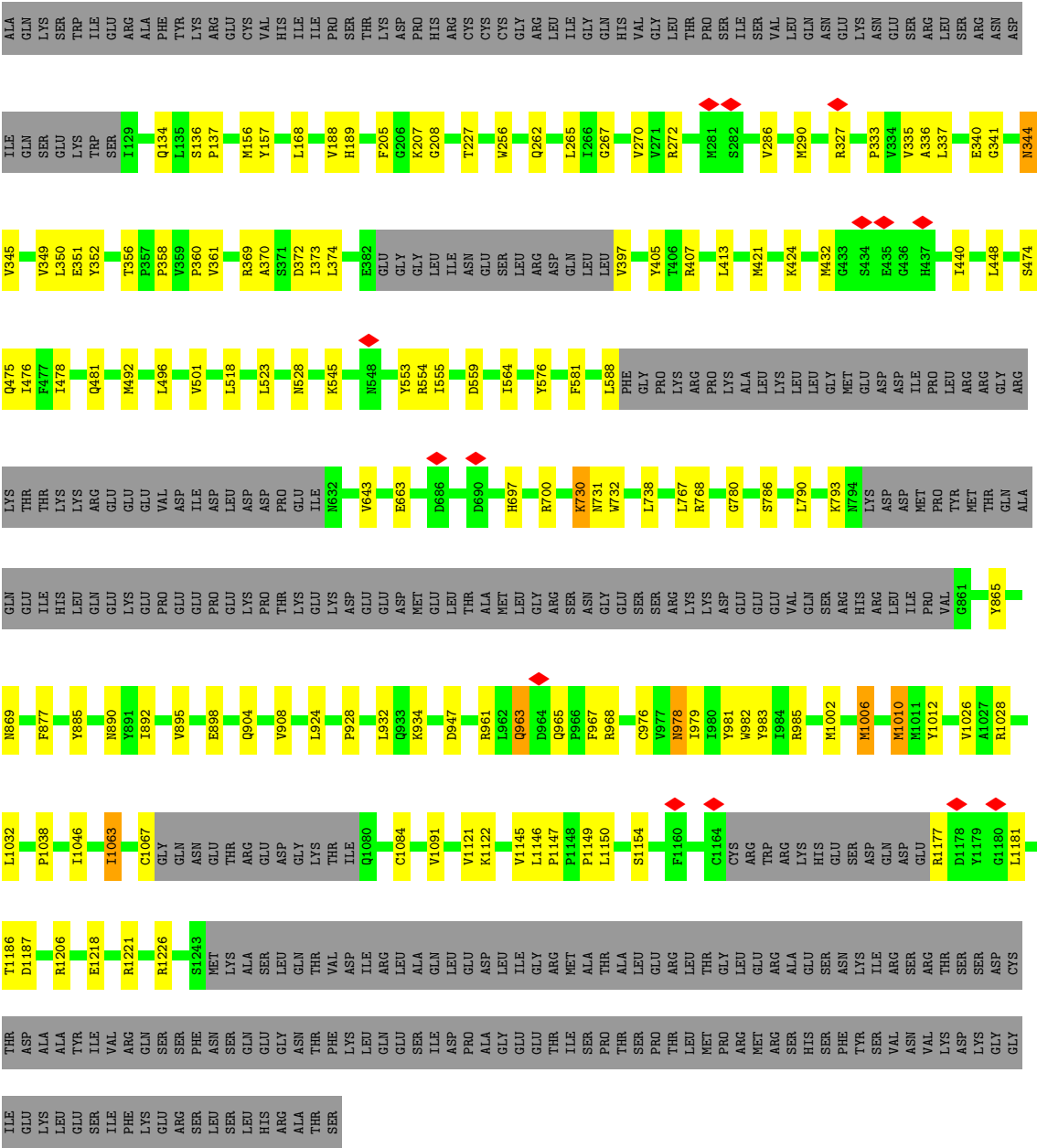
Chain B:



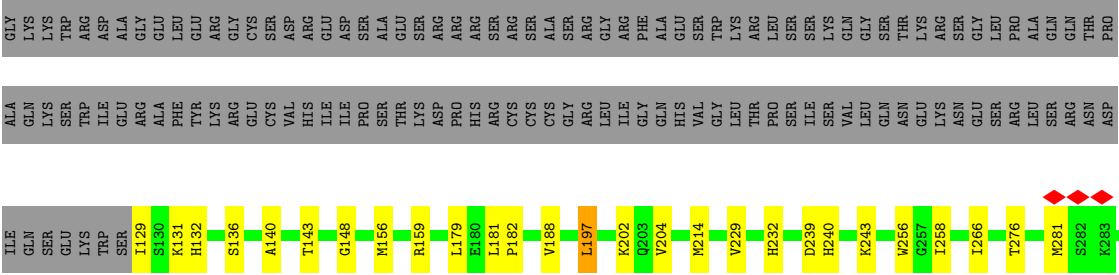
- Molecule 1: Transient receptor potential cation channel, subfamily M, member 3

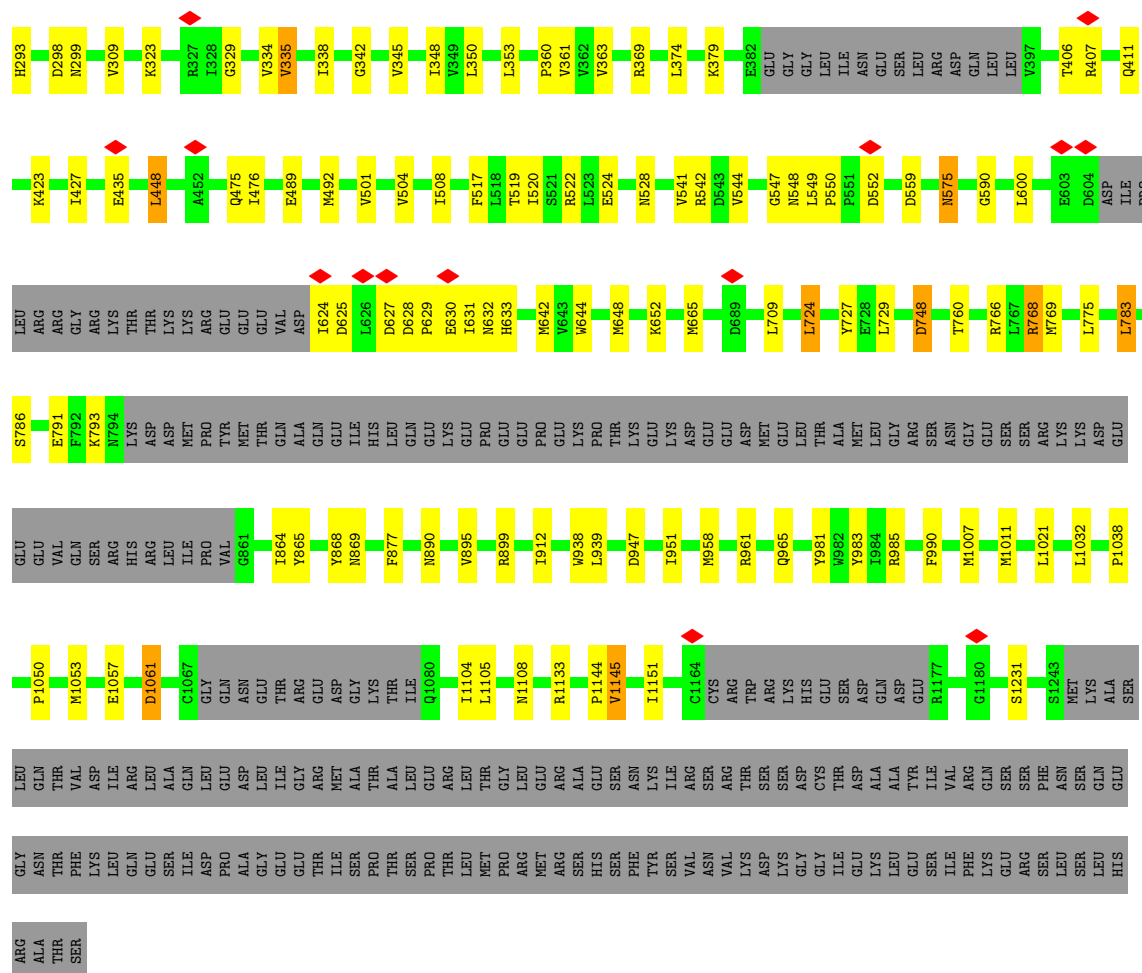
Chain C:





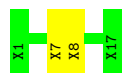
• Molecule 1: Transient receptor potential cation channel, subfamily M, member 3





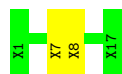
- Molecule 2: Unidentified segment at the N-terminus of TRPM3

Chain E: 88% 12%



- Molecule 2: Unidentified segment at the N-terminus of TRPM3

Chain F: 88% 12%



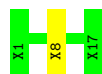
- Molecule 2: Unidentified segment at the N-terminus of TRPM3

Chain G: 82% 18%



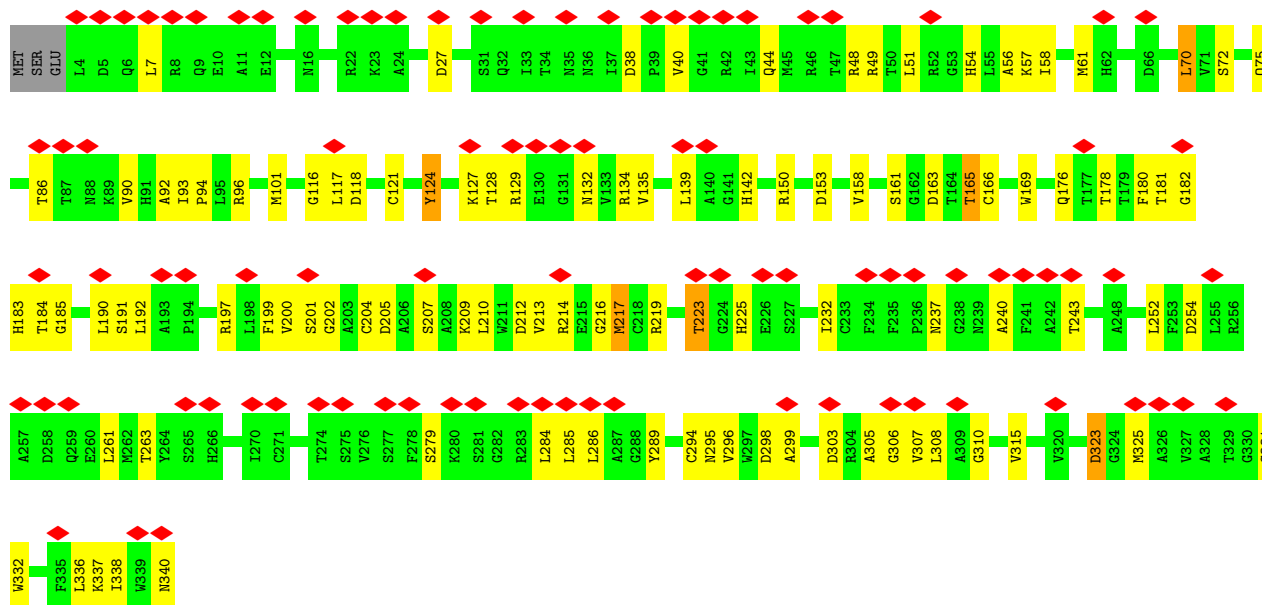
- Molecule 2: Unidentified segment at the N-terminus of TRPM3

Chain H:  94% 6%



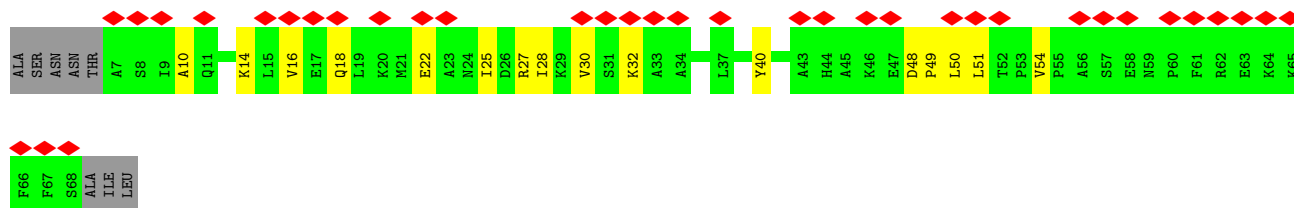
- Molecule 3: Guanine nucleotide-binding protein G(I)/G(S)/G(T) subunit beta-1

Chain I:  27% 68% 29% ..



- Molecule 4: Guanine nucleotide-binding protein G(I)/G(S)/G(O) subunit gamma-2

Chain J:  51% 66% 23% 11%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	47972	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 ($e^-/\text{\AA}^2$)	51.4	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.081	Depositor
Minimum map value	-0.041	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.012	Depositor
Map size (Å)	414.72003, 414.72003, 414.72003	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	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: 3PH, 9Z9, NA, PIO

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	0.15	0/8023	0.31	0/10848
1	B	0.14	0/8023	0.30	0/10848
1	C	0.15	0/8023	0.31	0/10848
1	D	0.16	0/8216	0.32	0/11108
3	I	0.09	0/2651	0.28	0/3593
4	J	0.10	0/492	0.29	0/661
All	All	0.15	0/35428	0.31	0/47906

There are no bond length outliers.

There are no bond angle outliers.

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	7841	0	7955	67	0
1	B	7841	0	7955	86	0
1	C	7841	0	7955	78	0
1	D	8030	0	8151	85	0
2	E	85	0	23	2	0
2	F	85	0	23	2	0
2	G	85	0	24	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	85	0	24	1	0
3	I	2598	0	2509	66	0
4	J	484	0	496	12	0
5	A	45	0	66	1	0
5	B	45	0	66	3	0
5	C	45	0	66	1	0
5	D	45	0	66	1	0
6	A	47	0	44	1	0
6	B	47	0	44	1	0
6	C	47	0	44	1	0
6	D	47	0	44	1	0
7	A	39	0	0	14	0
7	B	39	0	0	15	0
7	C	78	0	0	29	0
8	C	1	0	0	0	0
All	All	35500	0	35555	435	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 6.

All (435) 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)
7:A:1403:9Z9:C73	7:A:1403:9Z9:C74	1.81	1.58
7:C:1403:9Z9:C73	7:C:1403:9Z9:C74	1.81	1.57
7:A:1403:9Z9:C07	7:A:1403:9Z9:C06	1.84	1.56
7:B:1401:9Z9:C73	7:B:1401:9Z9:C74	1.81	1.54
7:B:1401:9Z9:C06	7:B:1401:9Z9:C07	1.84	1.53
7:C:1401:9Z9:C73	7:C:1401:9Z9:C74	1.81	1.53
7:C:1403:9Z9:C06	7:C:1403:9Z9:C07	1.84	1.52
7:C:1401:9Z9:C06	7:C:1401:9Z9:C07	1.84	1.51
7:C:1401:9Z9:C74	7:C:1401:9Z9:C75	1.90	1.48
7:C:1403:9Z9:C74	7:C:1403:9Z9:C75	1.90	1.47
7:B:1401:9Z9:C74	7:B:1401:9Z9:C75	1.90	1.47
7:A:1403:9Z9:C74	7:A:1403:9Z9:C75	1.90	1.46
7:C:1401:9Z9:C75	7:C:1401:9Z9:C03	2.30	1.08
7:A:1403:9Z9:C75	7:A:1403:9Z9:C03	2.32	1.06
7:C:1403:9Z9:C75	7:C:1403:9Z9:C03	2.34	1.06
7:B:1401:9Z9:C75	7:B:1401:9Z9:C03	2.34	1.03
7:A:1403:9Z9:C06	7:A:1403:9Z9:C02	1.78	1.02
7:C:1403:9Z9:C06	7:C:1403:9Z9:C02	1.78	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:C:1401:9Z9:C06	7:C:1401:9Z9:C02	1.78	0.98
7:B:1401:9Z9:C06	7:B:1401:9Z9:C02	1.78	0.95
7:B:1401:9Z9:C74	7:B:1401:9Z9:O80	2.14	0.95
7:C:1403:9Z9:C74	7:C:1403:9Z9:O80	2.14	0.95
7:C:1401:9Z9:C74	7:C:1401:9Z9:O80	2.15	0.93
7:A:1403:9Z9:C74	7:A:1403:9Z9:O80	2.15	0.92
1:D:520:ILE:HG13	1:D:632:ASN:ND2	1.89	0.88
7:B:1401:9Z9:C07	7:B:1401:9Z9:C05	2.51	0.84
7:C:1401:9Z9:C07	7:C:1401:9Z9:C05	2.52	0.83
1:D:520:ILE:HG13	1:D:632:ASN:HD21	1.43	0.81
7:A:1403:9Z9:C07	7:A:1403:9Z9:C05	2.49	0.81
7:A:1403:9Z9:C06	7:A:1403:9Z9:C15	2.61	0.79
7:C:1401:9Z9:C06	7:C:1401:9Z9:C15	2.61	0.79
1:B:407:ARG:HG3	1:B:411:GLN:HB3	1.65	0.79
7:B:1401:9Z9:C06	7:B:1401:9Z9:C15	2.61	0.79
7:C:1403:9Z9:C06	7:C:1403:9Z9:C15	2.62	0.78
7:B:1401:9Z9:C06	7:B:1401:9Z9:C10	2.63	0.77
7:C:1403:9Z9:C07	7:C:1403:9Z9:C05	2.52	0.76
7:A:1403:9Z9:C06	7:A:1403:9Z9:C10	2.63	0.75
1:D:520:ILE:H	1:D:632:ASN:HD21	1.33	0.74
7:C:1403:9Z9:C06	7:C:1403:9Z9:C10	2.65	0.74
7:C:1401:9Z9:C06	7:C:1401:9Z9:C10	2.66	0.73
3:I:212:ASP:HB3	3:I:216:GLY:H	1.54	0.72
1:A:1028:ARG:HH11	1:A:1063:ILE:HD13	1.56	0.70
1:D:520:ILE:N	1:D:632:ASN:HD21	1.90	0.70
7:C:1401:9Z9:C75	7:C:1401:9Z9:O80	2.40	0.69
7:B:1401:9Z9:C75	7:B:1401:9Z9:O80	2.40	0.69
7:A:1403:9Z9:C75	7:A:1403:9Z9:O80	2.41	0.69
1:C:189:HIS:HE2	1:C:352:TYR:HH	1.40	0.69
7:C:1403:9Z9:C75	7:C:1403:9Z9:O80	2.40	0.69
1:B:559:ASP:OD1	1:B:559:ASP:N	2.25	0.69
1:B:931:LEU:HA	1:B:934:LYS:HE2	1.73	0.69
7:C:1403:9Z9:C06	7:C:1403:9Z9:C08	2.70	0.68
7:A:1403:9Z9:C06	7:A:1403:9Z9:C08	2.70	0.68
7:C:1403:9Z9:C06	7:C:1403:9Z9:C01	2.72	0.67
1:D:628:ASP:O	1:D:629:PRO:C	2.38	0.67
1:B:208:GLY:HA3	1:B:440:ILE:HB	1.77	0.67
7:C:1401:9Z9:C06	7:C:1401:9Z9:C08	2.71	0.66
1:D:407:ARG:HH21	1:D:411:GLN:HE21	1.42	0.66
1:D:520:ILE:H	1:D:632:ASN:ND2	1.93	0.66
7:B:1401:9Z9:C06	7:B:1401:9Z9:C08	2.71	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:548:ASN:ND2	3:I:92:ALA:O	2.28	0.65
1:D:766:ARG:HE	1:D:793:LYS:HD2	1.60	0.65
3:I:48:ARG:HG3	3:I:340:ASN:HB3	1.77	0.65
3:I:254:ASP:HB3	3:I:261:LEU:HD11	1.79	0.65
1:B:256:TRP:HB3	1:B:298:ASP:HB3	1.77	0.65
7:B:1401:9Z9:C06	7:B:1401:9Z9:C01	2.73	0.65
1:B:137:PRO:HA	2:F:7:UNK:HA	1.79	0.64
1:D:353:LEU:HD11	1:D:427:ILE:HD11	1.78	0.64
1:D:1061:ASP:OD1	1:D:1061:ASP:N	2.30	0.64
1:B:185:LEU:HB3	1:B:334:VAL:HG23	1.80	0.64
7:C:1401:9Z9:C06	7:C:1401:9Z9:C01	2.74	0.64
1:D:524:GLU:O	1:D:528:ASN:ND2	2.28	0.64
1:D:1144:PRO:HB2	1:D:1151:ILE:HD11	1.79	0.63
7:A:1403:9Z9:C06	7:A:1403:9Z9:C01	2.73	0.63
1:B:369:ARG:NH1	1:B:405:TYR:OH	2.31	0.63
1:B:1177:ARG:HG3	1:B:1181:LEU:HD22	1.81	0.63
1:D:552:ASP:O	3:I:96:ARG:NH1	2.32	0.63
1:D:590:GLY:HA2	3:I:117:LEU:HD13	1.82	0.62
3:I:323:ASP:OD1	3:I:323:ASP:N	2.31	0.62
1:B:1002:MET:HE2	1:B:1122:LYS:HA	1.82	0.62
1:A:326:THR:OG1	1:A:330:GLN:NE2	2.34	0.61
1:B:965:GLN:HG3	1:B:968:ARG:HH21	1.65	0.61
1:C:1028:ARG:HD2	1:C:1063:ILE:HD13	1.81	0.61
1:C:1177:ARG:HG2	1:C:1181:LEU:HD12	1.82	0.61
1:A:470:ASP:OD1	1:A:470:ASP:N	2.34	0.61
1:B:492:MET:HE1	1:B:504:VAL:HG13	1.82	0.61
3:I:121:CYS:HB3	3:I:139:LEU:HB2	1.82	0.61
3:I:127:LYS:O	3:I:129:ARG:NH1	2.34	0.61
1:B:963:GLN:HG2	1:B:967:PHE:HD2	1.64	0.61
1:C:344:ASN:N	1:C:344:ASN:OD1	2.33	0.61
1:B:138:THR:O	1:B:268:ARG:NH2	2.34	0.61
1:D:632:ASN:C	1:D:632:ASN:HD22	2.09	0.60
1:C:349:VAL:HG23	1:C:361:VAL:HG11	1.83	0.60
1:C:1006:MET:SD	1:C:1006:MET:N	2.75	0.60
1:B:489:GLU:OE2	1:B:522:ARG:NH1	2.34	0.60
1:C:205:PHE:HA	1:C:440:ILE:HG13	1.83	0.60
1:C:963:GLN:HG2	1:C:967:PHE:HD2	1.66	0.59
3:I:51:LEU:HB2	3:I:336:LEU:HB3	1.85	0.59
3:I:232:ILE:HG22	3:I:243:THR:HA	1.85	0.59
1:B:335:VAL:HG23	1:B:360:PRO:HG2	1.85	0.58
1:A:559:ASP:N	1:A:559:ASP:OD1	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:692:SER:O	1:B:696:ASN:ND2	2.34	0.58
1:C:341:GLY:HA3	1:C:345:VAL:HG21	1.86	0.58
3:I:207:SER:HB3	3:I:223:THR:HG23	1.84	0.58
3:I:128:THR:OG1	3:I:132:ASN:O	2.21	0.58
3:I:279:SER:OG	4:J:48:ASP:OD2	2.22	0.58
1:A:524:GLU:O	1:A:528:ASN:ND2	2.36	0.57
1:C:890:ASN:ND2	1:C:983:TYR:OH	2.36	0.57
1:C:1002:MET:HE2	1:C:1122:LYS:HA	1.86	0.57
1:B:676:LYS:NZ	1:B:744:ALA:O	2.32	0.57
1:D:890:ASN:ND2	1:D:983:TYR:OH	2.37	0.57
1:A:313:ARG:HG3	1:A:317:LYS:HE2	1.86	0.57
1:A:136:SER:O	2:E:8:UNK:N	2.38	0.57
1:A:208:GLY:HA3	1:A:440:ILE:HB	1.86	0.57
1:C:205:PHE:HE1	1:C:337:LEU:HD21	1.69	0.57
1:A:130:SER:O	1:A:132:HIS:ND1	2.34	0.56
1:A:325:ASN:O	1:A:327:ARG:NH1	2.38	0.56
1:C:947:ASP:OD2	1:C:985:ARG:NH2	2.38	0.56
1:D:338:ILE:HB	1:D:363:VAL:HG12	1.85	0.56
1:D:644:TRP:NE1	1:D:648:MET:SD	2.78	0.56
1:C:545:LYS:NZ	1:C:553:TYR:OH	2.38	0.56
1:B:524:GLU:O	1:B:528:ASN:ND2	2.39	0.56
1:B:767:LEU:HD23	1:B:769:MET:HE2	1.87	0.56
1:A:323:LYS:NZ	1:A:329:GLY:O	2.38	0.56
1:B:947:ASP:OD1	1:B:981:TYR:OH	2.21	0.56
1:C:965:GLN:OE1	1:C:968:ARG:NH2	2.39	0.56
1:C:981:TYR:HB2	5:C:1402:3PH:H3C2	1.87	0.56
1:B:947:ASP:OD2	1:B:985:ARG:NH2	2.36	0.55
3:I:296:VAL:O	3:I:305:ALA:N	2.38	0.55
1:A:492:MET:HE1	1:A:508:ILE:HG12	1.88	0.55
1:B:1186:THR:OG1	1:B:1187:ASP:N	2.37	0.55
1:A:152:SER:OG	1:A:154:LYS:NZ	2.37	0.55
1:C:136:SER:O	2:G:8:UNK:N	2.39	0.55
1:A:432:MET:SD	1:A:432:MET:N	2.77	0.55
1:A:1186:THR:OG1	1:A:1187:ASP:N	2.39	0.55
1:D:899:ARG:O	1:D:961:ARG:NH1	2.39	0.55
1:D:323:LYS:NZ	1:D:329:GLY:O	2.38	0.55
4:J:18:GLN:NE2	4:J:22:GLU:OE1	2.39	0.55
1:A:260:GLU:HG2	1:A:286:VAL:HG12	1.88	0.55
1:C:663:GLU:OE2	1:C:731:ASN:ND2	2.39	0.55
1:D:865:TYR:O	1:D:869:ASN:ND2	2.32	0.55
1:D:256:TRP:HB3	1:D:298:ASP:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:964:ASP:OD1	1:A:964:ASP:N	2.38	0.54
1:B:644:TRP:NE1	1:B:648:MET:SD	2.81	0.54
3:I:325:MET:O	3:I:340:ASN:ND2	2.34	0.54
4:J:25:ILE:O	4:J:27:ARG:NH1	2.40	0.54
3:I:296:VAL:N	3:I:306:GLY:O	2.40	0.54
1:D:630:GLU:O	1:D:631:ILE:C	2.51	0.54
3:I:56:ALA:HB1	3:I:75:GLN:HB3	1.90	0.54
1:B:145:GLU:OE1	1:B:145:GLU:N	2.41	0.54
3:I:40:VAL:HG11	4:J:51:LEU:HD11	1.90	0.54
1:A:760:THR:OG1	1:A:1133:ARG:NH2	2.41	0.53
1:B:467:ASN:HB2	1:B:500:ARG:HH21	1.74	0.53
1:B:498:LEU:HB3	1:B:500:ARG:HH11	1.74	0.53
1:D:148:GLY:HA3	1:D:276:THR:HG22	1.89	0.53
1:D:239:ASP:O	1:D:243:LYS:NZ	2.40	0.53
1:D:760:THR:OG1	1:D:1133:ARG:NH2	2.41	0.53
1:A:981:TYR:HB2	5:A:1401:3PH:H3C1	1.91	0.53
1:B:1105:LEU:HB2	1:C:1010:MET:HE2	1.90	0.53
3:I:166:CYS:HB2	3:I:180:PHE:HB2	1.90	0.53
1:D:136:SER:O	2:H:8:UNK:N	2.41	0.53
3:I:134:ARG:NH1	3:I:135:VAL:O	2.42	0.53
1:A:247:LYS:NZ	1:A:248:ILE:O	2.42	0.52
3:I:118:ASP:OD1	3:I:118:ASP:N	2.40	0.52
1:A:139:ASP:OD1	1:A:139:ASP:N	2.42	0.52
1:C:476:ILE:O	1:C:481:GLN:NE2	2.42	0.52
3:I:169:TRP:HA	3:I:176:GLN:HA	1.91	0.52
1:D:748:ASP:OD1	1:D:748:ASP:N	2.41	0.52
3:I:150:ARG:HH22	3:I:190:LEU:HD12	1.75	0.52
1:A:644:TRP:NE1	1:A:648:MET:SD	2.83	0.52
1:B:244:SER:O	1:C:1206:ARG:NH1	2.43	0.52
1:D:281:MET:SD	1:D:281:MET:N	2.83	0.52
1:C:270:VAL:HG13	1:C:272:ARG:HD3	1.92	0.51
3:I:217:MET:SD	3:I:217:MET:N	2.78	0.51
1:B:926:SER:HB3	1:B:934:LYS:HD2	1.91	0.51
3:I:7:LEU:HD22	4:J:16:VAL:HG11	1.92	0.51
1:C:898:GLU:O	1:C:961:ARG:NH2	2.43	0.51
1:A:169:LEU:HD11	1:A:311:LEU:HD11	1.93	0.51
3:I:240:ALA:HA	3:I:254:ASP:HA	1.93	0.51
1:D:1007:MET:HG2	1:D:1011[A]:MET:HE2	1.93	0.51
1:A:885:TYR:HE1	1:A:908:VAL:HG13	1.76	0.50
1:C:256:TRP:O	1:C:262:GLN:NE2	2.45	0.50
1:C:421:MET:HE3	1:C:424:LYS:HD2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1104:ILE:O	1:D:1108:ASN:ND2	2.44	0.50
3:I:252:LEU:HD22	3:I:261:LEU:HD12	1.92	0.50
1:D:632:ASN:ND2	1:D:632:ASN:O	2.44	0.50
3:I:315:VAL:HA	3:I:331:SER:HA	1.94	0.50
1:B:435:GLU:O	1:B:438:GLN:NE2	2.45	0.49
3:I:161:SER:OG	3:I:163:ASP:OD1	2.29	0.49
1:B:528:ASN:OD1	1:B:554:ARG:NH2	2.45	0.49
1:C:767:LEU:O	1:C:793:LYS:NZ	2.45	0.49
1:D:181:LEU:HD12	1:D:182:PRO:HD2	1.94	0.49
1:A:1067:CYS:HB3	1:A:1084:CYS:HA	1.95	0.49
1:B:1005:LYS:HD2	1:B:1121:VAL:HG13	1.95	0.49
1:D:197:LEU:HD11	1:D:202:LYS:HB3	1.94	0.49
3:I:286:LEU:HD23	3:I:296:VAL:HG13	1.94	0.49
1:D:489:GLU:HG3	1:D:517:PHE:HD1	1.78	0.49
3:I:61:MET:HB2	3:I:70:LEU:HD12	1.94	0.49
1:B:587:ASN:OD1	1:B:587:ASN:N	2.45	0.49
1:D:501:VAL:HG11	1:D:652:LYS:HE3	1.94	0.49
3:I:49:ARG:O	3:I:338:ILE:N	2.43	0.49
1:A:962:LEU:O	1:A:963:GLN:NE2	2.46	0.48
1:C:208:GLY:HA3	1:C:440:ILE:HB	1.94	0.48
3:I:57:LYS:HD2	3:I:332:TRP:HB3	1.95	0.48
1:A:964:ASP:O	1:A:968:ARG:NH2	2.45	0.48
1:C:413:LEU:HD23	1:C:413:LEU:H	1.77	0.48
1:A:963:GLN:HB2	1:A:967:PHE:HB2	1.96	0.48
1:A:1230:MET:HE2	1:D:1231:SER:HB2	1.95	0.48
1:D:624:ILE:O	1:D:625:ASP:C	2.55	0.48
1:B:911:TYR:HD2	1:B:1149:PRO:HG3	1.79	0.48
1:B:1024:PHE:HD1	1:B:1094:ILE:HG22	1.78	0.48
7:B:1401:9Z9:C06	7:B:1401:9Z9:C09	2.91	0.48
1:B:502:ASP:OD1	1:B:502:ASP:N	2.47	0.48
7:A:1403:9Z9:C07	7:A:1403:9Z9:C02	2.89	0.48
1:B:1109:LEU:HD13	1:C:1006:MET:HG2	1.96	0.48
1:D:665:MET:HG3	1:D:709:LEU:HD23	1.94	0.48
3:I:184:THR:HG22	3:I:205:ASP:HB3	1.94	0.48
1:D:129:ILE:HD11	1:D:132:HIS:HB2	1.95	0.48
1:D:229:VAL:HA	1:D:232:HIS:HD2	1.78	0.48
1:D:345:VAL:HA	1:D:348:ILE:HG22	1.95	0.48
1:A:478:ILE:HG22	1:A:479:TYR:H	1.78	0.48
7:A:1403:9Z9:C06	7:A:1403:9Z9:C09	2.91	0.48
1:B:192:LEU:HD23	1:B:192:LEU:H	1.78	0.48
1:B:198:GLN:NE2	1:B:366:GLY:O	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:356:THR:O	1:C:358:PRO:HD3	2.13	0.47
1:C:1218:GLU:OE2	1:C:1221:ARG:NH2	2.39	0.47
7:C:1401:9Z9:C07	7:C:1401:9Z9:C02	2.91	0.47
1:B:768:ARG:HE	1:B:791:GLU:HG2	1.78	0.47
3:I:38:ASP:OD1	3:I:38:ASP:N	2.47	0.47
1:B:890:ASN:ND2	1:B:983:TYR:OH	2.38	0.47
1:B:984:ILE:HD11	5:B:1402:3PH:H2D1	1.97	0.47
1:C:576:TYR:HA	1:C:581:PHE:HD2	1.80	0.47
1:D:492:MET:HE1	1:D:504:VAL:HG13	1.96	0.47
3:I:212:ASP:OD1	3:I:213:VAL:N	2.48	0.47
1:C:1146:LEU:HD12	1:C:1154:SER:HB2	1.96	0.47
3:I:197:ARG:HG2	3:I:214:ARG:HH22	1.79	0.47
1:A:568:MET:HE1	1:A:670:VAL:HG13	1.95	0.47
1:C:189:HIS:NE2	1:C:352:TYR:OH	2.41	0.47
1:D:143:THR:OG1	1:D:156:MET:SD	2.70	0.47
3:I:27:ASP:H	4:J:30:VAL:HG12	1.78	0.47
3:I:165:THR:HG22	3:I:181:THR:HG22	1.95	0.47
1:A:1028:ARG:HH22	1:A:1049:MET:HB3	1.78	0.47
1:B:981:TYR:HD1	5:B:1402:3PH:H371	1.79	0.47
1:C:928:PRO:HD2	1:C:934:LYS:HG2	1.96	0.47
7:C:1403:9Z9:C06	7:C:1403:9Z9:C09	2.92	0.47
1:D:542:ARG:NE	1:D:547:GLY:O	2.48	0.47
1:A:253:ILE:HG12	1:A:295:ILE:HB	1.96	0.47
1:B:221:PHE:O	1:B:312:ARG:NH2	2.47	0.47
1:B:989:ILE:HG22	6:B:1403:PIO:H5B	1.96	0.47
1:A:137:PRO:HA	2:E:7:UNK:HA	1.97	0.47
1:B:964:ASP:OD1	1:B:964:ASP:N	2.48	0.47
1:C:1038:PRO:HD3	1:D:895:VAL:HG13	1.96	0.47
3:I:51:LEU:N	3:I:336:LEU:O	2.46	0.47
3:I:323:ASP:HB2	3:I:325:MET:HE1	1.96	0.47
1:A:421:MET:HE3	1:A:424:LYS:HE3	1.97	0.46
1:B:279:ASN:HB3	1:B:282:SER:HB3	1.96	0.46
1:C:697:HIS:CD2	1:C:700:ARG:HH21	2.32	0.46
1:C:892:ILE:HD11	1:C:904:GLN:HB2	1.97	0.46
7:C:1401:9Z9:C06	7:C:1401:9Z9:C09	2.93	0.46
1:D:990:PHE:HE1	6:D:1402:PIO:H6B	1.79	0.46
7:C:1401:9Z9:C05	7:C:1401:9Z9:C01	2.86	0.46
1:A:256:TRP:HB3	1:A:298:ASP:HB3	1.97	0.46
1:B:519:THR:HG23	1:B:522:ARG:H	1.81	0.46
1:D:630:GLU:C	1:D:632:ASN:N	2.71	0.46
3:I:153:ASP:OD1	3:I:153:ASP:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:947:ASP:OD1	1:D:985:ARG:NH1	2.49	0.46
1:A:926:SER:HB3	1:A:934:LYS:HG2	1.98	0.46
1:B:454:ALA:O	1:B:459:GLN:NE2	2.46	0.46
5:D:1401:3PH:H3A2	5:D:1401:3PH:H3D2	1.79	0.46
3:I:284:LEU:HD23	4:J:50:LEU:HD13	1.97	0.46
1:B:256:TRP:O	1:B:262:GLN:NE2	2.48	0.46
7:B:1401:9Z9:C05	7:B:1401:9Z9:C01	2.86	0.46
1:B:210:ILE:HD13	1:B:240:HIS:HB2	1.97	0.46
1:B:256:TRP:NE1	1:B:262:GLN:OE1	2.49	0.46
3:I:289:TYR:HE2	3:I:295:ASN:H	1.63	0.46
1:A:1032:LEU:O	1:A:1085:LYS:NZ	2.45	0.46
1:B:768:ARG:NH2	1:B:792:PHE:O	2.45	0.46
1:C:156:MET:HE3	1:C:156:MET:HB2	1.85	0.46
1:C:327:ARG:HA	1:C:327:ARG:HD3	1.77	0.46
1:D:179:LEU:HD11	1:D:293:HIS:CE1	2.51	0.46
1:D:600:LEU:HD13	3:I:101:MET:HE2	1.98	0.46
1:B:136:SER:O	2:F:8:UNK:N	2.49	0.45
1:C:528:ASN:HD21	1:C:555:ILE:H	1.64	0.45
4:J:49:PRO:HB3	4:J:54:VAL:HG23	1.98	0.45
1:B:1053:MET:HE1	1:B:1063:ILE:HD13	1.98	0.45
1:C:137:PRO:HA	2:G:7:UNK:HA	1.99	0.45
1:C:474:SER:OG	1:C:475:GLN:OE1	2.25	0.45
1:C:134:GLN:N	2:G:10:UNK:O	2.38	0.45
1:B:441:ASP:HA	1:B:444:ILE:HG22	1.99	0.45
1:C:790:LEU:HD23	1:C:790:LEU:HA	1.82	0.45
1:D:140:ALA:HB3	1:D:159:ARG:HD2	1.98	0.45
1:C:424:LYS:HE3	1:C:424:LYS:HB2	1.83	0.45
1:B:131:LYS:HD3	1:B:131:LYS:HA	1.88	0.45
3:I:294[A]:CYS:HB2	3:I:308:LEU:HB2	1.98	0.45
3:I:298:ASP:N	3:I:303:ASP:O	2.47	0.45
1:C:370:ALA:HA	1:C:373:ILE:HB	1.97	0.45
1:C:978:ASN:HD22	1:C:982:TRP:HE1	1.65	0.45
3:I:183:HIS:ND1	3:I:205:ASP:OD2	2.38	0.45
1:B:1147:PRO:HB2	1:B:1149:PRO:HD2	1.99	0.45
1:C:207:LYS:HA	1:C:207:LYS:HD3	1.75	0.45
1:A:1003:ILE:HG23	1:D:1105:LEU:HD21	1.97	0.45
1:C:865:TYR:O	1:C:869:ASN:ND2	2.35	0.45
1:A:713:SER:HG	1:A:1200:CYS:HG	1.65	0.44
1:C:1002:MET:HG2	1:C:1121:VAL:HG12	1.98	0.44
3:I:210:LEU:HB3	3:I:219:ARG:HB2	1.98	0.44
1:A:679:ALA:HB2	1:A:695:LEU:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:768:ARG:HG2	1:D:791:GLU:HG3	1.99	0.44
1:D:958:MET:HE3	1:D:958:MET:HB3	1.93	0.44
1:B:549:LEU:HD21	1:B:553:TYR:HD2	1.81	0.44
1:D:868:TYR:O	1:D:1145:VAL:HG11	2.18	0.44
3:I:191:SER:O	3:I:199:PHE:HB2	2.18	0.44
1:A:773:SER:OG	6:A:1402:PIO:O11	2.28	0.44
1:C:924:LEU:O	1:C:934:LYS:NZ	2.42	0.44
1:B:346:ILE:HD11	1:B:370:ALA:HB1	2.00	0.44
1:D:520:ILE:CG1	1:D:632:ASN:HD21	2.23	0.44
1:B:682:ALA:HB1	1:B:687:MET:HG2	1.99	0.44
1:D:627:ASP:O	1:D:628:ASP:C	2.61	0.44
3:I:101:MET:HB2	3:I:116:GLY:HA2	2.00	0.44
3:I:294[B]:CYS:HB3	3:I:308:LEU:HB2	1.99	0.44
4:J:10:ALA:O	4:J:14:LYS:NZ	2.45	0.44
3:I:44:GLN:HE22	3:I:307:VAL:H	1.66	0.44
1:A:370:ALA:HA	1:A:373:ILE:HB	1.99	0.44
1:B:349:VAL:HG23	1:B:361:VAL:HG11	1.99	0.44
1:B:557:LEU:HA	1:B:560:ILE:HD12	1.99	0.44
1:A:299:ASN:OD1	1:A:299:ASN:N	2.51	0.44
1:B:353:LEU:HD11	1:B:427:ILE:HD11	1.99	0.44
1:B:649:LYS:HE3	1:B:649:LYS:HB2	1.88	0.44
1:C:157:TYR:OH	1:C:265:LEU:O	2.35	0.44
1:C:335:VAL:HG23	1:C:360:PRO:HG2	1.99	0.44
1:C:336:ALA:HB3	1:C:361:VAL:HG12	1.99	0.44
1:C:976:CYS:O	1:C:979:ILE:HG22	2.18	0.43
3:I:285:LEU:HG	3:I:299:ALA:HB2	2.00	0.43
1:A:1065:PRO:O	1:A:1067:CYS:N	2.51	0.43
1:B:345:VAL:HA	1:B:348:ILE:HG22	2.01	0.43
7:B:1401:9Z9:C07	7:B:1401:9Z9:C02	2.91	0.43
1:D:1021:LEU:HG	1:D:1050:PRO:HB2	1.99	0.43
1:B:165:LYS:HB2	1:B:168:LEU:HD23	2.01	0.43
1:D:541:VAL:HA	1:D:544:VAL:HG12	2.00	0.43
1:A:1063:ILE:HD11	1:A:1091:VAL:HG11	2.01	0.43
1:B:1063:ILE:HD11	1:B:1095:MET:SD	2.58	0.43
1:C:256:TRP:NE1	1:C:262:GLN:OE1	2.52	0.43
3:I:58:ILE:HD11	3:I:72:SER:HB2	1.99	0.43
3:I:310:GLY:O	3:I:337:LYS:NZ	2.46	0.43
1:A:324:ILE:HG23	1:A:326:THR:H	1.84	0.43
1:B:179:LEU:HD11	1:B:293:HIS:CE1	2.53	0.43
6:C:1404:PIO:H2AA	6:C:1404:PIO:H2C	1.73	0.43
1:D:353:LEU:HB3	1:D:423:LYS:HD3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:947:ASP:OD1	1:D:981:TYR:OH	2.30	0.43
1:B:324:ILE:HG23	1:B:326:THR:H	1.83	0.43
1:A:892:ILE:HG12	1:A:905:GLU:HB2	2.01	0.43
1:B:341:GLY:HA3	1:B:345:VAL:HG11	2.01	0.43
1:B:518:LEU:HD23	1:B:518:LEU:HA	1.86	0.43
1:B:712:GLN:NE2	1:B:1203:GLU:OE1	2.52	0.43
1:C:780:GLY:HA2	1:C:786:SER:HB3	2.01	0.43
3:I:237:ASN:HB3	4:J:40:TYR:CZ	2.54	0.43
1:A:488:LEU:HD23	1:A:488:LEU:HA	1.85	0.42
1:A:1081:LEU:HD23	1:A:1081:LEU:H	1.84	0.42
1:B:519:THR:HG22	1:B:522:ARG:HB2	2.01	0.42
1:C:1026:VAL:HG22	1:C:1046:ILE:HD13	2.01	0.42
3:I:201:SER:OG	3:I:202:GLY:N	2.52	0.42
1:B:407:ARG:H	1:B:407:ARG:HG2	1.45	0.42
1:A:919:LYS:HE3	1:A:938:TRP:HZ2	1.84	0.42
1:D:549:LEU:HD12	1:D:550:PRO:HD2	2.01	0.42
4:J:28:ILE:HB	4:J:32:LYS:HD3	2.01	0.42
1:D:575:ASN:N	1:D:575:ASN:OD1	2.50	0.42
1:B:550:PRO:HA	1:B:551:PRO:HD3	1.92	0.42
1:C:333:PRO:HB2	1:C:448:LEU:HD11	2.01	0.42
7:C:1403:9Z9:C07	7:C:1403:9Z9:C02	2.91	0.42
1:D:379:LYS:HB3	1:D:379:LYS:HE3	1.81	0.42
3:I:200:VAL:HG23	3:I:232:ILE:HD12	2.01	0.42
1:C:369:ARG:HD3	1:C:369:ARG:HA	1.79	0.42
1:D:214:MET:HE2	1:D:214:MET:HB3	1.98	0.42
1:A:345:VAL:HA	1:A:348:ILE:HG12	2.02	0.42
5:B:1402:3PH:H242	5:B:1402:3PH:H271	1.70	0.42
1:C:492:MET:HE3	1:C:492:MET:HB3	1.95	0.42
1:A:279:ASN:HB2	1:A:281:MET:SD	2.59	0.42
1:C:1146:LEU:HA	1:C:1147:PRO:HD3	1.92	0.42
1:D:519:THR:OG1	1:D:520:ILE:N	2.52	0.42
1:A:960:LEU:HD13	1:A:960:LEU:HA	1.93	0.42
1:D:724:LEU:HD23	1:D:727:TYR:HE1	1.84	0.42
1:A:200:LYS:HE2	1:A:200:LYS:HB2	1.93	0.42
1:A:369:ARG:HD3	1:A:369:ARG:HA	1.93	0.42
1:C:478:ILE:HG12	1:C:481:GLN:HE21	1.85	0.41
1:C:1147:PRO:HB2	1:C:1149:PRO:HD2	2.02	0.41
1:D:369:ARG:HD3	1:D:369:ARG:HA	1.82	0.41
1:B:336:ALA:HB3	1:B:361:VAL:HG12	2.01	0.41
1:C:369:ARG:O	1:C:372:ASP:N	2.53	0.41
1:C:1186:THR:OG1	1:C:1187:ASP:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:356:THR:HB	1:B:357:PRO:HD3	2.01	0.41
1:B:909:ILE:HG12	1:B:958:MET:HE3	2.00	0.41
1:B:560:ILE:HA	1:B:563:VAL:HG12	2.02	0.41
1:C:1187:ASP:N	1:C:1187:ASP:OD1	2.53	0.41
1:D:335:VAL:HG23	1:D:360:PRO:HG2	2.02	0.41
1:D:342:GLY:H	1:D:345:VAL:HG22	1.85	0.41
3:I:58:ILE:HD12	3:I:58:ILE:HA	1.79	0.41
1:A:964:ASP:HA	1:A:968:ARG:HH21	1.86	0.41
1:B:139:ASP:N	1:B:139:ASP:OD1	2.53	0.41
1:B:1099:LEU:HD23	1:B:1099:LEU:HA	1.95	0.41
1:C:730:LYS:H	1:C:730:LYS:NZ	2.18	0.41
1:A:556:SER:N	1:A:559:ASP:OD2	2.48	0.41
1:A:1231:SER:HB2	1:B:1230:MET:HE2	2.02	0.41
1:D:628:ASP:HB2	1:D:631:ILE:HB	2.03	0.41
1:C:663:GLU:OE1	1:C:732:TRP:NE1	2.51	0.41
1:D:489:GLU:HG3	1:D:517:PHE:CD1	2.55	0.41
1:A:348:ILE:O	1:A:351:GLU:HG3	2.20	0.41
1:B:1228:GLU:HB2	1:C:1226:ARG:NH2	2.35	0.41
1:C:885:TYR:HE1	1:C:908:VAL:HG13	1.85	0.41
1:C:1067:CYS:HB2	1:C:1084:CYS:HA	2.02	0.41
1:D:548:ASN:O	3:I:94:PRO:HG3	2.21	0.41
1:D:783:LEU:HD12	1:D:786:SER:HB2	2.03	0.41
3:I:27:ASP:OD1	3:I:27:ASP:N	2.54	0.41
4:J:25:ILE:HD12	4:J:25:ILE:HA	1.96	0.41
1:D:1053:MET:HE2	1:D:1053:MET:HB3	1.84	0.41
3:I:124:TYR:HD1	3:I:124:TYR:HA	1.73	0.41
3:I:185:GLY:HA3	3:I:204:CYS:HB2	2.03	0.41
1:A:568:MET:HG2	1:A:572:TYR:CG	2.56	0.40
1:C:518:LEU:HD12	1:C:518:LEU:HA	1.91	0.40
1:D:775:LEU:HD23	1:D:775:LEU:H	1.86	0.40
1:D:912:ILE:HG12	1:D:951:ILE:HG13	2.02	0.40
1:D:965:GLN:HE21	1:D:965:GLN:HB2	1.66	0.40
3:I:182:GLY:HA3	3:I:209:LYS:NZ	2.36	0.40
1:A:927:GLU:HB3	1:A:928:PRO:HD3	2.03	0.40
1:C:432:MET:HE2	1:C:432:MET:HB3	1.88	0.40
1:C:564:ILE:HD11	1:C:643:VAL:HG11	2.04	0.40
1:D:131:LYS:NZ	1:D:299:ASN:O	2.53	0.40
1:D:448:LEU:HD13	1:D:448:LEU:HA	1.89	0.40
1:A:345:VAL:HG23	1:A:348:ILE:HD11	2.04	0.40
1:A:422:LYS:HE3	1:A:422:LYS:HB3	1.89	0.40
1:A:496:LEU:HD23	1:A:526:LEU:HD11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:933:GLN:N	1:A:933:GLN:OE1	2.54	0.40
1:C:738:LEU:HD23	1:C:738:LEU:HA	1.89	0.40
1:D:938:TRP:HE3	1:D:939:LEU:HD12	1.86	0.40
1:D:1053:MET:HA	1:D:1057:GLU:O	2.22	0.40
1:C:267:GLY:HA3	1:C:272:ARG:HD3	2.02	0.40
1:A:895:VAL:HG13	1:D:1038:PRO:HD3	2.02	0.40
1:B:738:LEU:HD23	1:B:738:LEU:HA	1.89	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	A	956/1370 (70%)	914 (96%)	42 (4%)	0	100	100
1	B	956/1370 (70%)	913 (96%)	43 (4%)	0	100	100
1	C	956/1370 (70%)	907 (95%)	49 (5%)	0	100	100
1	D	980/1370 (72%)	933 (95%)	47 (5%)	0	100	100
3	I	337/340 (99%)	323 (96%)	14 (4%)	0	100	100
4	J	60/70 (86%)	58 (97%)	2 (3%)	0	100	100
All	All	4245/5890 (72%)	4048 (95%)	197 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	A	859/1218 (70%)	832 (97%)	27 (3%)	35	57
1	B	858/1218 (70%)	825 (96%)	33 (4%)	29	53
1	C	859/1218 (70%)	825 (96%)	34 (4%)	28	52
1	D	880/1218 (72%)	846 (96%)	34 (4%)	28	52
3	I	282/283 (100%)	266 (94%)	16 (6%)	18	44
4	J	51/57 (90%)	51 (100%)	0	100	100
All	All	3789/5212 (73%)	3645 (96%)	144 (4%)	30	53

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

Mol	Chain	Res	Type
1	A	170	LEU
1	A	204	VAL
1	A	240	HIS
1	A	258	ILE
1	A	260	GLU
1	A	265	LEU
1	A	276	THR
1	A	304	LYS
1	A	332	VAL
1	A	352	TYR
1	A	356	THR
1	A	405	TYR
1	A	421	MET
1	A	485	VAL
1	A	496	LEU
1	A	559	ASP
1	A	586	HIS
1	A	588	LEU
1	A	710	LEU
1	A	761	ASP
1	A	877	PHE
1	A	948	LEU
1	A	960	LEU
1	A	1145	VAL
1	A	1155	HIS
1	A	1186	THR
1	A	1197	GLU
1	B	151	HIS

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Mol	Chain	Res	Type
1	B	164	THR
1	B	172	LEU
1	B	188	VAL
1	B	204	VAL
1	B	258	ILE
1	B	291	HIS
1	B	345	VAL
1	B	351	GLU
1	B	352	TYR
1	B	374	LEU
1	B	407	ARG
1	B	412	HIS
1	B	416	ILE
1	B	445	LEU
1	B	448	LEU
1	B	458	ASP
1	B	462	LEU
1	B	501	VAL
1	B	528	ASN
1	B	537	LEU
1	B	559	ASP
1	B	577	THR
1	B	711	ASP
1	B	726	THR
1	B	779	LEU
1	B	877	PHE
1	B	1021	LEU
1	B	1048	TYR
1	B	1081	LEU
1	B	1145	VAL
1	B	1186	THR
1	B	1242	HIS
1	C	168	LEU
1	C	188	VAL
1	C	227	THR
1	C	286	VAL
1	C	290	MET
1	C	340	GLU
1	C	344	ASN
1	C	350	LEU
1	C	351	GLU
1	C	374	LEU

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Mol	Chain	Res	Type
1	C	397	VAL
1	C	405	TYR
1	C	407	ARG
1	C	496	LEU
1	C	501	VAL
1	C	523	LEU
1	C	554	ARG
1	C	559	ASP
1	C	588	LEU
1	C	730	LYS
1	C	768	ARG
1	C	877	PHE
1	C	895	VAL
1	C	932	LEU
1	C	963	GLN
1	C	978	ASN
1	C	1006	MET
1	C	1010	MET
1	C	1012	TYR
1	C	1032	LEU
1	C	1063	ILE
1	C	1091	VAL
1	C	1145	VAL
1	C	1150	LEU
1	D	188	VAL
1	D	197	LEU
1	D	204	VAL
1	D	240	HIS
1	D	258	ILE
1	D	266	ILE
1	D	309	VAL
1	D	334	VAL
1	D	335	VAL
1	D	350	LEU
1	D	361	VAL
1	D	374	LEU
1	D	406	THR
1	D	435	GLU
1	D	448	LEU
1	D	475	GLN
1	D	476	ILE
1	D	508	ILE

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Mol	Chain	Res	Type
1	D	522	ARG
1	D	559	ASP
1	D	575	ASN
1	D	633	HIS
1	D	642	MET
1	D	724	LEU
1	D	729	LEU
1	D	748	ASP
1	D	768	ARG
1	D	769	MET
1	D	783	LEU
1	D	864	ILE
1	D	877	PHE
1	D	1032	LEU
1	D	1061	ASP
1	D	1145	VAL
3	I	54	HIS
3	I	70	LEU
3	I	86	THR
3	I	90	VAL
3	I	93	ILE
3	I	124	TYR
3	I	142	HIS
3	I	158	VAL
3	I	165	THR
3	I	178	THR
3	I	192	LEU
3	I	217	MET
3	I	223	THR
3	I	225	HIS
3	I	263	THR
3	I	323	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (67) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	232	HIS
1	A	293	HIS
1	A	314	GLN
1	A	378	HIS
1	A	481	GLN
1	A	535	ASN

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Mol	Chain	Res	Type
1	A	719	GLN
1	A	963	GLN
1	A	993	ASN
1	A	1127	GLN
1	A	1213	ASN
1	A	1229	ASN
1	A	1238	ASN
1	B	151	HIS
1	B	153	ASN
1	B	314	GLN
1	B	330	GLN
1	B	411	GLN
1	B	481	GLN
1	B	490	GLN
1	B	639	HIS
1	B	680	HIS
1	B	933	GLN
1	B	978	ASN
1	B	1115	ASN
1	B	1116	ASN
1	B	1194	HIS
1	C	378	HIS
1	C	411	GLN
1	C	481	GLN
1	C	697	HIS
1	C	719	GLN
1	C	890	ASN
1	C	978	ASN
1	C	1115	ASN
1	C	1116	ASN
1	C	1229	ASN
1	C	1242	HIS
1	D	153	ASN
1	D	171	HIS
1	D	232	HIS
1	D	288	ASN
1	D	291	HIS
1	D	314	GLN
1	D	378	HIS
1	D	411	GLN
1	D	632	ASN
1	D	712	GLN

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Mol	Chain	Res	Type
1	D	716	GLN
1	D	746	HIS
1	D	965	GLN
1	D	978	ASN
1	D	1029	GLN
1	D	1062	GLN
1	D	1115	ASN
1	D	1116	ASN
1	D	1126	ASN
1	D	1127	GLN
1	D	1135	GLN
1	D	1161	GLN
1	D	1194	HIS
1	D	1199	GLN
1	D	1213	ASN
3	I	44	GLN
3	I	110	ASN
3	I	230	ASN
3	I	239	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 [i](#)

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 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	$\# Z > 2$
5	3PH	C	1402	-	44,44,47	1.16	3 (6%)	47,49,52	1.18	2 (4%)
7	9Z9	C	1403	-	44,44,44	9.37	22 (50%)	64,68,68	4.96	24 (37%)
6	PIO	C	1404	-	47,47,47	1.20	6 (12%)	62,65,65	1.05	3 (4%)
6	PIO	A	1402	-	47,47,47	1.19	6 (12%)	62,65,65	1.02	3 (4%)
6	PIO	B	1403	-	47,47,47	1.22	6 (12%)	62,65,65	1.06	3 (4%)
6	PIO	D	1402	-	47,47,47	1.22	6 (12%)	62,65,65	1.19	5 (8%)
7	9Z9	C	1401	-	44,44,44	9.37	22 (50%)	64,68,68	4.99	24 (37%)
7	9Z9	B	1401	-	44,44,44	9.37	22 (50%)	64,68,68	4.98	23 (35%)
5	3PH	B	1402	-	44,44,47	1.16	3 (6%)	47,49,52	1.20	2 (4%)
5	3PH	D	1401	-	44,44,47	1.17	3 (6%)	47,49,52	1.16	2 (4%)
5	3PH	A	1401	-	44,44,47	1.16	3 (6%)	47,49,52	1.19	2 (4%)
7	9Z9	A	1403	-	44,44,44	9.37	22 (50%)	64,68,68	4.99	26 (40%)

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	3PH	C	1402	-	-	24/46/46/49	-
7	9Z9	C	1403	-	-	0/12/100/100	0/6/6/6
6	PIO	C	1404	-	-	24/44/68/68	0/1/1/1
6	PIO	A	1402	-	-	14/44/68/68	0/1/1/1
6	PIO	B	1403	-	-	14/44/68/68	0/1/1/1
6	PIO	D	1402	-	-	26/44/68/68	0/1/1/1
7	9Z9	C	1401	-	-	2/12/100/100	0/6/6/6
7	9Z9	B	1401	-	-	1/12/100/100	0/6/6/6
5	3PH	B	1402	-	-	28/46/46/49	-
5	3PH	D	1401	-	-	22/46/46/49	-
5	3PH	A	1401	-	-	27/46/46/49	-
7	9Z9	A	1403	-	-	5/12/100/100	0/6/6/6

All (124) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	1401	9Z9	C02-C03	-29.63	1.01	1.56
7	A	1403	9Z9	C02-C03	-29.62	1.01	1.56
7	C	1403	9Z9	C02-C03	-29.59	1.01	1.56
7	B	1401	9Z9	C02-C03	-29.55	1.01	1.56
7	C	1401	9Z9	C03-C74	-25.02	0.83	1.54
7	A	1403	9Z9	C03-C74	-24.94	0.83	1.54
7	C	1403	9Z9	C03-C74	-24.86	0.84	1.54
7	B	1401	9Z9	C03-C74	-24.82	0.84	1.54
7	A	1403	9Z9	C05-C06	-23.29	1.04	1.54
7	B	1401	9Z9	C05-C06	-23.17	1.04	1.54
7	C	1403	9Z9	C05-C06	-23.10	1.04	1.54
7	C	1401	9Z9	C05-C06	-23.10	1.04	1.54
7	A	1403	9Z9	O72-C73	-18.67	1.03	1.42
7	B	1401	9Z9	O72-C73	-18.59	1.03	1.42
7	C	1403	9Z9	O72-C73	-18.57	1.03	1.42
7	C	1401	9Z9	O72-C73	-18.46	1.03	1.42
7	B	1401	9Z9	C75-C74	18.09	1.90	1.53
7	C	1403	9Z9	C75-C74	18.02	1.90	1.53
7	C	1401	9Z9	C75-C74	17.99	1.90	1.53
7	A	1403	9Z9	C75-C74	17.83	1.90	1.53
7	C	1403	9Z9	C07-C06	16.51	1.84	1.53
7	B	1401	9Z9	C07-C06	16.42	1.84	1.53
7	C	1401	9Z9	C07-C06	16.36	1.84	1.53
7	A	1403	9Z9	C07-C06	16.17	1.84	1.53
7	C	1401	9Z9	C73-C74	14.29	1.81	1.53
7	A	1403	9Z9	C73-C74	14.22	1.81	1.53
7	B	1401	9Z9	C73-C74	14.22	1.81	1.53
7	C	1403	9Z9	C73-C74	14.14	1.81	1.53
7	C	1401	9Z9	C02-C06	12.92	1.78	1.55
7	C	1403	9Z9	C02-C06	12.85	1.78	1.55
7	B	1401	9Z9	C02-C06	12.83	1.78	1.55
7	A	1403	9Z9	C02-C06	12.76	1.78	1.55
7	C	1403	9Z9	C10-C09	-12.18	1.29	1.53
7	B	1401	9Z9	C10-C09	-12.11	1.29	1.53
7	C	1401	9Z9	C10-C09	-12.08	1.29	1.53
7	A	1403	9Z9	C10-C09	-12.01	1.29	1.53
7	A	1403	9Z9	C15-C07	-9.99	1.37	1.53
7	C	1401	9Z9	C15-C07	-9.98	1.37	1.53
7	B	1401	9Z9	C15-C07	-9.89	1.37	1.53
7	C	1403	9Z9	C15-C07	-9.86	1.37	1.53
7	A	1403	9Z9	C10-C02	9.43	1.70	1.54
7	C	1401	9Z9	C10-C02	9.38	1.70	1.54
7	C	1403	9Z9	C10-C02	9.37	1.70	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	1401	9Z9	C10-C02	9.35	1.70	1.54
7	A	1403	9Z9	C03-C04	-5.87	1.43	1.54
7	C	1401	9Z9	C03-C04	-5.86	1.43	1.54
7	B	1401	9Z9	C14-C13	5.79	1.44	1.33
7	C	1403	9Z9	C14-C13	5.79	1.44	1.33
7	B	1401	9Z9	C03-C04	-5.77	1.43	1.54
7	C	1403	9Z9	C03-C04	-5.77	1.43	1.54
7	C	1401	9Z9	C14-C13	5.75	1.44	1.33
7	A	1403	9Z9	C14-C13	5.67	1.44	1.33
7	C	1403	9Z9	C09-C08	-4.84	1.45	1.53
7	B	1401	9Z9	C09-C08	-4.77	1.45	1.53
7	C	1401	9Z9	C09-C08	-4.73	1.46	1.53
7	A	1403	9Z9	C09-C08	-4.63	1.46	1.53
7	C	1403	9Z9	C15-C14	4.28	1.58	1.50
7	B	1401	9Z9	C15-C14	4.23	1.58	1.50
7	C	1401	9Z9	C15-C14	4.12	1.58	1.50
7	A	1403	9Z9	C15-C14	4.04	1.58	1.50
7	C	1401	9Z9	C19-C18	3.82	1.61	1.53
7	A	1403	9Z9	C19-C18	3.82	1.61	1.53
7	B	1401	9Z9	C19-C18	3.63	1.60	1.53
7	C	1403	9Z9	C19-C18	3.61	1.60	1.53
6	D	1402	PIO	P4-O4	3.53	1.65	1.59
6	B	1403	PIO	P4-O4	3.52	1.65	1.59
7	A	1403	9Z9	C11-C08	3.39	1.61	1.56
6	D	1402	PIO	P5-O5	3.38	1.65	1.59
7	B	1401	9Z9	C01-C02	3.30	1.59	1.54
6	A	1402	PIO	P5-O5	3.27	1.65	1.59
7	C	1403	9Z9	C01-C02	3.27	1.59	1.54
6	B	1403	PIO	P5-O5	3.25	1.65	1.59
7	C	1401	9Z9	C11-C08	3.23	1.61	1.56
6	C	1404	PIO	P4-O4	3.23	1.65	1.59
6	A	1402	PIO	P4-O4	3.22	1.65	1.59
7	C	1401	9Z9	C01-C02	3.21	1.59	1.54
6	C	1404	PIO	P5-O5	3.21	1.65	1.59
7	B	1401	9Z9	C11-C08	3.19	1.61	1.56
7	A	1403	9Z9	C01-C02	3.13	1.59	1.54
7	C	1403	9Z9	C11-C08	2.98	1.60	1.56
7	C	1403	9Z9	C07-C08	-2.94	1.48	1.53
5	C	1402	3PH	O21-C21	2.93	1.42	1.34
7	A	1403	9Z9	C07-C08	-2.93	1.48	1.53
5	D	1401	3PH	O31-C31	2.93	1.41	1.33
5	A	1401	3PH	O31-C31	2.93	1.41	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1402	3PH	O31-C31	2.92	1.41	1.33
5	A	1401	3PH	O21-C21	2.92	1.42	1.34
7	B	1401	9Z9	C07-C08	-2.91	1.48	1.53
7	B	1401	9Z9	C19-C11	-2.90	1.48	1.54
7	A	1403	9Z9	C19-C11	-2.88	1.48	1.54
7	C	1403	9Z9	C19-C11	-2.88	1.48	1.54
5	C	1402	3PH	O31-C31	2.88	1.41	1.33
5	D	1401	3PH	O21-C21	2.87	1.42	1.34
7	C	1401	9Z9	C19-C11	-2.83	1.49	1.54
5	B	1402	3PH	O21-C21	2.83	1.42	1.34
7	C	1401	9Z9	C07-C08	-2.79	1.48	1.53
7	A	1403	9Z9	C76-C73	2.75	1.56	1.51
7	C	1401	9Z9	C18-C17	2.72	1.58	1.51
7	A	1403	9Z9	C18-C17	2.72	1.58	1.51
7	B	1401	9Z9	C18-C17	2.71	1.58	1.51
6	B	1403	PIO	O2C-C2C	-2.66	1.40	1.46
7	C	1403	9Z9	C18-C17	2.66	1.58	1.51
5	B	1402	3PH	O21-C2	-2.64	1.40	1.46
6	C	1404	PIO	O2C-C2C	-2.61	1.40	1.46
7	B	1401	9Z9	C76-C73	2.61	1.55	1.51
5	C	1402	3PH	O21-C2	-2.60	1.40	1.46
5	D	1401	3PH	O21-C2	-2.60	1.40	1.46
7	C	1403	9Z9	C76-C73	2.56	1.55	1.51
5	A	1401	3PH	O21-C2	-2.55	1.40	1.46
7	C	1401	9Z9	C76-C73	2.55	1.55	1.51
6	D	1402	PIO	O2C-C2C	-2.52	1.40	1.46
6	B	1403	PIO	O3C-C1B	2.45	1.40	1.33
6	A	1402	PIO	O3C-C1B	2.38	1.40	1.33
6	A	1402	PIO	O2C-C1A	2.34	1.40	1.34
6	C	1404	PIO	O3C-C1B	2.33	1.40	1.33
6	D	1402	PIO	O3C-C1B	2.31	1.40	1.33
6	C	1404	PIO	O2C-C1A	2.27	1.40	1.34
6	D	1402	PIO	O2C-C1A	2.26	1.40	1.34
6	A	1402	PIO	O2C-C2C	-2.25	1.41	1.46
6	C	1404	PIO	O3C-C3C	-2.24	1.40	1.45
6	D	1402	PIO	O3C-C3C	-2.23	1.40	1.45
6	B	1403	PIO	O2C-C1A	2.22	1.40	1.34
6	A	1402	PIO	O3C-C3C	-2.19	1.40	1.45
6	B	1403	PIO	O3C-C3C	-2.18	1.40	1.45

All (119) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	1401	9Z9	C76-C73-C74	27.62	165.69	115.66
7	C	1403	9Z9	C76-C73-C74	27.46	165.39	115.66
7	C	1401	9Z9	C76-C73-C74	27.40	165.30	115.66
7	A	1403	9Z9	C76-C73-C74	27.40	165.29	115.66
7	C	1401	9Z9	O72-C73-C76	-18.87	70.23	108.54
7	A	1403	9Z9	O72-C73-C76	-18.80	70.37	108.54
7	B	1401	9Z9	O72-C73-C76	-18.79	70.39	108.54
7	C	1403	9Z9	O72-C73-C76	-18.65	70.66	108.54
7	C	1401	9Z9	O80-C73-O72	11.97	138.35	109.88
7	B	1401	9Z9	O80-C73-O72	11.49	137.21	109.88
7	C	1403	9Z9	O80-C73-O72	11.48	137.19	109.88
7	A	1403	9Z9	O80-C73-O72	11.25	136.64	109.88
7	B	1401	9Z9	O80-C73-C74	-8.47	82.01	107.26
7	C	1403	9Z9	O80-C73-C74	-8.36	82.34	107.26
7	C	1401	9Z9	O80-C73-C74	-8.30	82.52	107.26
7	A	1403	9Z9	O80-C73-C74	-8.22	82.74	107.26
7	A	1403	9Z9	C10-C02-C06	-6.06	98.18	107.25
7	B	1401	9Z9	C10-C02-C06	-6.01	98.25	107.25
7	A	1403	9Z9	C02-C06-C07	-5.69	106.33	114.41
7	C	1403	9Z9	C02-C06-C07	-5.49	106.62	114.41
7	C	1403	9Z9	C10-C02-C06	-5.46	99.08	107.25
7	A	1403	9Z9	C07-C15-C14	-5.41	105.27	112.76
7	B	1401	9Z9	C02-C06-C07	-5.33	106.85	114.41
7	C	1401	9Z9	C02-C06-C07	-5.20	107.03	114.41
7	C	1401	9Z9	C10-C02-C06	-4.98	99.80	107.25
7	C	1403	9Z9	C09-C08-C11	-4.73	107.25	113.08
7	C	1401	9Z9	C07-C15-C14	-4.73	106.21	112.76
7	C	1401	9Z9	C06-C05-C04	-4.62	94.35	102.40
6	C	1404	PIO	O2C-C1A-C2A	4.55	121.32	111.48
7	B	1401	9Z9	C06-C05-C04	-4.38	94.76	102.40
7	C	1403	9Z9	C06-C05-C04	-4.31	94.89	102.40
7	A	1403	9Z9	C06-C05-C04	-4.27	94.96	102.40
6	A	1402	PIO	O2C-C1A-C2A	4.24	120.66	111.48
7	B	1401	9Z9	C07-C15-C14	-4.23	106.91	112.76
6	D	1402	PIO	O2C-C1A-C2A	4.14	120.43	111.48
5	B	1402	3PH	O21-C21-C22	4.08	120.30	111.48
6	B	1403	PIO	O2C-C1A-C2A	4.07	120.29	111.48
7	B	1401	9Z9	C09-C08-C11	-4.07	108.07	113.08
5	A	1401	3PH	O21-C21-C22	4.02	120.17	111.48
7	B	1401	9Z9	C09-C10-C02	-4.00	106.00	112.74
5	D	1401	3PH	O21-C21-C22	3.99	120.12	111.48
5	C	1402	3PH	O21-C21-C22	3.91	119.94	111.48
7	C	1403	9Z9	C09-C10-C02	-3.91	106.15	112.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1401	9Z9	C09-C08-C11	-3.90	108.27	113.08
7	C	1401	9Z9	C09-C10-C02	-3.83	106.28	112.74
7	C	1403	9Z9	C07-C15-C14	-3.66	107.70	112.76
7	A	1403	9Z9	C15-C14-C13	-3.61	118.92	125.02
7	A	1403	9Z9	O80-C79-C78	-3.61	107.53	112.17
7	B	1401	9Z9	C73-C74-C03	3.60	108.67	103.37
7	C	1401	9Z9	C01-C02-C03	3.60	119.40	111.58
7	A	1403	9Z9	C73-C74-C03	3.56	108.62	103.37
7	C	1403	9Z9	C73-C74-C03	3.55	108.59	103.37
7	C	1403	9Z9	O80-C79-C78	-3.52	107.65	112.17
7	A	1403	9Z9	C09-C10-C02	-3.51	106.82	112.74
7	C	1403	9Z9	C01-C02-C03	3.47	119.14	111.58
7	A	1403	9Z9	C01-C02-C03	3.45	119.10	111.58
7	C	1401	9Z9	C73-C74-C03	3.43	108.43	103.37
7	B	1401	9Z9	C01-C02-C03	3.39	118.96	111.58
6	D	1402	PIO	C6-C1-C2	3.37	115.54	110.86
7	A	1403	9Z9	C09-C08-C11	-3.37	108.93	113.08
7	C	1401	9Z9	C74-C03-C04	3.33	109.64	104.27
7	B	1401	9Z9	O80-C79-C78	-3.17	108.10	112.17
7	C	1401	9Z9	O80-C79-C78	-3.14	108.14	112.17
7	C	1401	9Z9	C15-C14-C13	-3.07	119.84	125.02
7	A	1403	9Z9	C10-C09-C08	3.01	118.25	113.14
6	B	1403	PIO	O3C-C1B-C2B	2.97	120.88	111.83
7	C	1401	9Z9	C75-C74-C03	-2.95	108.59	114.50
6	D	1402	PIO	C5-C6-C1	2.92	115.03	109.11
7	B	1401	9Z9	C16-C13-C11	2.92	120.16	116.42
7	C	1403	9Z9	C74-C03-C04	2.91	108.96	104.27
5	A	1401	3PH	O31-C31-C32	2.89	120.66	111.83
7	C	1403	9Z9	C16-C13-C11	2.86	120.08	116.42
7	B	1401	9Z9	C15-C14-C13	-2.86	120.20	125.02
5	C	1402	3PH	O31-C31-C32	2.85	120.54	111.83
7	B	1401	9Z9	C74-C03-C04	2.85	108.86	104.27
5	B	1402	3PH	O31-C31-C32	2.85	120.52	111.83
7	C	1403	9Z9	C11-C08-C07	-2.84	108.57	112.71
6	D	1402	PIO	O3C-C1B-C2B	2.81	120.41	111.83
7	C	1403	9Z9	C15-C14-C13	-2.80	120.30	125.02
5	D	1401	3PH	O31-C31-C32	2.76	120.24	111.83
7	A	1403	9Z9	C11-C08-C07	-2.75	108.69	112.71
7	A	1403	9Z9	O72-C73-C74	-2.74	100.92	104.56
7	B	1401	9Z9	C11-C08-C07	-2.72	108.74	112.71
7	C	1401	9Z9	C16-C13-C11	2.71	119.89	116.42
7	A	1403	9Z9	C74-C03-C04	2.69	108.60	104.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1403	9Z9	C10-C09-C08	2.66	117.65	113.14
7	B	1401	9Z9	C10-C09-C08	2.64	117.63	113.14
7	C	1401	9Z9	C11-C08-C07	-2.61	108.90	112.71
7	C	1401	9Z9	C10-C09-C08	2.60	117.55	113.14
6	C	1404	PIO	O3C-C1B-C2B	2.59	119.74	111.83
7	A	1403	9Z9	C73-O72-C04	-2.59	100.98	108.14
6	A	1402	PIO	O3C-C1B-C2B	2.58	119.72	111.83
7	C	1401	9Z9	O72-C04-C03	2.56	108.60	105.12
7	B	1401	9Z9	O72-C73-C74	-2.52	101.21	104.56
7	B	1401	9Z9	O72-C04-C03	2.47	108.47	105.12
7	C	1403	9Z9	O72-C73-C74	-2.41	101.35	104.56
7	C	1403	9Z9	C73-O72-C04	-2.39	101.53	108.14
7	A	1403	9Z9	O72-C04-C03	2.38	108.35	105.12
7	A	1403	9Z9	C10-C02-C03	2.37	118.74	115.36
7	B	1401	9Z9	C73-O72-C04	-2.37	101.59	108.14
7	C	1403	9Z9	O72-C04-C03	2.37	108.33	105.12
7	A	1403	9Z9	C17-C16-C13	-2.32	108.03	111.45
7	C	1403	9Z9	C11-C13-C14	-2.29	119.59	122.93
7	C	1401	9Z9	C02-C03-C04	2.27	106.65	104.20
6	B	1403	PIO	C5-C6-C1	2.27	113.71	109.11
7	C	1403	9Z9	C01-C02-C06	-2.27	107.56	111.68
7	C	1401	9Z9	C10-C02-C03	2.26	118.57	115.36
7	A	1403	9Z9	C16-C13-C11	2.23	119.28	116.42
7	A	1403	9Z9	C75-C74-C03	-2.21	110.07	114.50
7	C	1401	9Z9	C15-C07-C06	-2.16	107.88	110.93
6	D	1402	PIO	C3-C2-C1	2.11	114.46	109.68
7	B	1401	9Z9	C79-O80-C73	2.10	117.31	113.69
7	B	1401	9Z9	C01-C02-C06	-2.05	107.97	111.68
6	A	1402	PIO	P4-O4-C4	-2.04	117.98	123.43
7	C	1403	9Z9	C79-O80-C73	2.04	117.21	113.69
7	A	1403	9Z9	C01-C02-C06	-2.03	107.99	111.68
7	C	1401	9Z9	C79-O80-C73	2.03	117.18	113.69
7	A	1403	9Z9	C79-O80-C73	2.02	117.17	113.69
6	C	1404	PIO	P5-O5-C5	-2.00	118.08	123.43

There are no chirality outliers.

All (187) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1401	3PH	C22-C21-O21-C2
5	B	1402	3PH	C1-O11-P-O13
5	B	1402	3PH	C1-O11-P-O14

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Mol	Chain	Res	Type	Atoms
5	B	1402	3PH	C1-O11-P-O12
5	C	1402	3PH	C1-O11-P-O13
5	C	1402	3PH	C1-O11-P-O14
5	C	1402	3PH	C1-O11-P-O12
5	C	1402	3PH	C22-C21-O21-C2
5	D	1401	3PH	O22-C21-O21-C2
5	D	1401	3PH	C22-C21-O21-C2
6	A	1402	PIO	C1C-O13-P1-O12
6	A	1402	PIO	C2A-C1A-O2C-C2C
6	B	1403	PIO	C3-C4-O4-P4
6	B	1403	PIO	C2A-C1A-O2C-C2C
6	C	1404	PIO	C1C-O13-P1-O1
6	C	1404	PIO	C1C-O13-P1-O11
6	C	1404	PIO	C1C-O13-P1-O12
6	C	1404	PIO	O1A-C1A-O2C-C2C
6	C	1404	PIO	C2A-C1A-O2C-C2C
6	D	1402	PIO	C1-O1-P1-O11
6	D	1402	PIO	C1-O1-P1-O13
6	D	1402	PIO	C5-C4-O4-P4
6	D	1402	PIO	O1A-C1A-O2C-C2C
6	D	1402	PIO	C2A-C1A-O2C-C2C
5	A	1401	3PH	O32-C31-O31-C3
6	B	1403	PIO	O1B-C1B-O3C-C3C
5	C	1402	3PH	O22-C21-O21-C2
6	A	1402	PIO	O1A-C1A-O2C-C2C
6	B	1403	PIO	O1A-C1A-O2C-C2C
5	A	1401	3PH	C32-C31-O31-C3
6	B	1403	PIO	C2B-C1B-O3C-C3C
5	A	1401	3PH	O22-C21-O21-C2
6	D	1402	PIO	C2B-C1B-O3C-C3C
6	D	1402	PIO	O1B-C1B-O3C-C3C
5	D	1401	3PH	C32-C31-O31-C3
5	D	1401	3PH	C31-C32-C33-C34
6	C	1404	PIO	C1A-C2A-C3A-C4A
5	A	1401	3PH	C31-C32-C33-C34
6	C	1404	PIO	C1B-C2B-C3B-C4B
5	B	1402	3PH	C32-C31-O31-C3
5	C	1402	3PH	C32-C31-O31-C3
5	D	1401	3PH	O32-C31-O31-C3
5	C	1402	3PH	O32-C31-O31-C3
6	A	1402	PIO	C1C-C2C-O2C-C1A
7	C	1401	9Z9	C23-C48-O49-C50

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Mol	Chain	Res	Type	Atoms
5	B	1402	3PH	O32-C31-O31-C3
5	C	1402	3PH	C2A-C2B-C2C-C2D
5	C	1402	3PH	C24-C25-C26-C27
5	C	1402	3PH	C37-C38-C39-C3A
5	A	1401	3PH	C24-C25-C26-C27
6	D	1402	PIO	C3A-C4A-C5A-C6A
5	B	1402	3PH	C3A-C3B-C3C-C3D
5	B	1402	3PH	C2A-C2B-C2C-C2D
5	A	1401	3PH	C29-C2A-C2B-C2C
5	D	1401	3PH	C3E-C3F-C3G-C3H
5	A	1401	3PH	C21-C22-C23-C24
5	D	1401	3PH	C21-C22-C23-C24
5	B	1402	3PH	C23-C24-C25-C26
5	B	1402	3PH	C33-C34-C35-C36
5	C	1402	3PH	C21-C22-C23-C24
6	A	1402	PIO	C1B-C2B-C3B-C4B
5	A	1401	3PH	C3C-C3D-C3E-C3F
5	A	1401	3PH	C28-C29-C2A-C2B
5	B	1402	3PH	C32-C33-C34-C35
5	A	1401	3PH	C36-C37-C38-C39
5	D	1401	3PH	C32-C33-C34-C35
5	B	1402	3PH	C26-C27-C28-C29
5	D	1401	3PH	C35-C36-C37-C38
5	B	1402	3PH	C22-C21-O21-C2
5	C	1402	3PH	C32-C33-C34-C35
5	B	1402	3PH	C38-C39-C3A-C3B
5	A	1401	3PH	C35-C36-C37-C38
5	A	1401	3PH	C32-C33-C34-C35
5	A	1401	3PH	C3B-C3C-C3D-C3E
5	C	1402	3PH	C36-C37-C38-C39
5	B	1402	3PH	C29-C2A-C2B-C2C
5	D	1401	3PH	C25-C26-C27-C28
5	C	1402	3PH	C2B-C2C-C2D-C2E
5	A	1401	3PH	O11-C1-C2-C3
5	B	1402	3PH	O22-C21-O21-C2
5	C	1402	3PH	C38-C39-C3A-C3B
6	D	1402	PIO	C3B-C4B-C5B-C6B
6	A	1402	PIO	C1C-C2C-C3C-O3C
5	C	1402	3PH	C3C-C3D-C3E-C3F
5	A	1401	3PH	C2-C1-O11-P
5	B	1402	3PH	C3C-C3D-C3E-C3F
6	C	1404	PIO	C4A-C5A-C6A-C7A

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Mol	Chain	Res	Type	Atoms
5	A	1401	3PH	C26-C27-C28-C29
6	B	1403	PIO	C2A-C3A-C4A-C5A
5	B	1402	3PH	C35-C36-C37-C38
5	A	1401	3PH	O21-C2-C3-O31
5	B	1402	3PH	O21-C2-C3-O31
5	D	1401	3PH	C3D-C3E-C3F-C3G
5	C	1402	3PH	C33-C34-C35-C36
7	B	1401	9Z9	C23-C48-O49-C50
6	D	1402	PIO	C4A-C5A-C6A-C7A
5	D	1401	3PH	C2C-C2D-C2E-C2F
5	C	1402	3PH	C1-C2-C3-O31
5	D	1401	3PH	C1-C2-C3-O31
6	C	1404	PIO	C1C-C2C-C3C-O3C
5	A	1401	3PH	C37-C38-C39-C3A
5	C	1402	3PH	C3A-C3B-C3C-C3D
6	A	1402	PIO	C2C-C1C-O13-P1
5	C	1402	3PH	O21-C2-C3-O31
6	C	1404	PIO	O2C-C2C-C3C-O3C
5	C	1402	3PH	C35-C36-C37-C38
5	B	1402	3PH	C24-C25-C26-C27
5	B	1402	3PH	C28-C29-C2A-C2B
6	C	1404	PIO	C4B-C5B-C6B-C7B
5	C	1402	3PH	O11-C1-C2-C3
5	D	1401	3PH	O11-C1-C2-C3
5	B	1402	3PH	C34-C35-C36-C37
5	D	1401	3PH	C1-O11-P-O13
6	D	1402	PIO	C6-C1-O1-P1
7	A	1403	9Z9	C16-C17-O20-C21
7	A	1403	9Z9	C18-C17-O20-C21
5	A	1401	3PH	O11-C1-C2-O21
5	D	1401	3PH	O11-C1-C2-O21
6	B	1403	PIO	C1B-C2B-C3B-C4B
5	D	1401	3PH	O21-C2-C3-O31
5	A	1401	3PH	C38-C39-C3A-C3B
5	D	1401	3PH	C23-C24-C25-C26
7	A	1403	9Z9	C23-C24-O25-C26
7	A	1403	9Z9	C23-C48-O49-C50
5	B	1402	3PH	C2B-C2C-C2D-C2E
6	D	1402	PIO	C1-O1-P1-O12
6	D	1402	PIO	C1B-C2B-C3B-C4B
5	B	1402	3PH	O11-C1-C2-C3
6	A	1402	PIO	O13-C1C-C2C-C3C

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Mol	Chain	Res	Type	Atoms
6	D	1402	PIO	C5B-C6B-C7B-C8B
5	D	1401	3PH	C2-C1-O11-P
5	B	1402	3PH	O11-C1-C2-O21
5	C	1402	3PH	O11-C1-C2-O21
6	B	1403	PIO	O13-C1C-C2C-O2C
6	D	1402	PIO	O2C-C2C-C3C-O3C
5	A	1401	3PH	C1-C2-C3-O31
5	B	1402	3PH	C1-C2-C3-O31
6	D	1402	PIO	C1C-C2C-C3C-O3C
5	B	1402	3PH	C27-C28-C29-C2A
6	A	1402	PIO	C1C-O13-P1-O1
6	A	1402	PIO	C1C-O13-P1-O11
6	B	1403	PIO	C2C-C1C-O13-P1
6	C	1404	PIO	C2C-C1C-O13-P1
5	A	1401	3PH	C1-O11-P-O12
5	D	1401	3PH	C26-C27-C28-C29
5	B	1402	3PH	C25-C26-C27-C28
7	C	1401	9Z9	C23-C24-O25-C26
6	C	1404	PIO	C3A-C4A-C5A-C6A
6	D	1402	PIO	C5-O5-P5-O52
5	A	1401	3PH	C2A-C2B-C2C-C2D
5	A	1401	3PH	C2B-C2C-C2D-C2E
6	D	1402	PIO	C1C-C2C-O2C-C1A
6	D	1402	PIO	C3C-C2C-O2C-C1A
5	D	1401	3PH	C39-C3A-C3B-C3C
6	B	1403	PIO	O13-C1C-C2C-C3C
6	D	1402	PIO	O13-C1C-C2C-C3C
5	A	1401	3PH	C22-C23-C24-C25
6	A	1402	PIO	C5-O5-P5-O53
6	C	1404	PIO	C5-O5-P5-O53
6	D	1402	PIO	C4-O4-P4-O42
6	B	1403	PIO	C5B-C6B-C7B-C8B
6	C	1404	PIO	C2A-C3A-C4A-C5A
6	D	1402	PIO	C2-C1-O1-P1
6	C	1404	PIO	C5A-C6A-C7A-C8A
6	B	1403	PIO	C4B-C5B-C6B-C7B
6	C	1404	PIO	O3C-C1B-C2B-C3B
7	A	1403	9Z9	C22-C21-O20-C17
6	A	1402	PIO	C1-O1-P1-O11
6	C	1404	PIO	C1-O1-P1-O11
5	A	1401	3PH	C2-C3-O31-C31
6	D	1402	PIO	O2C-C1A-C2A-C3A

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Mol	Chain	Res	Type	Atoms
6	D	1402	PIO	C3-C4-O4-P4
6	C	1404	PIO	O2C-C1A-C2A-C3A
6	A	1402	PIO	C4-O4-P4-O41
6	A	1402	PIO	C5-O5-P5-O51
6	C	1404	PIO	C5-O5-P5-O52
5	C	1402	3PH	C26-C27-C28-C29
5	D	1401	3PH	C33-C34-C35-C36
6	C	1404	PIO	C2B-C1B-O3C-C3C
6	B	1403	PIO	C4-O4-P4-O42
6	B	1403	PIO	C5-O5-P5-O53
6	C	1404	PIO	O1B-C1B-O3C-C3C
6	C	1404	PIO	O1A-C1A-C2A-C3A
6	C	1404	PIO	O1B-C1B-C2B-C3B
6	D	1402	PIO	O3C-C1B-C2B-C3B
6	D	1402	PIO	O1A-C1A-C2A-C3A
5	B	1402	3PH	C3D-C3E-C3F-C3G

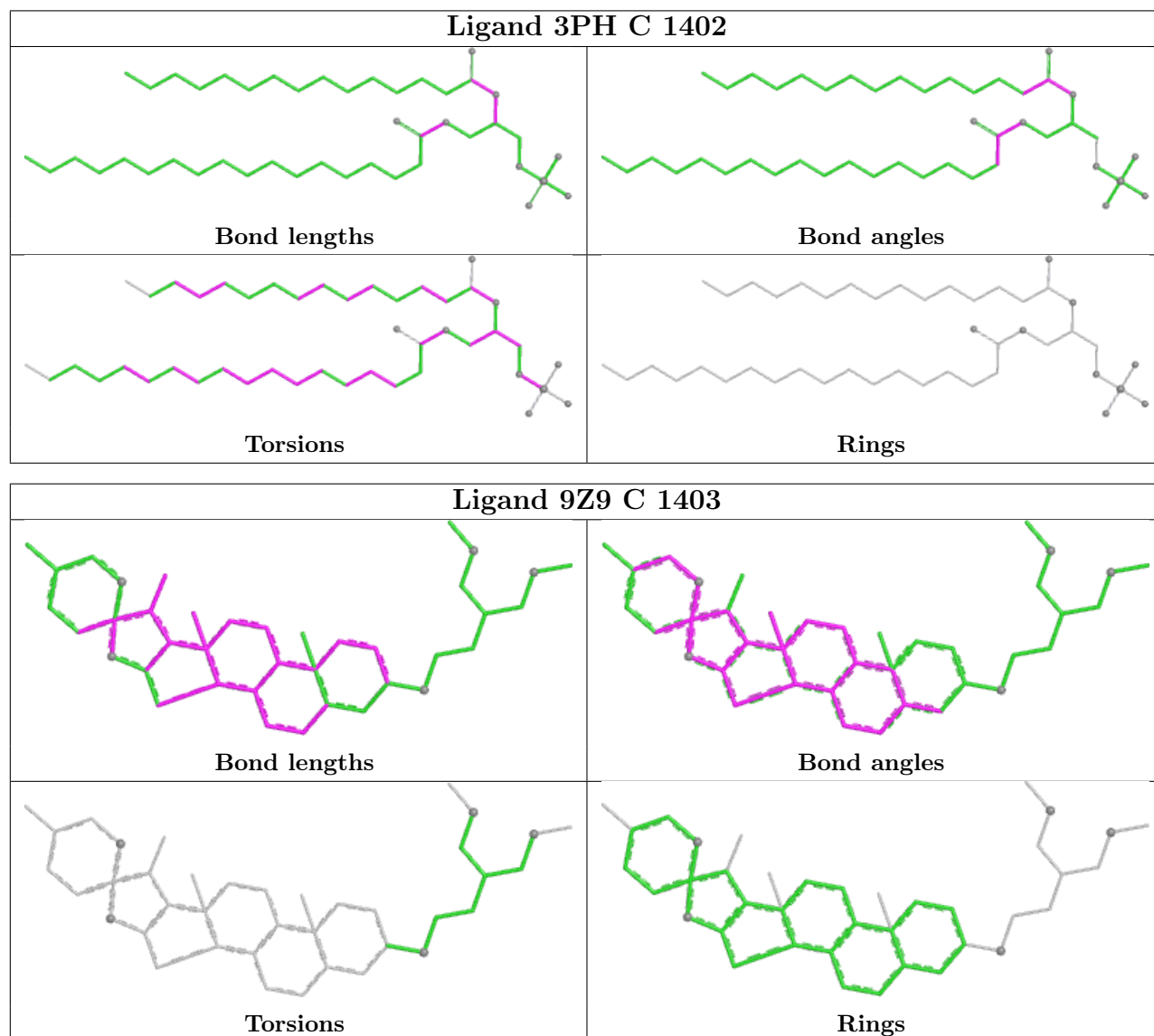
There are no ring outliers.

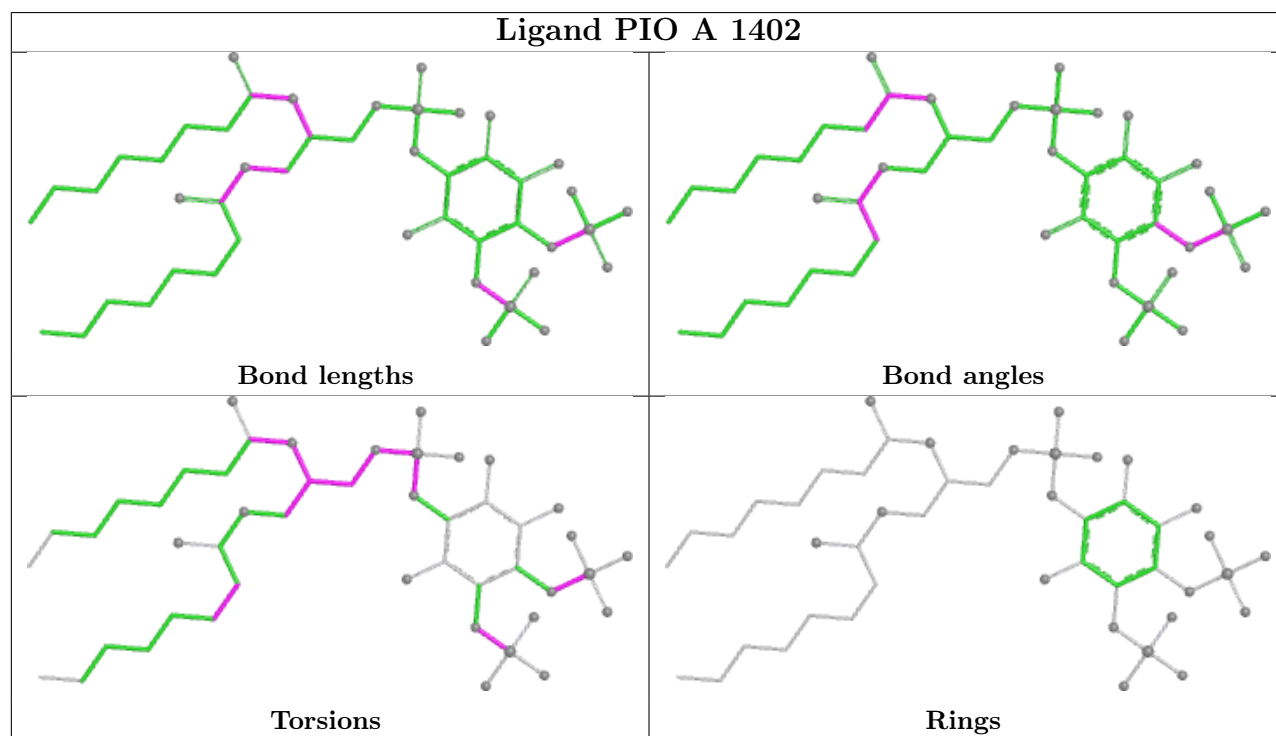
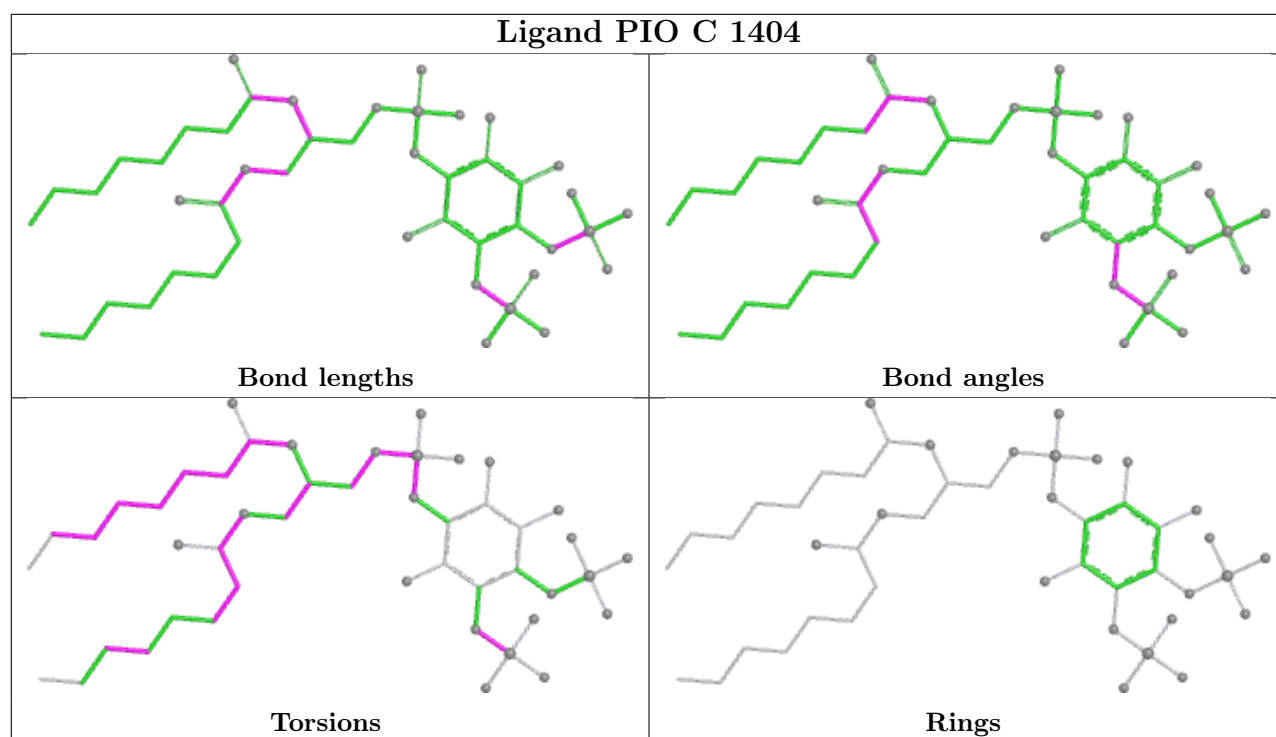
12 monomers are involved in 68 short contacts:

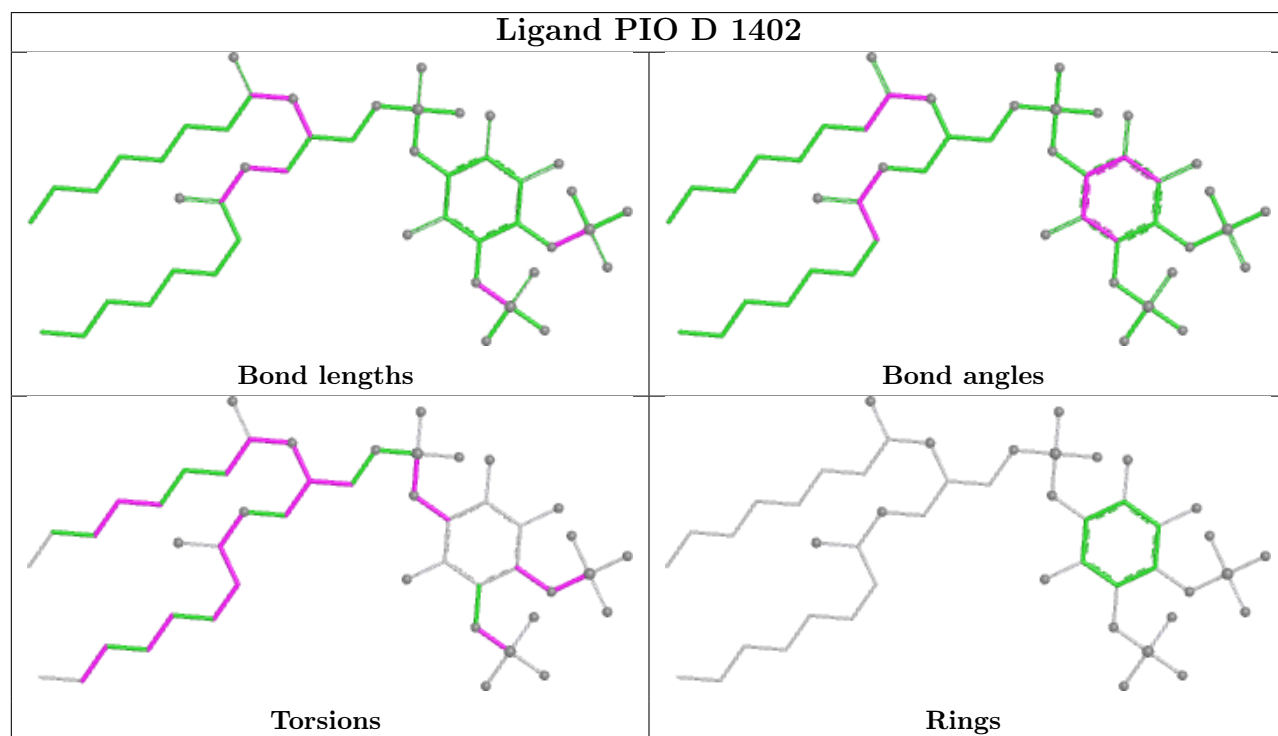
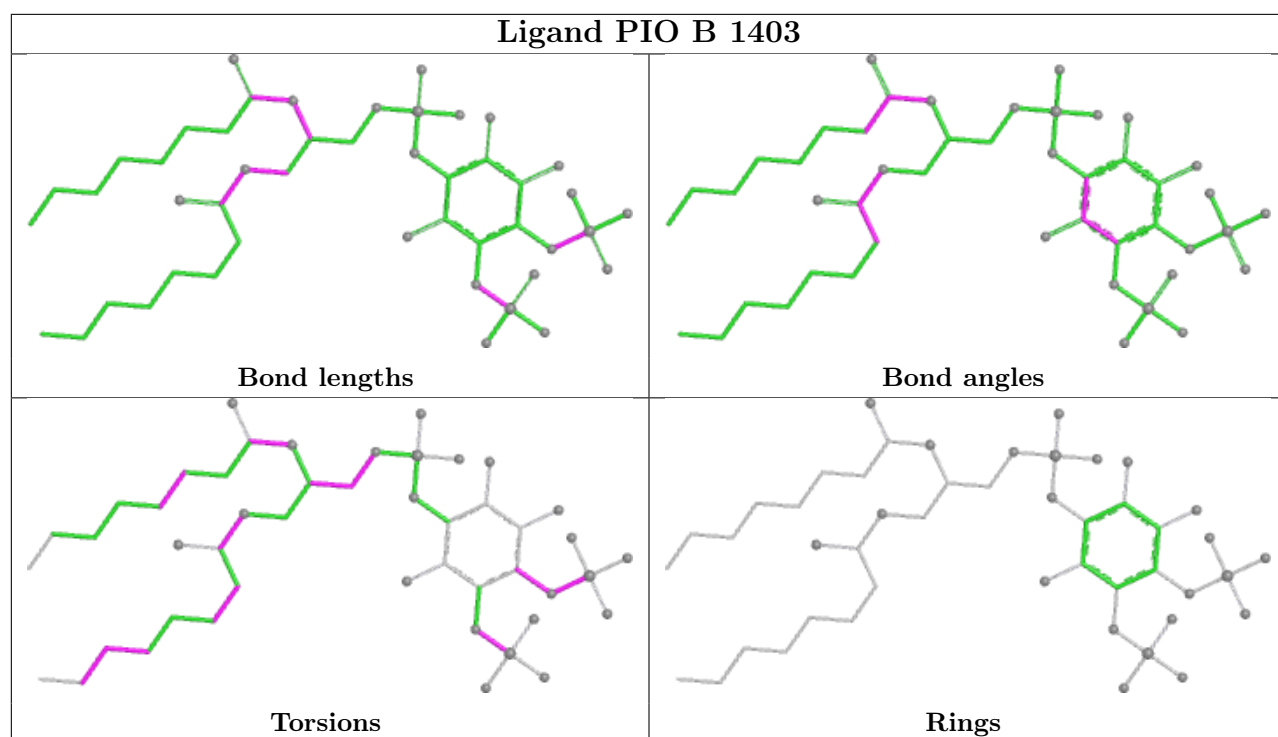
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	1402	3PH	1	0
7	C	1403	9Z9	14	0
6	C	1404	PIO	1	0
6	A	1402	PIO	1	0
6	B	1403	PIO	1	0
6	D	1402	PIO	1	0
7	C	1401	9Z9	15	0
7	B	1401	9Z9	15	0
5	B	1402	3PH	3	0
5	D	1401	3PH	1	0
5	A	1401	3PH	1	0
7	A	1403	9Z9	14	0

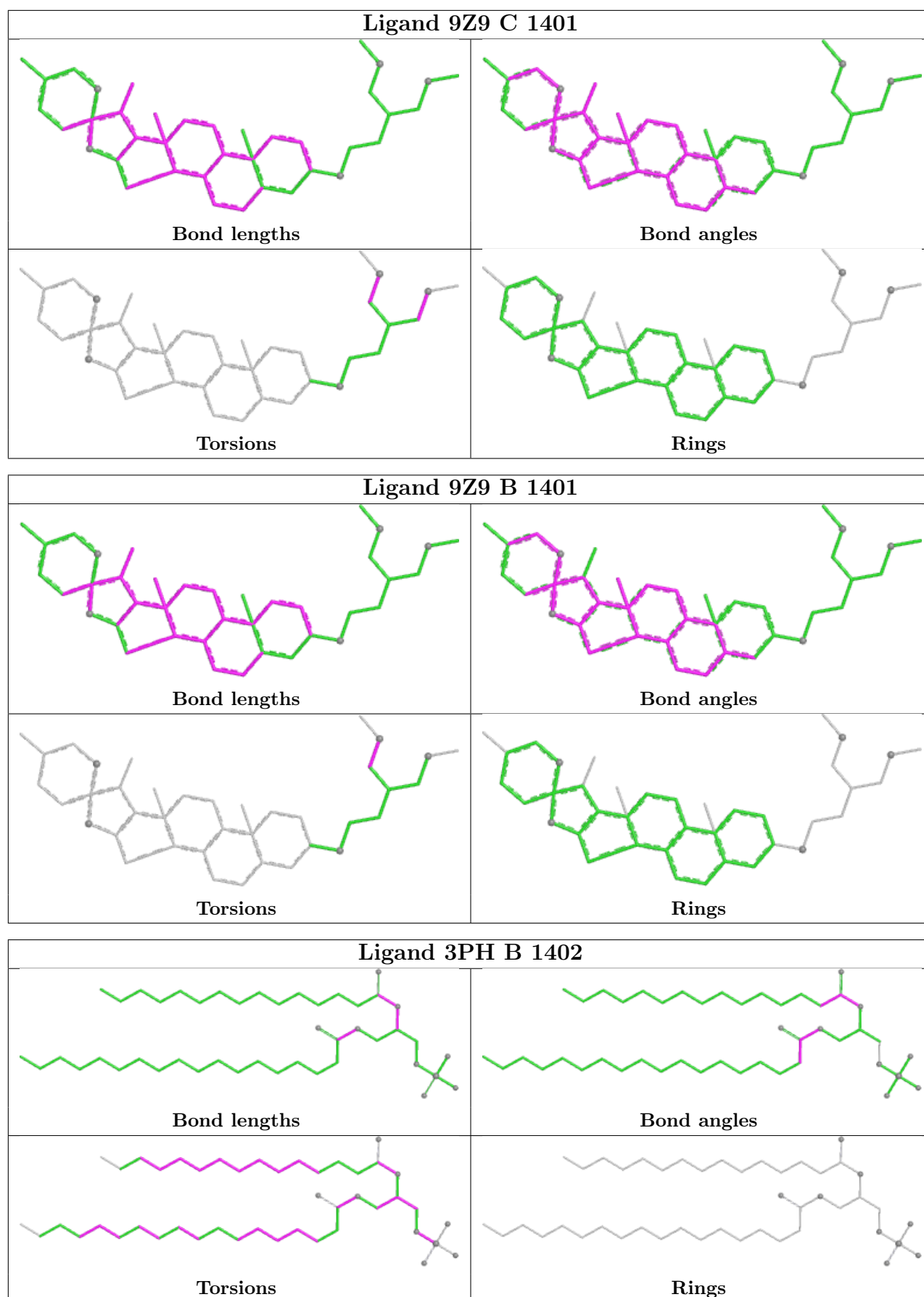
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

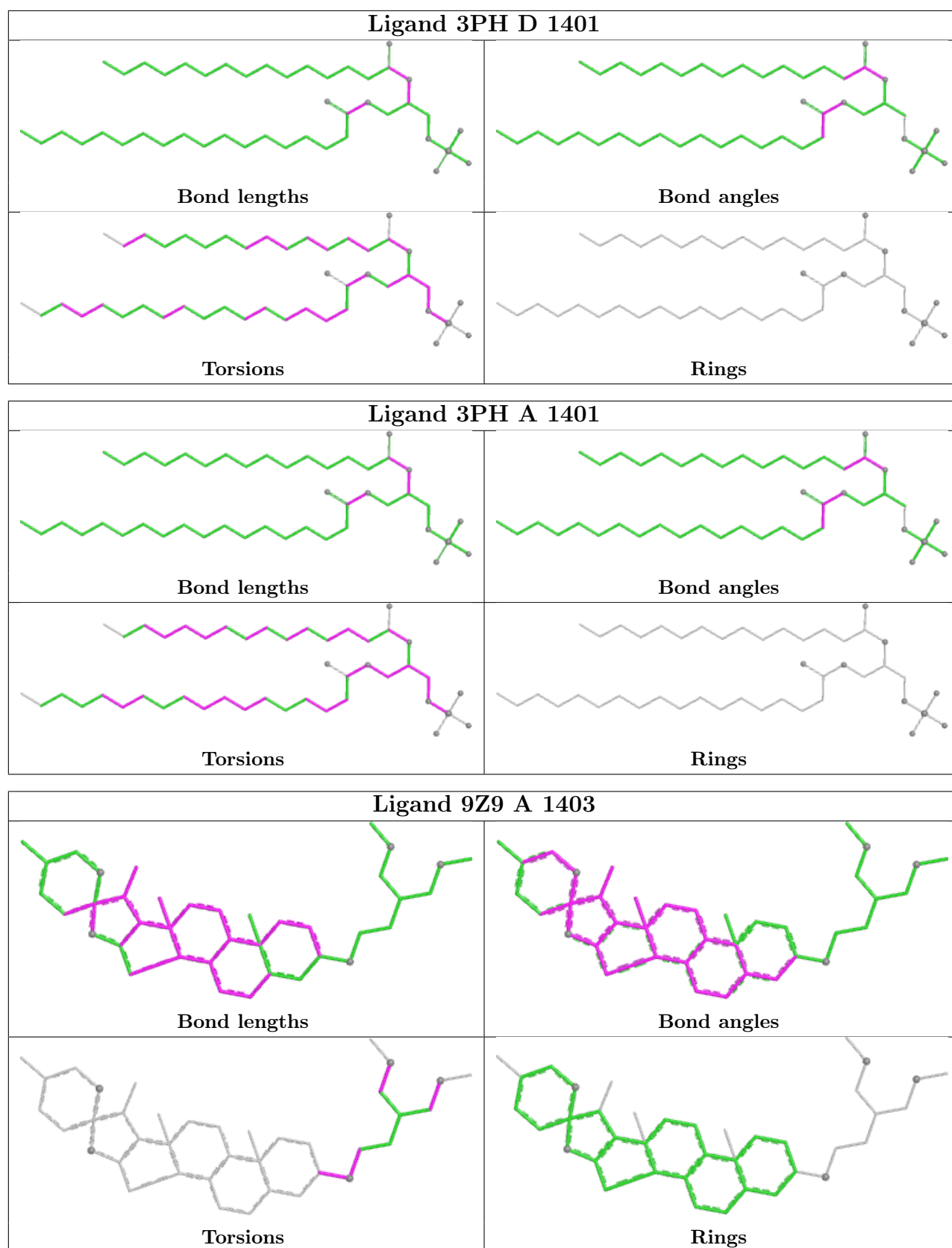
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

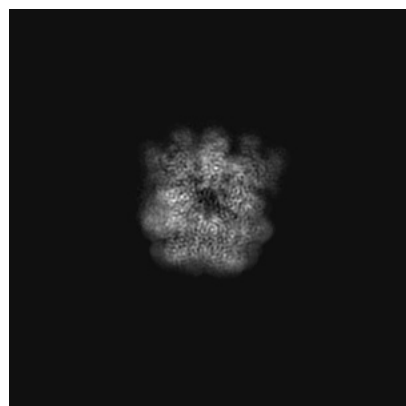
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-27345. These allow visual inspection of the internal detail of the map and identification of artifacts.

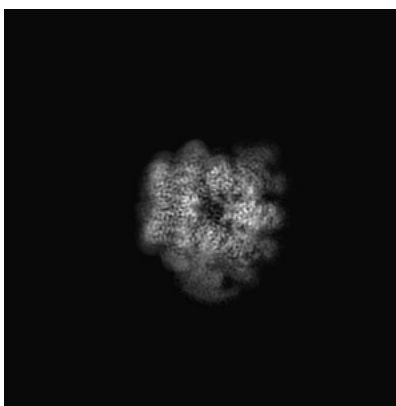
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

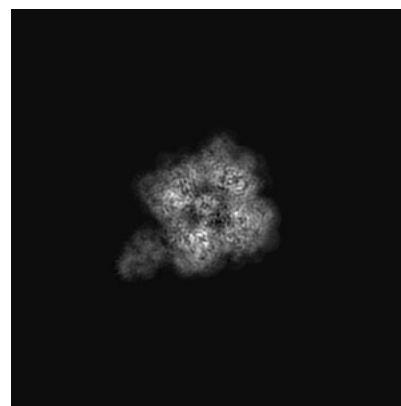
6.1.1 Primary map



X

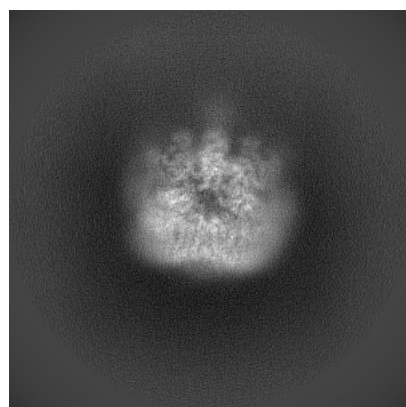


Y

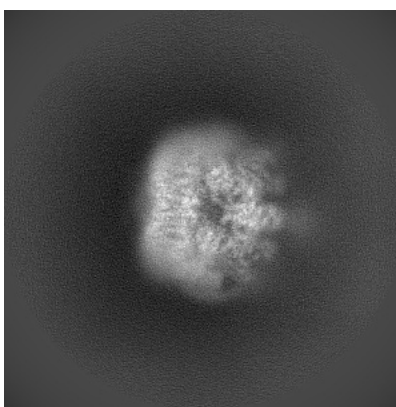


Z

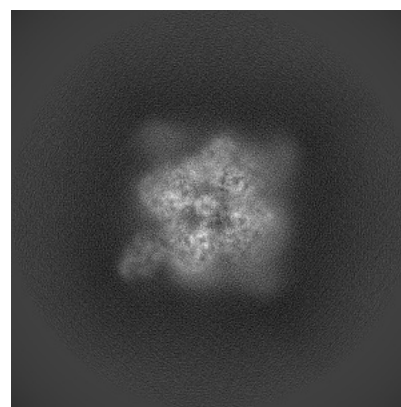
6.1.2 Raw map



X



Y

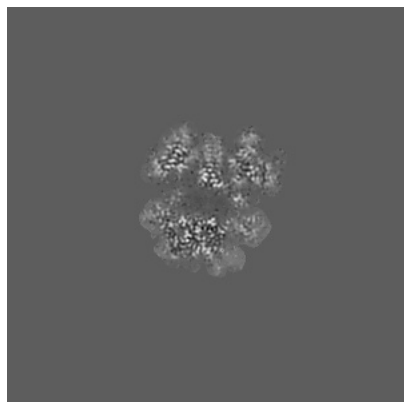


Z

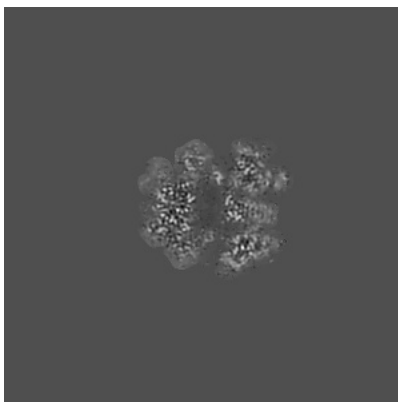
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

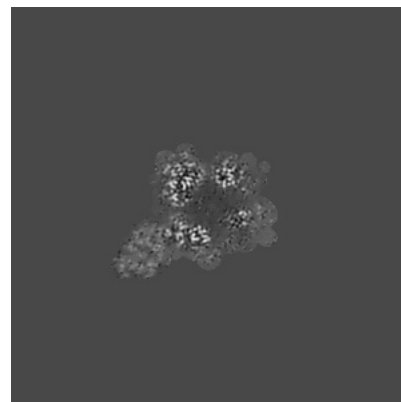
6.2.1 Primary map



X Index: 192

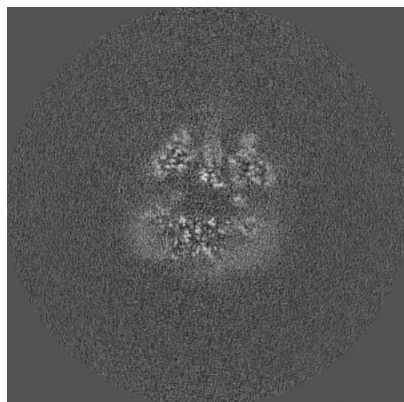


Y Index: 192

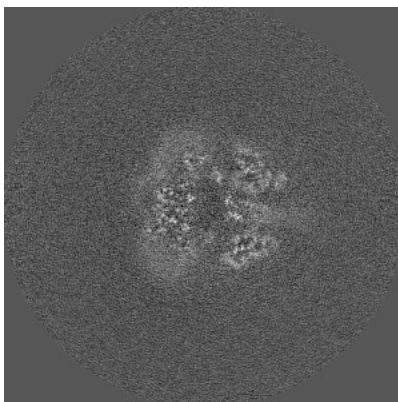


Z Index: 192

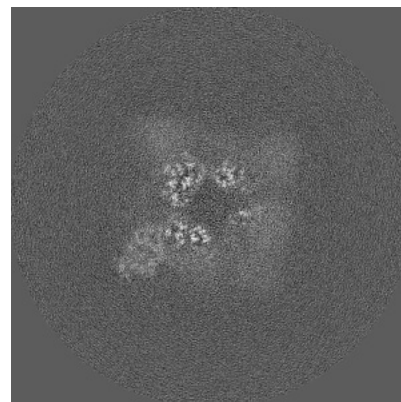
6.2.2 Raw map



X Index: 192



Y Index: 192

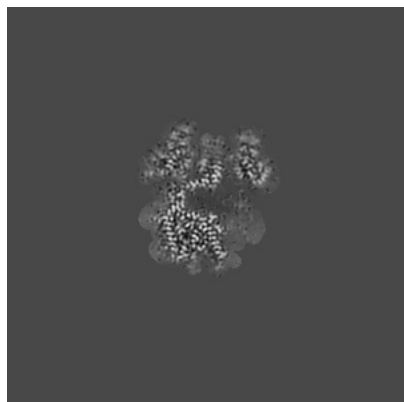


Z Index: 192

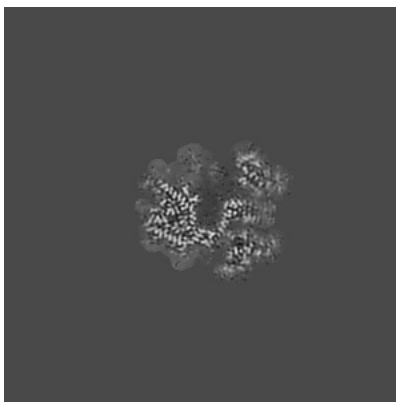
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

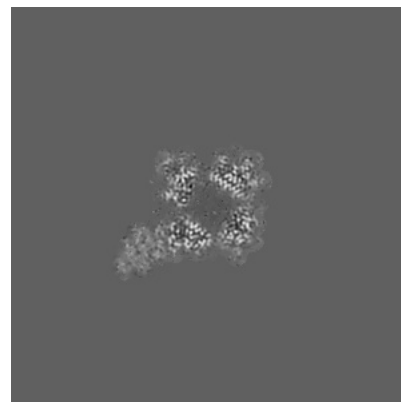
6.3.1 Primary map



X Index: 186

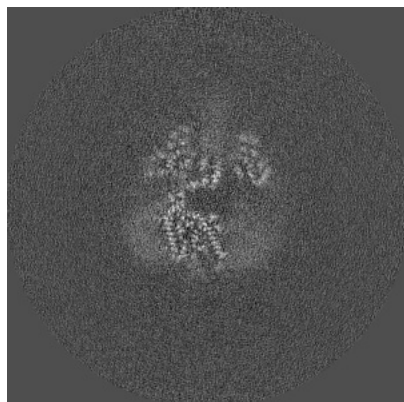


Y Index: 198

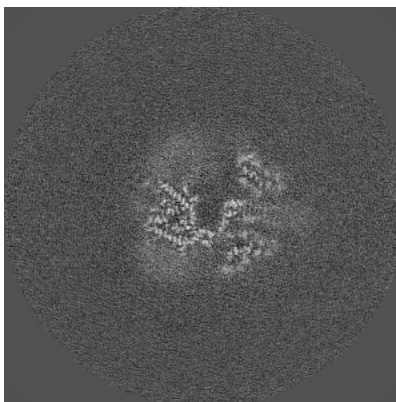


Z Index: 200

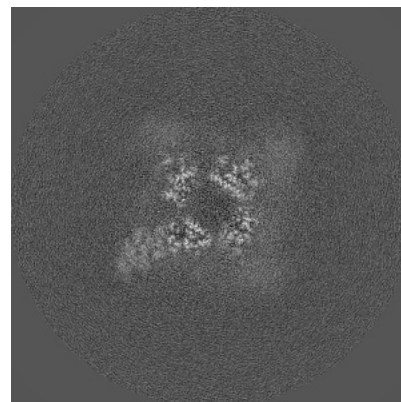
6.3.2 Raw map



X Index: 186



Y Index: 198

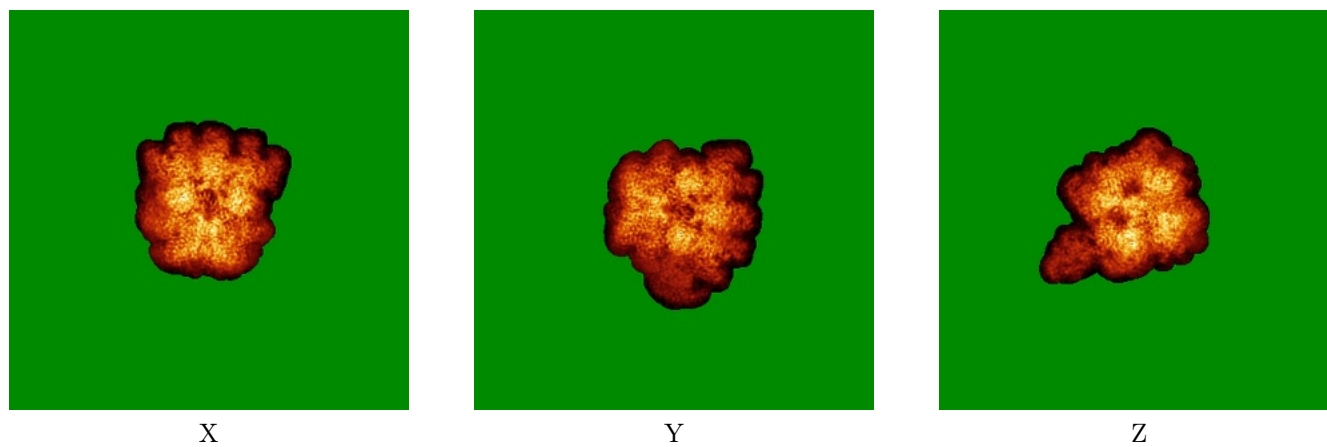


Z Index: 200

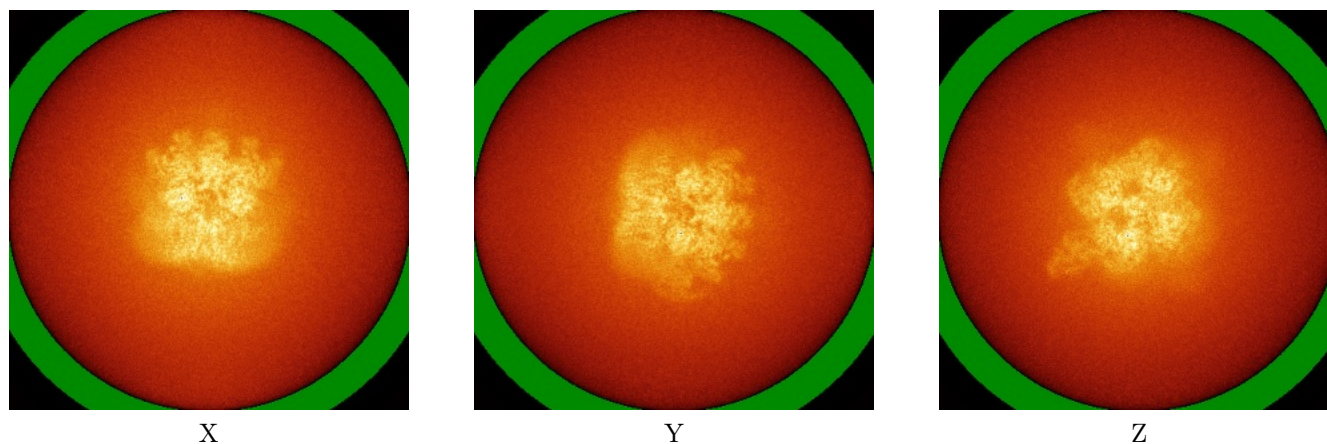
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

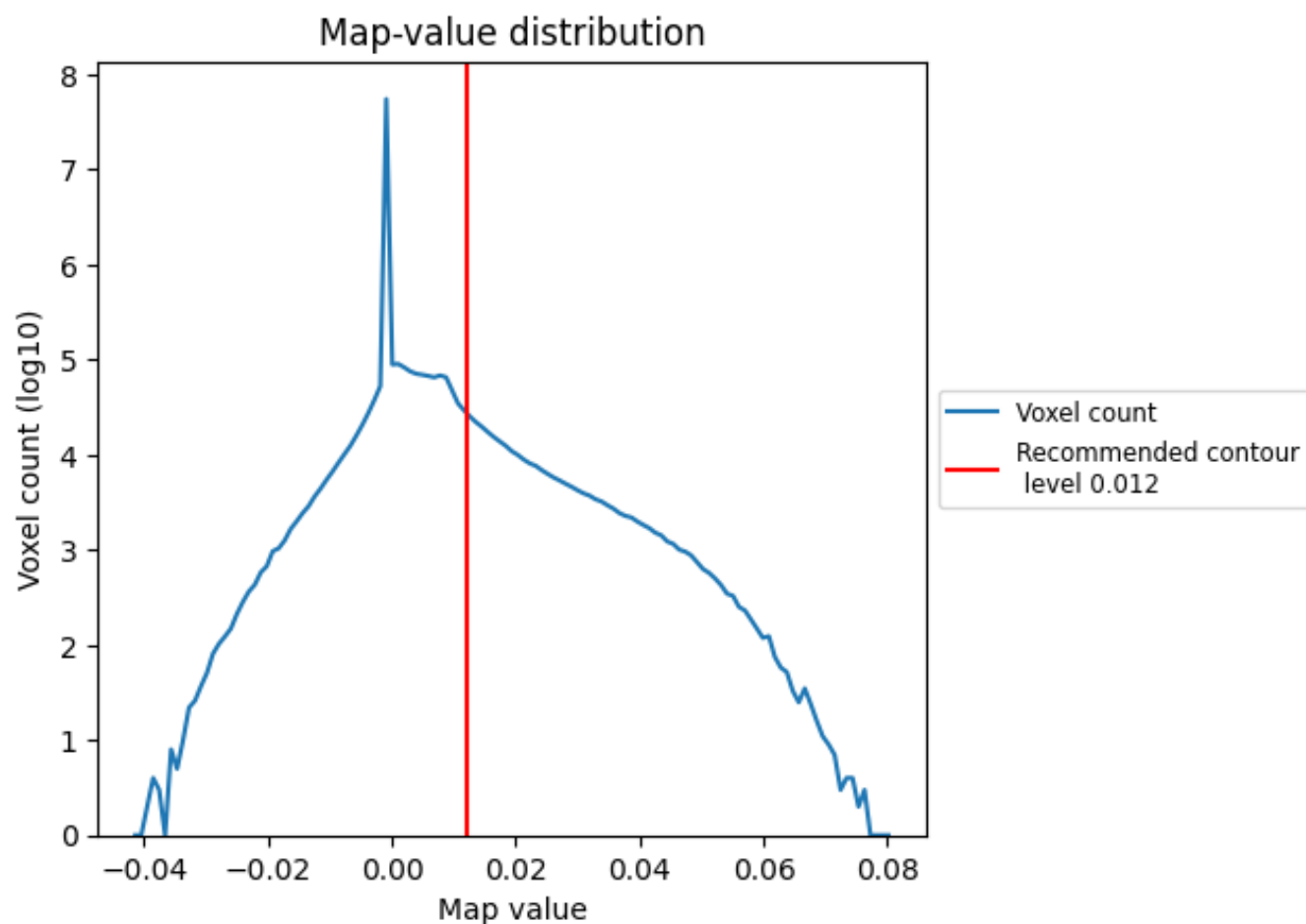
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

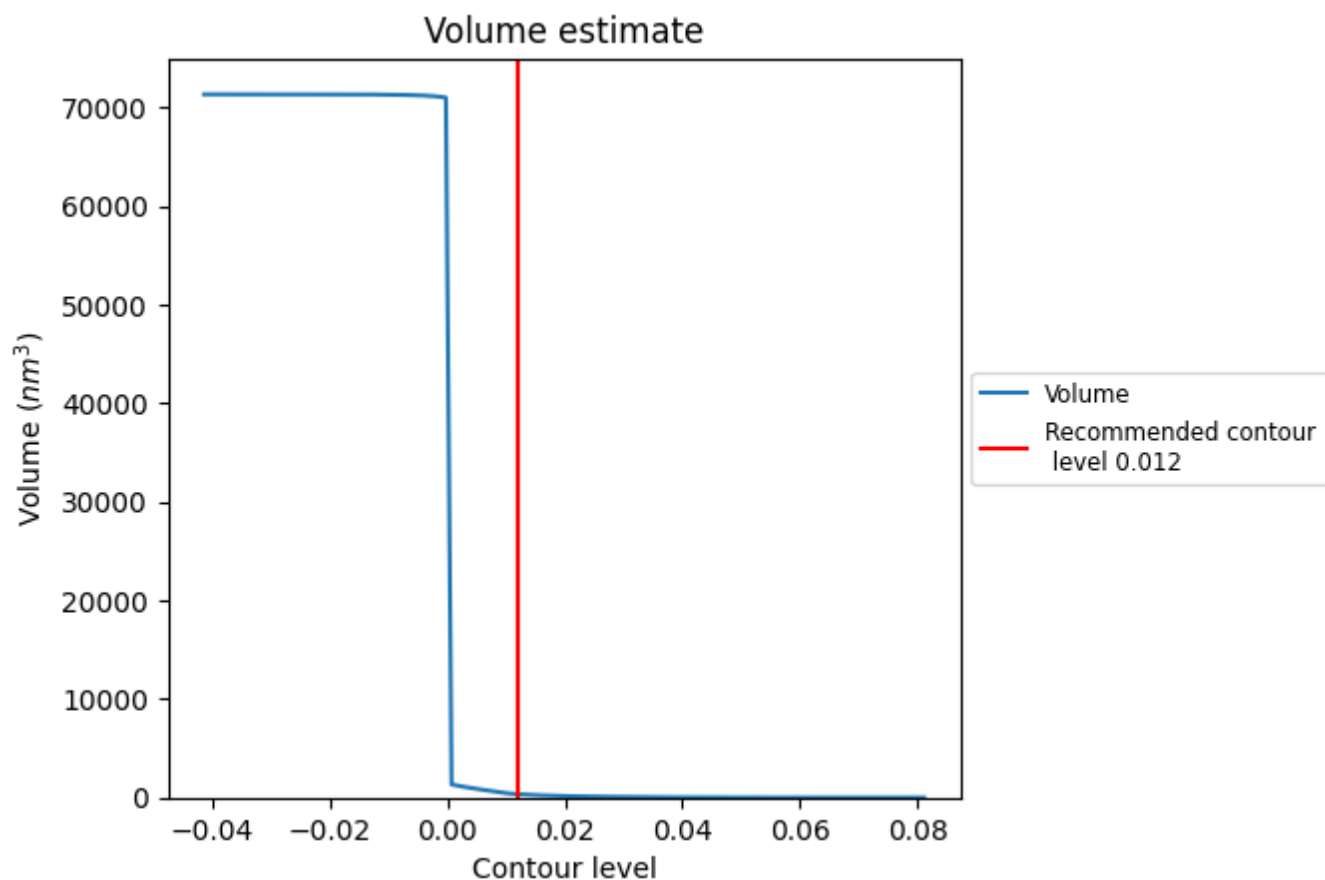
This section contains the results of statistical analysis of the map.

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.

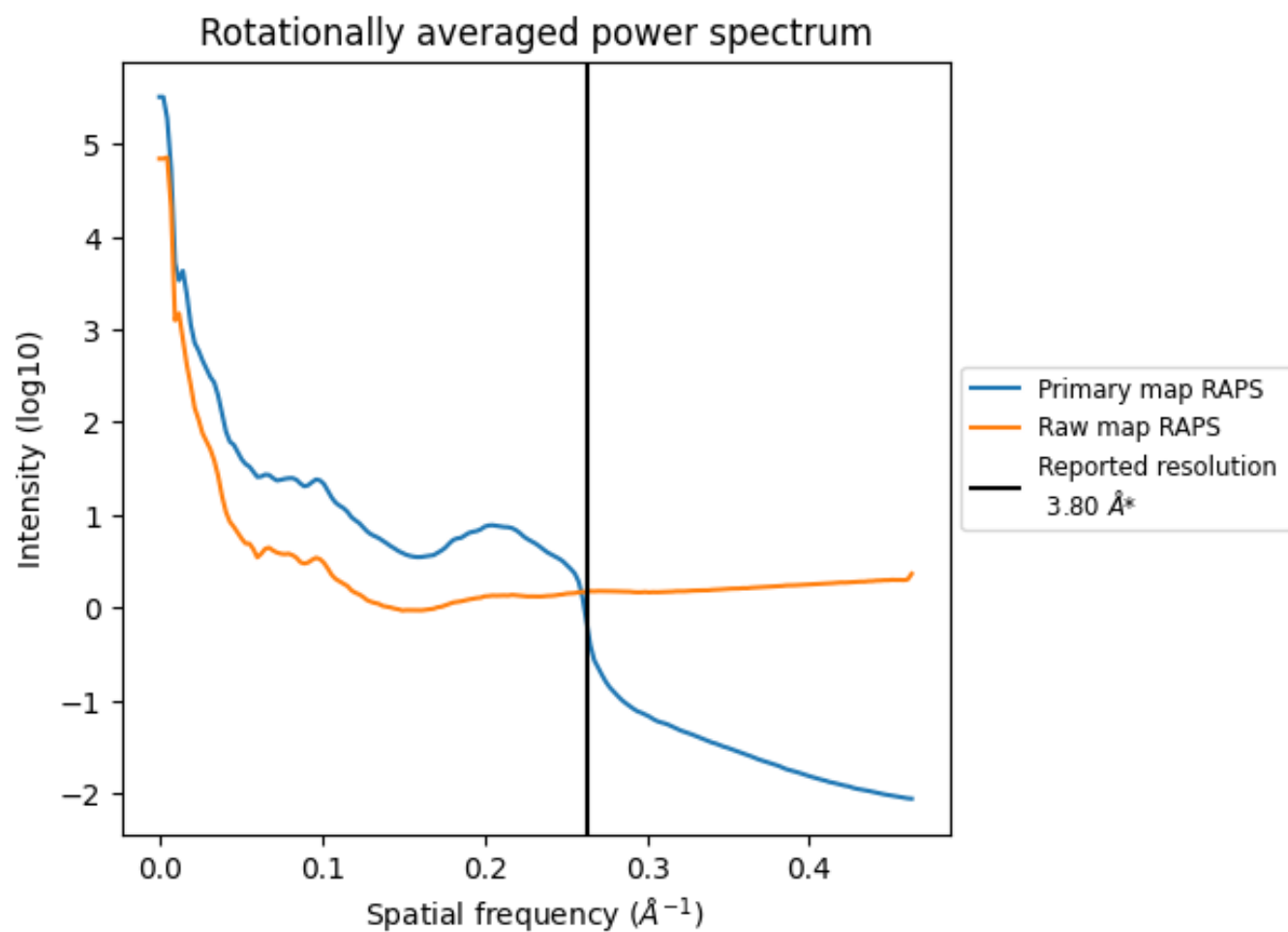
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 345 nm^3 ; this corresponds to an approximate mass of 312 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 ⓘ

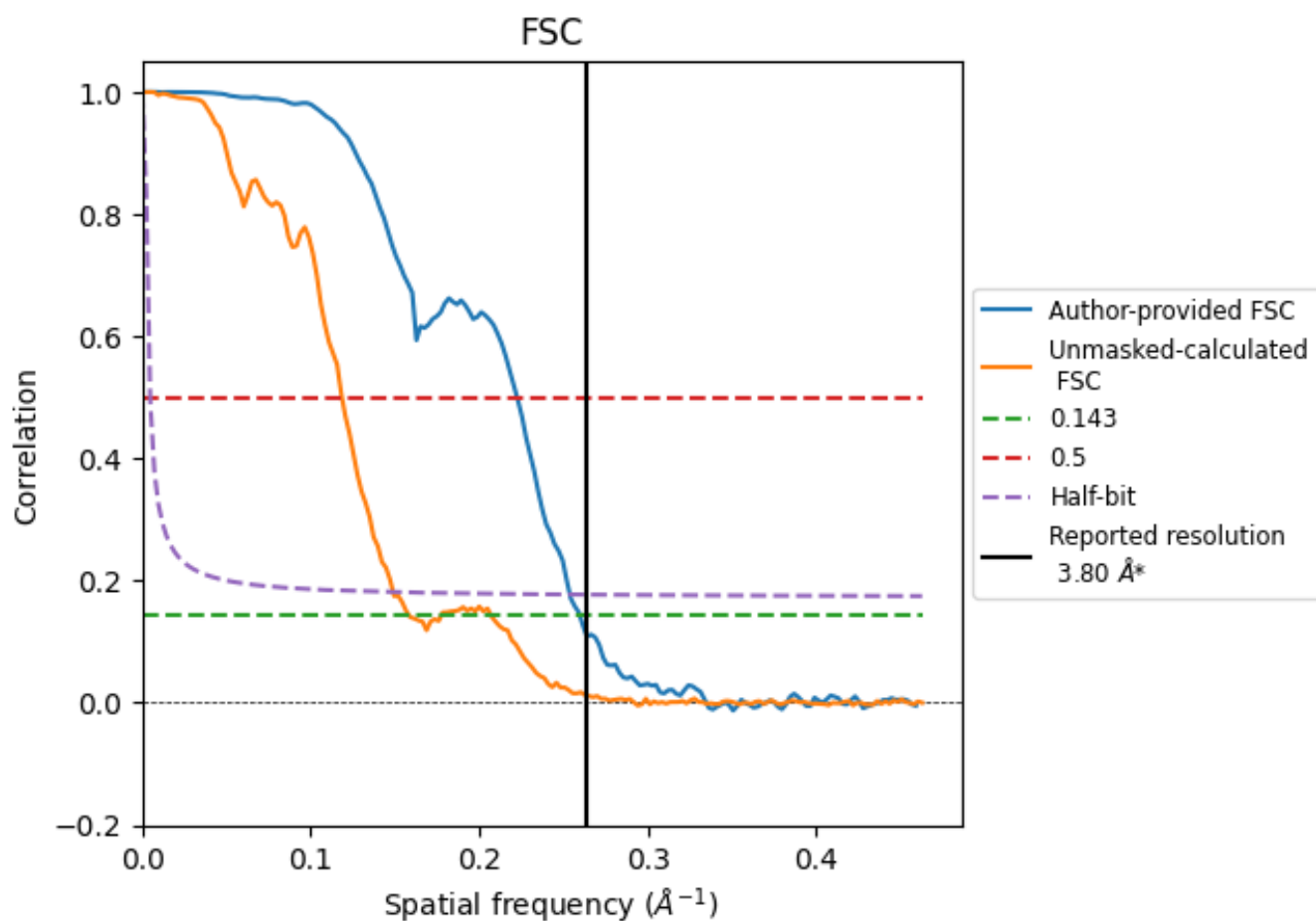


*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

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 0.263 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

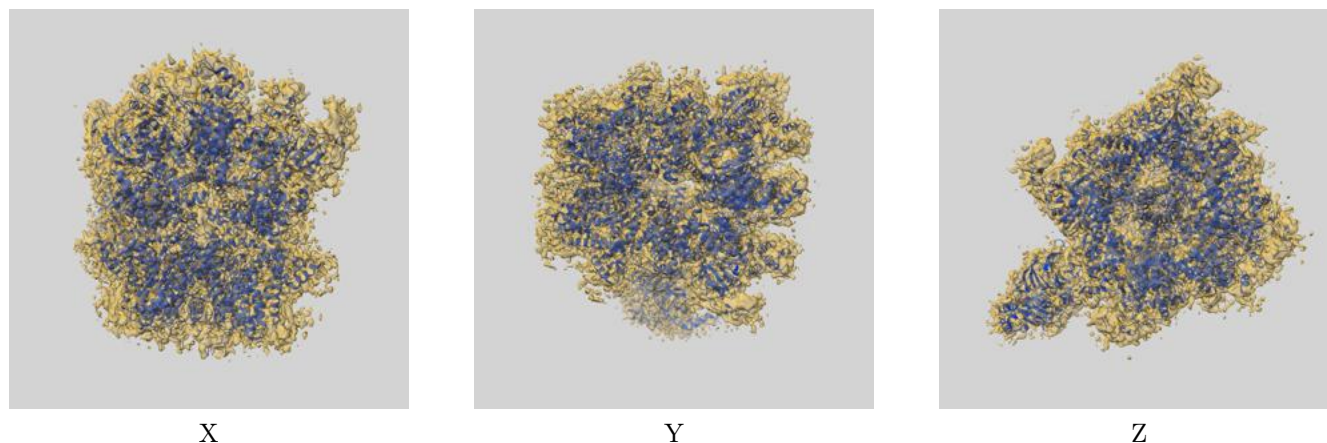
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.85	4.49	3.94
Unmasked-calculated*	6.31	8.44	6.72

*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 6.31 differs from the reported value 3.8 by more than 10 %

9 Map-model fit [i](#)

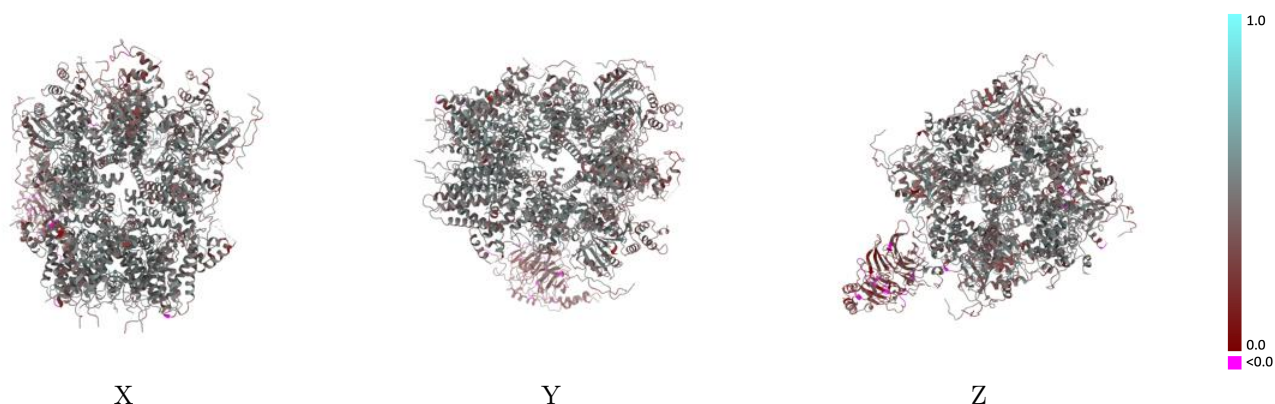
This section contains information regarding the fit between EMDB map EMD-27345 and PDB model 8DDX. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

9.1 Map-model overlay [i](#)



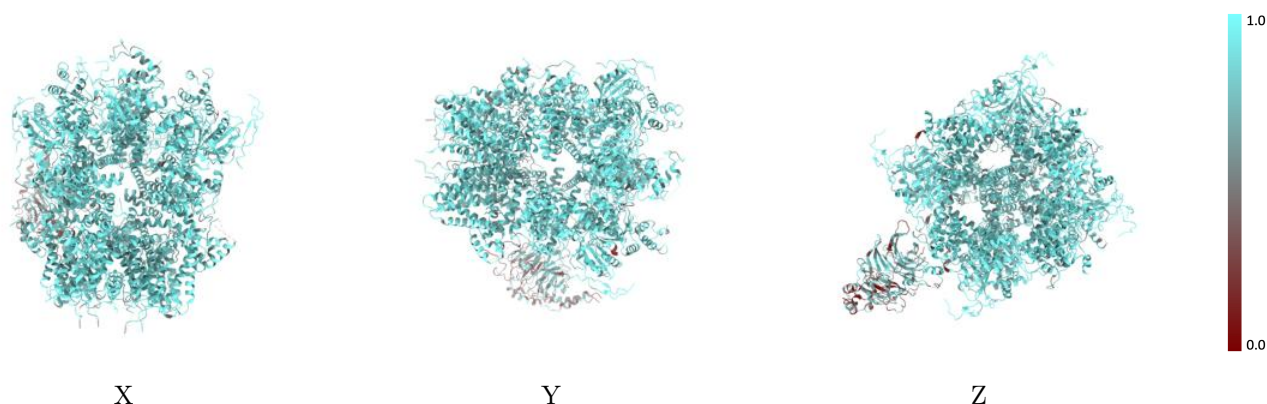
The images above show the 3D surface view of the map at the recommended contour level 0.012 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



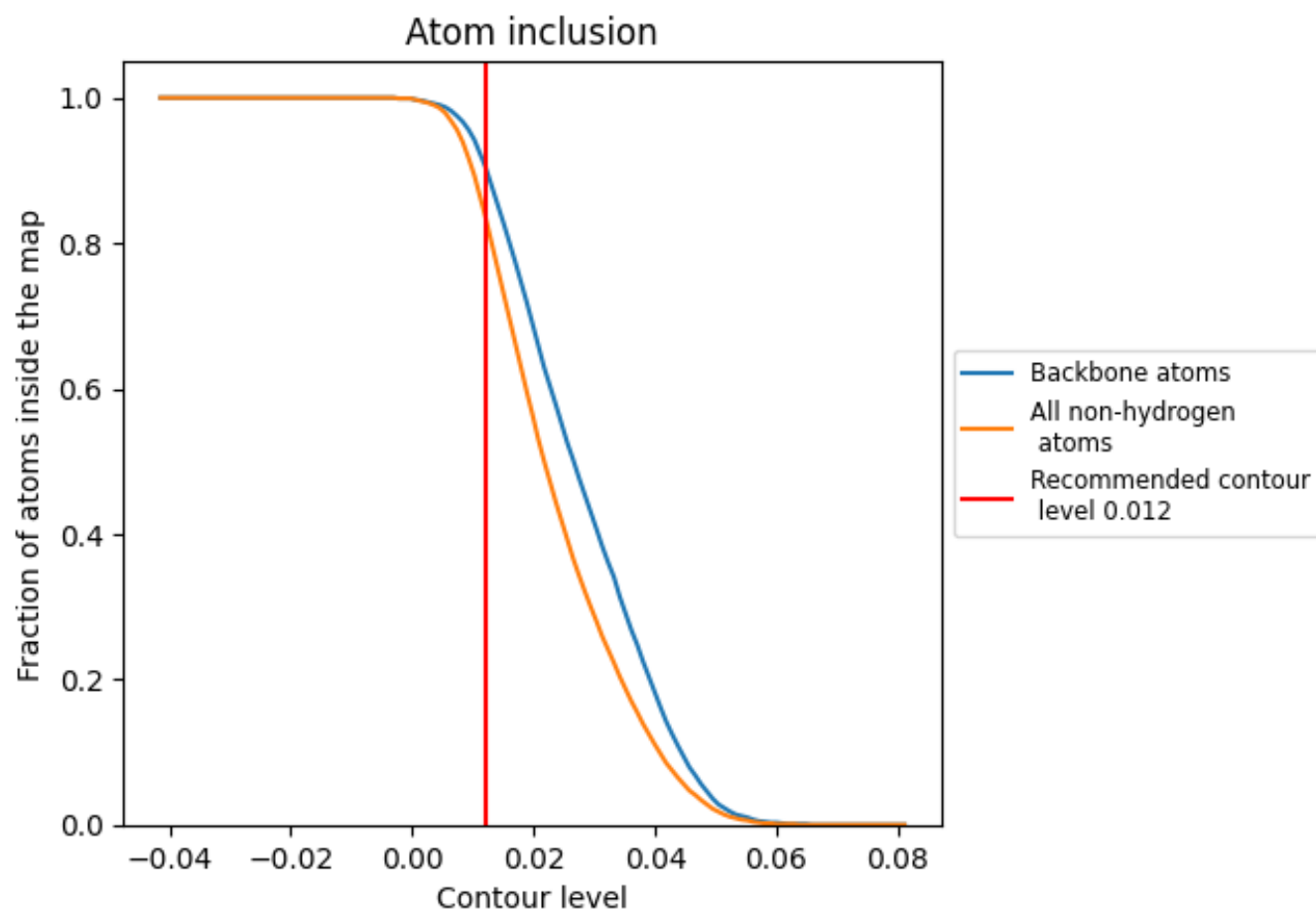
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.012).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 84% 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 (0.012) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8390	0.4300
A	0.8700	0.4520
B	0.8640	0.4460
C	0.8670	0.4510
D	0.8610	0.4500
E	0.9530	0.4120
F	0.9530	0.4390
G	0.9770	0.4470
H	0.9650	0.4370
I	0.5810	0.2320
J	0.4090	0.2100

1.0

0.0

<0.0