



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 03:53 PM UTC

PDB ID : 7SAC / pdb_00007sac
EMDB ID : EMD-24948
Title : S-(+)-ketamine bound GluN1a-GluN2B NMDA receptors at 3.69 Angstrom resolution
Authors : Chou, T.-H.; Furukawa, H.
Deposited on : 2021-09-22
Resolution : 3.69 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:

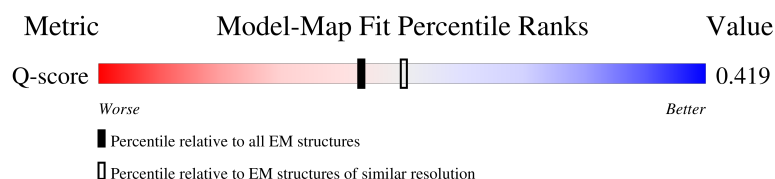
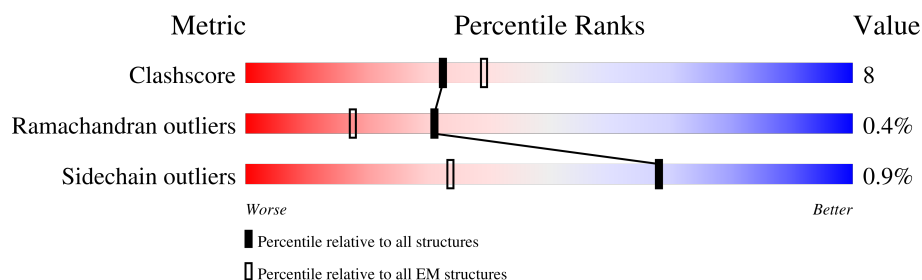
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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 3.69 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	11284 (3.19 - 4.18)

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	847	6% 81% 12% 6%
1	C	847	6% 81% 12% 6%
2	B	883	7% 74% 13% 11%
2	D	883	7% 70% 17% 11%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	E	2	50% 100%
3	F	2	50% 50%
3	G	2	100%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 23474 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 Glutamate receptor ionotropic, NMDA 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	795	Total	C	N	O	S	0	0
			5888	3765	993	1099	31		
1	C	795	Total	C	N	O	S	0	0
			5870	3755	994	1093	28		

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	SER	CYS	conflict	UNP P35439
A	61	GLN	ASN	conflict	UNP P35439
A	239	ASP	ASN	conflict	UNP P35439
A	350	GLN	ASN	conflict	UNP P35439
A	471	GLN	ASN	conflict	UNP P35439
A	491	GLN	ASN	conflict	UNP P35439
A	771	GLN	ASN	conflict	UNP P35439
A	844	ASN	ARG	conflict	UNP P35439
A	845	GLY	ARG	conflict	UNP P35439
A	846	ALA	LYS	conflict	UNP P35439
C	22	SER	CYS	conflict	UNP P35439
C	61	GLN	ASN	conflict	UNP P35439
C	239	ASP	ASN	conflict	UNP P35439
C	350	GLN	ASN	conflict	UNP P35439
C	471	GLN	ASN	conflict	UNP P35439
C	491	GLN	ASN	conflict	UNP P35439
C	771	GLN	ASN	conflict	UNP P35439
C	844	ASN	ARG	conflict	UNP P35439
C	845	GLY	ARG	conflict	UNP P35439
C	846	ALA	LYS	conflict	UNP P35439

- Molecule 2 is a protein called Glutamate receptor ionotropic, NMDA 2B.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	785	Total	C	N	O	S	0	0
			5747	3705	937	1072	33		
2	D	785	Total	C	N	O	S	0	0
			5709	3685	928	1064	32		

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-30	MET	-	expression tag	UNP Q00960
B	-29	GLY	-	expression tag	UNP Q00960
B	-28	THR	-	expression tag	UNP Q00960
B	-27	MET	-	expression tag	UNP Q00960
B	-26	ARG	-	expression tag	UNP Q00960
B	-25	LEU	-	expression tag	UNP Q00960
B	-24	PHE	-	expression tag	UNP Q00960
B	-23	LEU	-	expression tag	UNP Q00960
B	-22	LEU	-	expression tag	UNP Q00960
B	-21	ALA	-	expression tag	UNP Q00960
B	-20	VAL	-	expression tag	UNP Q00960
B	-19	LEU	-	expression tag	UNP Q00960
B	-18	PHE	-	expression tag	UNP Q00960
B	-17	LEU	-	expression tag	UNP Q00960
B	-16	PHE	-	expression tag	UNP Q00960
B	-15	SER	-	expression tag	UNP Q00960
B	-14	PHE	-	expression tag	UNP Q00960
B	-13	ALA	-	expression tag	UNP Q00960
B	-12	ARG	-	expression tag	UNP Q00960
B	-11	ALA	-	expression tag	UNP Q00960
B	-10	THR	-	expression tag	UNP Q00960
B	-9	GLY	-	expression tag	UNP Q00960
B	-8	TRP	-	expression tag	UNP Q00960
B	-7	SER	-	expression tag	UNP Q00960
B	-6	HIS	-	expression tag	UNP Q00960
B	-5	PRO	-	expression tag	UNP Q00960
B	-4	GLN	-	expression tag	UNP Q00960
B	-3	PHE	-	expression tag	UNP Q00960
B	-2	GLU	-	expression tag	UNP Q00960
B	-1	LYS	-	expression tag	UNP Q00960
B	0	GLY	-	expression tag	UNP Q00960
B	1	GLY	-	expression tag	UNP Q00960
B	2	GLY	-	expression tag	UNP Q00960
B	3	SER	-	expression tag	UNP Q00960
B	4	GLY	-	expression tag	UNP Q00960

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	5	GLY	-	expression tag	UNP Q00960
B	6	GLY	-	expression tag	UNP Q00960
B	7	SER	-	expression tag	UNP Q00960
B	8	GLY	-	expression tag	UNP Q00960
B	9	GLY	-	expression tag	UNP Q00960
B	10	SER	-	expression tag	UNP Q00960
B	11	ALA	-	expression tag	UNP Q00960
B	12	TRP	-	expression tag	UNP Q00960
B	13	SER	-	expression tag	UNP Q00960
B	14	HIS	-	expression tag	UNP Q00960
B	15	PRO	-	expression tag	UNP Q00960
B	16	GLN	-	expression tag	UNP Q00960
B	17	PHE	-	expression tag	UNP Q00960
B	18	GLU	-	expression tag	UNP Q00960
B	19	LYS	-	expression tag	UNP Q00960
B	20	GLY	-	expression tag	UNP Q00960
B	21	ALA	-	expression tag	UNP Q00960
B	22	LEU	-	expression tag	UNP Q00960
B	23	VAL	-	expression tag	UNP Q00960
B	24	PRO	-	expression tag	UNP Q00960
B	25	ARG	-	expression tag	UNP Q00960
B	26	GLY	-	expression tag	UNP Q00960
B	849	SER	CYS	conflict	UNP Q00960
D	-30	MET	-	expression tag	UNP Q00960
D	-29	GLY	-	expression tag	UNP Q00960
D	-28	THR	-	expression tag	UNP Q00960
D	-27	MET	-	expression tag	UNP Q00960
D	-26	ARG	-	expression tag	UNP Q00960
D	-25	LEU	-	expression tag	UNP Q00960
D	-24	PHE	-	expression tag	UNP Q00960
D	-23	LEU	-	expression tag	UNP Q00960
D	-22	LEU	-	expression tag	UNP Q00960
D	-21	ALA	-	expression tag	UNP Q00960
D	-20	VAL	-	expression tag	UNP Q00960
D	-19	LEU	-	expression tag	UNP Q00960
D	-18	PHE	-	expression tag	UNP Q00960
D	-17	LEU	-	expression tag	UNP Q00960
D	-16	PHE	-	expression tag	UNP Q00960
D	-15	SER	-	expression tag	UNP Q00960
D	-14	PHE	-	expression tag	UNP Q00960
D	-13	ALA	-	expression tag	UNP Q00960
D	-12	ARG	-	expression tag	UNP Q00960

Continued on next page...

Continued from previous page...

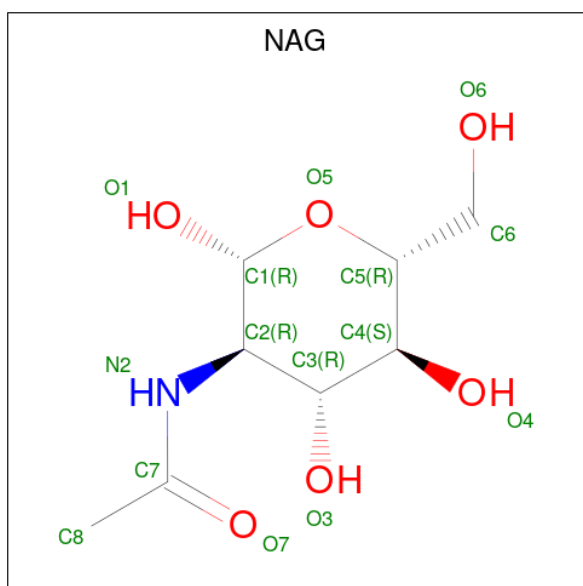
Chain	Residue	Modelled	Actual	Comment	Reference
D	-11	ALA	-	expression tag	UNP Q00960
D	-10	THR	-	expression tag	UNP Q00960
D	-9	GLY	-	expression tag	UNP Q00960
D	-8	TRP	-	expression tag	UNP Q00960
D	-7	SER	-	expression tag	UNP Q00960
D	-6	HIS	-	expression tag	UNP Q00960
D	-5	PRO	-	expression tag	UNP Q00960
D	-4	GLN	-	expression tag	UNP Q00960
D	-3	PHE	-	expression tag	UNP Q00960
D	-2	GLU	-	expression tag	UNP Q00960
D	-1	LYS	-	expression tag	UNP Q00960
D	0	GLY	-	expression tag	UNP Q00960
D	1	GLY	-	expression tag	UNP Q00960
D	2	GLY	-	expression tag	UNP Q00960
D	3	SER	-	expression tag	UNP Q00960
D	4	GLY	-	expression tag	UNP Q00960
D	5	GLY	-	expression tag	UNP Q00960
D	6	GLY	-	expression tag	UNP Q00960
D	7	SER	-	expression tag	UNP Q00960
D	8	GLY	-	expression tag	UNP Q00960
D	9	GLY	-	expression tag	UNP Q00960
D	10	SER	-	expression tag	UNP Q00960
D	11	ALA	-	expression tag	UNP Q00960
D	12	TRP	-	expression tag	UNP Q00960
D	13	SER	-	expression tag	UNP Q00960
D	14	HIS	-	expression tag	UNP Q00960
D	15	PRO	-	expression tag	UNP Q00960
D	16	GLN	-	expression tag	UNP Q00960
D	17	PHE	-	expression tag	UNP Q00960
D	18	GLU	-	expression tag	UNP Q00960
D	19	LYS	-	expression tag	UNP Q00960
D	20	GLY	-	expression tag	UNP Q00960
D	21	ALA	-	expression tag	UNP Q00960
D	22	LEU	-	expression tag	UNP Q00960
D	23	VAL	-	expression tag	UNP Q00960
D	24	PRO	-	expression tag	UNP Q00960
D	25	ARG	-	expression tag	UNP Q00960
D	26	GLY	-	expression tag	UNP Q00960
D	849	SER	CYS	conflict	UNP Q00960

- Molecule 3 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
3	E	2	Total	C	N	O	0	0
			28	16	2	10		
3	F	2	Total	C	N	O	0	0
			28	16	2	10		
3	G	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



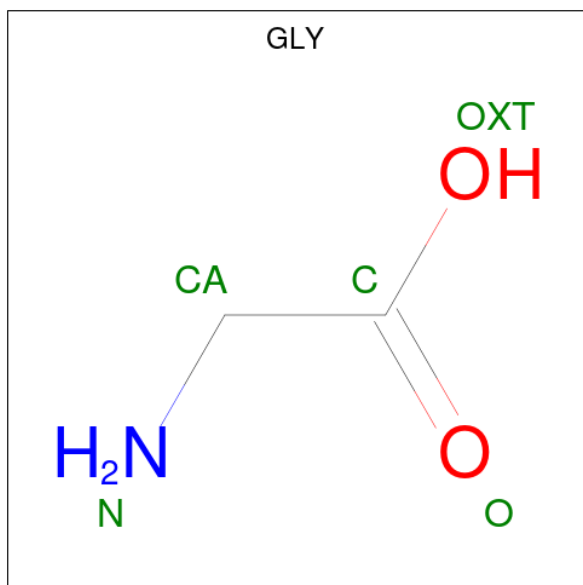
Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	D	1	Total	C	N	O	0
			14	8	1	5	

Continued on next page...

Continued from previous page...

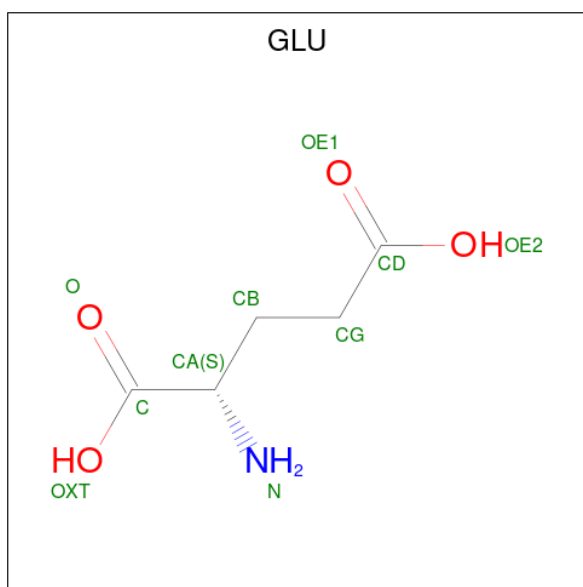
Mol	Chain	Residues	Atoms				AltConf
4	D	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 5 is GLYCINE (CCD ID: GLY) (formula: $C_2H_5NO_2$) (labeled as "Ligand of Interest" by depositor).



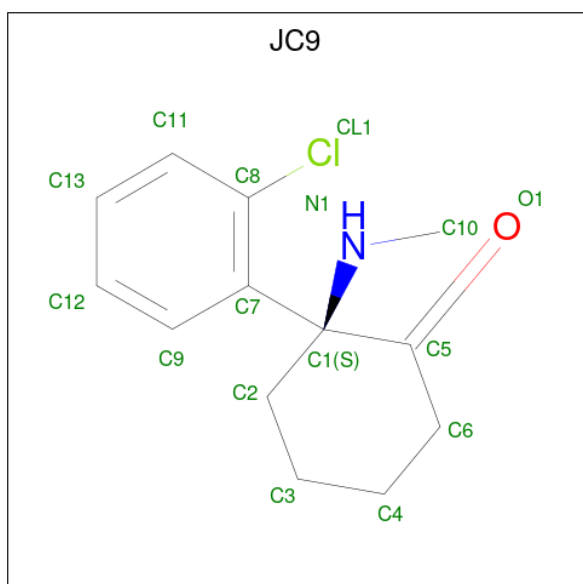
Mol	Chain	Residues	Atoms				AltConf
5	A	1	Total	C	N	O	0
			5	2	1	2	
5	C	1	Total	C	N	O	0
			5	2	1	2	

- Molecule 6 is GLUTAMIC ACID (CCD ID: GLU) (formula: $C_5H_9NO_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
6	B	1	Total	C	N	O	0
			10	5	1	4	
6	D	1	Total	C	N	O	0
			10	5	1	4	

- Molecule 7 is (2 {S})-2-(2-chlorophenyl)-2-(methylamino)cyclohexan-1-one (CCD ID: JC9) (formula: C₁₃H₁₆ClNO) (labeled as "Ligand of Interest" by depositor).



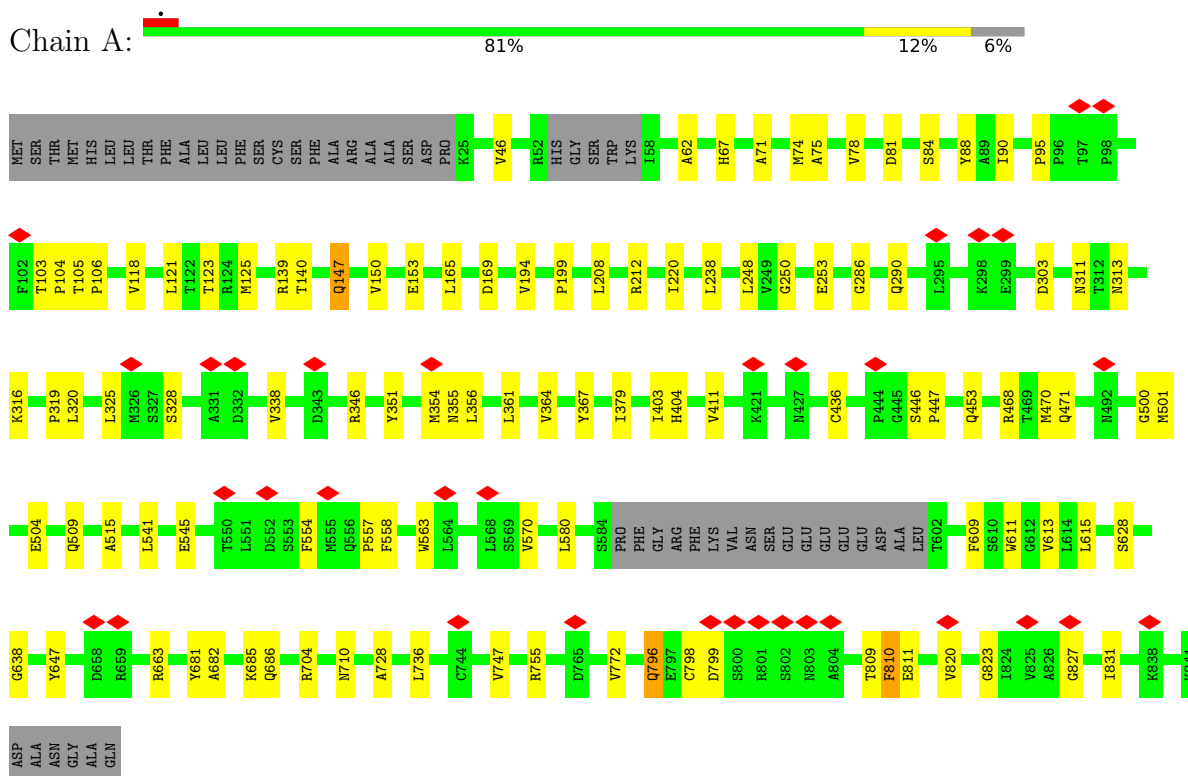
Mol	Chain	Residues	Atoms					AltConf
7	B	1	Total	C	Cl	N	O	1
			48	39	3	3	3	

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.

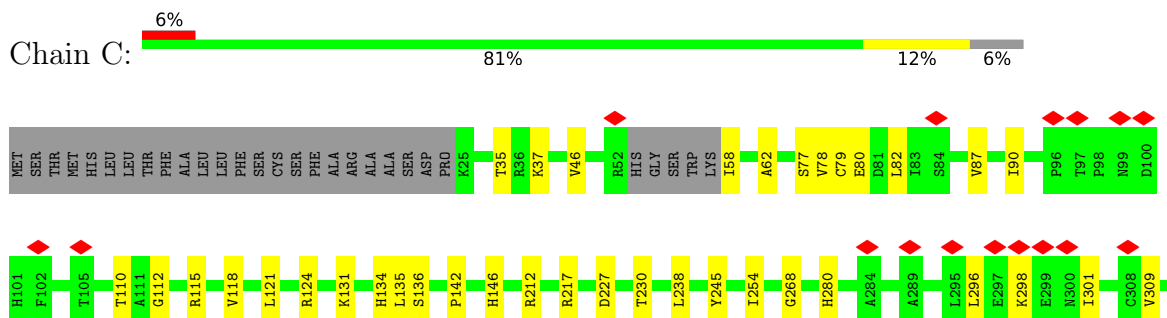
- Molecule 1: Glutamate receptor ionotropic, NMDA 1

Chain A:



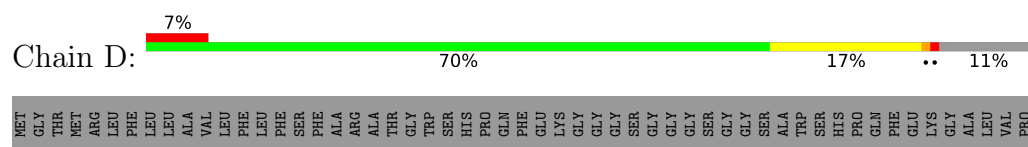
- Molecule 1: Glutamate receptor ionotropic, NMDA 1

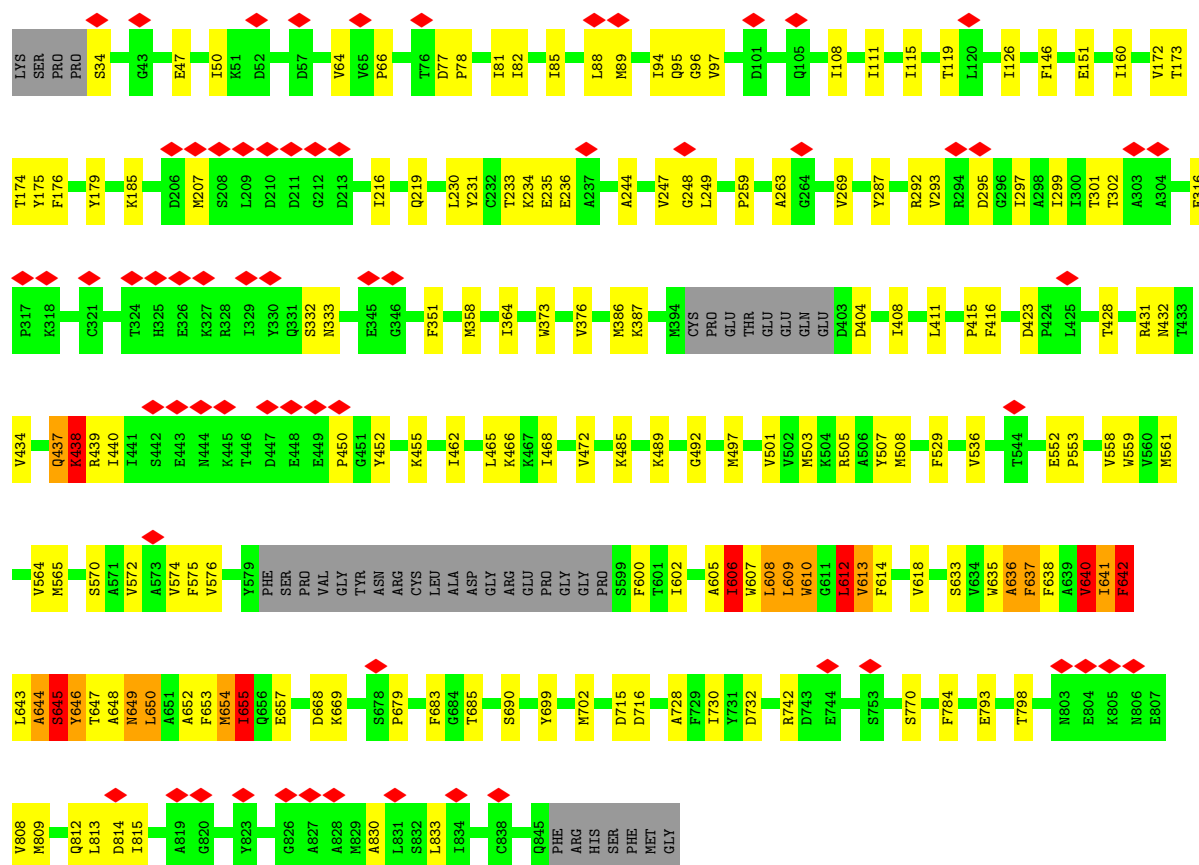
Chain C:



- Molecule 2: Glutamate receptor ionotropic, NMDA 2B

- Molecule 2: Glutamate receptor ionotropic, NMDA 2B





- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 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	434625	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	57.6	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	12.536	Depositor
Minimum map value	-5.454	Depositor
Average map value	0.004	Depositor
Map value standard deviation	0.352	Depositor
Recommended contour level	1.6	Depositor
Map size ($\text{\AA}$)	342.4, 342.4, 342.4	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.856, 0.856, 0.856	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: JC9, NAG

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	0/6015	0.34	0/8207
1	C	0.37	7/5995 (0.1%)	0.40	2/8183 (0.0%)
2	B	0.57	22/5866 (0.4%)	0.50	7/8009 (0.1%)
2	D	0.73	32/5829 (0.5%)	0.59	11/7965 (0.1%)
All	All	0.50	61/23705 (0.3%)	0.47	20/32364 (0.1%)

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
2	B	0	2

All (61) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	438	LYS	CA-C	-9.99	1.40	1.52
2	B	642	PHE	CA-C	-8.56	1.42	1.52
2	B	437	GLN	C-O	-8.28	1.14	1.24
2	D	636	ALA	CA-C	-7.93	1.42	1.52
2	D	610	TRP	C-O	-7.76	1.15	1.24
2	D	640	VAL	CA-C	-7.67	1.41	1.52
2	D	646	TYR	CA-C	-7.50	1.43	1.52
2	D	637	PHE	C-O	-7.41	1.15	1.24
2	D	613	VAL	C-O	-7.41	1.15	1.24
2	D	646	TYR	C-O	-7.25	1.15	1.24
1	C	647	TYR	N-CA	-7.22	1.37	1.46
2	D	645	SER	CA-C	-7.15	1.43	1.52
2	D	610	TRP	CA-C	-7.08	1.43	1.52
2	D	647	THR	C-O	-6.96	1.16	1.24

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	648	ALA	CA-C	-6.72	1.44	1.52
2	D	650	LEU	C-O	-6.71	1.15	1.24
2	D	649	ASN	C-O	-6.69	1.16	1.24
2	D	612	LEU	C-O	-6.65	1.15	1.24
2	D	637	PHE	CA-C	-6.58	1.44	1.52
2	B	646	TYR	CA-C	-6.57	1.44	1.52
2	B	644	ALA	C-O	-6.54	1.16	1.24
2	B	436	CYS	C-O	-6.45	1.16	1.23
2	B	648	ALA	C-N	-6.40	1.25	1.33
2	B	647	THR	CA-C	-6.32	1.44	1.52
2	B	642	PHE	C-O	-6.30	1.16	1.24
2	B	644	ALA	CA-C	-6.25	1.44	1.52
2	B	437	GLN	CA-CB	-6.25	1.44	1.53
1	C	647	TYR	CA-C	-6.19	1.44	1.52
2	B	653	PHE	C-O	-6.14	1.17	1.24
2	D	636	ALA	C-O	-6.12	1.17	1.24
2	B	642	PHE	N-CA	-6.09	1.39	1.46
2	B	438	LYS	C-O	-6.02	1.16	1.24
2	D	640	VAL	C-O	-5.98	1.16	1.24
2	B	647	THR	C-O	-5.89	1.17	1.24
2	D	437	GLN	CA-C	-5.80	1.45	1.52
2	B	649	ASN	C-O	-5.80	1.17	1.24
2	B	645	SER	C-O	-5.72	1.17	1.24
2	B	650	LEU	C-O	-5.71	1.17	1.24
2	D	646	TYR	CA-CB	-5.69	1.44	1.53
2	D	642	PHE	C-O	-5.67	1.17	1.24
2	B	652	ALA	C-O	-5.61	1.17	1.24
1	C	647	TYR	C-O	-5.58	1.17	1.24
2	D	645	SER	C-O	-5.57	1.17	1.24
2	D	641	ILE	C-O	-5.55	1.17	1.24
2	B	655	ILE	CA-C	-5.52	1.45	1.52
1	C	655	LEU	C-O	-5.45	1.17	1.24
2	D	655	ILE	CA-CB	-5.43	1.48	1.54
1	C	648	THR	C-O	-5.43	1.17	1.24
2	D	437	GLN	CA-CB	-5.34	1.45	1.53
2	D	652	ALA	C-O	-5.34	1.17	1.24
2	D	641	ILE	CA-C	-5.25	1.46	1.52
2	B	655	ILE	C-O	-5.25	1.18	1.24
2	D	437	GLN	C-O	-5.24	1.18	1.24
2	D	612	LEU	CA-CB	-5.24	1.44	1.53
1	C	646	SER	C-O	-5.23	1.18	1.24
2	D	647	THR	CA-C	-5.17	1.46	1.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	608	LEU	CA-C	-5.12	1.46	1.52
2	D	645	SER	N-CA	-5.04	1.40	1.46
2	B	617	SER	N-CA	-5.04	1.40	1.46
2	B	649	ASN	CA-C	-5.03	1.46	1.52
1	C	646	SER	CA-C	-5.00	1.46	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	606	ILE	CB-CA-C	-12.06	95.86	112.14
2	D	613	VAL	CB-CA-C	-9.33	99.54	112.14
1	C	655	LEU	N-CA-C	8.05	119.84	111.14
2	D	437	GLN	N-CA-C	7.70	121.49	108.02
2	D	612	LEU	CA-CB-CG	-7.31	90.73	116.30
2	B	642	PHE	CB-CA-C	-6.70	100.65	110.96
2	D	642	PHE	CB-CA-C	6.60	121.33	110.90
2	B	653	PHE	CB-CA-C	-6.57	99.88	110.79
2	D	613	VAL	N-CA-C	6.42	117.17	110.62
2	D	640	VAL	N-CA-CB	6.32	122.20	110.77
2	D	640	VAL	CA-C-O	-6.31	113.65	120.47
2	D	647	THR	CB-CA-C	-5.99	100.85	110.79
2	D	644	ALA	CA-C-N	-5.94	112.71	120.44
2	D	644	ALA	C-N-CA	-5.94	112.71	120.44
2	B	648	ALA	CA-C-O	5.87	126.98	120.82
2	B	616	ASN	N-CA-C	5.53	119.02	111.39
2	B	648	ALA	CA-C-N	-5.33	113.08	120.38
2	B	648	ALA	C-N-CA	-5.33	113.08	120.38
1	C	648	THR	N-CA-CB	5.19	117.84	110.16
2	B	643	LEU	N-CA-C	5.13	116.87	111.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	356	TYR	Peptide
2	B	435	PRO	Mainchain

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	5888	0	5599	73	0
1	C	5870	0	5580	83	0
2	B	5747	0	5437	91	0
2	D	5709	0	5368	164	0
3	E	28	0	25	0	0
3	F	28	0	25	1	0
3	G	28	0	25	0	0
4	A	28	0	26	0	0
4	B	14	0	13	0	0
4	C	28	0	26	0	0
4	D	28	0	26	0	0
5	A	5	0	2	0	0
5	C	5	0	2	0	0
6	B	10	0	5	0	0
6	D	10	0	5	2	0
7	B	48	0	0	5	0
All	All	23474	0	22164	387	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 8.

All (387) 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)
2:D:609:LEU:CD1	2:D:635:TRP:NE1	1.78	1.46
1:C:650:ASN:ND2	2:D:813:LEU:HD11	1.54	1.21
2:D:609:LEU:HD11	2:D:635:TRP:CE2	1.76	1.19
2:D:609:LEU:CD1	2:D:635:TRP:CD1	2.31	1.14
2:D:609:LEU:HD11	2:D:635:TRP:CD1	1.80	1.14
2:D:609:LEU:HD11	2:D:635:TRP:NE1	1.54	1.10
2:D:609:LEU:CD1	2:D:635:TRP:CE2	2.34	1.10
2:D:607:TRP:CZ2	2:D:618:VAL:HG11	1.89	1.06
2:D:438:LYS:HD3	2:D:455:LYS:NZ	1.69	1.05
7:B:903[C]:JC9:N1	7:B:903[C]:JC9:CL1	2.24	1.05
2:D:609:LEU:HD12	2:D:635:TRP:NE1	1.79	0.95
1:C:650:ASN:HD22	2:D:813:LEU:HD11	1.21	0.91
2:D:438:LYS:HD3	2:D:455:LYS:HZ3	1.27	0.90
2:D:609:LEU:HD13	2:D:635:TRP:NE1	1.85	0.89
1:C:810:PHE:CE2	1:C:813:MET:CB	2.56	0.89

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:609:LEU:HD13	2:D:635:TRP:CD1	2.09	0.87
2:B:650:LEU:HD12	2:B:650:LEU:O	1.76	0.86
2:D:386:MET:HE3	2:D:387:LYS:H	1.41	0.85
1:C:810:PHE:CD2	1:C:813:MET:CB	2.60	0.85
2:D:609:LEU:HD11	2:D:635:TRP:CD2	2.12	0.84
2:D:613:VAL:HG13	2:D:614:PHE:CD1	2.13	0.83
2:B:615:ASN:HD22	2:B:615:ASN:C	1.87	0.82
2:D:607:TRP:HZ2	2:D:618:VAL:HG11	1.40	0.81
2:D:609:LEU:HD11	2:D:635:TRP:CG	2.15	0.81
2:B:553:PRO:CB	2:B:649:ASN:HD22	1.95	0.80
2:D:609:LEU:CD1	2:D:635:TRP:HE1	1.88	0.78
2:D:609:LEU:HD12	2:D:635:TRP:CE2	2.17	0.78
1:A:554:PHE:HB3	1:A:563:TRP:HH2	1.49	0.77
2:B:553:PRO:HB3	2:B:649:ASN:ND2	1.99	0.77
2:D:439:ARG:HD2	2:D:439:ARG:O	1.85	0.76
1:C:558:PHE:CE2	1:C:647:TYR:HD2	2.03	0.76
2:D:600:PHE:CZ	2:D:605:ALA:HB2	2.22	0.75
1:C:650:ASN:HD22	2:D:813:LEU:CD1	1.98	0.75
2:D:503:MET:SD	2:D:505:ARG:NH2	2.58	0.75
1:C:634:MET:HE3	2:D:610:TRP:CD2	2.23	0.74
2:D:438:LYS:HD3	2:D:455:LYS:HZ1	1.50	0.74
1:A:125:MET:O	1:A:139:ARG:NH2	2.22	0.72
2:D:173:THR:HG23	2:D:231:TYR:HB3	1.71	0.72
1:C:650:ASN:HD21	2:D:813:LEU:HD11	1.47	0.72
2:B:101:ASP:OD1	2:B:102:ASP:N	2.21	0.72
2:B:650:LEU:HD12	2:B:650:LEU:C	2.13	0.71
2:D:650:LEU:C	2:D:650:LEU:HD23	2.15	0.71
2:D:438:LYS:CD	2:D:455:LYS:NZ	2.52	0.70
2:D:376:VAL:HG12	2:D:386:MET:HE1	1.73	0.70
2:B:648:ALA:O	2:B:649:ASN:C	2.33	0.69
1:A:647:TYR:OH	2:D:641:ILE:HD11	1.92	0.69
2:B:553:PRO:CB	2:B:649:ASN:ND2	2.56	0.69
1:C:651:LEU:HD23	1:C:651:LEU:O	1.93	0.69
1:A:123:THR:O	1:A:139:ARG:NH1	2.26	0.69
1:C:634:MET:CE	2:D:610:TRP:CE3	2.76	0.68
2:D:558:VAL:HG23	2:D:559:TRP:HD1	1.59	0.68
1:C:653:ALA:HB2	2:D:654:MET:HB3	1.76	0.67
2:B:553:PRO:HB2	2:B:649:ASN:HD22	1.60	0.67
2:D:151:GLU:OE2	2:D:185:LYS:NZ	2.20	0.67
1:A:468:ARG:O	1:A:471:GLN:NE2	2.28	0.67
2:D:606:ILE:HD13	2:D:606:ILE:N	2.09	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:654:MET:SD	2:D:654:MET:N	2.65	0.66
1:C:124:ARG:NH1	1:C:142:PRO:O	2.28	0.66
1:C:87:VAL:HG21	1:C:90:ILE:HD11	1.75	0.66
1:A:46:VAL:HG11	1:A:62:ALA:HB2	1.78	0.66
1:C:90:ILE:HB	1:C:118:VAL:HG22	1.77	0.66
1:C:484:PHE:HA	1:C:501:MET:HE3	1.77	0.66
2:D:572:VAL:HA	2:D:575:PHE:CE1	2.31	0.66
2:B:64:VAL:HG23	2:B:66:PRO:HD3	1.77	0.66
2:D:50:ILE:HD11	2:D:293:VAL:HG21	1.77	0.65
2:D:64:VAL:HG23	2:D:66:PRO:HD3	1.77	0.65
1:C:634:MET:HE3	2:D:610:TRP:CE3	2.32	0.65
2:B:553:PRO:HB3	2:B:649:ASN:HD22	1.58	0.65
2:D:608:LEU:C	2:D:608:LEU:HD23	2.21	0.65
2:B:507:TYR:HA	2:B:767:GLN:HG3	1.79	0.64
1:A:554:PHE:HB3	1:A:563:TRP:CH2	2.32	0.64
1:C:309:VAL:HA	2:D:77:ASP:HB3	1.80	0.64
2:D:231:TYR:HD1	2:D:259:PRO:HG3	1.63	0.64
2:D:607:TRP:HZ2	2:D:618:VAL:CG1	2.09	0.64
2:D:613:VAL:HG13	2:D:614:PHE:N	2.11	0.64
1:A:647:TYR:CZ	2:D:641:ILE:HD11	2.33	0.64
1:A:95:PRO:HG3	1:A:103:THR:HG21	1.80	0.63
1:C:651:LEU:HD23	1:C:651:LEU:C	2.23	0.63
1:A:286:GLY:O	1:A:290:GLN:NE2	2.31	0.63
2:D:613:VAL:HG22	2:D:613:VAL:O	1.99	0.63
2:D:633:SER:O	2:D:636:ALA:HB3	1.98	0.62
2:B:86:CYS:HA	2:B:89:MET:HE2	1.80	0.62
2:D:78:PRO:HA	2:D:81:ILE:HD12	1.82	0.62
2:D:613:VAL:CG1	2:D:614:PHE:N	2.62	0.61
2:B:552:GLU:CB	2:B:653:PHE:HE2	2.12	0.61
2:B:389:TYR:HE2	2:B:452:TYR:HB2	1.66	0.60
2:B:648:ALA:O	2:B:651:ALA:N	2.33	0.60
1:C:634:MET:CE	2:D:610:TRP:CD2	2.84	0.60
1:C:456:TYR:HA	1:C:460:ILE:HD11	1.83	0.60
1:A:403:ILE:HD12	1:A:501:MET:HE3	1.84	0.60
1:C:810:PHE:O	1:C:813:MET:N	2.35	0.59
2:D:507:TYR:OH	2:D:770:SER:OG	2.20	0.59
2:D:600:PHE:CE1	2:D:605:ALA:HB2	2.37	0.59
1:A:220:ILE:HG12	1:A:248:LEU:HB2	1.84	0.59
1:A:663:ARG:HH21	1:A:747:VAL:HG11	1.68	0.59
2:D:612:LEU:O	2:D:612:LEU:HD23	2.02	0.58
1:C:463:LEU:HD13	1:C:514:VAL:HG21	1.85	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:903[C]:JC9:CL1	7:B:903[C]:JC9:C10	2.88	0.58
7:B:903[B]:JC9:C5	7:B:903[B]:JC9:CL1	2.88	0.58
2:D:655:ILE:HG22	2:D:655:ILE:O	2.02	0.58
2:B:497:MET:HE3	2:B:509:ALA:HB1	1.85	0.58
2:D:438:LYS:O	2:D:438:LYS:HG2	2.04	0.58
2:B:552:GLU:CB	2:B:653:PHE:CE2	2.87	0.57
2:D:646:TYR:CD1	2:D:646:TYR:C	2.81	0.57
2:D:297:ILE:O	2:D:301:THR:HG23	2.05	0.57
2:D:690:SER:N	6:D:903:GLU:OE1	2.37	0.57
1:A:67:HIS:HA	1:A:74:MET:HE1	1.86	0.57
2:D:642:PHE:CD1	2:D:642:PHE:C	2.83	0.57
2:B:414:ALA:HB1	2:B:415:PRO:HD2	1.87	0.57
2:D:507:TYR:HH	2:D:770:SER:HG	1.51	0.57
2:D:216:ILE:HA	2:D:219:GLN:HG2	1.86	0.56
1:C:647:TYR:CD1	1:C:647:TYR:C	2.83	0.56
1:A:436:CYS:N	1:A:453:GLN:O	2.38	0.56
2:D:640:VAL:HG22	2:D:640:VAL:O	2.04	0.56
2:D:683:PHE:HB3	2:D:728:ALA:HB3	1.87	0.56
1:C:400:ILE:HB	1:C:476:VAL:HG12	1.87	0.56
2:B:234:LYS:NZ	2:B:263:ALA:O	2.32	0.56
2:D:600:PHE:CZ	2:D:605:ALA:CB	2.89	0.56
2:D:82:ILE:HA	2:D:85:ILE:HG22	1.88	0.56
2:B:460:PHE:HE1	2:B:792:LEU:HB3	1.70	0.56
1:A:250:GLY:N	1:A:253:GLU:OE2	2.38	0.56
2:B:438:LYS:O	2:B:438:LYS:HG2	2.05	0.56
2:D:94:ILE:HG22	2:D:96:GLY:H	1.72	0.55
2:D:610:TRP:CD1	2:D:610:TRP:C	2.84	0.55
2:B:646:TYR:CD1	2:B:646:TYR:C	2.85	0.55
2:D:287:TYR:OH	2:D:295:ASP:OD2	2.24	0.54
1:A:88:TYR:OH	1:A:303:ASP:OD1	2.24	0.54
2:D:642:PHE:HD1	2:D:643:LEU:N	2.06	0.54
2:D:439:ARG:HB2	2:D:452:TYR:CE1	2.43	0.54
1:A:364:VAL:HG11	1:A:379:ILE:HG23	1.90	0.54
1:C:35:THR:HG22	1:C:37:LYS:H	1.72	0.54
2:D:234:LYS:NZ	2:D:263:ALA:O	2.39	0.54
2:D:642:PHE:O	2:D:643:LEU:C	2.45	0.54
1:A:404:HIS:CD2	1:A:411:VAL:HG22	2.42	0.54
2:D:613:VAL:HG13	2:D:614:PHE:CG	2.42	0.54
2:B:360:PRO:HG2	2:B:362:LEU:HD11	1.89	0.54
1:A:169:ASP:HB2	1:A:199:PRO:HD3	1.91	0.53
2:B:653:PHE:CZ	1:C:807:THR:HG22	2.44	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:408:ILE:HD13	2:D:465:LEU:HD21	1.90	0.53
1:C:77:SER:HA	1:C:80:GLU:HB2	1.91	0.53
2:D:438:LYS:CD	2:D:455:LYS:HZ3	2.09	0.53
1:C:254:ILE:HD11	1:C:268:GLY:HA3	1.91	0.53
2:B:437:GLN:O	2:B:437:GLN:HG3	2.07	0.53
2:D:638:PHE:CD1	2:D:638:PHE:C	2.85	0.52
1:C:830:LEU:HA	1:C:833:ILE:HD12	1.90	0.52
2:D:434:VAL:HG13	2:D:466:LYS:HE2	1.91	0.52
1:A:504:GLU:OE1	1:A:509:GLN:NE2	2.42	0.52
2:B:358:MET:O	2:B:359:HIS:ND1	2.43	0.52
1:C:654:PHE:CD1	1:C:654:PHE:C	2.87	0.52
2:D:715:ASP:OD1	2:D:716:ASP:N	2.43	0.52
2:B:642:PHE:CD1	2:B:642:PHE:C	2.88	0.52
2:D:439:ARG:HD2	2:D:439:ARG:C	2.33	0.52
2:B:134:MET:SD	2:B:135:ALA:N	2.82	0.52
1:C:131:LYS:HG2	1:C:135:LEU:HD13	1.92	0.52
1:C:355:ASN:OD1	1:C:356:LEU:N	2.38	0.52
2:D:650:LEU:HD23	2:D:650:LEU:O	2.09	0.52
2:D:653:PHE:N	2:D:653:PHE:CD1	2.76	0.52
1:A:208:LEU:HD23	1:A:238:LEU:HD12	1.90	0.52
1:C:668:ASN:OD1	1:C:673:ARG:NH2	2.43	0.52
1:C:559:GLN:N	2:D:812:GLN:O	2.42	0.52
1:C:649:ALA:O	1:C:650:ASN:C	2.48	0.52
2:D:607:TRP:CZ2	2:D:618:VAL:CG1	2.77	0.52
2:D:742:ARG:NH2	2:D:798:THR:O	2.43	0.52
1:C:112:GLY:O	1:C:115:ARG:NH1	2.38	0.51
2:D:561:MET:HA	2:D:564:VAL:HG22	1.92	0.51
1:A:446:SER:HB2	1:A:447:PRO:HD3	1.91	0.51
2:D:64:VAL:HG21	2:D:301:THR:HB	1.91	0.51
1:A:820:VAL:HG13	2:D:565:MET:HE1	1.93	0.51
2:D:732:ASP:OD2	6:D:903:GLU:N	2.43	0.51
1:C:651:LEU:C	1:C:651:LEU:CD2	2.84	0.51
2:B:47:GLU:HA	2:B:50:ILE:HD12	1.92	0.50
1:A:796:GLN:NE2	1:A:799:ASP:HB3	2.27	0.50
2:B:501:VAL:HA	2:B:506:ALA:HB3	1.94	0.50
1:A:311:ASN:HD21	1:A:313:ASN:HB2	1.77	0.50
2:D:431:ARG:HG2	2:D:432:ASN:H	1.77	0.50
2:D:536:VAL:HG22	2:D:730:ILE:HG12	1.93	0.50
1:A:354:MET:HE2	1:A:361:LEU:HB3	1.93	0.50
2:D:50:ILE:HD11	2:D:293:VAL:CG2	2.41	0.50
2:D:172:VAL:O	2:D:230:LEU:HA	2.12	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:247:VAL:HG13	2:D:249:LEU:HD23	1.93	0.50
2:D:699:TYR:HB2	2:D:702:MET:HB2	1.93	0.50
2:D:207:MET:HE3	2:D:236:GLU:HB3	1.93	0.50
1:C:541:LEU:HD21	1:C:746:LEU:HD13	1.94	0.49
2:D:78:PRO:HB3	2:D:111:ILE:HG12	1.93	0.49
1:A:755:ARG:NH1	2:D:793:GLU:OE1	2.45	0.49
2:B:311:HIS:O	2:B:311:HIS:ND1	2.44	0.49
2:B:467:LYS:NZ	2:B:788:GLU:HG2	2.27	0.49
1:C:681:TYR:HB3	1:C:728:ALA:HB3	1.93	0.49
2:D:437:GLN:O	2:D:437:GLN:HG3	2.12	0.49
2:B:685:THR:HG22	2:B:730:ILE:HB	1.93	0.49
1:C:541:LEU:HB2	1:C:736:LEU:HD13	1.94	0.49
2:D:85:ILE:HD11	2:D:97:VAL:HG21	1.93	0.49
2:D:438:LYS:CD	2:D:455:LYS:HZ1	2.19	0.49
2:B:663:VAL:HG13	2:B:751:ILE:HD11	1.93	0.49
2:D:34:SER:N	2:D:64:VAL:O	2.45	0.49
2:B:615:ASN:O	2:B:615:ASN:ND2	2.44	0.48
2:D:115:ILE:O	2:D:119:THR:OG1	2.24	0.48
2:D:608:LEU:C	2:D:608:LEU:CD2	2.85	0.48
1:C:646:SER:OG	1:C:647:TYR:N	2.45	0.48
2:D:175:TYR:O	2:D:179:TYR:HB2	2.13	0.48
2:D:439:ARG:HG2	2:D:440:ILE:N	2.29	0.48
2:B:507:TYR:HD1	2:B:767:GLN:HE21	1.61	0.48
2:D:653:PHE:N	2:D:653:PHE:HD1	2.10	0.48
2:D:160:ILE:HD11	2:D:364:ILE:HD13	1.96	0.48
1:A:75:ALA:HA	1:A:78:VAL:HG12	1.95	0.48
1:C:577:LEU:HD22	1:C:632:LEU:HD23	1.96	0.48
1:A:325:LEU:O	1:A:328:SER:OG	2.30	0.48
2:D:613:VAL:CG1	2:D:614:PHE:CD1	2.93	0.48
1:A:810:PHE:HZ	2:D:561:MET:HG2	1.79	0.48
1:C:634:MET:HE1	2:D:610:TRP:CE3	2.48	0.48
1:C:711:TYR:HE2	1:C:726:LEU:HD22	1.79	0.47
2:D:146:PHE:HE1	2:D:351:PHE:HZ	1.61	0.47
2:B:35:ILE:HD11	2:B:314:ILE:HD11	1.95	0.47
2:B:258:VAL:HB	2:B:262:VAL:HG11	1.96	0.47
2:B:343:THR:OG1	2:B:347:ARG:O	2.31	0.47
2:B:750:THR:O	2:B:751:ILE:HG22	2.14	0.47
2:D:358:MET:HE3	2:D:358:MET:HA	1.96	0.47
1:A:809:THR:HG23	1:A:811:GLU:HG3	1.97	0.47
2:B:818:MET:O	2:B:822:PHE:HD1	1.97	0.47
1:C:634:MET:HE3	2:D:610:TRP:CE2	2.49	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:439:ARG:C	2:D:439:ARG:CD	2.87	0.47
2:D:552:GLU:CB	2:D:653:PHE:HE2	2.28	0.47
1:C:82:LEU:HD13	1:C:87:VAL:HG22	1.96	0.47
2:D:332:SER:OG	2:D:333:ASN:N	2.47	0.47
1:A:351:TYR:O	1:A:367:TYR:N	2.45	0.47
2:B:389:TYR:CE2	2:B:452:TYR:HB2	2.49	0.47
2:D:415:PRO:HG2	2:D:416:PHE:CD1	2.50	0.47
2:B:565:MET:HE1	1:C:817:PHE:CG	2.50	0.47
1:C:368:ASN:HB3	3:F:1:NAG:O5	2.15	0.47
1:A:81:ASP:O	1:A:84:SER:OG	2.29	0.46
1:A:355:ASN:OD1	1:A:356:LEU:N	2.48	0.46
2:B:101:ASP:OD1	2:B:103:THR:N	2.48	0.46
2:B:212:GLY:O	2:B:216:ILE:HG12	2.15	0.46
2:B:467:LYS:HZ3	2:B:788:GLU:HG2	1.80	0.46
2:B:699:TYR:HB3	2:B:702:MET:HG2	1.97	0.46
1:A:570:VAL:HG11	1:A:611:TRP:HD1	1.81	0.46
2:B:358:MET:HE2	2:B:358:MET:HA	1.95	0.46
1:C:682:ALA:HB1	1:C:710:ASN:HA	1.96	0.46
2:D:126:ILE:O	2:D:292:ARG:NH1	2.46	0.46
1:A:796:GLN:HE21	1:A:798:CYS:C	2.23	0.46
1:C:79:CYS:HA	1:C:82:LEU:HB3	1.97	0.46
2:D:269:VAL:HG22	2:D:373:TRP:HE1	1.81	0.46
1:A:354:MET:CE	1:A:361:LEU:HB3	2.46	0.46
2:D:570:SER:O	2:D:574:VAL:HG23	2.16	0.46
2:D:644:ALA:O	2:D:645:SER:C	2.56	0.46
2:D:830:ALA:O	2:D:833:LEU:HG	2.16	0.46
2:D:95:GLN:HE22	2:D:316:GLU:HA	1.81	0.46
2:D:808:VAL:HG13	2:D:809:MET:N	2.30	0.46
2:B:820:GLY:O	2:B:824:MET:HG2	2.16	0.45
1:A:105:THR:OG1	1:A:106:PRO:HD3	2.16	0.45
1:A:501:MET:HE1	1:A:515:ALA:HB3	1.98	0.45
1:C:810:PHE:O	1:C:811:GLU:C	2.58	0.45
1:A:103:THR:OG1	1:A:104:PRO:HD3	2.16	0.45
1:C:631:ILE:HA	1:C:634:MET:HB2	1.97	0.45
1:C:657:LEU:HD11	2:D:655:ILE:HG23	1.97	0.45
1:C:810:PHE:CD2	1:C:810:PHE:O	2.70	0.45
1:A:165:LEU:HD11	1:A:194:VAL:HG12	1.99	0.45
2:B:357:GLN:O	2:B:358:MET:HB2	2.17	0.45
2:B:639:ALA:O	2:B:642:PHE:HB3	2.17	0.45
2:B:643:LEU:O	2:B:643:LEU:HG	2.14	0.45
2:B:230:LEU:HD21	2:B:232:CYS:HB2	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:903[C]:JC9:CL1	7:B:903[C]:JC9:O1	2.72	0.45
2:D:85:ILE:O	2:D:89:MET:HG2	2.17	0.45
2:B:615:ASN:C	2:B:615:ASN:ND2	2.59	0.45
1:C:406:GLU:HB3	1:C:407:PRO:HD3	1.99	0.45
1:A:71:ALA:HB1	1:A:106:PRO:HG3	1.98	0.44
2:B:565:MET:HE1	1:C:817:PHE:HB3	1.98	0.44
1:C:217:ARG:HD3	1:C:245:TYR:CE1	2.52	0.44
2:D:553:PRO:HB3	2:D:649:ASN:ND2	2.32	0.44
1:A:580:LEU:HD12	1:A:628:SER:HB2	1.99	0.44
2:B:141:SER:O	2:B:142:MET:HE2	2.17	0.44
2:B:544:THR:HG23	2:B:810:SER:HB2	1.99	0.44
2:B:714:VAL:HG11	2:B:739:MET:HE3	1.99	0.44
2:D:173:THR:HG22	2:D:174:THR:O	2.17	0.44
2:B:204:LEU:HD23	2:B:204:LEU:O	2.18	0.44
1:C:78:VAL:HG11	1:C:110:THR:HG21	1.99	0.44
1:C:364:VAL:HG12	1:C:364:VAL:O	2.17	0.44
1:C:722:ARG:NH1	1:C:743:LYS:HD2	2.33	0.44
1:A:147:GLN:HA	1:A:150:VAL:HG12	2.00	0.44
2:B:755:LYS:HB3	2:B:755:LYS:HE2	1.73	0.44
1:C:507:SER:O	1:C:509:GLN:HG3	2.17	0.44
1:C:810:PHE:O	1:C:810:PHE:HD2	2.00	0.44
1:A:809:THR:HG23	1:A:811:GLU:H	1.83	0.44
1:A:541:LEU:HD22	1:A:736:LEU:HD22	1.99	0.44
2:B:298:ALA:HA	2:B:301:THR:HG1	1.82	0.44
1:C:398:LEU:HD21	1:C:768:TRP:HE1	1.83	0.44
1:A:663:ARG:HA	1:A:663:ARG:HD2	1.89	0.44
1:A:638:GLY:HA2	2:B:614:PHE:CE2	2.53	0.43
2:B:107:ALA:O	2:B:111:ILE:HG13	2.18	0.43
1:C:46:VAL:HG11	1:C:62:ALA:HB2	1.99	0.43
1:A:558:PHE:HD2	1:A:563:TRP:HE1	1.66	0.43
2:B:137:LYS:NZ	2:B:143:PHE:O	2.50	0.43
2:B:498:ILE:HA	2:B:501:VAL:HG12	1.99	0.43
2:B:603:GLY:HA2	2:B:606:ILE:HG22	1.99	0.43
1:C:663:ARG:HE	1:C:749:THR:HG22	1.84	0.43
2:B:749:VAL:HG22	2:B:750:THR:H	1.83	0.43
1:C:580:LEU:HD21	1:C:628:SER:OG	2.19	0.43
2:D:411:LEU:HD22	2:D:485:LYS:O	2.19	0.43
1:A:316:LYS:O	1:A:319:PRO:HD2	2.19	0.43
1:A:470:MET:HE1	1:A:772:VAL:HA	2.01	0.43
2:B:654:MET:HB2	2:B:654:MET:HE3	1.44	0.43
1:C:58:ILE:HG22	1:C:296:LEU:HD22	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:287:TYR:HE2	2:D:292:ARG:HG2	1.84	0.43
1:A:681:TYR:HB3	1:A:728:ALA:HB3	2.01	0.43
2:B:545:VAL:HG22	2:B:547:PRO:HD3	2.01	0.43
2:D:108:ILE:HA	2:D:111:ILE:HG22	2.01	0.43
1:A:823:GLY:O	1:A:827:GLY:N	2.52	0.43
2:D:47:GLU:O	2:D:50:ILE:HG22	2.18	0.43
2:D:423:ASP:HB3	2:D:428:THR:O	2.19	0.43
2:D:815:ILE:HG22	2:D:815:ILE:O	2.18	0.43
2:B:537:MET:SD	2:B:736:LEU:HD22	2.59	0.43
2:B:750:THR:O	2:B:750:THR:HG22	2.19	0.43
2:B:788:GLU:OE2	2:B:788:GLU:N	2.50	0.43
1:C:559:GLN:HG3	1:C:561:THR:H	1.84	0.43
1:A:404:HIS:HD2	1:A:411:VAL:HG22	1.83	0.42
1:C:134:HIS:C	1:C:136:SER:H	2.27	0.42
2:D:66:PRO:HG3	2:D:301:THR:HG21	1.99	0.42
1:A:704:ARG:CZ	2:B:430:MET:HE2	2.48	0.42
2:B:558:VAL:HG23	2:B:559:TRP:HD1	1.83	0.42
1:C:408:PHE:HB3	1:C:459:CYS:SG	2.58	0.42
1:C:212:ARG:NH2	1:C:238:LEU:O	2.52	0.42
2:D:462:ILE:HG22	2:D:466:LYS:HE3	2.00	0.42
2:D:508:MET:HE3	2:D:508:MET:HB2	1.90	0.42
2:B:653:PHE:CZ	1:C:807:THR:CG2	3.02	0.42
1:A:355:ASN:HD22	1:A:379:ILE:HG22	1.85	0.42
1:A:609:PHE:O	1:A:613:VAL:HG23	2.20	0.42
1:A:827:GLY:O	1:A:831:ILE:HG12	2.19	0.42
2:D:552:GLU:CB	2:D:653:PHE:CE2	3.02	0.42
1:A:338:VAL:HG22	1:A:346:ARG:HE	1.85	0.42
1:C:227:ASP:O	1:C:230:THR:OG1	2.35	0.42
2:D:111:ILE:O	2:D:115:ILE:HG13	2.19	0.42
2:D:287:TYR:CE2	2:D:292:ARG:HG2	2.55	0.42
2:D:295:ASP:O	2:D:299:ILE:HG23	2.20	0.42
1:A:103:THR:O	1:A:106:PRO:HD2	2.19	0.42
1:A:682:ALA:HB1	1:A:710:ASN:HA	2.02	0.42
2:B:644:ALA:HB1	1:C:648:THR:OG1	2.20	0.42
2:D:233:THR:OG1	2:D:236:GLU:HG2	2.18	0.42
1:A:570:VAL:HG11	1:A:611:TRP:CD1	2.55	0.42
1:A:685:LYS:HG2	1:A:686:GLN:HG3	2.02	0.42
2:B:187:ARG:O	2:B:191:GLU:HG2	2.20	0.42
2:B:249:LEU:HB2	2:B:254:TYR:CE2	2.55	0.42
1:C:487:GLN:CD	1:C:496:LYS:HE2	2.45	0.42
2:D:299:ILE:HA	2:D:302:THR:HG22	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:244:ALA:O	2:D:248:GLY:N	2.49	0.42
2:D:529:PHE:HA	2:D:784:PHE:HZ	1.85	0.42
1:A:545:GLU:N	1:A:545:GLU:OE1	2.53	0.41
1:C:479:VAL:HG13	1:C:482:GLY:H	1.84	0.41
1:C:558:PHE:HE2	1:C:647:TYR:HD2	1.63	0.41
1:A:90:ILE:HB	1:A:118:VAL:HG22	2.02	0.41
2:B:362:LEU:HD13	2:B:379:TRP:CZ2	2.55	0.41
1:A:212:ARG:HH12	1:A:238:LEU:HB3	1.85	0.41
1:A:557:PRO:O	2:B:812:GLN:HA	2.21	0.41
1:C:539:THR:HG23	1:C:736:LEU:HD12	2.01	0.41
7:B:903[A]:JC9:C5	7:B:903[A]:JC9:CL1	3.06	0.41
2:D:685:THR:HG22	2:D:730:ILE:HB	2.02	0.41
1:A:153:GLU:HA	1:A:153:GLU:OE1	2.20	0.41
2:B:249:LEU:HD23	2:B:249:LEU:H	1.85	0.41
2:B:554:PHE:HE1	1:C:808:LEU:HD21	1.85	0.41
2:B:646:TYR:O	2:B:647:THR:C	2.61	0.41
2:D:235:GLU:HG2	2:D:236:GLU:N	2.35	0.41
2:D:814:ASP:N	2:D:814:ASP:OD1	2.52	0.41
2:B:540:ARG:NH2	2:B:746:CYS:O	2.52	0.41
2:B:430:MET:SD	2:B:430:MET:N	2.91	0.41
2:D:468:ILE:O	2:D:472:VAL:HG22	2.20	0.41
2:B:162:GLU:HB3	2:B:197:TRP:HH2	1.86	0.41
2:B:439:ARG:HD2	2:B:440:ILE:N	2.36	0.41
1:C:121:LEU:HD22	1:C:280:HIS:HB3	2.03	0.41
2:D:497:MET:O	2:D:501:VAL:HG23	2.20	0.41
2:D:602:ILE:CD1	2:D:606:ILE:HD11	2.51	0.41
2:D:668:ASP:OD1	2:D:669:LYS:N	2.54	0.41
1:A:320:LEU:HD23	1:A:320:LEU:H	1.87	0.41
2:B:269:VAL:HG11	2:B:371:ARG:HB3	2.03	0.41
2:D:404:ASP:OD1	2:D:404:ASP:N	2.53	0.41
1:A:500:GLY:O	1:A:504:GLU:HG2	2.21	0.40
1:C:298:LYS:H	1:C:301:ILE:HD11	1.86	0.40
1:C:751:GLU:CD	1:C:751:GLU:H	2.29	0.40
1:A:615:LEU:HD11	2:D:637:PHE:HD1	1.87	0.40
2:B:683:PHE:HE1	2:B:706:MET:HE2	1.86	0.40
2:D:88:LEU:HB3	2:D:94:ILE:HD12	2.02	0.40
2:D:612:LEU:HD23	2:D:612:LEU:C	2.47	0.40
1:A:121:LEU:HA	1:A:140:THR:HB	2.04	0.40
2:D:115:ILE:HG13	2:D:115:ILE:H	1.71	0.40
2:D:489:LYS:HZ3	2:D:492:GLY:HA2	1.86	0.40
2:D:572:VAL:O	2:D:576:VAL:HG23	2.22	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:640:VAL:O	2:D:640:VAL:HG13	2.21	0.40
2:B:121:THR:HA	2:B:122:PRO:HD3	1.90	0.40
1:C:507:SER:OG	1:C:509:GLN:OE1	2.35	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	789/847 (93%)	748 (95%)	40 (5%)	1 (0%)	48	78
1	C	789/847 (93%)	761 (96%)	27 (3%)	1 (0%)	48	78
2	B	779/883 (88%)	734 (94%)	40 (5%)	5 (1%)	21	52
2	D	779/883 (88%)	725 (93%)	50 (6%)	4 (0%)	24	56
All	All	3136/3460 (91%)	2968 (95%)	157 (5%)	11 (0%)	31	60

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	289	LEU
2	B	450	PRO
2	B	679	PRO
1	C	664	ILE
2	D	450	PRO
2	D	679	PRO
2	B	414	ALA
2	B	751	ILE
2	D	657	GLU
1	A	796	GLN
2	D	176	PHE

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	595/728 (82%)	593 (100%)	2 (0%)	86	83
1	C	589/728 (81%)	584 (99%)	5 (1%)	73	76
2	B	577/762 (76%)	572 (99%)	5 (1%)	70	75
2	D	568/762 (74%)	559 (98%)	9 (2%)	55	68
All	All	2329/2980 (78%)	2308 (99%)	21 (1%)	68	75

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

Mol	Chain	Res	Type
1	A	147	GLN
1	A	810	PHE
2	B	347	ARG
2	B	647	THR
2	B	650	LEU
2	B	653	PHE
2	B	655	ILE
1	C	146	HIS
1	C	647	TYR
1	C	648	THR
1	C	651	LEU
1	C	655	LEU
2	D	438	LYS
2	D	606	ILE
2	D	609	LEU
2	D	612	LEU
2	D	640	VAL
2	D	642	PHE
2	D	645	SER
2	D	654	MET
2	D	655	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	A	276	ASN
1	A	311	ASN
1	A	357	GLN
1	A	405	GLN
1	A	650	ASN
1	A	724	ASN
1	A	796	GLN
2	B	159	ASN
2	B	192	ASN
2	B	437	GLN
2	B	615	ASN
2	B	649	ASN
2	B	782	GLN
1	C	177	GLN
1	C	393	GLN
1	C	686	GLN
2	D	153	GLN
2	D	219	GLN
2	D	656	GLN
2	D	775	GLN

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](#)

6 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
3	NAG	E	1	2,3	14,14,15	0.28	0	17,19,21	0.46	0
3	NAG	E	2	3	14,14,15	0.23	0	17,19,21	0.45	0
3	NAG	F	1	1,3	14,14,15	0.59	0	17,19,21	0.68	1 (5%)
3	NAG	F	2	3	14,14,15	0.28	0	17,19,21	0.48	0
3	NAG	G	1	2,3	14,14,15	0.34	0	17,19,21	0.55	0
3	NAG	G	2	3	14,14,15	0.25	0	17,19,21	0.40	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
3	NAG	E	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	NAG	G	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	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(°)
3	F	1	NAG	C1-O5-C5	2.06	114.94	112.19

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	F	2	NAG	C4-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
3	F	1	NAG	C1-C2-N2-C7

Continued on next page...

Continued from previous page...

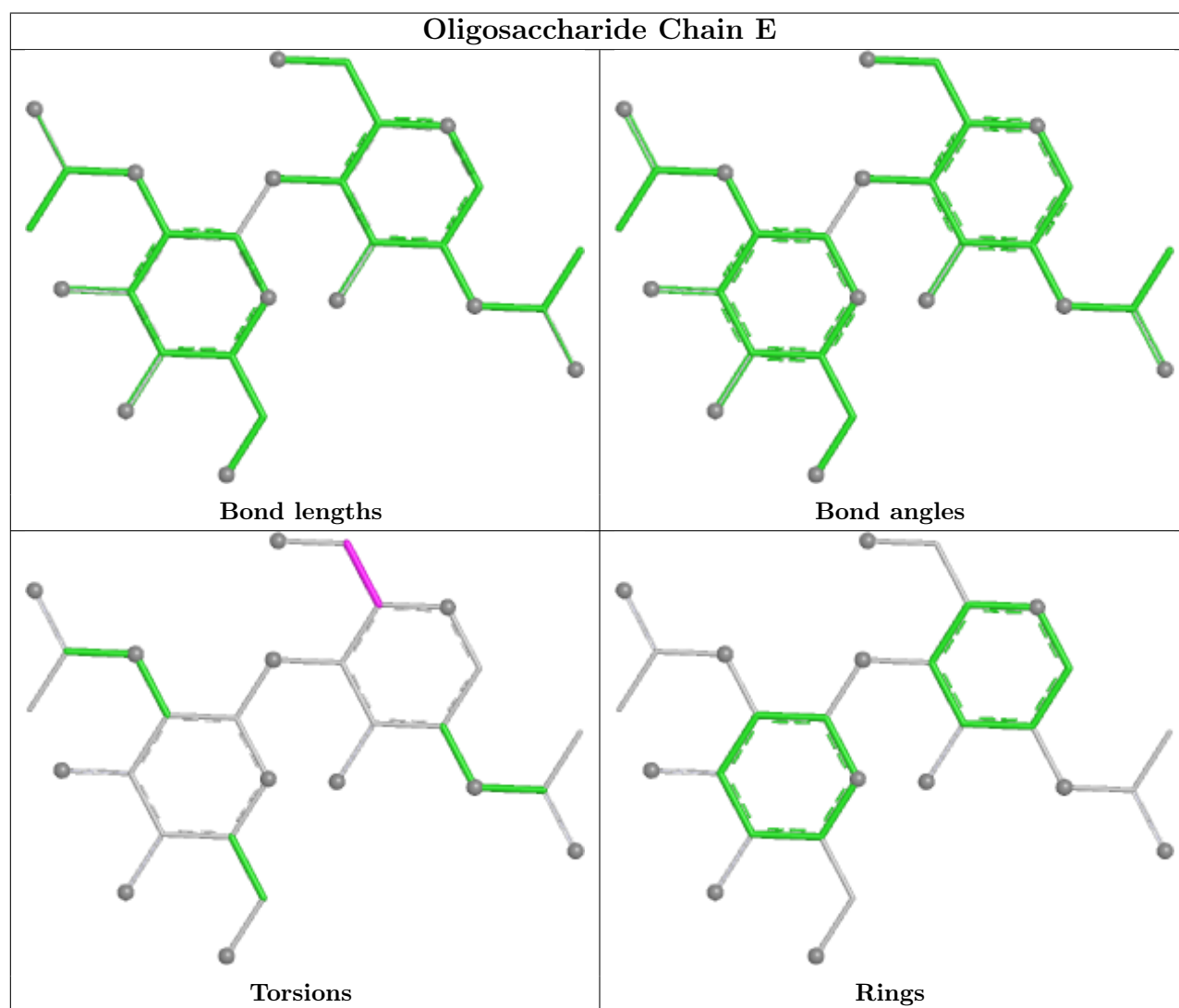
Mol	Chain	Res	Type	Atoms
3	F	1	NAG	C3-C2-N2-C7

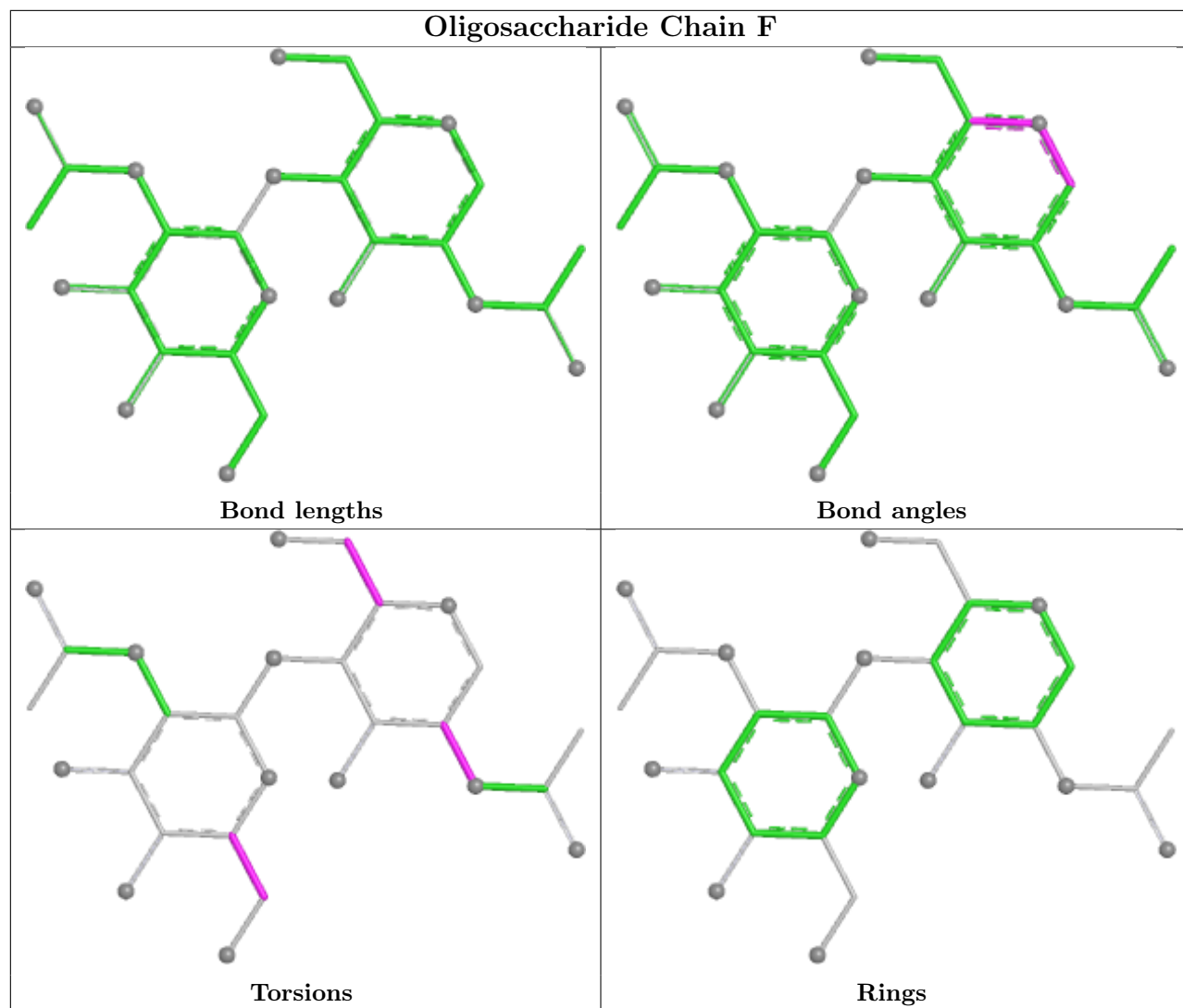
There are no ring outliers.

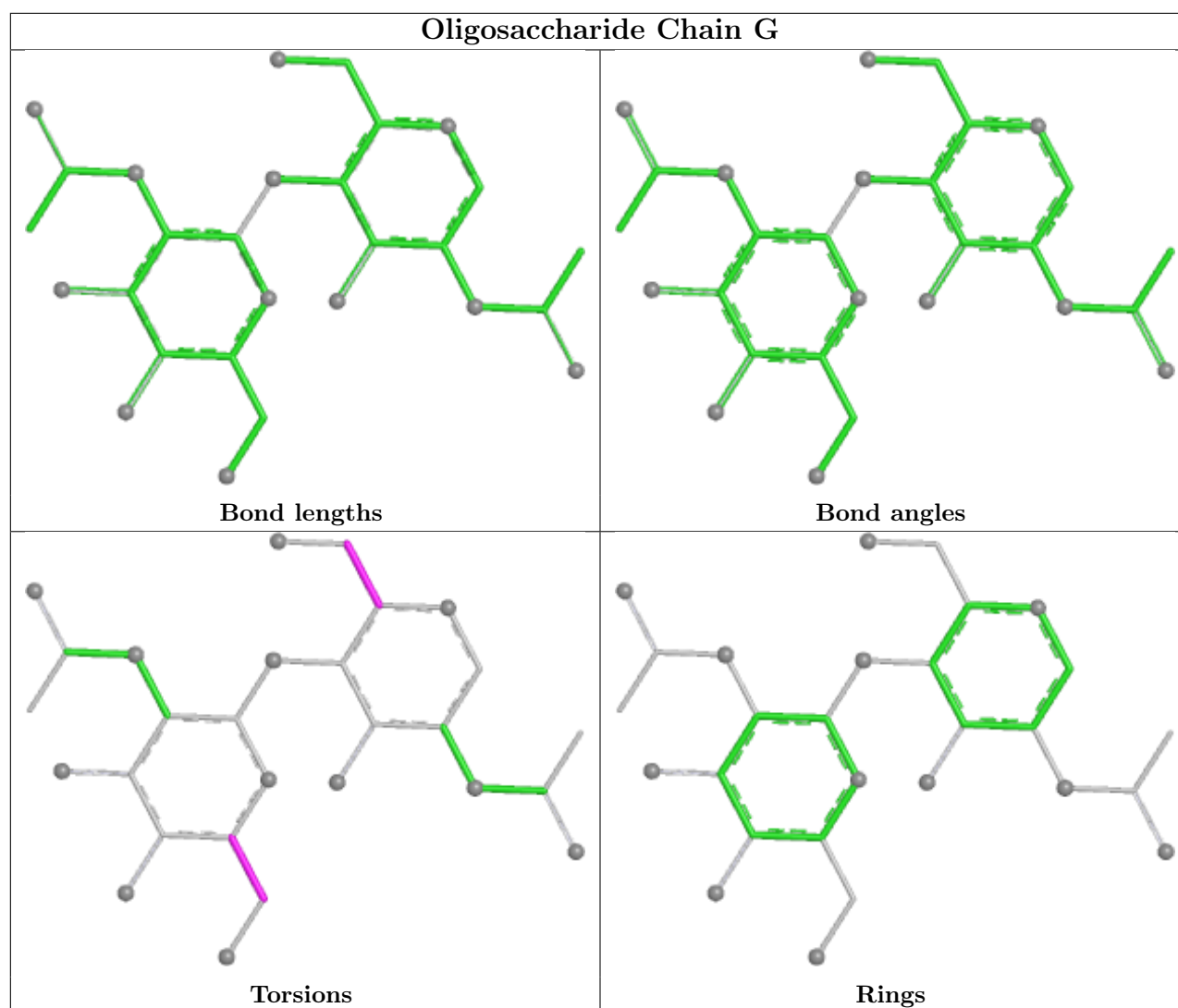
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1	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](#)

14 ligands 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	$\# Z > 2$
6	GLU	D	903	-	8,9,9	1.09	1 (12%)	8,11,11	1.24	1 (12%)
5	GLY	A	903	-	4,4,4	1.15	1 (25%)	3,4,4	1.64	1 (33%)
5	GLY	C	903	-	4,4,4	1.16	1 (25%)	3,4,4	1.63	1 (33%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GLU	B	902	-	8,9,9	1.10	1 (12%)	8,11,11	1.19	1 (12%)
4	NAG	A	902	1	14,14,15	0.28	0	17,19,21	0.59	0
4	NAG	A	901	-	14,14,15	0.20	0	17,19,21	0.44	0
7	JC9	B	903[A]	-	15,17,17	1.36	2 (13%)	15,24,24	2.05	5 (33%)
7	JC9	B	903[C]	-	15,17,17	2.19	6 (40%)	15,24,24	2.84	6 (40%)
4	NAG	B	901	2	14,14,15	0.95	1 (7%)	17,19,21	1.46	1 (5%)
4	NAG	C	902	-	14,14,15	0.18	0	17,19,21	0.43	0
4	NAG	D	902	-	14,14,15	0.20	0	17,19,21	0.43	0
4	NAG	D	901	-	14,14,15	0.42	0	17,19,21	0.77	1 (5%)
7	JC9	B	903[B]	-	15,17,17	2.23	6 (40%)	15,24,24	3.00	5 (33%)
4	NAG	C	901	-	14,14,15	0.20	0	17,19,21	0.44	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	GLU	D	903	-	-	0/9/9/9	-
5	GLY	A	903	-	-	0/2/2/2	-
5	GLY	C	903	-	-	0/2/2/2	-
6	GLU	B	902	-	-	0/9/9/9	-
4	NAG	A	902	1	-	4/6/23/26	0/1/1/1
4	NAG	A	901	-	-	2/6/23/26	0/1/1/1
7	JC9	B	903[A]	-	-	0/6/23/23	0/2/2/2
7	JC9	B	903[C]	-	-	3/6/23/23	0/2/2/2
4	NAG	B	901	2	-	0/6/23/26	0/1/1/1
4	NAG	C	902	-	-	0/6/23/26	0/1/1/1
4	NAG	D	902	-	-	2/6/23/26	0/1/1/1
4	NAG	D	901	-	-	0/6/23/26	0/1/1/1
7	JC9	B	903[B]	-	-	4/6/23/23	0/2/2/2
4	NAG	C	901	-	-	2/6/23/26	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	903[C]	JC9	C8-C7	-4.54	1.33	1.39
7	B	903[C]	JC9	O1-C5	-4.38	1.14	1.21
7	B	903[A]	JC9	C6-C5	4.14	1.56	1.50

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	903[B]	JC9	C6-C5	4.00	1.56	1.50
7	B	903[B]	JC9	C9-C7	-3.59	1.34	1.39
7	B	903[B]	JC9	C1-C7	-3.52	1.50	1.53
4	B	901	NAG	O5-C1	3.48	1.49	1.43
7	B	903[B]	JC9	C2-C1	-2.78	1.50	1.55
7	B	903[B]	JC9	O1-C5	-2.66	1.17	1.21
7	B	903[C]	JC9	C6-C5	2.51	1.54	1.50
7	B	903[B]	JC9	C2-C3	-2.43	1.46	1.52
7	B	903[C]	JC9	C1-C7	-2.26	1.51	1.53
6	D	903	GLU	OXT-C	-2.25	1.23	1.30
6	B	902	GLU	OXT-C	-2.23	1.23	1.30
5	C	903	GLY	OXT-C	-2.22	1.23	1.30
5	A	903	GLY	OXT-C	-2.19	1.23	1.30
7	B	903[A]	JC9	C8-CL1	2.18	1.78	1.73
7	B	903[C]	JC9	C9-C7	-2.15	1.36	1.39
7	B	903[C]	JC9	C2-C3	-2.10	1.47	1.52

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	903[B]	JC9	C9-C7-C1	-7.02	113.25	120.56
7	B	903[C]	JC9	C7-C8-CL1	-6.59	113.74	121.70
7	B	903[C]	JC9	C1-C7-C8	-5.95	117.97	123.43
4	B	901	NAG	C1-O5-C5	5.83	119.99	112.19
7	B	903[B]	JC9	C1-C7-C8	5.56	128.52	123.43
7	B	903[B]	JC9	C7-C8-CL1	4.82	127.52	121.70
7	B	903[A]	JC9	C9-C7-C1	-4.61	115.76	120.56
7	B	903[C]	JC9	C9-C7-C1	4.11	124.84	120.56
7	B	903[B]	JC9	C11-C8-CL1	-3.93	110.69	118.42
7	B	903[A]	JC9	C2-C3-C4	3.65	117.04	111.35
4	D	901	NAG	C1-O5-C5	2.82	115.97	112.19
6	D	903	GLU	OXT-C-O	-2.73	117.88	124.08
6	B	902	GLU	OXT-C-O	-2.67	118.01	124.08
7	B	903[B]	JC9	O1-C5-C6	2.61	125.94	121.51
7	B	903[C]	JC9	C11-C8-C7	2.56	124.93	121.73
7	B	903[A]	JC9	C3-C2-C1	2.48	117.24	112.24
7	B	903[A]	JC9	C1-C7-C8	2.26	125.50	123.43
7	B	903[C]	JC9	C12-C13-C11	-2.26	117.45	120.24
7	B	903[C]	JC9	C3-C2-C1	2.25	116.78	112.24
7	B	903[A]	JC9	C9-C7-C8	2.10	118.74	116.51
5	A	903	GLY	OXT-C-O	-2.07	118.00	123.33
5	C	903	GLY	OXT-C-O	-2.06	118.03	123.33

There are no chirality outliers.

All (17) torsion outliers are listed below:

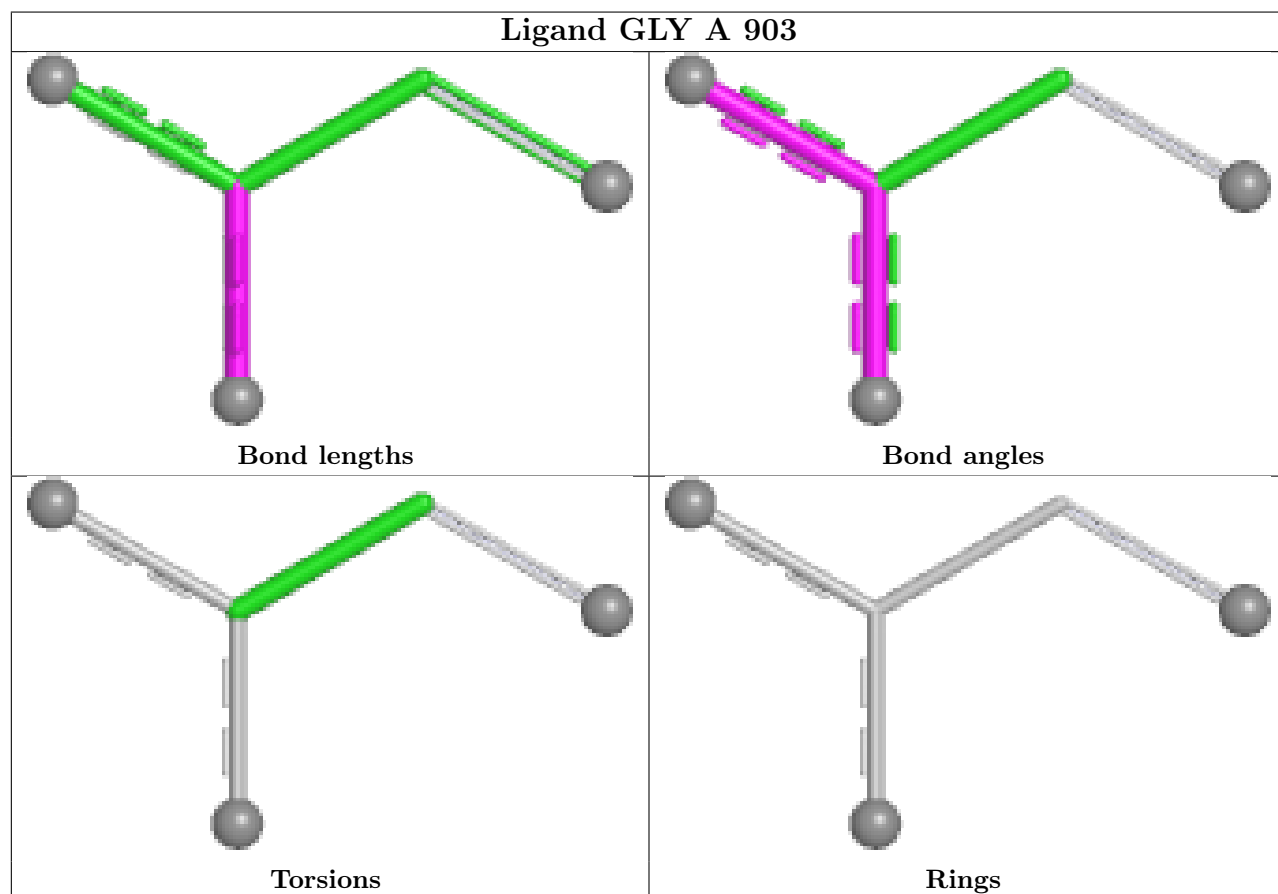
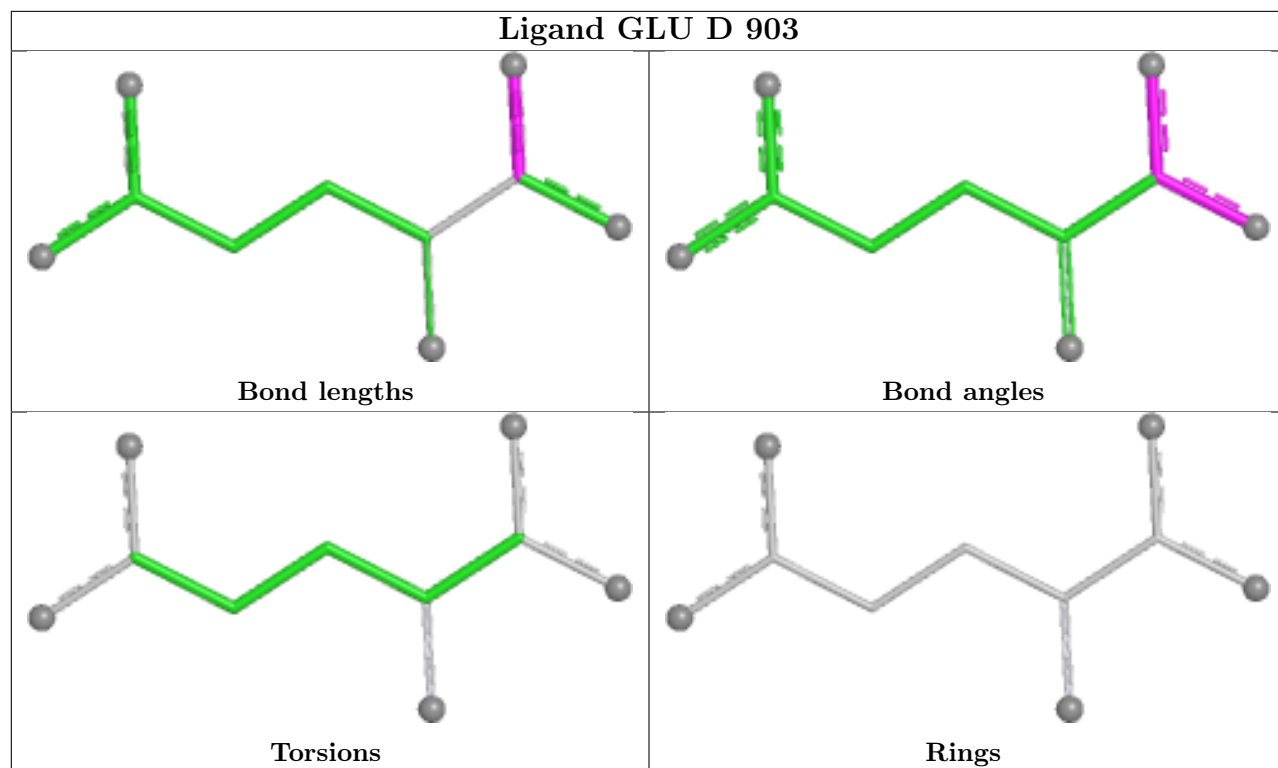
Mol	Chain	Res	Type	Atoms
7	B	903[B]	JC9	C5-C1-C7-C8
7	B	903[B]	JC9	C5-C1-C7-C9
7	B	903[C]	JC9	N1-C1-C7-C8
7	B	903[C]	JC9	C5-C1-C7-C8
7	B	903[C]	JC9	C5-C1-C7-C9
4	A	901	NAG	O5-C5-C6-O6
4	A	902	NAG	O5-C5-C6-O6
4	A	902	NAG	C4-C5-C6-O6
4	C	901	NAG	O5-C5-C6-O6
4	A	901	NAG	C4-C5-C6-O6
4	C	901	NAG	C4-C5-C6-O6
4	D	902	NAG	C4-C5-C6-O6
4	D	902	NAG	O5-C5-C6-O6
4	A	902	NAG	C3-C2-N2-C7
4	A	902	NAG	C1-C2-N2-C7
7	B	903[B]	JC9	N1-C1-C7-C8
7	B	903[B]	JC9	N1-C1-C7-C9

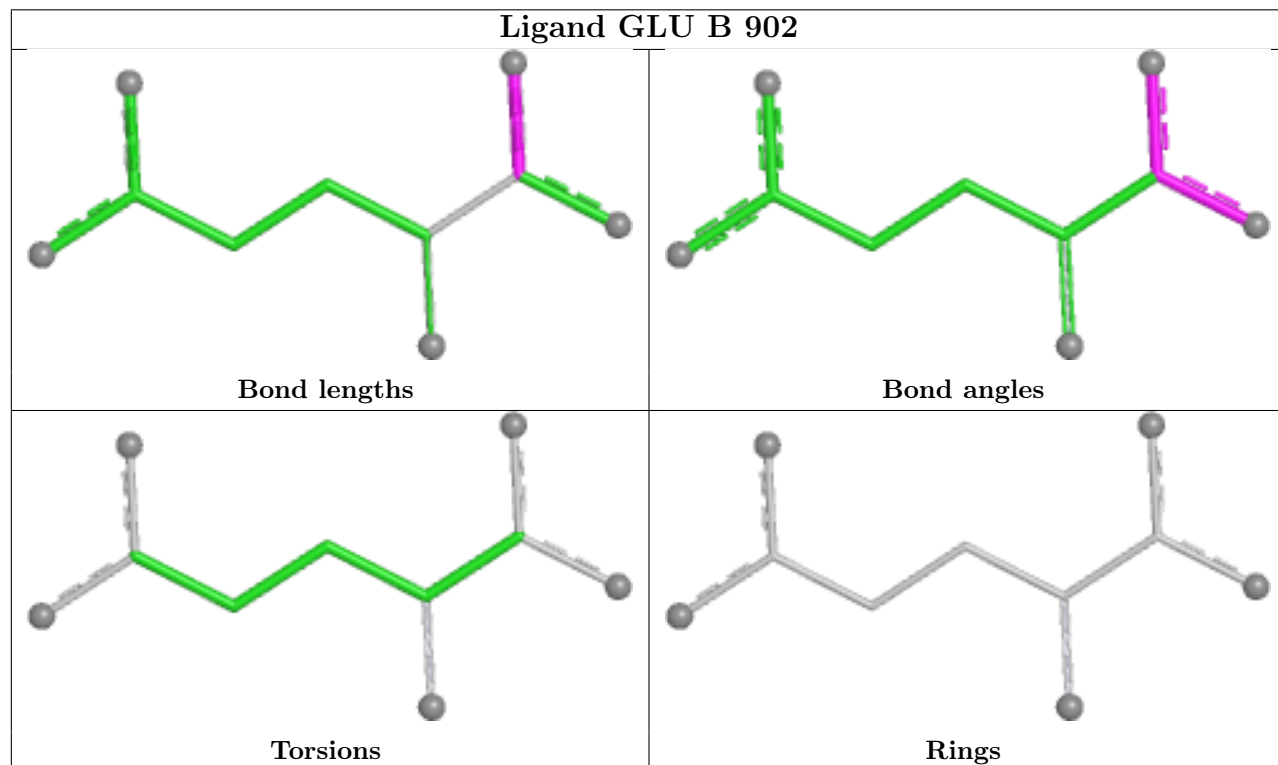
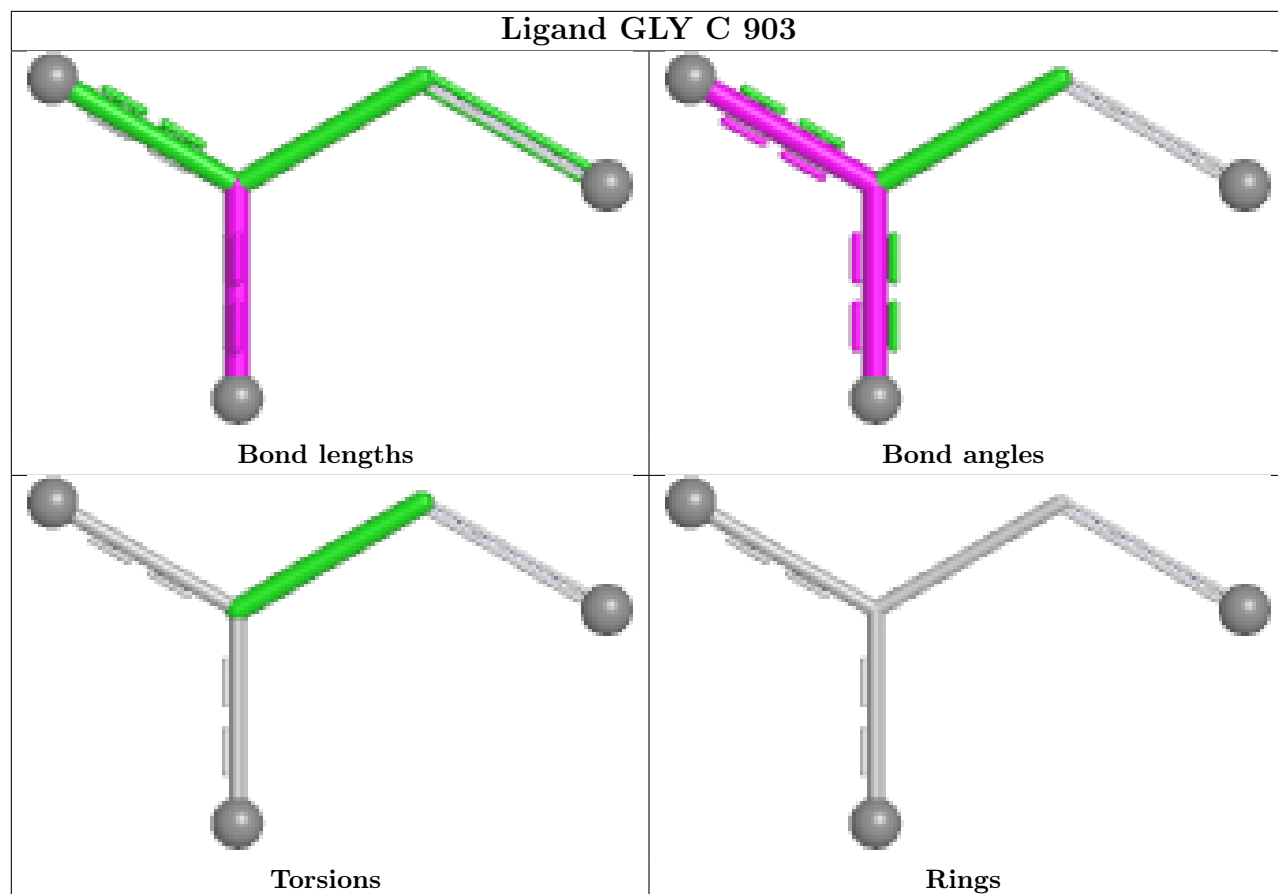
There are no ring outliers.

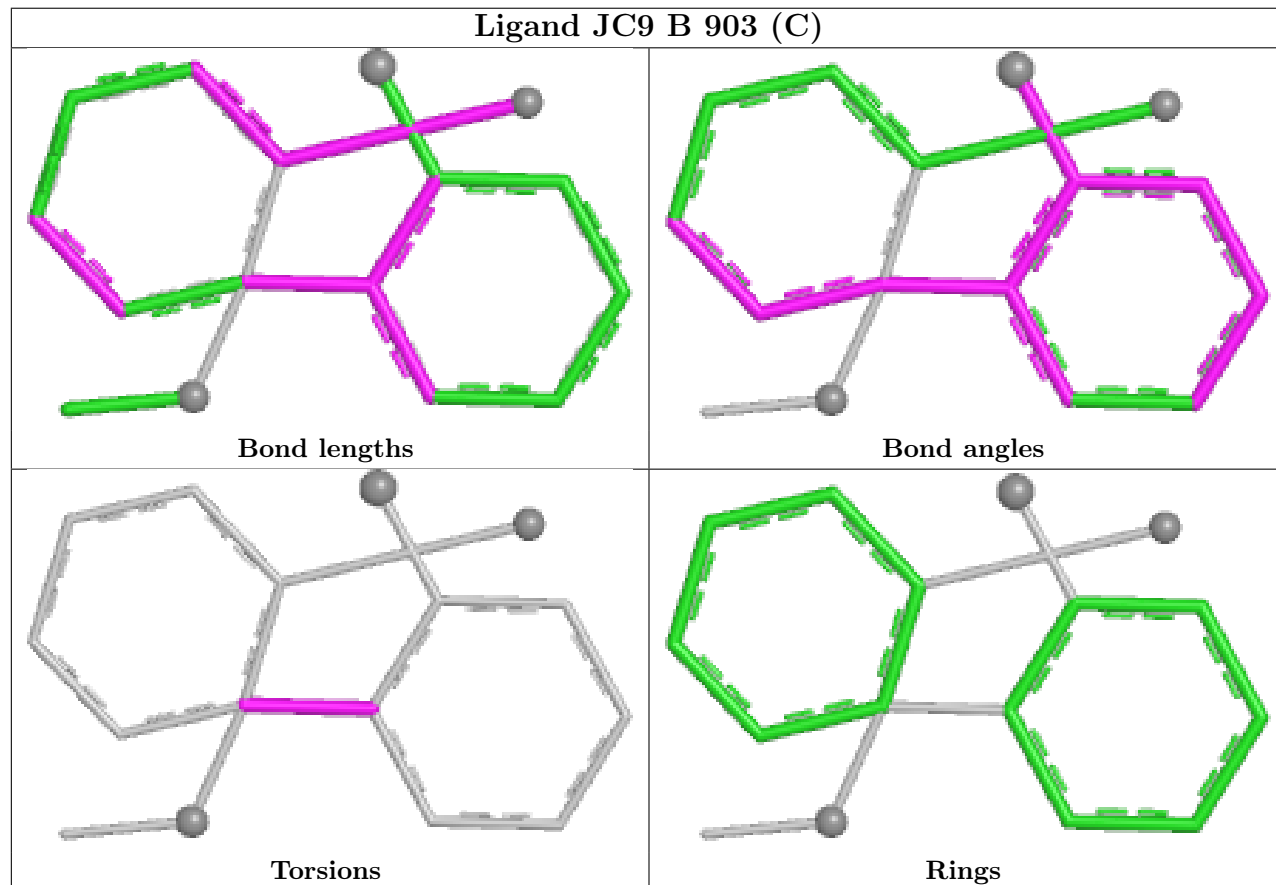
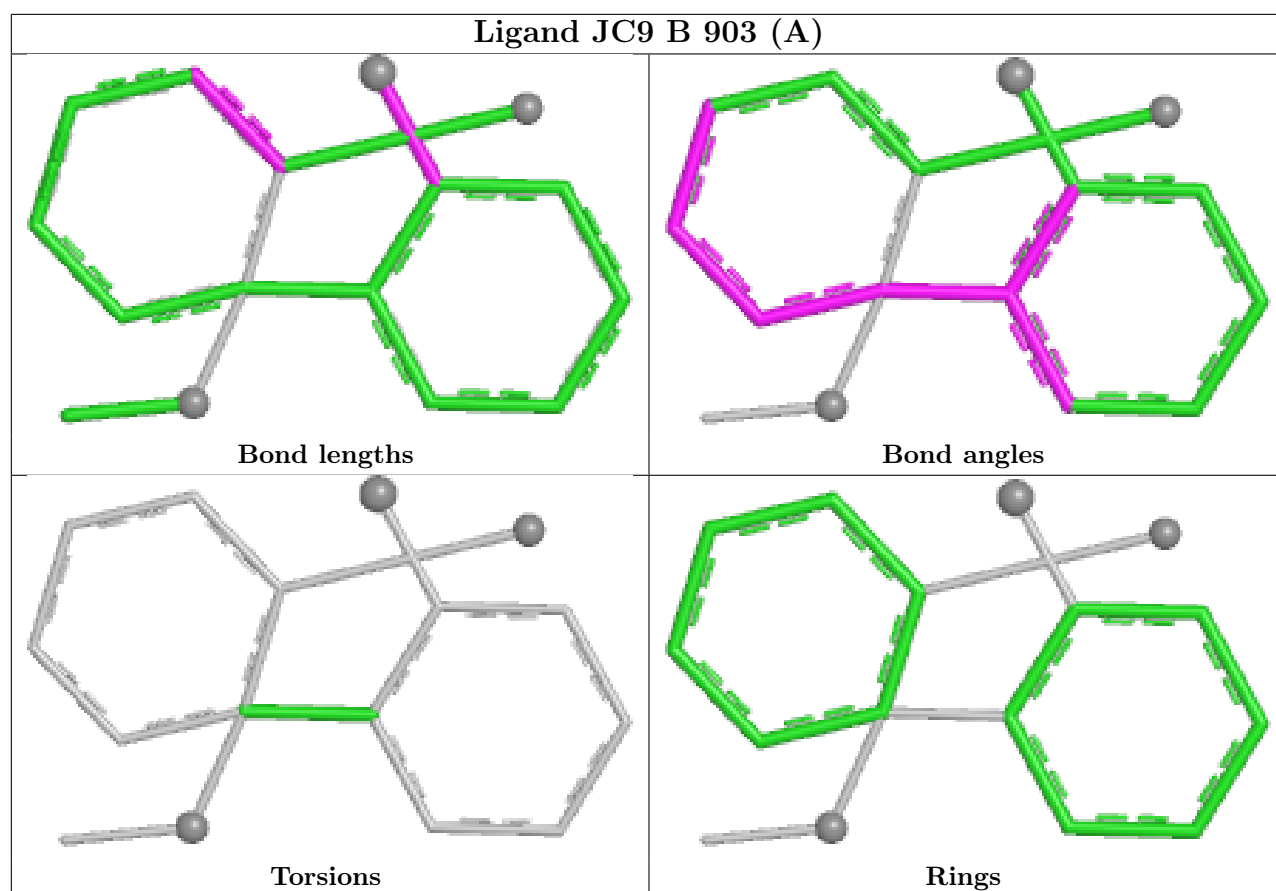
4 monomers are involved in 7 short contacts:

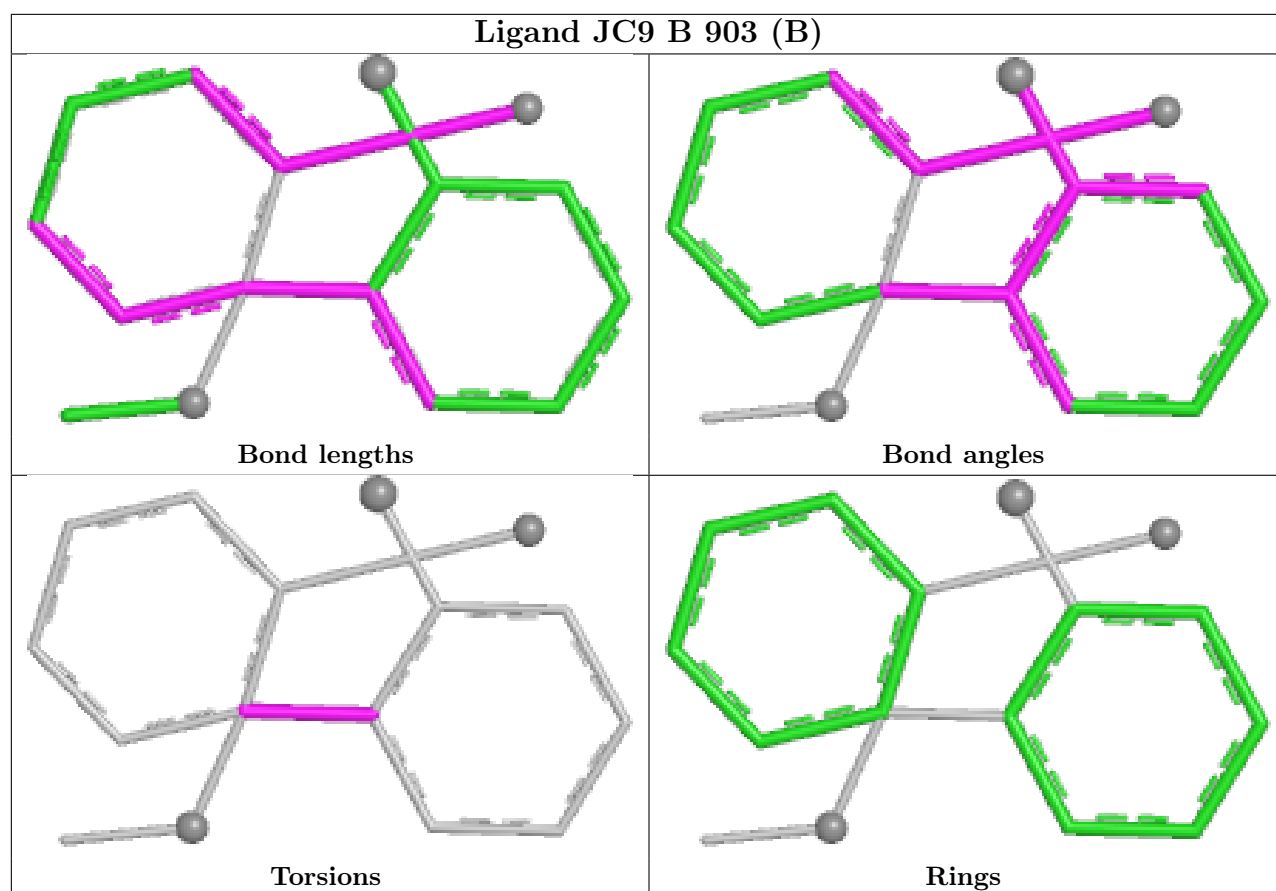
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	903	GLU	2	0
7	B	903[A]	JC9	1	0
7	B	903[C]	JC9	3	0
7	B	903[B]	JC9	1	0

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.

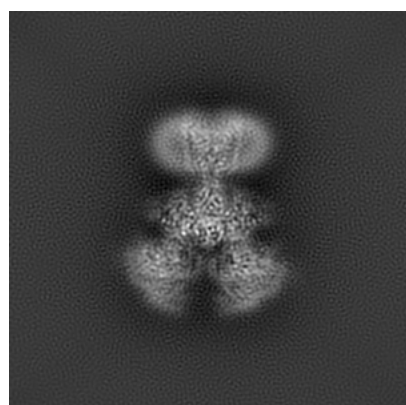
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24948. These allow visual inspection of the internal detail of the map and identification of artifacts.

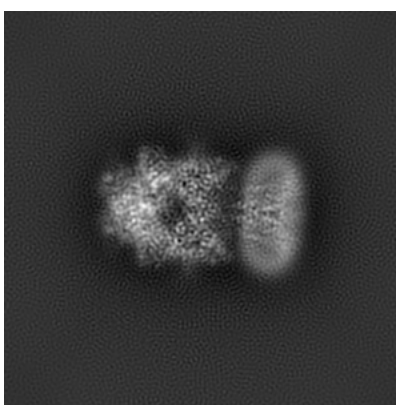
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

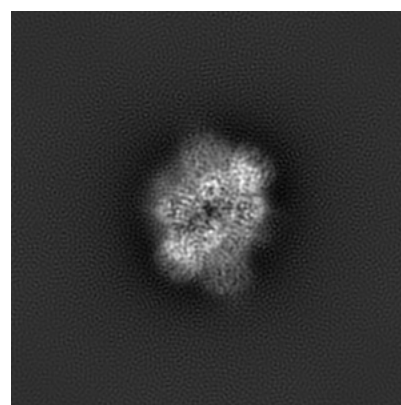
6.1.1 Primary map



X



Y

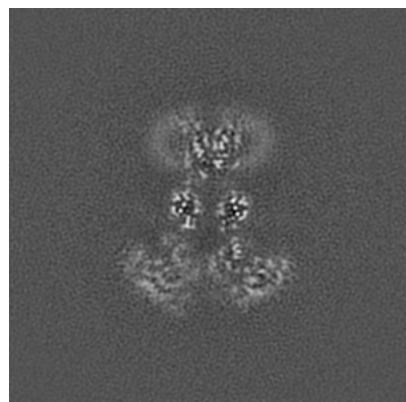


Z

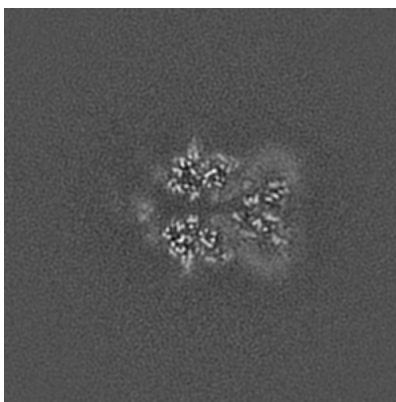
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

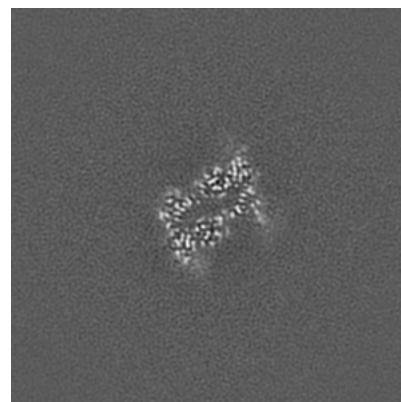
6.2.1 Primary map



X Index: 200



Y Index: 200

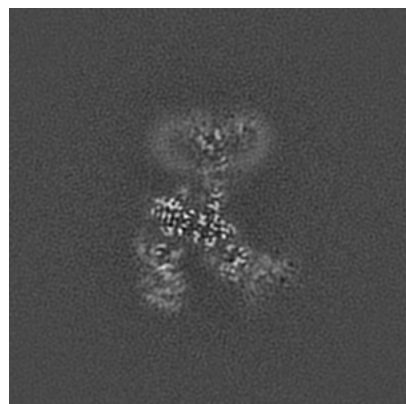


Z Index: 200

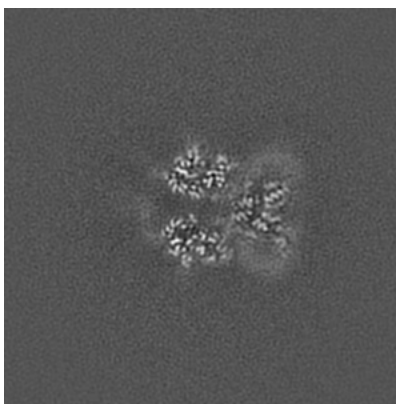
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

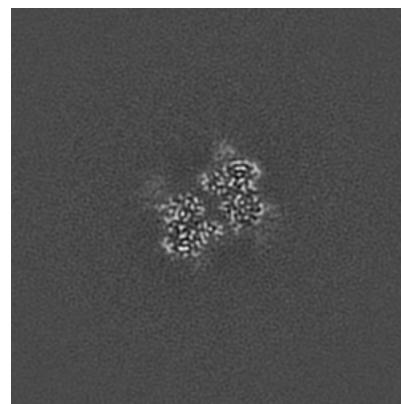
6.3.1 Primary map



X Index: 177



Y Index: 198

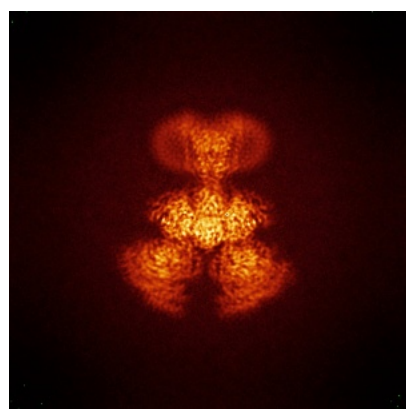


Z Index: 187

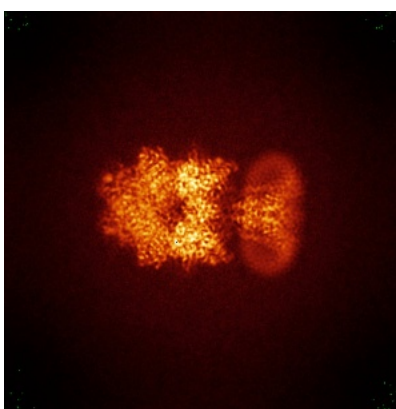
The images above show the largest variance slices of the map in three orthogonal directions.

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

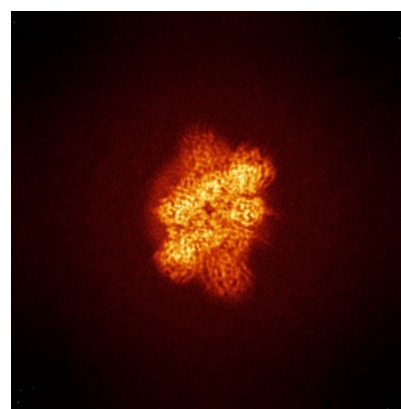
6.4.1 Primary map



X



Y

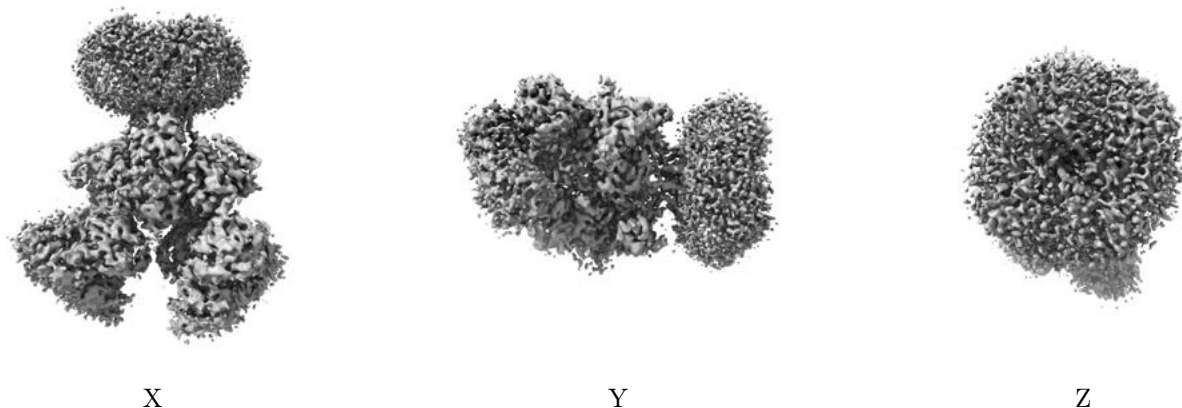


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 1.6. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

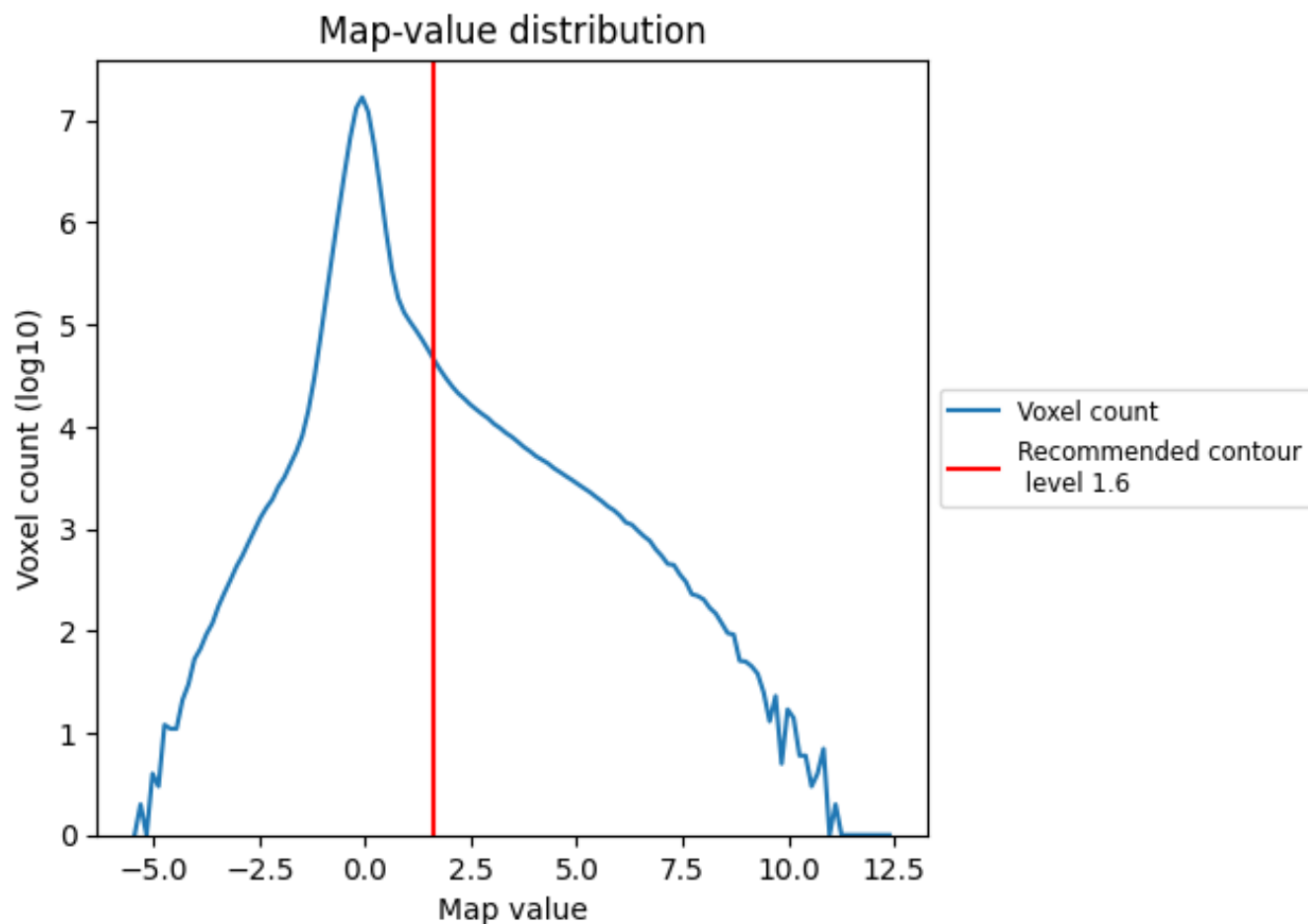
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

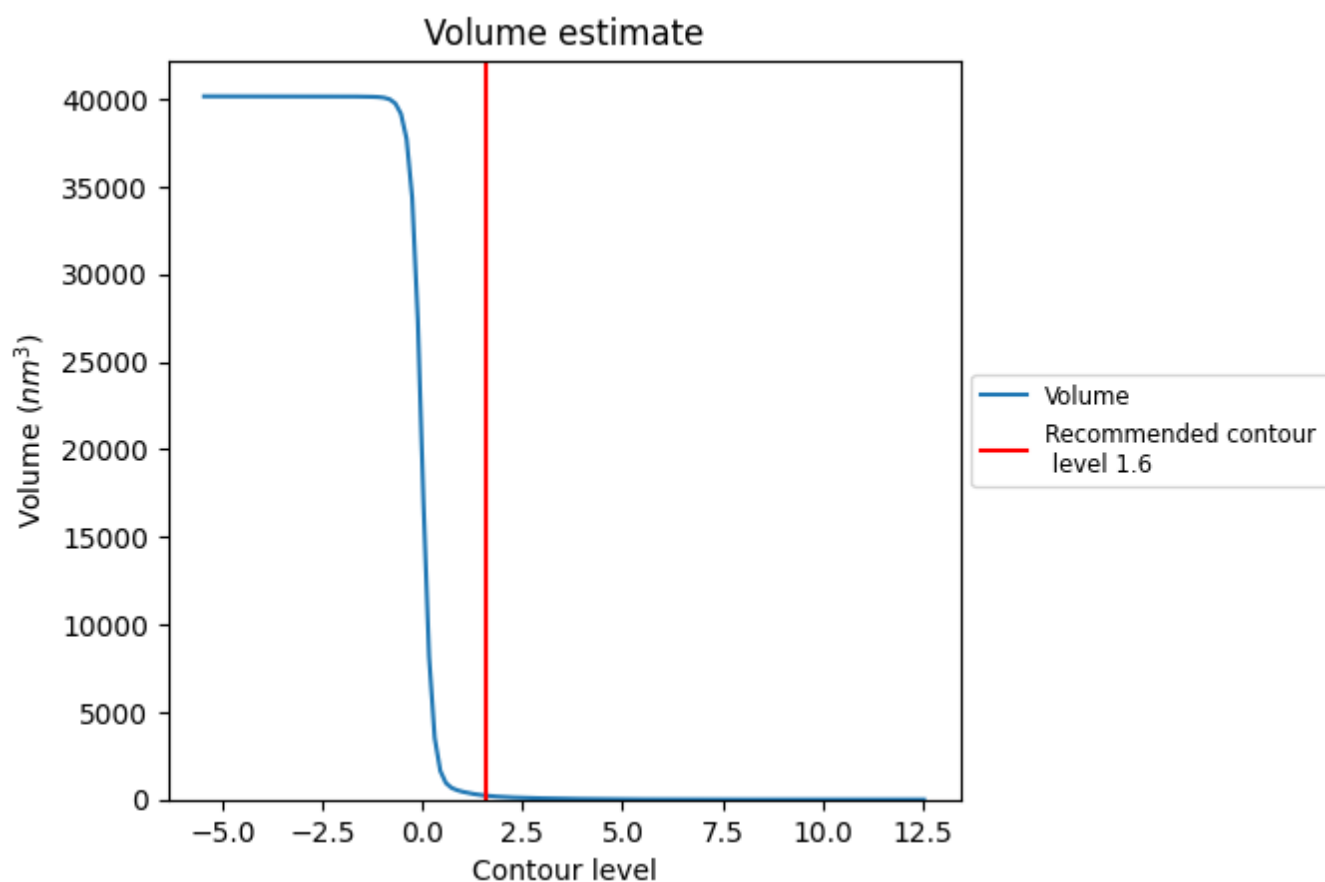
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

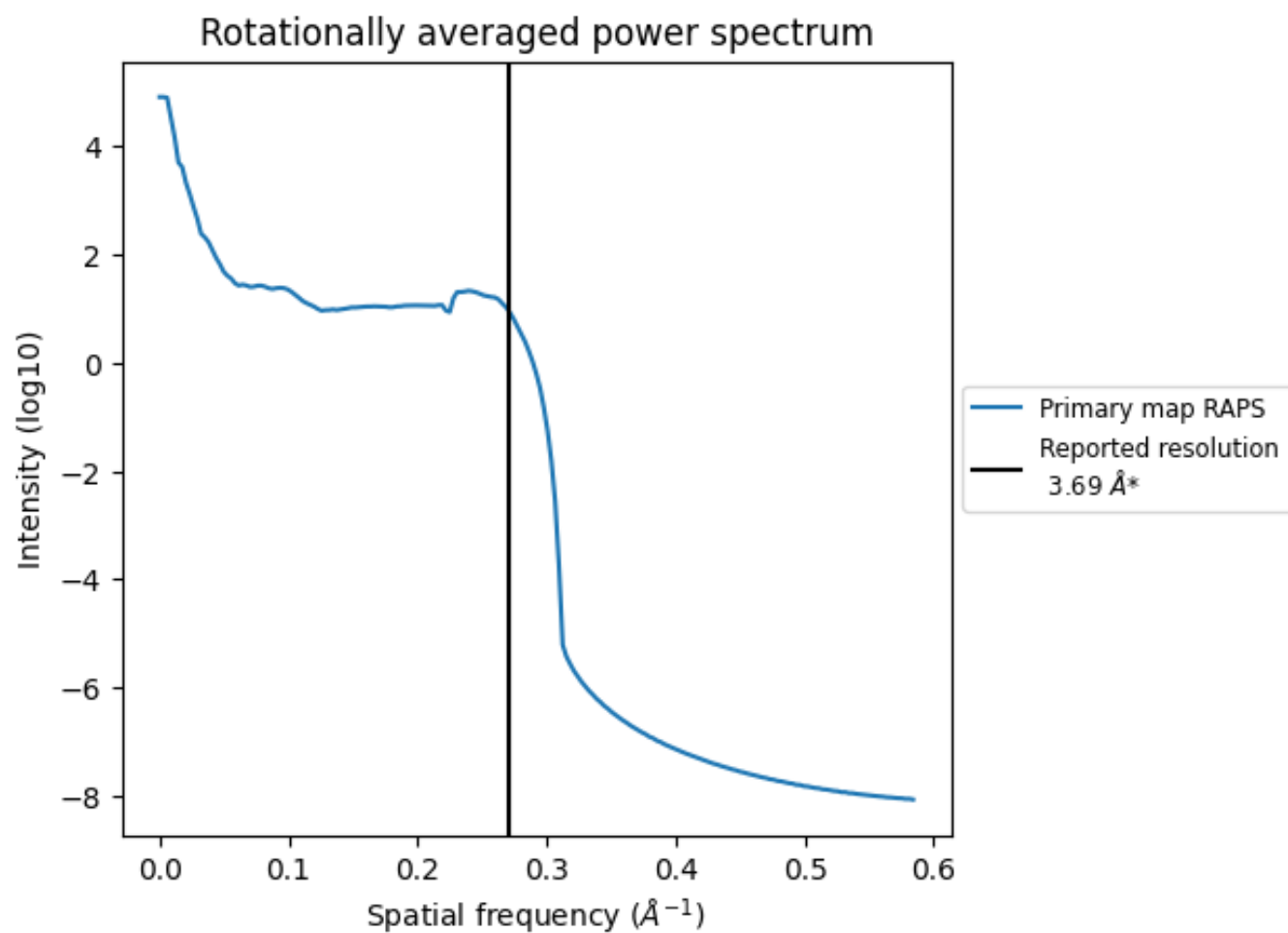
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 222 nm^3 ; this corresponds to an approximate mass of 201 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 0.271 Å⁻¹

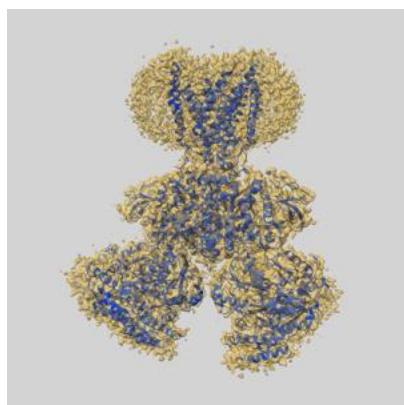
8 Fourier-Shell correlation

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

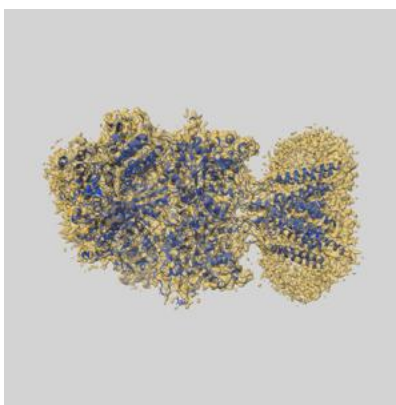
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-24948 and PDB model 7SAC. Per-residue inclusion information can be found in section [3](#) on page [11](#).

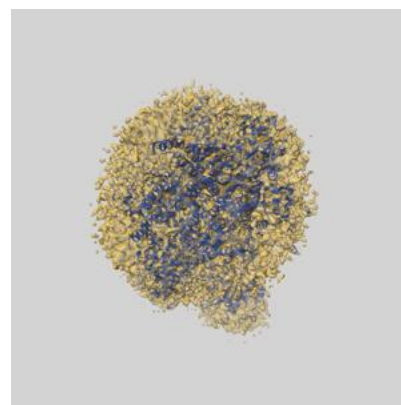
9.1 Map-model overlay [i](#)



X



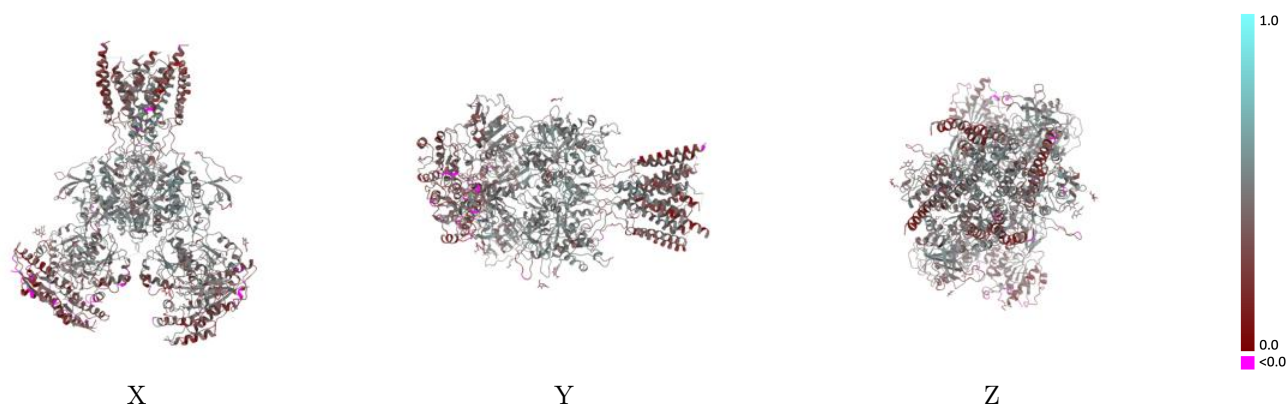
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 1.6 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](#)

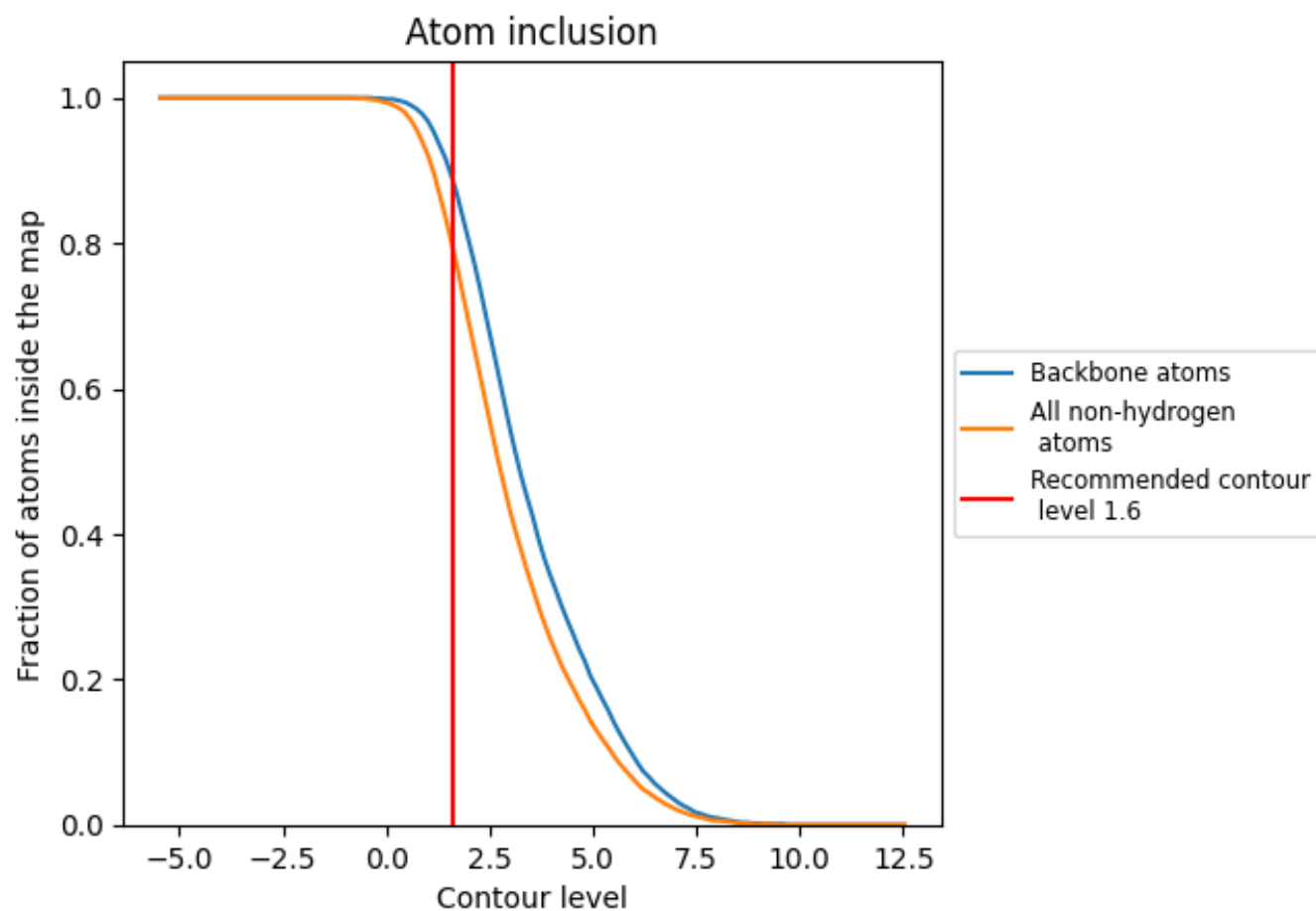


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 89% of all backbone atoms, 80% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (1.6) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7960	0.4190
A	0.8190	0.4340
B	0.7900	0.4220
C	0.7960	0.4120
D	0.7840	0.4100
E	0.5360	0.3680
F	0.3930	0.1800
G	0.6790	0.3960

1.0

0.0

<0.0