



Full wwPDB EM Validation Report ⓘ

Mar 10, 2026 – 10:17 AM UTC

PDB ID : 7VWJ / pdb_00007vwj
EMDB ID : EMD-32154
Title : Matrix arm of deactive state CI from rotenone-NADH dataset
Authors : Gu, J.K.; Yang, M.J.
Deposited on : 2021-11-10
Resolution : 2.60 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

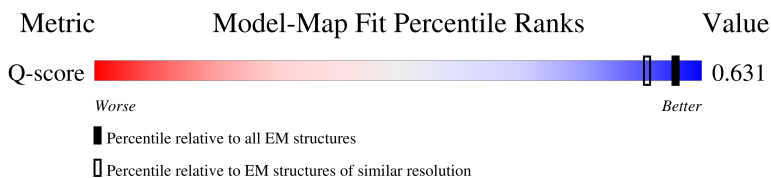
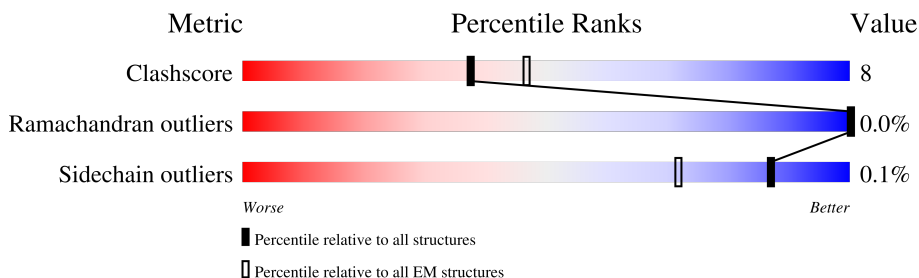
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.60 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	8728 (2.10 - 3.10)

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	431	8% 79% 21%
2	B	176	79% 20%
3	C	156	5% 87% 13%
4	E	115	11% 88% 12%

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Mol	Chain	Length	Quality of chain
5	F	86	
6	G	88	
7	H	112	
8	I	112	
9	J	341	
10	K	42	
11	L	125	
12	M	690	
13	N	144	
14	O	217	
15	P	208	
16	Q	385	
17	T	96	
18	W	29	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
19	SF4	C	301	-	-	X	-

2 Entry composition [i](#)

There are 30 unique types of molecules in this entry. The entry contains 28748 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 NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	431	Total	C	N	O	S	0	0
			3318	2095	591	612	20		

- Molecule 2 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	176	Total	C	N	O	S	0	0
			1412	887	243	269	13		

- Molecule 3 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	156	Total	C	N	O	S	0	0
			1248	794	227	213	14		

- Molecule 4 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	E	115	Total	C	N	O	S	0	0
			971	619	179	168	5		

- Molecule 5 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	F	86	Total	C	N	O	S	0	0
			691	434	129	126	2		

- Molecule 6 is a protein called Acyl carrier protein, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	G	88	Total	C	N	O	S	0	0
			693	447	102	139	5		

- Molecule 7 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	H	112	Total	C	N	O	S	0	0
			910	588	154	165	3		

- Molecule 8 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	I	97	Total	C	N	O	S	0	0
			780	491	147	139	3		

- Molecule 9 is a protein called NADH dehydrogenase ubiquinone 1 alpha subcomplex subunit 9, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	J	297	Total	C	N	O	S	0	0
			2359	1514	421	416	8		

- Molecule 10 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	K	42	Total	C	N	O	S	0	0
			355	219	67	68	1		

- Molecule 11 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	L	125	Total	C	N	O	S	0	0
			1016	642	181	190	3		

- Molecule 12 is a protein called NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	M	690	Total	C	N	O	S	0	0
			5296	3320	923	1014	39		

- Molecule 13 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	N	144	Total	C	N	O	S	0	0
			1204	770	218	212	4		

- Molecule 14 is a protein called NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	O	217	Total	C	N	O	S	0	0
			1671	1065	281	315	10		

- Molecule 15 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	P	208	Total	C	N	O	S	0	0
			1738	1124	298	314	2		

- Molecule 16 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	Q	379	Total	C	N	O	S	0	0
			3044	1945	522	554	23		

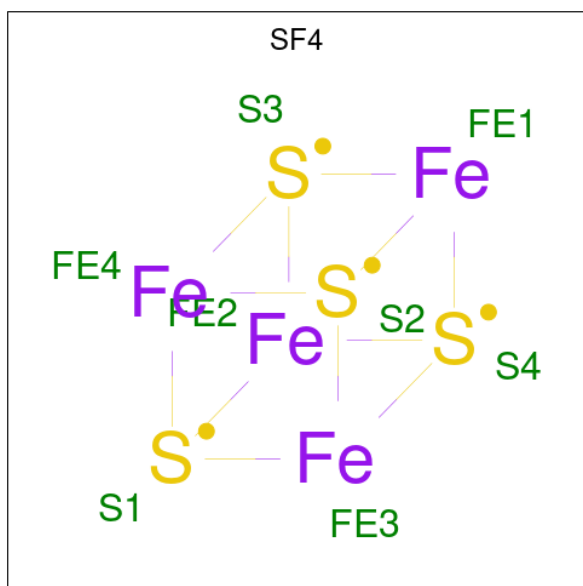
- Molecule 17 is a protein called NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	T	96	Total	C	N	O	S	0	0
			741	452	140	146	3		

- Molecule 18 is a protein called NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 13.

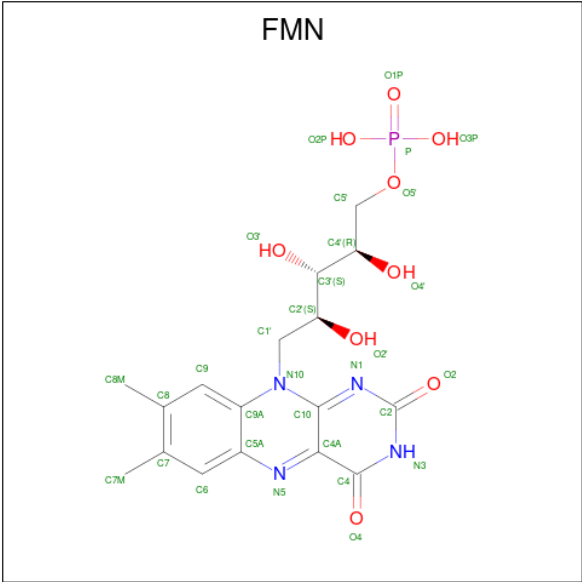
Mol	Chain	Residues	Atoms					AltConf	Trace
18	W	29	Total	C	N	O	S	0	0
			224	141	43	39	1		

- Molecule 19 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4) (labeled as "Ligand of Interest" by depositor).



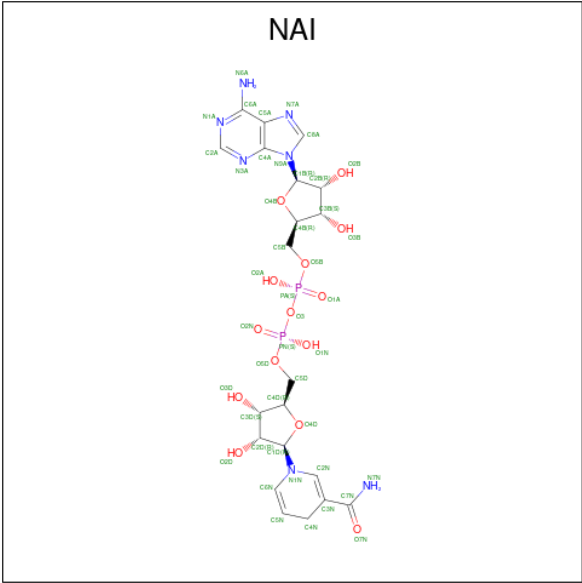
Mol	Chain	Residues	Atoms			AltConf
19	A	1	Total	Fe	S	0
			8	4	4	
19	B	1	Total	Fe	S	0
			8	4	4	
19	B	1	Total	Fe	S	0
			8	4	4	
19	C	1	Total	Fe	S	0
			8	4	4	
19	M	1	Total	Fe	S	0
			8	4	4	
19	M	1	Total	Fe	S	0
			8	4	4	

- Molecule 20 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $\text{C}_{17}\text{H}_{21}\text{N}_4\text{O}_9\text{P}$) (labeled as "Ligand of Interest" by depositor).



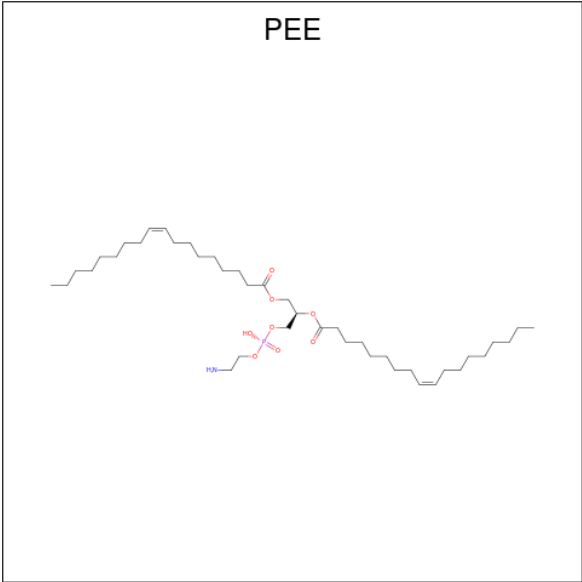
Mol	Chain	Residues	Atoms					AltConf
20	A	1	Total	C	N	O	P	0
			31	17	4	9	1	

- Molecule 21 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).



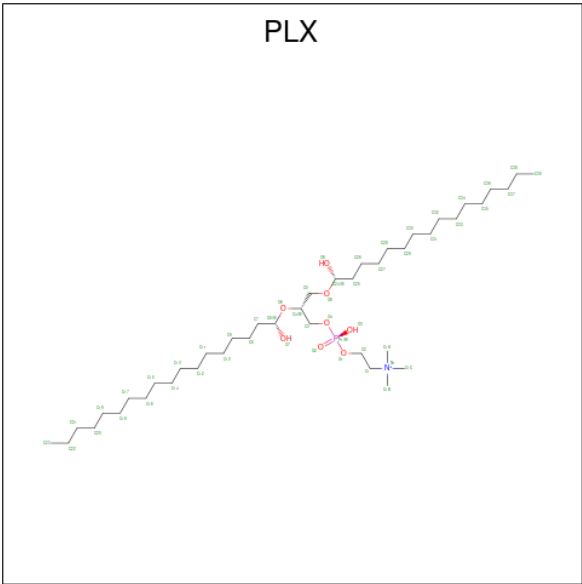
Mol	Chain	Residues	Atoms					AltConf
21	A	1	Total	C	N	O	P	0
			44	21	7	14	2	

- Molecule 22 is 1,2-dioleoyl-sn-glycero-3-phosphoethanolamine (CCD ID: PEE) (formula: $C_{41}H_{78}NO_8P$) (labeled as "Ligand of Interest" by depositor).



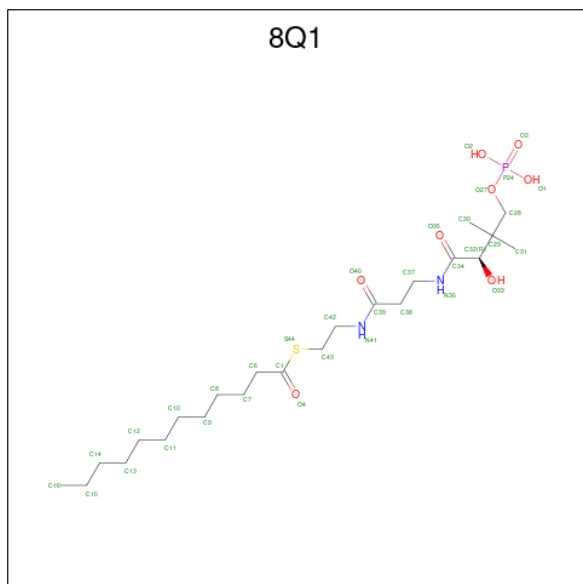
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
22	C	1	47	37	1	8	1	0

- Molecule 23 is (9R,11S)-9-([[(1S)-1-HYDROXYHEXADECYL]OXY}METHYL)-2,2-DIMETHYL-5,7,10-TRIOXA-2LAMBDA 5 -AZA-6LAMBDA 5 -PHOSPHAOCTACOSANE-6,6,11-TRIOL (CCD ID: PLX) (formula: C₄₂H₈₉NO₈P) (labeled as "Ligand of Interest" by depositor).



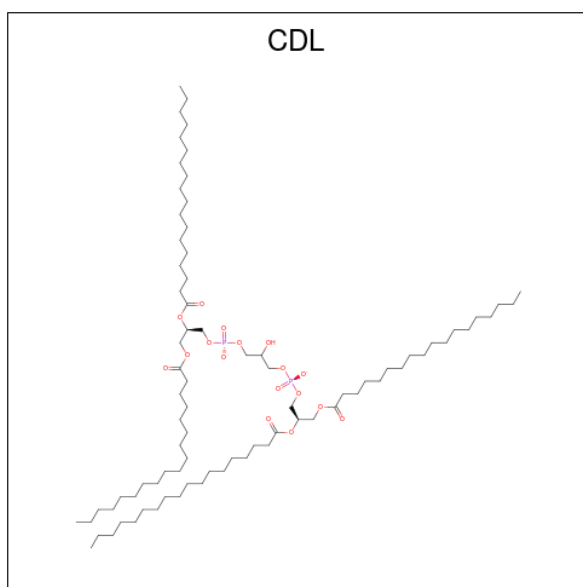
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
23	C	1	52	42	1	8	1	0

- Molecule 24 is S-[2-({N-[(2R)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-beta-alanyl}amino)ethyl] dodecanethioate (CCD ID: 8Q1) (formula: $C_{23}H_{45}N_2O_8PS$) (labeled as "Ligand of Interest" by depositor).



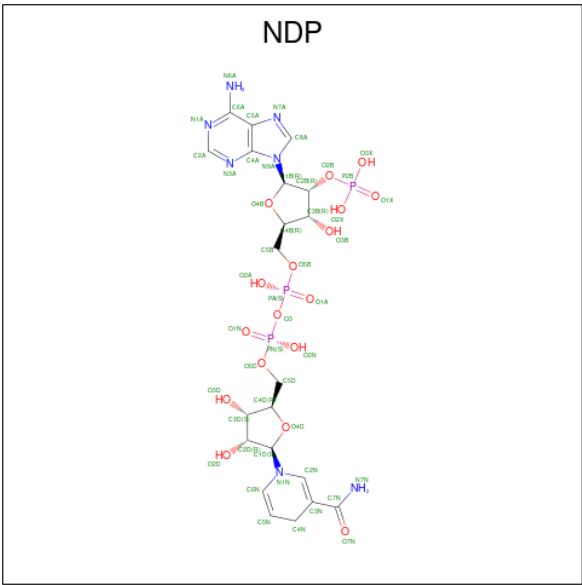
Mol	Chain	Residues	Atoms						AltConf
			Total	C	N	O	P	S	
24	G	1	35	23	2	8	1	1	0

- Molecule 25 is CARDIOLIPIN (CCD ID: CDL) (formula: $C_{81}H_{156}O_{17}P_2$) (labeled as "Ligand of Interest" by depositor).



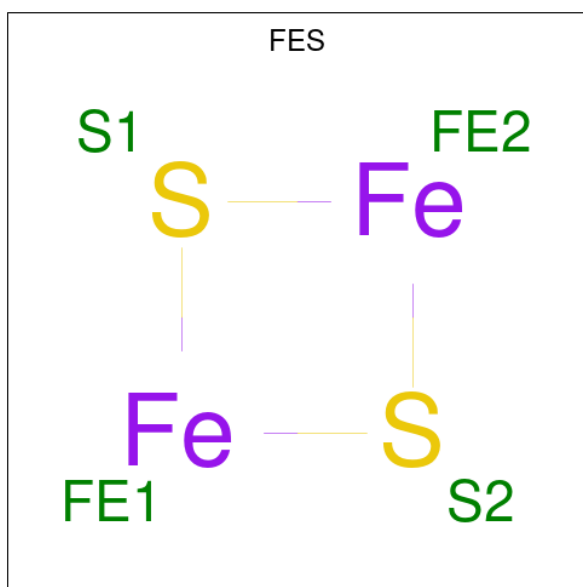
Mol	Chain	Residues	Atoms				AltConf
25	I	1	Total	C	O	P	0
			51	32	17	2	

- Molecule 26 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
26	J	1	Total	C	N	O	P	0
			48	21	7	17	3	

- Molecule 27 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe₂S₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
27	M	1	Total	Fe	S	0
			4	2	2	
27	O	1	Total	Fe	S	0
			4	2	2	

- Molecule 28 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
28	M	1	Total	Mg	0
			1	1	

- Molecule 29 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
29	T	1	Total	Zn	0
			1	1	

- Molecule 30 is water.

Mol	Chain	Residues	Atoms		AltConf
30	A	47	Total	O	0
			47	47	
30	B	76	Total	O	0
			76	76	

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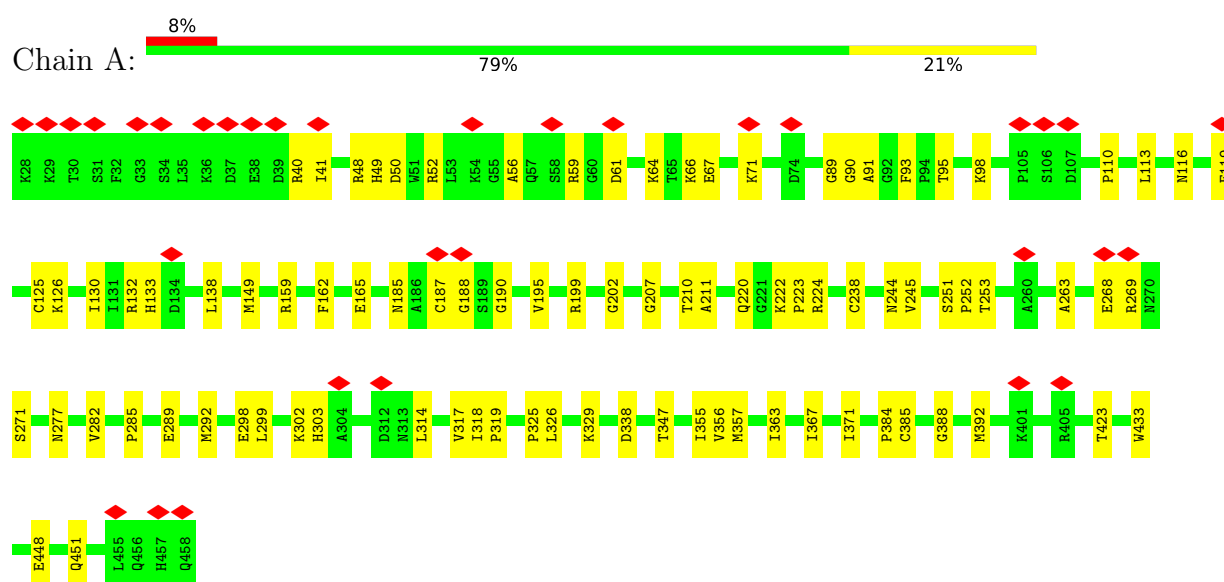
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Mol	Chain	Residues	Atoms		AltConf
30	C	54	Total 54	O 54	0
30	E	3	Total 3	O 3	0
30	F	1	Total 1	O 1	0
30	H	4	Total 4	O 4	0
30	I	12	Total 12	O 12	0
30	J	8	Total 8	O 8	0
30	K	4	Total 4	O 4	0
30	L	22	Total 22	O 22	0
30	M	189	Total 189	O 189	0
30	N	6	Total 6	O 6	0
30	O	16	Total 16	O 16	0
30	P	84	Total 84	O 84	0
30	Q	175	Total 175	O 175	0
30	T	9	Total 9	O 9	0
30	W	1	Total 1	O 1	0

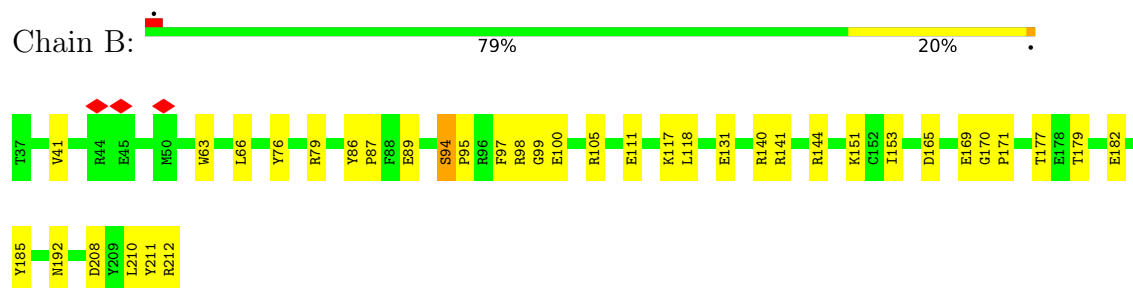
3 Residue-property plots

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

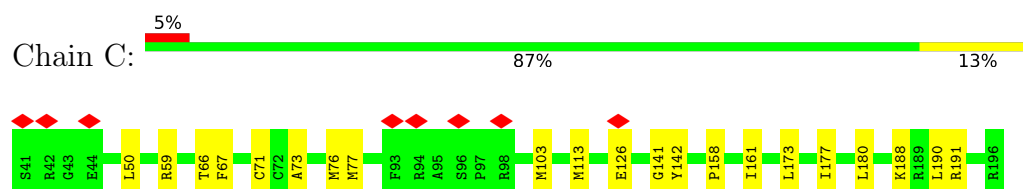
- Molecule 1: NADH dehydrogenase [ubiquinone] flavoprotein 1, mitochondrial



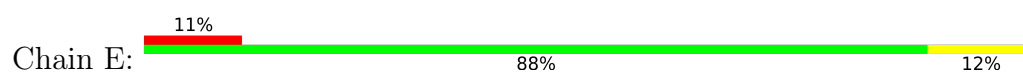
- Molecule 2: NADH dehydrogenase [ubiquinone] iron-sulfur protein 8, mitochondrial



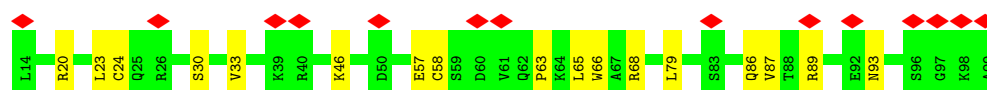
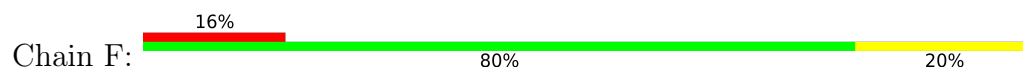
- Molecule 3: NADH dehydrogenase [ubiquinone] iron-sulfur protein 7, mitochondrial



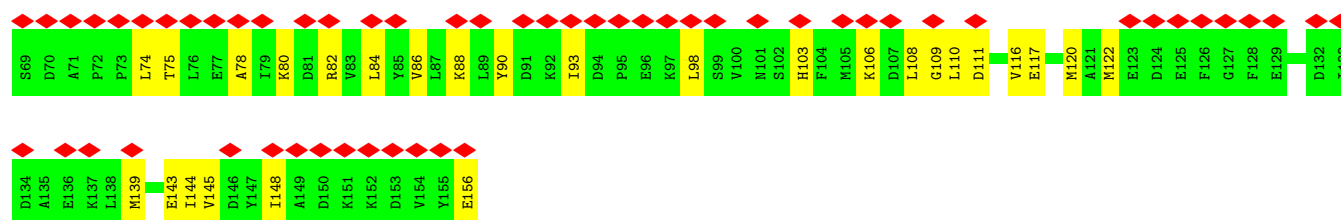
- Molecule 4: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 6



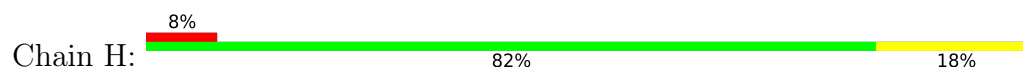
- Molecule 5: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 2



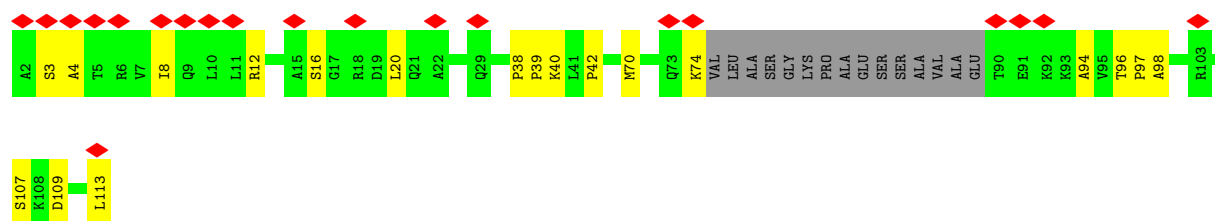
- Molecule 6: Acyl carrier protein, mitochondrial



- Molecule 7: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 5

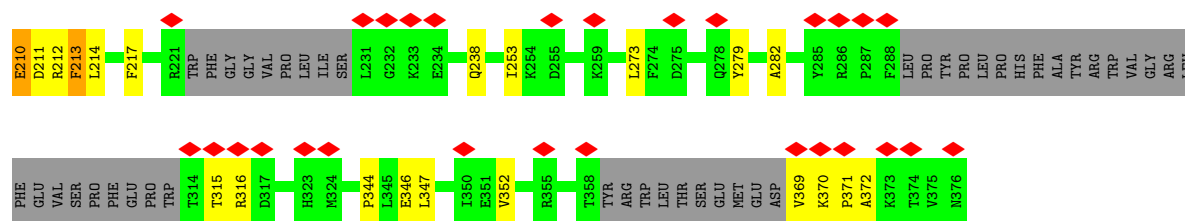


- Molecule 8: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 7

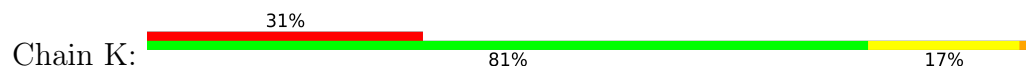


- Molecule 9: NADH dehydrogenase ubiquinone 1 alpha subcomplex subunit 9, mitochondrial

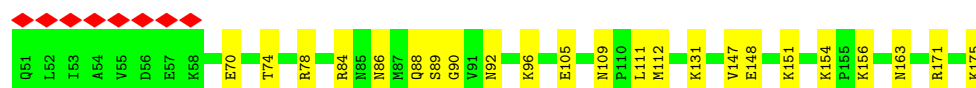
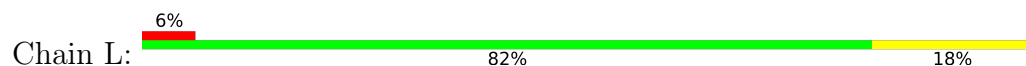




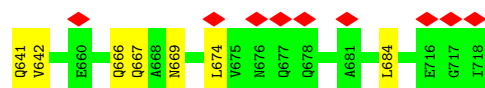
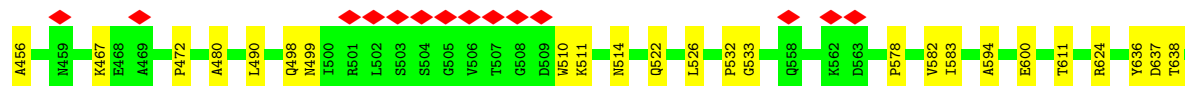
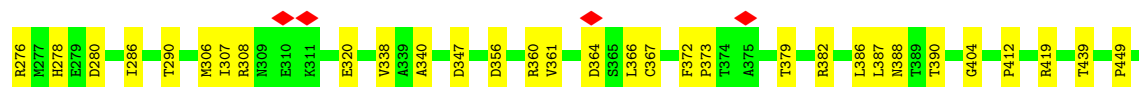
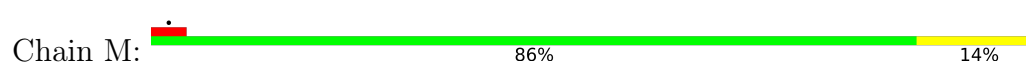
- Molecule 10: NADH dehydrogenase [ubiquinone] flavoprotein 3, mitochondrial



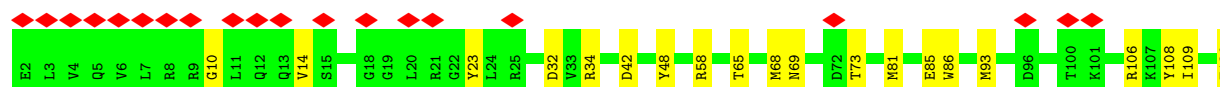
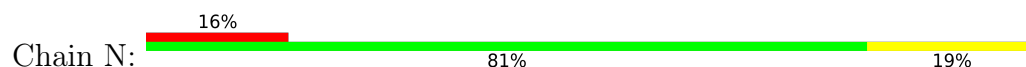
- Molecule 11: NADH dehydrogenase [ubiquinone] iron-sulfur protein 4, mitochondrial



- Molecule 12: NADH-ubiquinone oxidoreductase 75 kDa subunit, mitochondrial

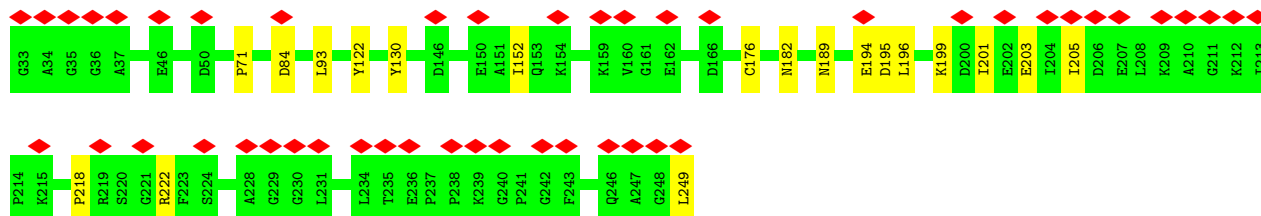
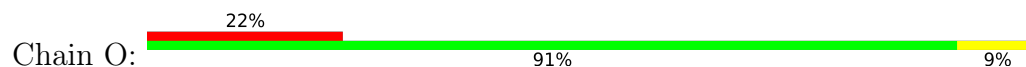


- Molecule 13: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 12

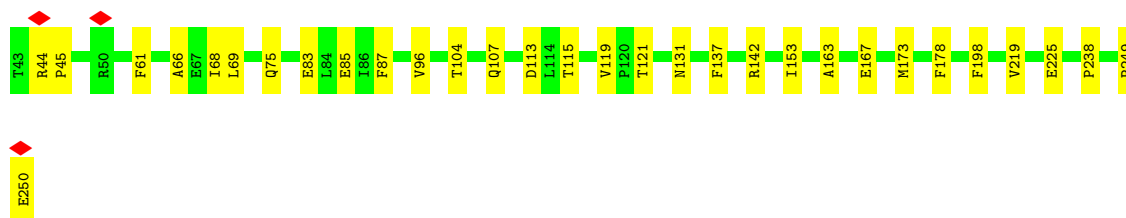
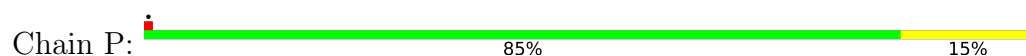




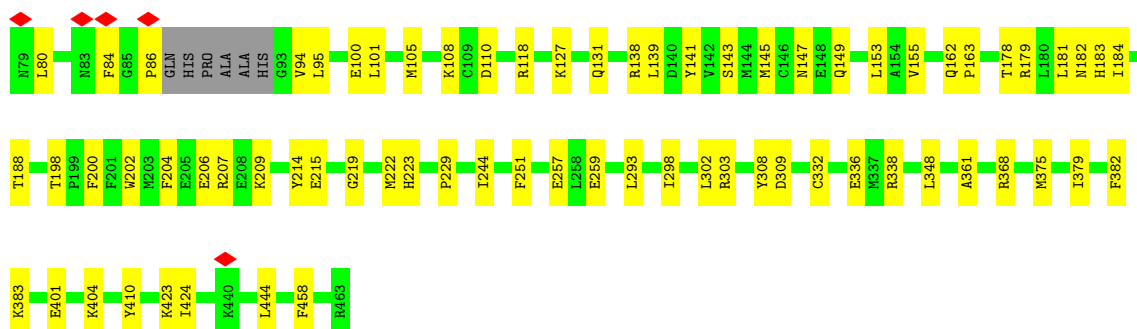
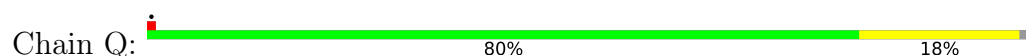
- Molecule 14: NADH dehydrogenase [ubiquinone] flavoprotein 2, mitochondrial



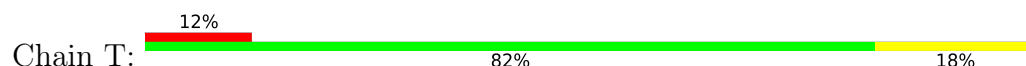
- Molecule 15: NADH dehydrogenase [ubiquinone] iron-sulfur protein 3, mitochondrial



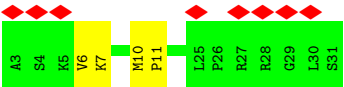
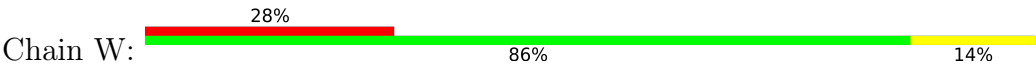
- Molecule 16: NADH dehydrogenase [ubiquinone] iron-sulfur protein 2, mitochondrial



- Molecule 17: NADH dehydrogenase [ubiquinone] iron-sulfur protein 6, mitochondrial



- Molecule 18: NADH dehydrogenase [ubiquinone] 1 alpha subcomplex subunit 13



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	169835	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1300	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.090	Depositor
Minimum map value	-0.047	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.0165	Depositor
Map size (Å)	274.9952, 274.9952, 274.9952	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.5371, 0.5371, 0.5371	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, CDL, SF4, NDP, FES, NAI, 2MR, 8Q1, MG, PEE, PLX, FMN

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	0.13	0/3393	0.32	0/4584
2	B	0.15	0/1443	0.39	2/1952 (0.1%)
3	C	0.14	0/1279	0.34	0/1730
4	E	0.13	0/995	0.29	0/1340
5	F	0.13	0/702	0.37	0/945
6	G	0.12	0/705	0.32	0/956
7	H	0.13	0/929	0.33	0/1258
8	I	0.18	0/798	0.42	0/1079
9	J	0.15	0/2411	0.37	1/3254 (0.0%)
10	K	0.10	0/365	0.30	0/493
11	L	0.11	0/1039	0.29	0/1403
12	M	0.11	0/5384	0.29	0/7295
13	N	0.11	0/1245	0.30	0/1694
14	O	0.13	0/1711	0.34	0/2328
15	P	0.13	0/1789	0.32	0/2436
16	Q	0.14	0/3101	0.32	0/4189
17	T	0.11	0/755	0.26	0/1018
18	W	0.11	0/230	0.34	0/309
All	All	0.13	0/28274	0.33	3/38263 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	J	213	PHE	N-CA-C	5.97	117.46	111.07
2	B	94	SER	CA-C-N	5.09	124.70	119.05
2	B	94	SER	C-N-CA	5.09	124.70	119.05

There are no chirality outliers.

There are no planarity outliers.

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	3318	0	3280	62	0
2	B	1412	0	1363	36	0
3	C	1248	0	1254	27	0
4	E	971	0	975	9	0
5	F	691	0	704	11	0
6	G	693	0	671	19	0
7	H	910	0	950	13	0
8	I	780	0	808	21	0
9	J	2359	0	2402	48	0
10	K	355	0	329	11	0
11	L	1016	0	1016	21	0
12	M	5296	0	5326	63	0
13	N	1204	0	1162	19	0
14	O	1671	0	1673	13	0
15	P	1738	0	1693	22	0
16	Q	3044	0	3018	55	0
17	T	741	0	702	12	0
18	W	224	0	230	3	0
19	A	8	0	0	1	0
19	B	16	0	0	0	0
19	C	8	0	0	2	0
19	M	16	0	0	0	0
20	A	31	0	19	3	0
21	A	44	0	27	7	0
22	C	47	0	71	10	0
23	C	52	0	88	6	0
24	G	35	0	0	1	0
25	I	51	0	46	19	0
26	J	48	0	24	2	0
27	M	4	0	0	0	0
27	O	4	0	0	0	0
28	M	1	0	0	0	0
29	T	1	0	0	0	0
30	A	47	0	0	5	0
30	B	76	0	0	2	0
30	C	54	0	0	0	0
30	E	3	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
30	F	1	0	0	0	0
30	H	4	0	0	0	0
30	I	12	0	0	0	0
30	J	8	0	0	0	0
30	K	4	0	0	0	0
30	L	22	0	0	2	0
30	M	189	0	0	2	0
30	N	6	0	0	0	0
30	O	16	0	0	0	0
30	P	84	0	0	1	0
30	Q	175	0	0	1	0
30	T	9	0	0	0	0
30	W	1	0	0	0	0
All	All	28748	0	27831	421	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 (421) 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)
26:J:401:NDP:O4D	26:J:401:NDP:C4D	1.69	1.20
21:A:503:NAI:C1B	21:A:503:NAI:O4B	1.64	1.16
8:I:3:SER:O	25:I:201:CDL:HA22	1.49	1.11
3:C:50:LEU:CD2	22:C:302:PEE:H27	1.89	1.01
25:I:201:CDL:H532	25:I:201:CDL:CA7	1.93	0.99
2:B:97:PHE:HB2	16:Q:215:GLU:OE2	1.63	0.99
8:I:3:SER:O	25:I:201:CDL:CA2	2.15	0.94
9:J:64:PHE:CD1	9:J:210:GLU:HG2	2.03	0.94
3:C:50:LEU:HD21	22:C:302:PEE:H27	1.50	0.94
1:A:89:GLY:O	21:A:503:NAI:H2N	1.71	0.90
9:J:133:GLU:OE1	9:J:212:ARG:CD	2.19	0.90
9:J:64:PHE:CE1	9:J:210:GLU:HG2	2.08	0.89
19:C:301:SF4:S4	16:Q:223:HIS:CD2	2.67	0.88
25:I:201:CDL:H532	25:I:201:CDL:C31	2.05	0.87
9:J:208:GLY:H	9:J:211:ASP:CG	1.81	0.86
3:C:50:LEU:HD22	22:C:302:PEE:H27	1.59	0.85
9:J:132:ARG:NH2	26:J:401:NDP:O2X	2.10	0.84
2:B:99:GLY:HA3	2:B:170:GLY:O	1.78	0.83
9:J:208:GLY:O	9:J:211:ASP:OD1	1.96	0.83
3:C:71:CYS:HB2	16:Q:222:MET:HE1	1.63	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:I:201:CDL:C55	25:I:201:CDL:H322	2.12	0.80
9:J:207:PHE:HA	9:J:211:ASP:OD2	1.83	0.79
21:A:503:NAI:O1N	30:A:601:HOH:O	2.01	0.78
12:M:149:ASP:HB2	16:Q:361:ALA:HB3	1.65	0.77
11:L:86:ASN:HD22	12:M:275:PRO:HG2	1.50	0.76
2:B:99:GLY:CA	2:B:170:GLY:O	2.33	0.75
17:T:79:VAL:HG21	17:T:84:ILE:HD13	1.69	0.75
1:A:210:THR:HB	1:A:224:ARG:HG2	1.70	0.73
13:N:68:MET:HG2	13:N:69:ASN:H	1.52	0.73
3:C:190:LEU:HD21	22:C:302:PEE:H14	1.71	0.72
11:L:112:MET:HE3	13:N:126:PRO:HB2	1.71	0.72
9:J:133:GLU:OE1	9:J:212:ARG:HD3	1.88	0.72
2:B:89:GLU:OE2	13:N:34:ARG:NH2	2.24	0.71
2:B:63:TRP:HB3	2:B:66:LEU:HD12	1.72	0.71
9:J:133:GLU:OE1	9:J:212:ARG:HD2	1.89	0.71
3:C:71:CYS:HB3	16:Q:141:TYR:CG	2.26	0.70
6:G:109:GLY:O	6:G:110:LEU:HD23	1.91	0.70
3:C:126:GLU:O	9:J:89:TYR:OH	2.10	0.70
16:Q:149:GLN:NE2	16:Q:309:ASP:OD2	2.24	0.69
13:N:32:ASP:OD2	13:N:58:ARG:NH2	2.25	0.68
1:A:285:PRO:O	14:O:222:ARG:NH2	2.26	0.68
9:J:192:ARG:NH1	9:J:198:ALA:O	2.27	0.68
1:A:282:VAL:HG13	1:A:356:VAL:HG21	1.74	0.68
3:C:50:LEU:CD2	22:C:302:PEE:C18	2.70	0.68
6:G:88:LYS:HD3	6:G:98:LEU:HD21	1.75	0.67
15:P:83:GLU:OE1	15:P:142:ARG:NH2	2.27	0.67
13:N:42:ASP:HB2	13:N:48:TYR:HE1	1.59	0.67
2:B:177:THR:HG21	2:B:182:GLU:HB2	1.77	0.67
9:J:214:LEU:HD12	9:J:352:VAL:HG11	1.75	0.67
2:B:165:ASP:OD1	16:Q:368:ARG:NH2	2.27	0.67
12:M:308:ARG:NH2	12:M:578:PRO:O	2.28	0.67
6:G:116:VAL:HG12	6:G:120:MET:HE2	1.76	0.66
16:Q:178:THR:OG1	16:Q:214:TYR:OH	2.10	0.66
8:I:3:SER:O	25:I:201:CDL:C1	2.43	0.66
8:I:40:LYS:HB3	18:W:7:LYS:H	1.60	0.66
2:B:177:THR:HG22	2:B:179:THR:H	1.61	0.66
25:I:201:CDL:H532	25:I:201:CDL:OA9	1.95	0.66
1:A:93:PHE:CD1	1:A:98:LYS:HG3	2.31	0.66
1:A:385:CYS:HB2	19:A:501:SF4:S4	2.36	0.65
5:F:65:LEU:HB2	5:F:79:LEU:HD11	1.77	0.65
10:K:78:HIS:CE1	10:K:79:HIS:HE1	2.15	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
16:Q:302:LEU:HB2	16:Q:401:GLU:HB2	1.79	0.65
11:L:154:LYS:NZ	30:L:201:HOH:O	2.30	0.64
3:C:50:LEU:HD21	22:C:302:PEE:C18	2.26	0.64
3:C:73:ALA:O	3:C:77:MET:HG3	1.97	0.64
16:Q:143:SER:O	16:Q:147:ASN:ND2	2.30	0.64
1:A:326:LEU:HD22	1:A:363:ILE:HD11	1.80	0.64
3:C:188:LYS:HD3	3:C:191:ARG:HD3	1.78	0.64
9:J:64:PHE:CD1	9:J:210:GLU:CG	2.79	0.63
4:E:22:SER:OG	4:E:31:ARG:NH1	2.30	0.63
8:I:70:MET:HE3	15:P:66:ALA:HB1	1.80	0.63
2:B:99:GLY:O	2:B:169:GLU:CG	2.46	0.63
12:M:128:CYS:SG	12:M:140:GLN:NE2	2.70	0.63
1:A:159:ARG:NH2	14:O:176:CYS:O	2.29	0.63
3:C:71:CYS:SG	16:Q:223:HIS:CE1	2.91	0.63
2:B:97:PHE:HA	16:Q:219:GLY:O	1.98	0.62
11:L:90:GLY:HA3	15:P:238:PRO:HB2	1.80	0.62
12:M:222:ILE:HA	12:M:225:ILE:HG12	1.80	0.62
9:J:176:SER:O	9:J:182:ARG:NH1	2.30	0.62
1:A:91:ALA:CB	21:A:503:NAI:H52A	2.28	0.62
10:K:78:HIS:CE1	10:K:79:HIS:CE1	2.87	0.62
3:C:59:ARG:NH1	23:C:303:PLX:O3	2.33	0.61
7:H:25:LYS:HE2	7:H:60:LYS:HG2	1.82	0.61
9:J:214:LEU:CD1	9:J:352:VAL:HG11	2.30	0.61
8:I:3:SER:O	25:I:201:CDL:H1	2.00	0.61
2:B:100:GLU:OE2	2:B:185:TYR:OH	2.17	0.61
25:I:201:CDL:H532	25:I:201:CDL:C32	2.30	0.61
7:H:56:LEU:HG	7:H:60:LYS:HE2	1.83	0.61
14:O:196:LEU:HD13	14:O:201:ILE:HD13	1.82	0.61
2:B:131:GLU:HB2	2:B:144:ARG:HB3	1.82	0.61
20:A:502:FMN:N5	21:A:503:NAI:H4N	2.16	0.61
17:T:47:ASP:O	17:T:52:ARG:NH1	2.34	0.60
3:C:190:LEU:CD2	22:C:302:PEE:H14	2.31	0.60
5:F:79:LEU:HD22	5:F:87:VAL:HG12	1.84	0.60
9:J:64:PHE:HD1	9:J:210:GLU:HG2	1.63	0.60
9:J:213:PHE:CE1	9:J:217:PHE:HE2	2.20	0.59
7:H:7:LYS:HA	7:H:16:VAL:HG21	1.83	0.59
14:O:182:ASN:HB3	14:O:194:GLU:HB3	1.84	0.59
2:B:99:GLY:O	2:B:169:GLU:HG2	2.03	0.59
12:M:419:ARG:NH1	12:M:439:THR:O	2.34	0.59
1:A:93:PHE:CE1	1:A:98:LYS:HG3	2.38	0.58
1:A:113:LEU:HD13	1:A:149:MET:HE1	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:E:64:ARG:NH1	6:G:117:GLU:OE2	2.36	0.58
9:J:64:PHE:HD1	9:J:210:GLU:CG	2.15	0.58
18:W:10:MET:HE3	18:W:11:PRO:HD2	1.85	0.58
2:B:117:LYS:NZ	30:B:1009:HOH:O	2.34	0.58
15:P:44:ARG:HE	15:P:45:PRO:HD3	1.69	0.58
15:P:104:THR:O	15:P:107:GLN:NE2	2.26	0.58
13:N:132:LYS:NZ	13:N:136:GLU:OE1	2.34	0.58
16:Q:86:PRO:HB3	16:Q:94:VAL:HG23	1.85	0.58
9:J:208:GLY:N	9:J:211:ASP:CG	2.59	0.58
1:A:50:ASP:O	1:A:59:ARG:NH2	2.35	0.58
13:N:127:TYR:OH	17:T:61:GLU:O	2.16	0.58
2:B:98:ARG:HB3	2:B:169:GLU:CD	2.28	0.57
25:I:201:CDL:H532	25:I:201:CDL:H322	1.86	0.57
2:B:94:SER:OG	2:B:95:PRO:HD2	2.04	0.57
4:E:39:TRP:O	4:E:43:VAL:HG23	2.04	0.57
15:P:85:GLU:OE2	15:P:142:ARG:NH1	2.33	0.57
17:T:78:GLU:OE2	17:T:119:ARG:NH1	2.35	0.57
12:M:278:HIS:HD2	12:M:280:ASP:H	1.51	0.57
12:M:456:ALA:O	12:M:499:ASN:ND2	2.38	0.57
12:M:137:CYS:HB3	12:M:140:GLN:HB2	1.87	0.56
1:A:90:GLY:HA3	21:A:503:NAI:H1D	1.87	0.56
2:B:105:ARG:HD2	13:N:108:TYR:CG	2.40	0.56
6:G:82:ARG:O	6:G:86:VAL:HG22	2.06	0.56
11:L:89:SER:O	12:M:59:GLN:NE2	2.33	0.56
11:L:156:LYS:HD2	12:M:67:GLU:HG2	1.87	0.56
1:A:64:LYS:NZ	14:O:249:LEU:O	2.31	0.55
1:A:41:ILE:HG12	1:A:253:THR:HG21	1.87	0.55
7:H:44:TYR:HB2	15:P:68:ILE:HG23	1.87	0.55
1:A:319:PRO:HG2	1:A:347:THR:HG21	1.89	0.55
3:C:158:PRO:HB2	9:J:92:MET:HE3	1.87	0.55
1:A:40:ARG:NH1	1:A:289:GLU:O	2.40	0.55
1:A:119:GLU:O	1:A:159:ARG:NH1	2.40	0.55
9:J:37:HIS:HD2	9:J:40:LEU:HG	1.72	0.55
4:E:100:ARG:NH1	30:E:202:HOH:O	2.40	0.54
9:J:50:SER:OG	15:P:225:GLU:OE2	2.23	0.54
12:M:307:ILE:HG22	12:M:582:VAL:HG22	1.89	0.54
17:T:69:ASP:O	17:T:73:GLU:HG3	2.08	0.54
2:B:99:GLY:O	2:B:169:GLU:HG3	2.08	0.54
12:M:275:PRO:HB3	12:M:286:ILE:HG23	1.88	0.54
12:M:472:PRO:O	12:M:510:TRP:NE1	2.33	0.54
22:C:302:PEE:H75	23:C:303:PLX:H172	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:K:100:GLN:HB3	14:O:71:PRO:HA	1.89	0.54
3:C:66:THR:HG21	3:C:76:MET:HE2	1.90	0.54
13:N:124:TYR:HD1	17:T:61:GLU:HG3	1.73	0.54
12:M:217:GLU:HG2	12:M:218:LEU:HG	1.90	0.54
1:A:95:THR:HG22	30:A:619:HOH:O	2.07	0.53
8:I:98:ALA:HB3	15:P:137:PHE:O	2.07	0.53
6:G:90:TYR:HD2	6:G:93:ILE:HG13	1.72	0.53
14:O:199:LYS:O	14:O:203:GLU:HG3	2.09	0.53
6:G:75:THR:HG23	6:G:78:ALA:H	1.74	0.53
1:A:185:ASN:OD1	1:A:190:GLY:N	2.36	0.53
9:J:208:GLY:N	9:J:211:ASP:HB3	2.24	0.53
16:Q:181:LEU:HD23	16:Q:207:ARG:HG2	1.90	0.53
1:A:244:ASN:ND2	20:A:502:FMN:O2	2.34	0.53
11:L:70:GLU:O	11:L:74:THR:OG1	2.27	0.53
6:G:139:MET:HE2	6:G:139:MET:HA	1.90	0.53
9:J:346:GLU:HG2	9:J:371:PRO:HB3	1.89	0.53
10:K:74:ARG:HG3	10:K:74:ARG:HH11	1.73	0.53
11:L:151:LYS:NZ	30:L:205:HOH:O	2.41	0.53
16:Q:206:GLU:OE2	16:Q:209:LYS:NZ	2.41	0.53
2:B:208:ASP:HB3	2:B:212:ARG:HH21	1.73	0.53
5:F:68:ARG:NH2	12:M:364:ASP:OD1	2.42	0.52
11:L:112:MET:CE	13:N:126:PRO:HB2	2.37	0.52
16:Q:101:LEU:HD11	16:Q:444:LEU:HD13	1.90	0.52
3:C:113:MET:HE1	16:Q:86:PRO:HB2	1.91	0.52
8:I:4:ALA:HA	25:I:201:CDL:OA2	2.09	0.52
7:H:51:ILE:HD11	8:I:94:ALA:HB3	1.90	0.52
13:N:65:THR:O	13:N:73:THR:OG1	2.25	0.52
1:A:268:GLU:HG3	1:A:269:ARG:HD3	1.91	0.52
5:F:46:LYS:HE3	12:M:674:LEU:HD21	1.91	0.52
9:J:84:TYR:HD2	9:J:86:CYS:O	1.92	0.52
3:C:67:PHE:HE2	3:C:103:MET:HE2	1.74	0.51
12:M:636:TYR:CD1	12:M:642:VAL:HG22	2.44	0.51
14:O:152:ILE:HG12	14:O:205:ILE:HD11	1.92	0.51
5:F:89:ARG:O	5:F:93:ASN:ND2	2.36	0.51
6:G:144:ILE:O	6:G:148:ILE:HD12	2.11	0.51
2:B:79:ARG:HG2	8:I:12:ARG:CZ	2.40	0.51
1:A:222:LYS:O	11:L:175:LYS:NZ	2.39	0.51
11:L:109:ASN:ND2	11:L:111:LEU:O	2.39	0.51
12:M:388:ASN:HB3	12:M:511:LYS:HE2	1.92	0.51
25:I:201:CDL:C55	25:I:201:CDL:C32	2.86	0.51
11:L:88:GLN:NE2	12:M:141:ASP:OD2	2.34	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:J:83:PRO:HG3	9:J:119:VAL:HG11	1.93	0.51
12:M:666:GLN:OE1	12:M:667:GLN:NE2	2.44	0.51
15:P:173:MET:HB3	15:P:198:PHE:HB2	1.92	0.50
9:J:163:LYS:NZ	9:J:253:ILE:O	2.42	0.50
12:M:217:GLU:HG3	12:M:412:PRO:HB3	1.93	0.50
3:C:50:LEU:HD11	22:C:302:PEE:H59	1.91	0.50
1:A:388:GLY:O	1:A:392:MET:HG3	2.12	0.50
12:M:624:ARG:NH2	12:M:637:ASP:OD1	2.44	0.50
16:Q:100:GLU:HG2	16:Q:108:LYS:HB3	1.94	0.50
4:E:25:MET:HE1	4:E:79:VAL:HG21	1.92	0.49
25:I:201:CDL:H322	25:I:201:CDL:C54	2.42	0.49
3:C:161:ILE:HG13	3:C:180:LEU:HB2	1.94	0.49
12:M:637:ASP:N	12:M:641:GLN:OE1	2.38	0.49
12:M:638:THR:O	12:M:642:VAL:HG23	2.12	0.49
1:A:56:ALA:HB1	1:A:61:ASP:HB2	1.93	0.49
1:A:448:GLU:HA	1:A:451:GLN:HG2	1.94	0.49
8:I:16:SER:OG	8:I:20:LEU:HD11	2.12	0.49
6:G:84:LEU:O	6:G:88:LYS:HG2	2.12	0.49
9:J:37:HIS:HB3	9:J:40:LEU:HD12	1.93	0.49
1:A:91:ALA:HB3	30:A:617:HOH:O	2.11	0.48
2:B:144:ARG:NH1	11:L:112:MET:O	2.46	0.48
12:M:76:ARG:NH2	30:M:935:HOH:O	2.44	0.48
16:Q:105:MET:HE2	16:Q:105:MET:HA	1.94	0.48
1:A:48:ARG:HD3	10:K:70:ASN:O	2.14	0.48
12:M:182:CYS:HA	12:M:225:ILE:HD11	1.95	0.48
25:I:201:CDL:C32	25:I:201:CDL:C54	2.92	0.48
9:J:64:PHE:CZ	9:J:211:ASP:HB3	2.48	0.48
16:Q:139:LEU:HD11	16:Q:424:ILE:HD13	1.95	0.48
1:A:126:LYS:O	1:A:130:ILE:HG12	2.13	0.48
1:A:132:ARG:HB3	1:A:165:GLU:HG3	1.96	0.48
1:A:125:CYS:H	1:A:277:ASN:ND2	2.12	0.48
9:J:214:LEU:HD12	9:J:352:VAL:HG21	1.96	0.48
5:F:86:GLN:HG2	5:F:89:ARG:HH21	1.79	0.48
8:I:96:THR:HB	8:I:97:PRO:HD2	1.96	0.48
9:J:87:GLU:CD	9:J:88:PRO:HD2	2.39	0.48
13:N:10:GLY:O	13:N:14:VAL:HG23	2.14	0.48
17:T:34:THR:OG1	17:T:47:ASP:OD1	2.32	0.48
1:A:357:MET:HE1	1:A:363:ILE:HA	1.94	0.48
2:B:211:TYR:CZ	8:I:39:PRO:HG3	2.48	0.48
9:J:207:PHE:HB2	9:J:214:LEU:HG	1.96	0.48
11:L:163:ASN:O	11:L:171:ARG:HA	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
21:A:503:NAI:H6N	21:A:503:NAI:H2D	1.74	0.47
5:F:24:CYS:N	5:F:58:CYS:SG	2.87	0.47
12:M:522:GLN:NE2	30:M:936:HOH:O	2.45	0.47
9:J:37:HIS:NE2	17:T:49:ASP:HA	2.29	0.47
12:M:306:MET:HE1	13:N:139:PRO:HG3	1.96	0.47
15:P:131:ASN:ND2	30:P:315:HOH:O	2.44	0.47
12:M:390:THR:HA	12:M:600:GLU:HG2	1.95	0.47
1:A:185:ASN:O	1:A:187:CYS:N	2.47	0.47
1:A:292:MET:HE3	1:A:338:ASP:HA	1.97	0.47
11:L:105:GLU:HA	12:M:611:THR:HG21	1.96	0.47
12:M:262:VAL:HG23	12:M:276:ARG:HB2	1.96	0.47
12:M:387:LEU:HD12	12:M:514:ASN:HB3	1.96	0.47
16:Q:332:CYS:O	16:Q:336:GLU:HG3	2.14	0.47
1:A:132:ARG:HG3	1:A:133:HIS:CD2	2.50	0.47
2:B:111:GLU:O	2:B:141:ARG:NH1	2.47	0.47
12:M:43:VAL:HG12	12:M:55:LYS:HD2	1.95	0.47
6:G:109:GLY:C	6:G:110:LEU:HD23	2.39	0.47
8:I:70:MET:O	8:I:70:MET:SD	2.73	0.47
9:J:135:GLU:OE2	9:J:179:ARG:NH2	2.47	0.46
6:G:86:VAL:HG23	6:G:122:MET:HE1	1.97	0.46
12:M:36:VAL:HG22	12:M:102:ILE:HD12	1.96	0.46
2:B:140:ARG:HG3	12:M:238:PHE:CG	2.50	0.46
7:H:93:LYS:HG2	7:H:96:ARG:NH2	2.30	0.46
8:I:8:ILE:HG21	25:I:201:CDL:OA4	2.16	0.46
12:M:340:ALA:HB3	12:M:366:LEU:HD23	1.96	0.46
12:M:449:PRO:HG3	12:M:684:LEU:HD21	1.97	0.46
13:N:68:MET:SD	13:N:81:MET:HG2	2.55	0.46
1:A:67:GLU:O	1:A:71:LYS:HG2	2.16	0.46
2:B:76:TYR:OH	16:Q:257:GLU:OE1	2.24	0.46
12:M:191:GLY:HA3	12:M:439:THR:HB	1.98	0.46
30:B:1004:HOH:O	17:T:114:CYS:SG	2.60	0.46
2:B:151:LYS:HA	3:C:141:GLY:HA2	1.98	0.46
23:C:303:PLX:H6	23:C:303:PLX:H52	1.76	0.46
12:M:347:ASP:CB	12:M:594:ALA:HB1	2.46	0.46
16:Q:404:LYS:NZ	30:Q:530:HOH:O	2.48	0.46
1:A:317:VAL:HG22	1:A:356:VAL:HG12	1.98	0.46
2:B:153:ILE:CG2	3:C:142:TYR:HB2	2.46	0.46
4:E:123:TYR:CZ	12:M:320:GLU:HG3	2.51	0.46
7:H:114:TRP:CD2	7:H:115:PRO:HA	2.51	0.46
7:H:116:ILE:HG21	11:L:96:LYS:NZ	2.31	0.46
16:Q:303:ARG:HG3	16:Q:401:GLU:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ARG:HH21	10:K:77:GLN:HG2	1.80	0.46
2:B:171:PRO:HB3	13:N:93:MET:HE1	1.98	0.46
11:L:86:ASN:ND2	12:M:275:PRO:HG2	2.23	0.46
1:A:220:GLN:NE2	30:A:606:HOH:O	2.33	0.45
16:Q:95:LEU:HB2	16:Q:458:PHE:CZ	2.50	0.45
7:H:90:LEU:O	7:H:94:MET:HG2	2.16	0.45
12:M:144:MET:HG3	16:Q:383:LYS:HG3	1.97	0.45
12:M:306:MET:HB2	12:M:583:ILE:HB	1.98	0.45
9:J:208:GLY:N	9:J:211:ASP:CB	2.80	0.45
11:L:84:ARG:NH1	11:L:88:GLN:O	2.46	0.45
1:A:185:ASN:C	1:A:187:CYS:H	2.25	0.45
7:H:64:ASP:HB3	7:H:67:LYS:HB2	1.98	0.45
1:A:49:HIS:HA	10:K:74:ARG:NH2	2.31	0.45
6:G:117:GLU:HA	6:G:120:MET:HE3	1.98	0.45
9:J:165:ILE:HG12	9:J:199:THR:HB	1.97	0.45
16:Q:184:ILE:HD11	16:Q:251:PHE:CZ	2.52	0.45
9:J:214:LEU:HD23	9:J:214:LEU:HA	1.80	0.45
1:A:195:VAL:O	10:K:97:ARG:NH1	2.40	0.45
10:K:74:ARG:HG3	10:K:74:ARG:NH1	2.31	0.45
2:B:86:TYR:CD1	2:B:87:PRO:HA	2.52	0.45
4:E:66:MET:SD	4:E:69:LYS:NZ	2.84	0.45
6:G:93:ILE:HG23	6:G:108:LEU:HD12	1.97	0.45
25:I:201:CDL:H322	25:I:201:CDL:C53	2.46	0.44
10:K:98:MET:HE3	10:K:98:MET:HB3	1.89	0.44
3:C:173:LEU:O	3:C:177:ILE:HG12	2.17	0.44
12:M:136:GLU:OE2	12:M:272:ARG:NH2	2.49	0.44
12:M:347:ASP:HB3	12:M:594:ALA:HB1	1.98	0.44
7:H:76:GLN:HB2	7:H:79:GLU:OE2	2.18	0.44
1:A:318:ILE:HB	1:A:355:ILE:HB	1.99	0.44
19:C:301:SF4:S2	16:Q:138:ARG:HG2	2.58	0.44
9:J:369:VAL:HG12	9:J:370:LYS:HG2	1.98	0.44
11:L:131:LYS:HD2	11:L:147:VAL:HG11	2.00	0.44
15:P:113:ASP:OD1	16:Q:423:LYS:NZ	2.40	0.44
14:O:195:ASP:OD2	14:O:222:ARG:HD3	2.17	0.44
15:P:249:PRO:O	15:P:250:GLU:HG3	2.18	0.44
5:F:20:ARG:HB2	5:F:66:TRP:HB2	1.99	0.44
1:A:314:LEU:HD23	1:A:329:LYS:HG3	2.00	0.44
8:I:74:LYS:HD2	8:I:74:LYS:HA	1.79	0.44
4:E:101:THR:OG1	15:P:219:VAL:O	2.24	0.44
9:J:279:TYR:HB2	9:J:372:ALA:HB2	2.00	0.44
9:J:282:ALA:HB1	9:J:370:LYS:HE2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:50:LEU:HD22	22:C:302:PEE:C18	2.36	0.43
8:I:42:PRO:HD3	18:W:6:VAL:HG13	2.00	0.43
12:M:382:ARG:HG2	12:M:386:LEU:HD11	1.99	0.43
16:Q:179:ARG:HG2	16:Q:183:HIS:HD2	1.83	0.43
12:M:36:VAL:HB	12:M:56:VAL:HG21	1.99	0.43
1:A:299:LEU:HD12	1:A:303:HIS:ND1	2.33	0.43
7:H:41:ASN:OD1	7:H:41:ASN:N	2.52	0.43
13:N:106:ARG:HB2	13:N:109:ILE:HG13	2.00	0.43
16:Q:182:ASN:OD1	16:Q:404:LYS:NZ	2.47	0.43
1:A:48:ARG:HB3	10:K:70:ASN:O	2.18	0.43
23:C:303:PLX:H181	23:C:303:PLX:H212	1.80	0.43
16:Q:155:VAL:HB	16:Q:229:PRO:HB3	2.00	0.43
1:A:185:ASN:ND2	1:A:187:CYS:O	2.52	0.43
1:A:263:ALA:HA	1:A:271:SER:HB2	2.00	0.43
16:Q:179:ARG:HG2	16:Q:183:HIS:CD2	2.52	0.43
16:Q:383:LYS:HA	16:Q:383:LYS:HD2	1.82	0.43
6:G:80:LYS:HG2	6:G:145:VAL:HG21	2.00	0.43
16:Q:127:LYS:HB3	16:Q:131:GLN:HB3	2.00	0.43
17:T:79:VAL:O	17:T:121:PRO:HD3	2.19	0.43
2:B:118:LEU:HD11	16:Q:382:PHE:CE2	2.54	0.43
23:C:303:PLX:H21	23:C:303:PLX:H1C3	1.78	0.43
25:I:201:CDL:H532	25:I:201:CDL:H312	1.95	0.43
9:J:344:PRO:HG2	9:J:347:LEU:HG	2.01	0.43
14:O:194:GLU:HB2	14:O:218:PRO:HA	2.00	0.43
15:P:163:ALA:O	15:P:167:GLU:HG3	2.18	0.43
16:Q:84:PHE:CZ	16:Q:444:LEU:HD11	2.54	0.43
16:Q:145:MET:HE3	16:Q:214:TYR:CE2	2.54	0.43
16:Q:153:LEU:HD11	16:Q:308:TYR:CD1	2.53	0.43
16:Q:293:LEU:HD22	16:Q:298:ILE:HD12	2.01	0.43
13:N:14:VAL:HA	13:N:23:TYR:CD1	2.54	0.42
16:Q:198:THR:HG22	16:Q:202:TRP:CE2	2.53	0.42
12:M:356:ASP:O	12:M:360:ARG:HG2	2.19	0.42
15:P:75:GLN:HB3	15:P:87:PHE:CD1	2.54	0.42
2:B:41:VAL:HA	8:I:113:LEU:HB3	2.01	0.42
9:J:127:ILE:HG22	9:J:129:LEU:HD13	2.01	0.42
9:J:217:PHE:CZ	9:J:273:LEU:HD11	2.54	0.42
2:B:97:PHE:HA	16:Q:219:GLY:CA	2.50	0.42
25:I:201:CDL:C31	25:I:201:CDL:C53	2.86	0.42
12:M:467:LYS:HB2	12:M:467:LYS:HE3	1.84	0.42
15:P:69:LEU:HD22	15:P:96:VAL:HG22	2.01	0.42
1:A:116:ASN:ND2	1:A:207:GLY:O	2.36	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:ASP:N	1:A:338:ASP:OD1	2.53	0.42
3:C:188:LYS:HG3	3:C:188:LYS:O	2.19	0.42
5:F:30:SER:OG	5:F:63:PRO:HG3	2.20	0.42
16:Q:184:ILE:HG21	16:Q:207:ARG:HG3	2.02	0.42
17:T:104:LYS:HD2	17:T:107:LYS:HE3	2.02	0.42
2:B:153:ILE:HG22	3:C:142:TYR:HB2	2.02	0.42
24:G:201:8Q1:O40	24:G:201:8Q1:N36	2.52	0.42
12:M:498:GLN:HG3	12:M:669:ASN:OD1	2.20	0.42
14:O:130:TYR:HA	14:O:189:ASN:HD21	1.84	0.42
16:Q:204:PHE:HA	16:Q:207:ARG:HB2	2.02	0.42
7:H:76:GLN:O	7:H:78:GLU:N	2.53	0.42
12:M:338:VAL:HG21	12:M:361:VAL:HG11	2.02	0.42
1:A:211:ALA:HB2	1:A:223:PRO:HG3	2.01	0.41
12:M:367:CYS:HB3	12:M:533:GLY:O	2.19	0.41
12:M:372:PHE:H	12:M:532:PRO:HB2	1.85	0.41
9:J:315:THR:C	9:J:316:ARG:HD2	2.46	0.41
16:Q:410:TYR:HB3	16:Q:423:LYS:HB3	2.02	0.41
1:A:162:PHE:HB3	1:A:165:GLU:HB2	2.02	0.41
4:E:121:LYS:HA	4:E:124:VAL:HG22	2.02	0.41
6:G:75:THR:OG1	6:G:156:GLU:OE1	2.39	0.41
13:N:85:GLU:HG2	13:N:86:TRP:N	2.35	0.41
2:B:192:ASN:OD1	17:T:63:ASN:ND2	2.52	0.41
6:G:139:MET:N	6:G:143:GLU:OE2	2.31	0.41
12:M:49:VAL:HG13	12:M:102:ILE:HD13	2.01	0.41
12:M:347:ASP:OD1	12:M:347:ASP:N	2.43	0.41
12:M:379:THR:HG21	12:M:526:LEU:HD22	2.03	0.41
13:N:120:THR:HG22	13:N:122:GLN:H	1.86	0.41
16:Q:188:THR:HB	16:Q:200:PHE:HA	2.02	0.41
1:A:138:LEU:HD13	1:A:245:VAL:HG13	2.03	0.41
16:Q:108:LYS:HD3	16:Q:110:ASP:HB2	2.02	0.41
1:A:325:PRO:HG3	1:A:433:TRP:HB3	2.02	0.41
12:M:404:GLY:HA3	12:M:480:ALA:HB2	2.02	0.41
1:A:251:SER:OG	1:A:252:PRO:HD3	2.21	0.41
1:A:298:GLU:OE2	1:A:302:LYS:HG3	2.20	0.41
20:A:502:FMN:H9	20:A:502:FMN:H1'1	1.78	0.41
11:L:78:ARG:NH2	11:L:148:GLU:OE1	2.51	0.41
11:L:92:ASN:HB3	15:P:238:PRO:HA	2.03	0.41
16:Q:162:GLN:HA	16:Q:163:PRO:HD3	1.96	0.41
1:A:66:LYS:HE3	1:A:188:GLY:O	2.21	0.41
1:A:199:ARG:NH1	14:O:84:ASP:OD2	2.54	0.41
6:G:103:HIS:ND1	6:G:106:LYS:HG2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:M:373:PRO:HG2	12:M:490:LEU:HD22	2.03	0.41
15:P:119:VAL:HG12	15:P:121:THR:HG22	2.03	0.41
1:A:367:ILE:O	1:A:371:ILE:HG12	2.21	0.41
25:I:201:CDL:C53	25:I:201:CDL:H312	2.51	0.41
9:J:83:PRO:HB2	9:J:108:TRP:CD1	2.55	0.41
9:J:173:ASP:OD1	9:J:176:SER:N	2.54	0.41
9:J:204:SER:OG	9:J:238:GLN:O	2.38	0.41
12:M:81:GLU:OE2	12:M:106:SER:OG	2.33	0.41
1:A:384:PRO:HB2	1:A:423:THR:HG22	2.02	0.41
2:B:99:GLY:N	2:B:170:GLY:O	2.54	0.41
14:O:93:LEU:HD12	14:O:122:TYR:HB3	2.03	0.41
16:Q:244:ILE:HG22	16:Q:348:LEU:HD11	2.02	0.41
2:B:210:LEU:HD12	8:I:38:PRO:HB3	2.03	0.40
5:F:23:LEU:O	5:F:57:GLU:HA	2.20	0.40
6:G:74:LEU:H	6:G:74:LEU:HD23	1.86	0.40
16:Q:259:GLU:OE1	16:Q:338:ARG:NH2	2.41	0.40
16:Q:375:MET:HE3	16:Q:379:ILE:HG13	2.03	0.40
1:A:202:GLY:N	30:A:621:HOH:O	2.54	0.40
23:C:303:PLX:H281	23:C:303:PLX:H311	1.63	0.40
8:I:107:SER:C	8:I:109:ASP:H	2.29	0.40
12:M:250:SER:OG	12:M:251:ILE:N	2.47	0.40
15:P:153:ILE:O	15:P:178:PHE:HA	2.21	0.40
1:A:110:PRO:O	1:A:238:CYS:HB3	2.21	0.40
3:C:71:CYS:HB2	16:Q:222:MET:CE	2.43	0.40
5:F:23:LEU:HD13	5:F:33:VAL:HG12	2.03	0.40
16:Q:139:LEU:HD13	16:Q:424:ILE:HG21	2.03	0.40
9:J:172:ALA:HA	9:J:181:LEU:HB3	2.04	0.40
12:M:133:GLN:O	12:M:133:GLN:HG2	2.21	0.40
15:P:115:THR:HB	16:Q:423:LYS:HE3	2.02	0.40
8:I:97:PRO:HD3	15:P:61:PHE:CE1	2.56	0.40
12:M:252:ASP:OD2	12:M:290:THR:OG1	2.39	0.40
16:Q:80:LEU:HD23	16:Q:80:LEU:H	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

entries.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	429/431 (100%)	415 (97%)	14 (3%)	0	100	100
2	B	174/176 (99%)	171 (98%)	3 (2%)	0	100	100
3	C	154/156 (99%)	147 (96%)	7 (4%)	0	100	100
4	E	113/115 (98%)	110 (97%)	3 (3%)	0	100	100
5	F	84/86 (98%)	84 (100%)	0	0	100	100
6	G	86/88 (98%)	86 (100%)	0	0	100	100
7	H	110/112 (98%)	103 (94%)	7 (6%)	0	100	100
8	I	93/112 (83%)	83 (89%)	10 (11%)	0	100	100
9	J	289/341 (85%)	280 (97%)	8 (3%)	1 (0%)	36	58
10	K	40/42 (95%)	40 (100%)	0	0	100	100
11	L	123/125 (98%)	121 (98%)	2 (2%)	0	100	100
12	M	688/690 (100%)	674 (98%)	14 (2%)	0	100	100
13	N	142/144 (99%)	139 (98%)	3 (2%)	0	100	100
14	O	215/217 (99%)	207 (96%)	8 (4%)	0	100	100
15	P	206/208 (99%)	198 (96%)	8 (4%)	0	100	100
16	Q	374/385 (97%)	365 (98%)	9 (2%)	0	100	100
17	T	94/96 (98%)	92 (98%)	2 (2%)	0	100	100
18	W	27/29 (93%)	24 (89%)	3 (11%)	0	100	100
All	All	3441/3553 (97%)	3339 (97%)	101 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
9	J	38	HIS

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	345/345 (100%)	345 (100%)	0	100	100
2	B	151/151 (100%)	151 (100%)	0	100	100
3	C	132/132 (100%)	132 (100%)	0	100	100
4	E	107/107 (100%)	107 (100%)	0	100	100
5	F	76/76 (100%)	76 (100%)	0	100	100
6	G	76/81 (94%)	75 (99%)	1 (1%)	61	82
7	H	99/99 (100%)	99 (100%)	0	100	100
8	I	87/97 (90%)	87 (100%)	0	100	100
9	J	255/295 (86%)	254 (100%)	1 (0%)	84	93
10	K	41/41 (100%)	40 (98%)	1 (2%)	43	70
11	L	113/113 (100%)	113 (100%)	0	100	100
12	M	580/580 (100%)	580 (100%)	0	100	100
13	N	130/130 (100%)	130 (100%)	0	100	100
14	O	183/183 (100%)	183 (100%)	0	100	100
15	P	190/190 (100%)	190 (100%)	0	100	100
16	Q	327/331 (99%)	327 (100%)	0	100	100
17	T	79/79 (100%)	78 (99%)	1 (1%)	61	82
18	W	24/24 (100%)	24 (100%)	0	100	100
All	All	2995/3054 (98%)	2991 (100%)	4 (0%)	87	96

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
6	G	111	ASP
9	J	210	GLU
10	K	79	HIS
17	T	59	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	HIS
1	A	164	ASN
1	A	270	ASN
1	A	283	ASN
1	A	303	HIS

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Mol	Chain	Res	Type
1	A	313	ASN
1	A	452	GLN
4	E	70	ASN
5	F	48	ASN
6	G	101	ASN
7	H	37	GLN
8	I	51	ASN
8	I	110	GLN
9	J	79	GLN
9	J	93	HIS
9	J	122	HIS
9	J	138	ASN
9	J	150	HIS
9	J	215	ASN
9	J	356	HIS
9	J	376	ASN
10	K	78	HIS
10	K	79	HIS
11	L	86	ASN
12	M	123	ASN
12	M	388	ASN
12	M	459	ASN
12	M	652	ASN
12	M	666	GLN
12	M	667	GLN
13	N	12	GLN
13	N	113	HIS
14	O	187	GLN
14	O	191	ASN
15	P	196	HIS
16	Q	183	HIS
17	T	63	ASN
18	W	8	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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$
16	2MR	Q	118	16	10,12,13	2.02	1 (10%)	5,13,15	6.08	3 (60%)

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
16	2MR	Q	118	16	-	2/10/13/15	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	Q	118	2MR	CZ-NE	5.74	1.46	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	Q	118	2MR	NE-CZ-NH2	12.35	130.81	119.48
16	Q	118	2MR	CD-NE-CZ	4.37	131.58	123.36
16	Q	118	2MR	CQ2-NH2-CZ	3.41	130.97	123.65

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
16	Q	118	2MR	NE-CD-CG-CB
16	Q	118	2MR	CA-CB-CG-CD

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 2 are monoatomic - leaving 15 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
20	FMN	A	502	-	33,33,33	1.05	2 (6%)	48,50,50	1.24	8 (16%)
19	SF4	B	302	2	0,12,12	-	-	-	-	-
26	NDP	J	401	-	51,52,52	4.26	25 (49%)	71,80,80	2.20	12 (16%)
19	SF4	A	501	1	0,12,12	-	-	-	-	-
19	SF4	M	802	12	0,12,12	-	-	-	-	-
19	SF4	M	801	12	0,12,12	-	-	-	-	-
22	PEE	C	302	-	46,46,50	1.24	6 (13%)	49,51,55	1.01	2 (4%)
27	FES	M	803	12	0,4,4	-	-	-	-	-
24	8Q1	G	201	-	32,34,34	2.30	7 (21%)	39,43,43	1.75	11 (28%)
19	SF4	C	301	3,16	0,12,12	-	-	-	-	-
21	NAI	A	503	-	47,48,48	4.01	22 (46%)	64,73,73	1.68	11 (17%)
19	SF4	B	301	2	0,12,12	-	-	-	-	-
27	FES	O	301	14	0,4,4	-	-	-	-	-
25	CDL	I	201	-	50,50,99	1.29	4 (8%)	56,62,111	1.32	5 (8%)
23	PLX	C	303	-	51,51,51	1.21	4 (7%)	53,59,59	0.62	1 (1%)

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
20	FMN	A	502	-	-	8/18/18/18	0/3/3/3
19	SF4	B	302	2	-	-	0/6/5/5
26	NDP	J	401	-	-	8/34/77/77	0/5/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
19	SF4	A	501	1	-	-	0/6/5/5
19	SF4	M	802	12	-	-	0/6/5/5
22	PEE	C	302	-	-	20/50/50/54	-
19	SF4	M	801	12	-	-	0/6/5/5
27	FES	M	803	12	-	-	0/1/1/1
24	8Q1	G	201	-	-	16/41/41/41	-
19	SF4	C	301	3,16	-	-	0/6/5/5
21	NAI	A	503	-	-	14/29/72/72	0/5/5/5
19	SF4	B	301	2	-	-	0/6/5/5
27	FES	O	301	14	-	-	0/1/1/1
25	CDL	I	201	-	-	21/61/61/110	-
23	PLX	C	303	-	-	25/55/55/55	-

All (70) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	J	401	NDP	C3B-C2B	-12.82	1.24	1.53
26	J	401	NDP	O4D-C4D	10.79	1.69	1.45
21	A	503	NAI	C3D-C4D	-10.34	1.26	1.53
26	J	401	NDP	C3D-C4D	-10.01	1.27	1.53
21	A	503	NAI	O4B-C1B	9.50	1.64	1.42
21	A	503	NAI	O4B-C4B	-8.24	1.26	1.45
24	G	201	8Q1	P24-O27	7.95	1.85	1.60
26	J	401	NDP	O4B-C4B	-7.94	1.27	1.45
21	A	503	NAI	C2D-C1D	-7.69	1.29	1.53
26	J	401	NDP	C2N-C3N	7.48	1.55	1.35
21	A	503	NAI	C2B-C1B	-7.28	1.30	1.53
26	J	401	NDP	C6N-C5N	7.27	1.55	1.33
21	A	503	NAI	O4D-C4D	6.96	1.60	1.45
26	J	401	NDP	PN-O3	6.57	1.66	1.59
21	A	503	NAI	PA-O3	6.47	1.66	1.59
21	A	503	NAI	C2D-C3D	5.95	1.69	1.53
26	J	401	NDP	P2B-O2B	5.83	1.69	1.59
24	G	201	8Q1	C28-C29	5.77	1.62	1.52
26	J	401	NDP	C6A-N6A	5.74	1.48	1.34
21	A	503	NAI	PN-O3	5.60	1.65	1.59
21	A	503	NAI	O4D-C1D	5.54	1.54	1.42
26	J	401	NDP	PA-O3	5.45	1.65	1.59
26	J	401	NDP	C3B-C4B	5.35	1.66	1.53
21	A	503	NAI	C7N-N7N	5.29	1.48	1.33
21	A	503	NAI	C4N-C3N	-5.23	1.40	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
21	A	503	NAI	C6A-N6A	5.08	1.47	1.34
26	J	401	NDP	C6N-N1N	5.08	1.49	1.37
26	J	401	NDP	O4D-C1D	-5.00	1.30	1.42
26	J	401	NDP	O4B-C1B	4.78	1.53	1.42
26	J	401	NDP	C1B-N9A	-4.57	1.33	1.46
21	A	503	NAI	O2B-C2B	4.36	1.53	1.43
25	I	201	CDL	OB8-CB7	4.28	1.45	1.33
25	I	201	CDL	OA8-CA7	4.28	1.45	1.33
25	I	201	CDL	OA6-CA5	4.16	1.46	1.34
25	I	201	CDL	OB6-CB5	4.06	1.45	1.34
26	J	401	NDP	O2D-C2D	-3.97	1.33	1.43
24	G	201	8Q1	C1-S44	3.85	1.85	1.76
22	C	302	PEE	C18-C19	3.82	1.53	1.31
26	J	401	NDP	C7N-N7N	3.81	1.44	1.33
22	C	302	PEE	C39-C38	3.73	1.52	1.31
21	A	503	NAI	C7N-C3N	3.55	1.56	1.48
20	A	502	FMN	C4A-N5	3.41	1.38	1.30
24	G	201	8Q1	O27-C28	-3.34	1.33	1.43
21	A	503	NAI	C4N-C5N	-3.34	1.40	1.49
24	G	201	8Q1	C34-N36	3.31	1.41	1.33
24	G	201	8Q1	C6-C1	3.25	1.54	1.50
26	J	401	NDP	C8A-N9A	-3.16	1.32	1.37
26	J	401	NDP	C5A-N7A	-3.00	1.33	1.39
26	J	401	NDP	C7N-C3N	2.92	1.55	1.48
26	J	401	NDP	O3D-C3D	2.90	1.50	1.43
23	C	303	PLX	O6-C4	-2.84	1.41	1.44
26	J	401	NDP	O2B-C2B	2.74	1.53	1.44
21	A	503	NAI	C8A-N9A	-2.65	1.33	1.37
24	G	201	8Q1	C39-N41	2.64	1.39	1.33
22	C	302	PEE	O2-C2	-2.59	1.40	1.46
21	A	503	NAI	PN-O5D	2.47	1.69	1.59
20	A	502	FMN	C10-N1	2.46	1.38	1.33
22	C	302	PEE	O3-C30	2.46	1.40	1.33
21	A	503	NAI	C6N-C5N	2.45	1.40	1.33
21	A	503	NAI	C5B-C4B	2.42	1.58	1.51
23	C	303	PLX	C7-C6	2.36	1.55	1.50
22	C	302	PEE	O2-C10	2.32	1.40	1.34
21	A	503	NAI	O3B-C3B	-2.31	1.37	1.43
26	J	401	NDP	C2D-C3D	2.30	1.59	1.53
23	C	303	PLX	P1-O4	2.19	1.68	1.59
26	J	401	NDP	O7N-C7N	-2.18	1.19	1.24
22	C	302	PEE	O3-C3	-2.15	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
26	J	401	NDP	PA-O5B	2.12	1.67	1.59
21	A	503	NAI	C5A-N7A	-2.10	1.35	1.39
23	C	303	PLX	P1-O1	2.09	1.67	1.59

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	J	401	NDP	C3N-C2N-N1N	-7.79	111.77	123.20
26	J	401	NDP	C1D-N1N-C2N	-7.16	109.35	121.14
26	J	401	NDP	C6N-N1N-C2N	-6.55	112.31	119.32
21	A	503	NAI	C5A-C4A-N3A	-5.52	119.12	126.72
26	J	401	NDP	C1D-N1N-C6N	-5.49	109.16	120.77
26	J	401	NDP	C5A-C4A-N3A	-5.47	119.19	126.72
24	G	201	8Q1	C6-C1-S44	5.05	119.42	113.40
21	A	503	NAI	N3A-C2A-N1A	-4.37	121.96	128.58
26	J	401	NDP	N3A-C2A-N1A	-4.25	122.14	128.58
25	I	201	CDL	OB6-CB5-C51	4.09	120.33	111.48
25	I	201	CDL	OA6-CA5-C11	4.07	120.28	111.48
22	C	302	PEE	O2-C10-C11	3.95	120.03	111.48
21	A	503	NAI	N3A-C4A-N9A	3.91	133.82	127.17
24	G	201	8Q1	C43-S44-C1	3.89	113.35	101.84
26	J	401	NDP	N3A-C4A-N9A	3.86	133.72	127.17
21	A	503	NAI	C2A-N3A-C4A	3.69	120.83	111.83
26	J	401	NDP	C2A-N3A-C4A	3.61	120.64	111.83
21	A	503	NAI	C4A-C5A-N7A	-3.40	106.70	110.58
21	A	503	NAI	C5A-N7A-C8A	3.38	108.77	103.45
24	G	201	8Q1	O35-C34-N36	-3.36	115.86	122.98
26	J	401	NDP	C4A-C5A-N7A	-3.25	106.87	110.58
21	A	503	NAI	N9A-C8A-N7A	-3.23	109.35	113.94
20	A	502	FMN	C4-N3-C2	-3.22	119.92	125.64
26	J	401	NDP	N9A-C8A-N7A	-3.17	109.43	113.94
26	J	401	NDP	C5A-N7A-C8A	3.07	108.28	103.45
21	A	503	NAI	C3D-C2D-C1D	3.05	107.23	101.46
25	I	201	CDL	OA8-CA7-C31	2.89	120.64	111.83
25	I	201	CDL	CB4-OB6-CB5	-2.85	110.97	117.80
25	I	201	CDL	OB8-CB7-C71	2.78	120.30	111.83
21	A	503	NAI	C2D-C3D-C4D	2.76	107.95	102.61
21	A	503	NAI	C4D-O4D-C1D	-2.75	103.39	109.47
22	C	302	PEE	O3-C30-C31	2.72	120.13	111.83
24	G	201	8Q1	O40-C39-N41	-2.67	117.78	123.03
24	G	201	8Q1	O4-C1-S44	-2.66	119.30	122.68
20	A	502	FMN	C4A-C4-N3	2.64	119.97	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	G	201	8Q1	O2-P24-O27	-2.64	99.80	106.67
24	G	201	8Q1	C32-C34-N36	2.51	121.25	116.48
20	A	502	FMN	O4-C4-C4A	-2.48	119.98	126.53
26	J	401	NDP	C4A-N9A-C8A	2.47	108.33	105.74
20	A	502	FMN	C9A-C5A-N5	-2.42	119.88	122.45
20	A	502	FMN	C4A-C10-N10	2.42	119.95	116.48
24	G	201	8Q1	C38-C39-N41	2.37	120.67	116.34
24	G	201	8Q1	O1-P24-O2	2.36	116.64	107.80
20	A	502	FMN	C5A-C9A-N10	2.34	120.09	117.97
23	C	303	PLX	C1A-N1-C1	2.34	119.22	109.91
24	G	201	8Q1	O4-C1-C6	-2.34	121.28	123.98
20	A	502	FMN	C10-C4A-N5	-2.33	120.06	124.81
21	A	503	NAI	C4A-N9A-C8A	2.27	108.12	105.74
24	G	201	8Q1	O27-P24-O3	-2.06	100.87	106.44
20	A	502	FMN	C4A-C10-N1	-2.05	119.56	124.59

There are no chirality outliers.

All (112) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
20	A	502	FMN	N10-C1'-C2'-O2'
20	A	502	FMN	N10-C1'-C2'-C3'
20	A	502	FMN	C1'-C2'-C3'-O3'
20	A	502	FMN	C1'-C2'-C3'-C4'
20	A	502	FMN	O2'-C2'-C3'-O3'
20	A	502	FMN	O2'-C2'-C3'-C4'
20	A	502	FMN	C3'-C4'-C5'-O5'
21	A	503	NAI	C5B-O5B-PA-O1A
21	A	503	NAI	C5D-O5D-PN-O3
21	A	503	NAI	C5D-O5D-PN-O1N
22	C	302	PEE	C1-O3P-P-O1P
22	C	302	PEE	C1-O3P-P-O4P
22	C	302	PEE	C4-O4P-P-O1P
23	C	303	PLX	C3-O4-P1-O1
23	C	303	PLX	C3-O4-P1-O3
23	C	303	PLX	N1-C1-C2-O1
24	G	201	8Q1	O4-C1-S44-C43
24	G	201	8Q1	C6-C1-S44-C43
24	G	201	8Q1	N36-C37-C38-C39
24	G	201	8Q1	C42-C43-S44-C1
24	G	201	8Q1	C28-O27-P24-O3
24	G	201	8Q1	C28-O27-P24-O2

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Mol	Chain	Res	Type	Atoms
24	G	201	8Q1	C28-O27-P24-O1
25	I	201	CDL	CA2-OA2-PA1-OA3
25	I	201	CDL	CA2-OA2-PA1-OA5
25	I	201	CDL	CA3-OA5-PA1-OA2
25	I	201	CDL	CA3-OA5-PA1-OA4
25	I	201	CDL	C11-CA5-OA6-CA4
25	I	201	CDL	CB2-OB2-PB2-OB4
25	I	201	CDL	CB2-OB2-PB2-OB5
25	I	201	CDL	CB3-OB5-PB2-OB2
25	I	201	CDL	CB3-OB5-PB2-OB4
26	J	401	NDP	C5D-O5D-PN-O1N
26	J	401	NDP	C2N-C3N-C7N-O7N
25	I	201	CDL	OA7-CA5-OA6-CA4
25	I	201	CDL	O1-C1-CB2-OB2
22	C	302	PEE	C31-C30-O3-C3
21	A	503	NAI	C3B-C4B-C5B-O5B
21	A	503	NAI	O4D-C4D-C5D-O5D
21	A	503	NAI	C3D-C4D-C5D-O5D
22	C	302	PEE	O5-C30-O3-C3
23	C	303	PLX	C28-C29-C30-C31
26	J	401	NDP	C2D-C1D-N1N-C6N
21	A	503	NAI	O4B-C4B-C5B-O5B
22	C	302	PEE	C30-C31-C32-C33
25	I	201	CDL	CA6-CA4-OA6-CA5
24	G	201	8Q1	C12-C13-C14-C15
24	G	201	8Q1	C10-C11-C12-C13
23	C	303	PLX	C7-C8-C9-C10
23	C	303	PLX	C10-C11-C12-C13
23	C	303	PLX	C30-C31-C32-C33
23	C	303	PLX	C13-C14-C15-C16
25	I	201	CDL	C51-CB5-OB6-CB4
22	C	302	PEE	C15-C16-C17-C18
23	C	303	PLX	C31-C32-C33-C34
22	C	302	PEE	C11-C10-O2-C2
24	G	201	8Q1	C11-C12-C13-C14
23	C	303	PLX	C9-C10-C11-C12
25	I	201	CDL	OB7-CB5-OB6-CB4
23	C	303	PLX	C11-C12-C13-C14
23	C	303	PLX	C27-C28-C29-C30
23	C	303	PLX	C33-C34-C35-C36
22	C	302	PEE	O4-C10-O2-C2
26	J	401	NDP	C2N-C3N-C7N-N7N

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Mol	Chain	Res	Type	Atoms
21	A	503	NAI	PN-O3-PA-O5B
23	C	303	PLX	C14-C15-C16-C17
23	C	303	PLX	C25-C26-C27-C28
22	C	302	PEE	C11-C12-C13-C14
25	I	201	CDL	C51-C52-C53-C54
25	I	201	CDL	O1-C1-CA2-OA2
23	C	303	PLX	O9-C24-C25-C26
22	C	302	PEE	C34-C35-C36-C37
20	A	502	FMN	O4'-C4'-C5'-O5'
24	G	201	8Q1	O27-C28-C29-C30
24	G	201	8Q1	O27-C28-C29-C31
23	C	303	PLX	O8-C24-C25-C26
22	C	302	PEE	C42-C43-C44-C45
25	I	201	CDL	CA2-C1-CB2-OB2
23	C	303	PLX	C12-C13-C14-C15
22	C	302	PEE	C43-C44-C45-C46
21	A	503	NAI	C5B-O5B-PA-O2A
21	A	503	NAI	C5B-O5B-PA-O3
23	C	303	PLX	C3-C4-O6-C6
23	C	303	PLX	C5-C4-O6-C6
23	C	303	PLX	C2-O1-P1-O3
25	I	201	CDL	CA2-OA2-PA1-OA4
26	J	401	NDP	C5D-O5D-PN-O3
26	J	401	NDP	C5D-O5D-PN-O2N
22	C	302	PEE	C44-C45-C46-C47
23	C	303	PLX	C18-C19-C20-C21
22	C	302	PEE	C38-C39-C40-C41
21	A	503	NAI	C2D-C1D-N1N-C2N
24	G	201	8Q1	C29-C32-C34-O35
23	C	303	PLX	O6-C6-C7-C8
24	G	201	8Q1	C13-C14-C15-C16
24	G	201	8Q1	C29-C32-C34-N36
21	A	503	NAI	O4D-C1D-N1N-C2N
22	C	302	PEE	C16-C17-C18-C19
22	C	302	PEE	C18-C19-C20-C21
23	C	303	PLX	C16-C17-C18-C19
24	G	201	8Q1	C1-C6-C7-C8
23	C	303	PLX	O7-C6-C7-C8
25	I	201	CDL	C1-CA2-OA2-PA1
22	C	302	PEE	C19-C20-C21-C22
21	A	503	NAI	C2D-C1D-N1N-C6N
22	C	302	PEE	O3P-C1-C2-O2

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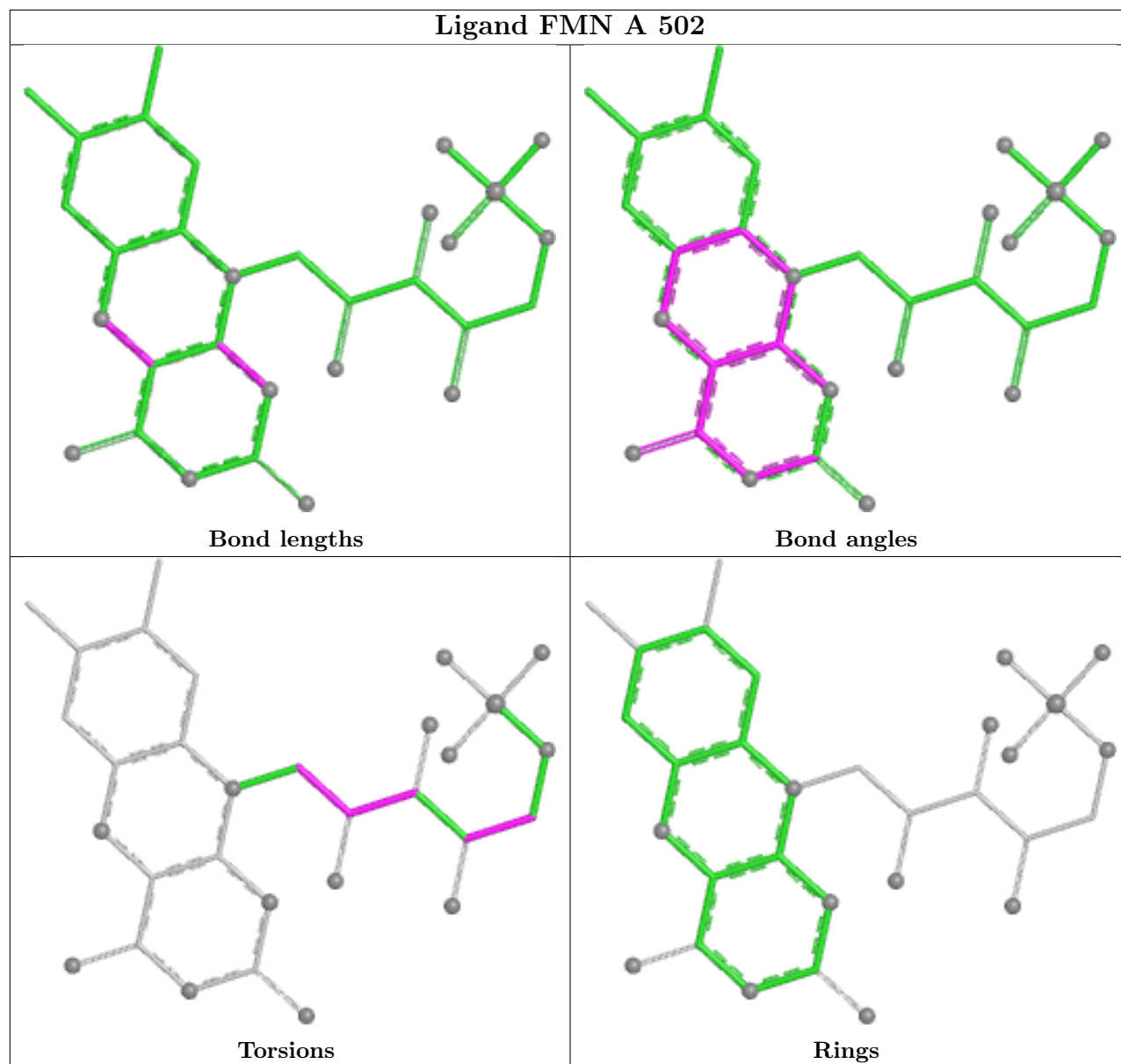
Mol	Chain	Res	Type	Atoms
25	I	201	CDL	C52-C51-CB5-OB6
26	J	401	NDP	O4D-C1D-N1N-C6N
22	C	302	PEE	O3-C30-C31-C32
26	J	401	NDP	C2B-O2B-P2B-O1X
25	I	201	CDL	C52-C51-CB5-OB7
21	A	503	NAI	O4D-C1D-N1N-C6N

There are no ring outliers.

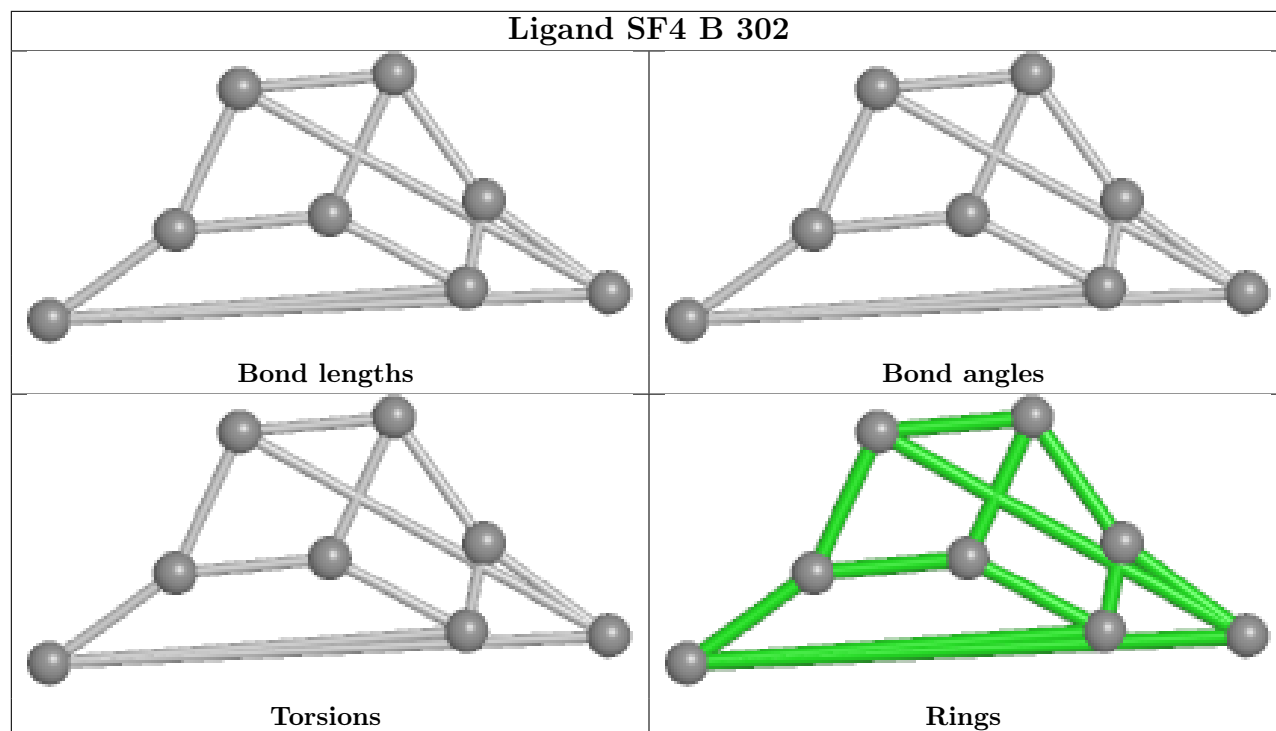
9 monomers are involved in 49 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	A	502	FMN	3	0
26	J	401	NDP	2	0
19	A	501	SF4	1	0
22	C	302	PEE	10	0
24	G	201	8Q1	1	0
19	C	301	SF4	2	0
21	A	503	NAI	7	0
25	I	201	CDL	19	0
23	C	303	PLX	6	0

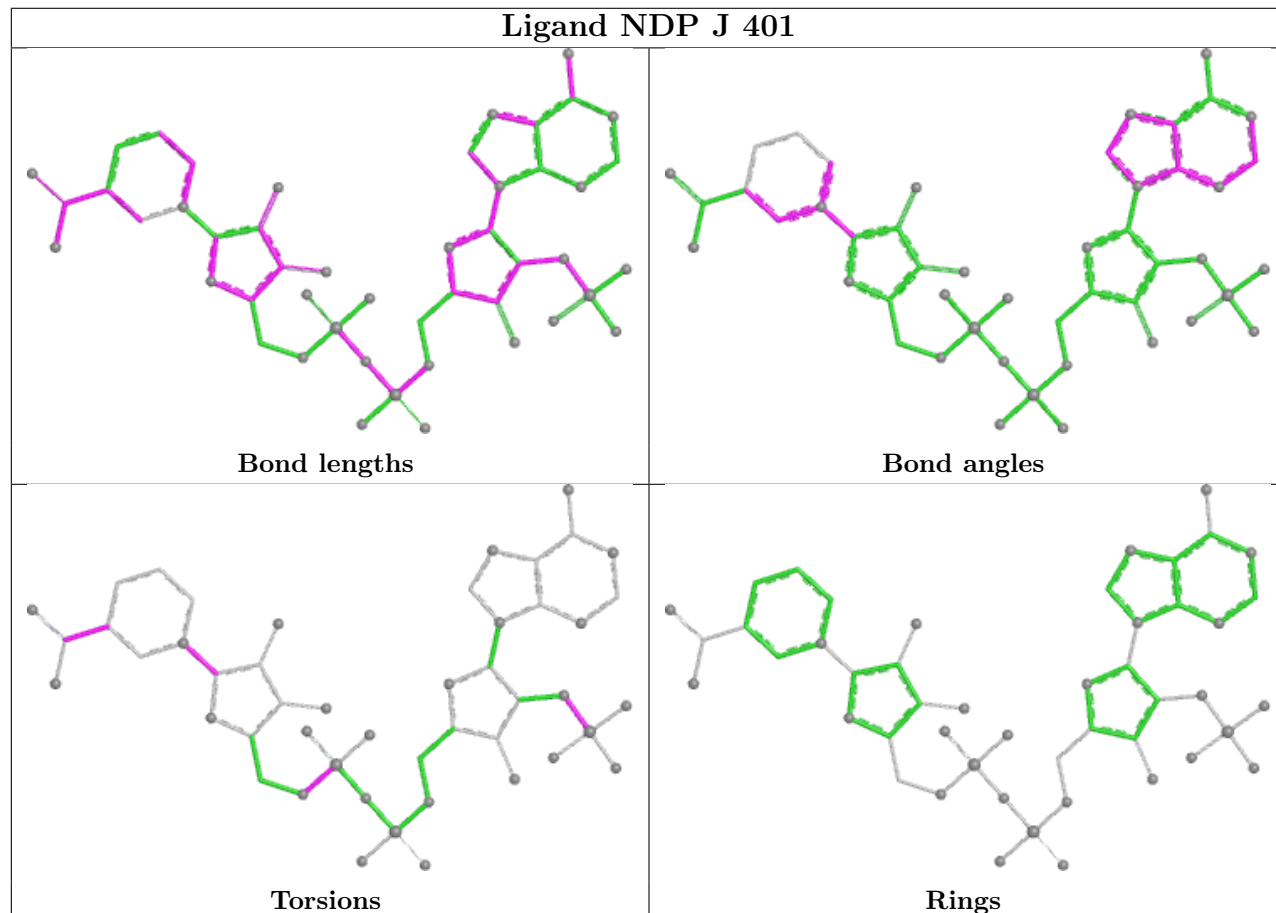
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

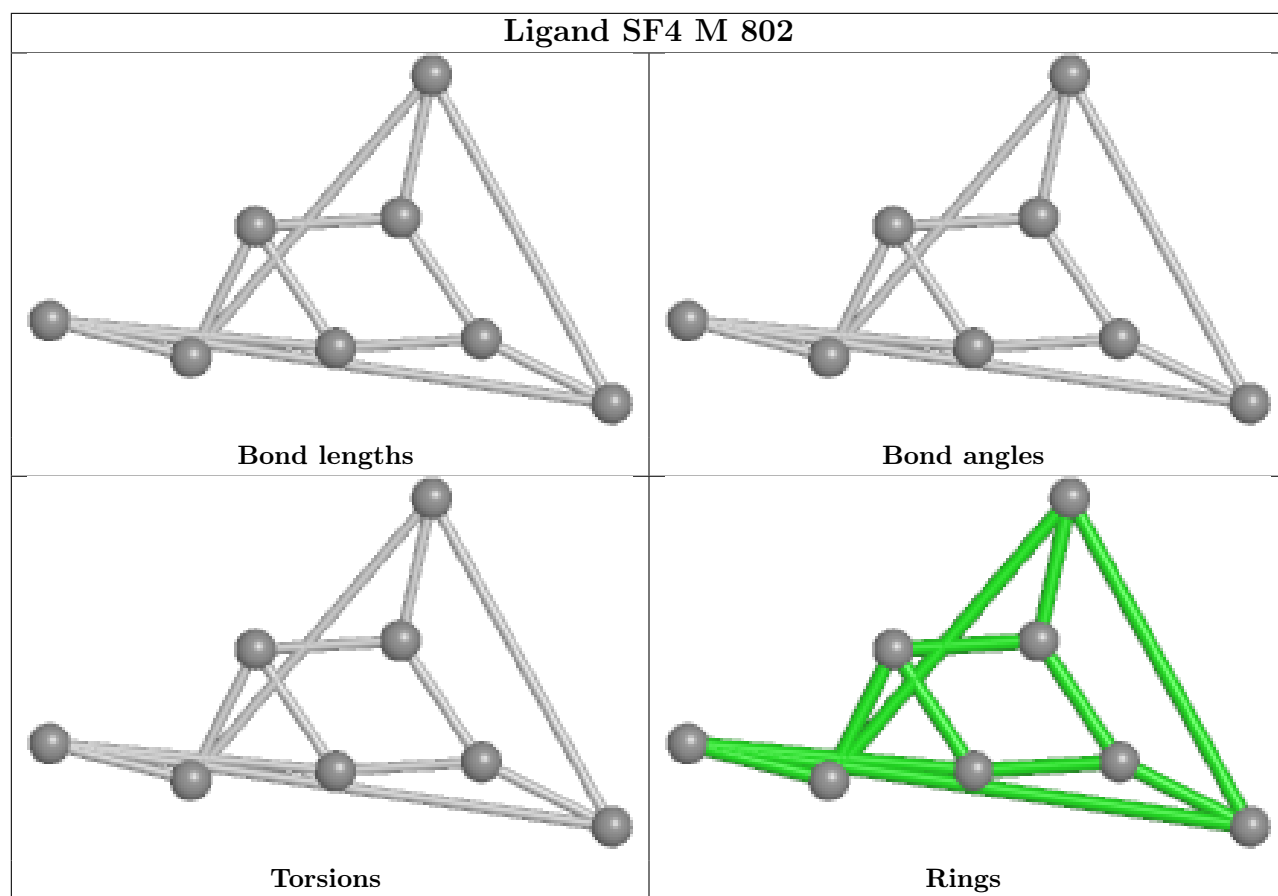
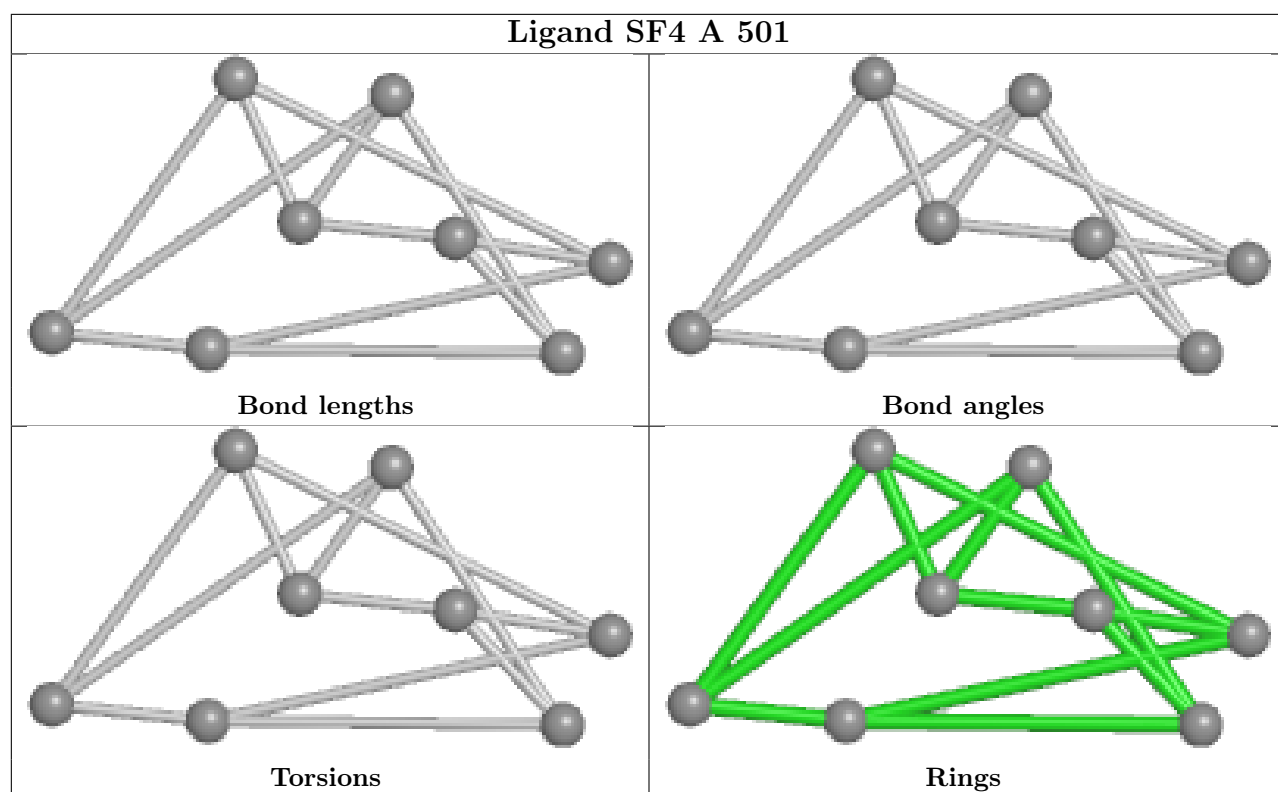


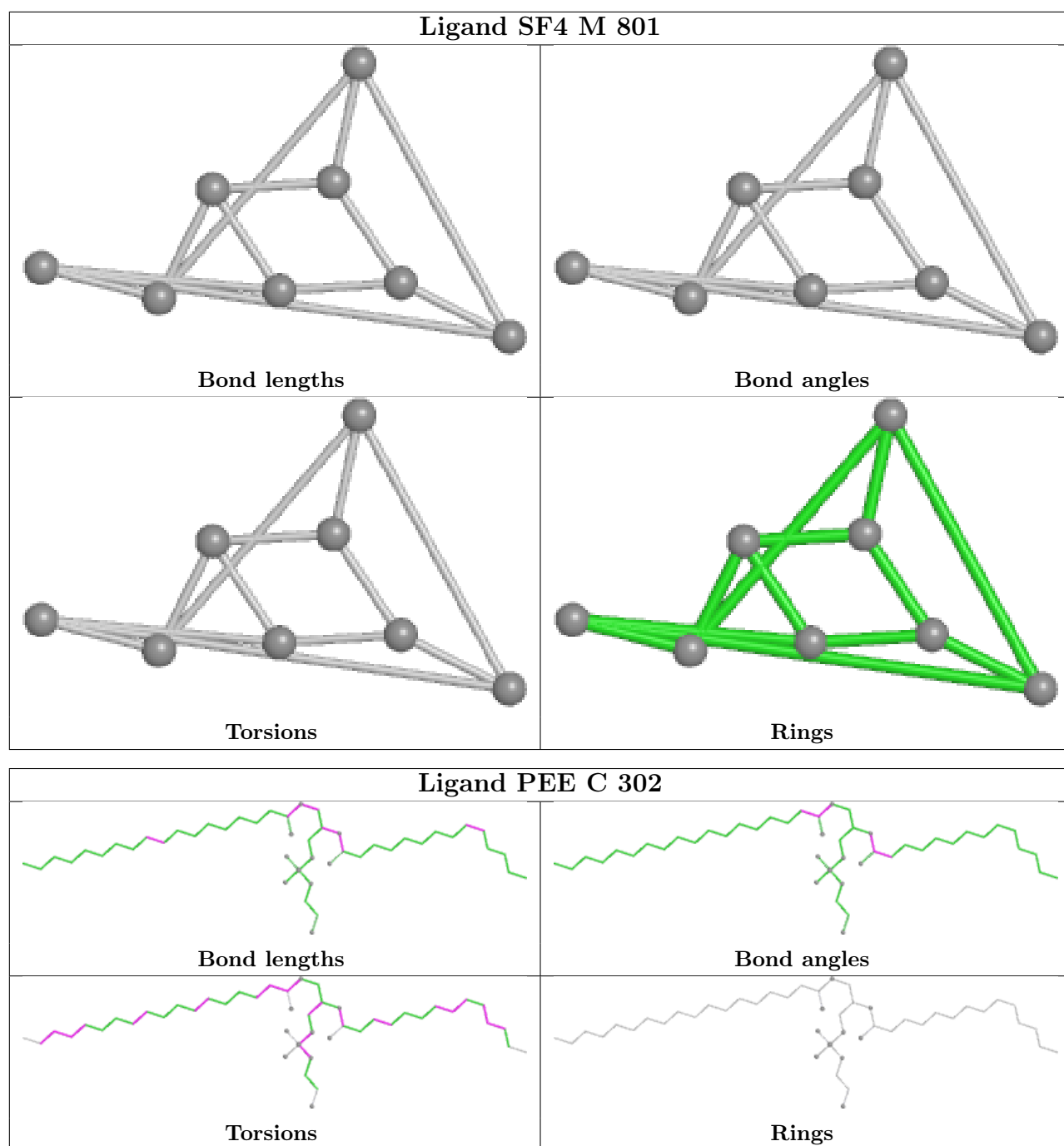
Ligand SF4 B 302

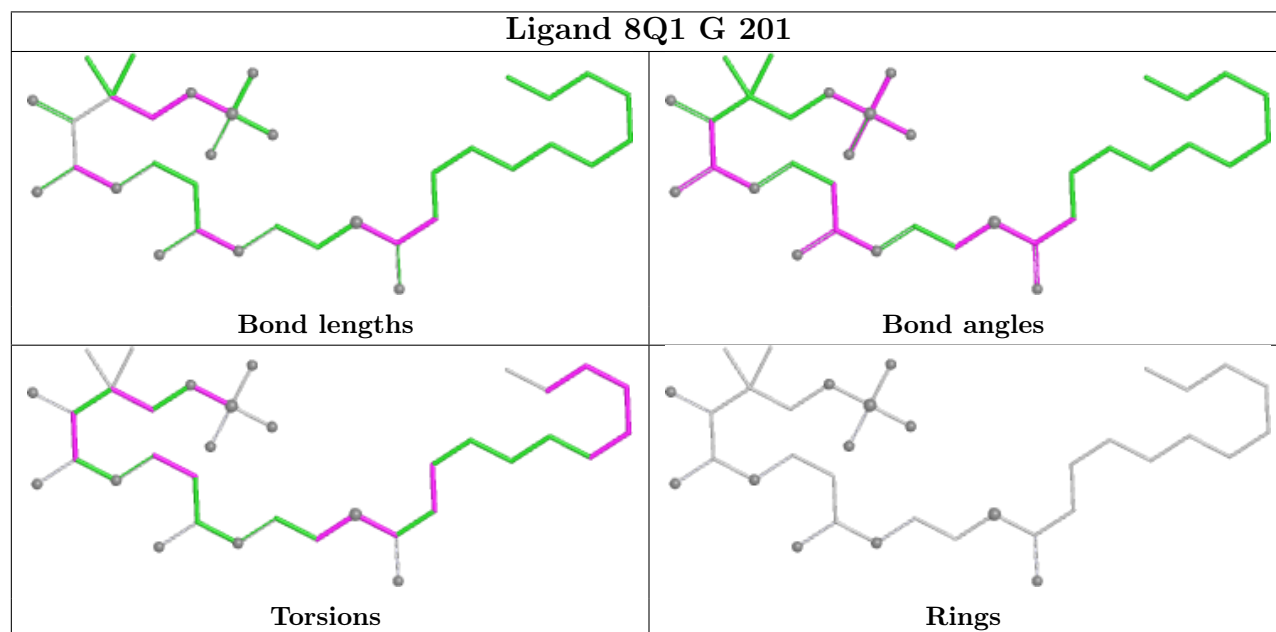
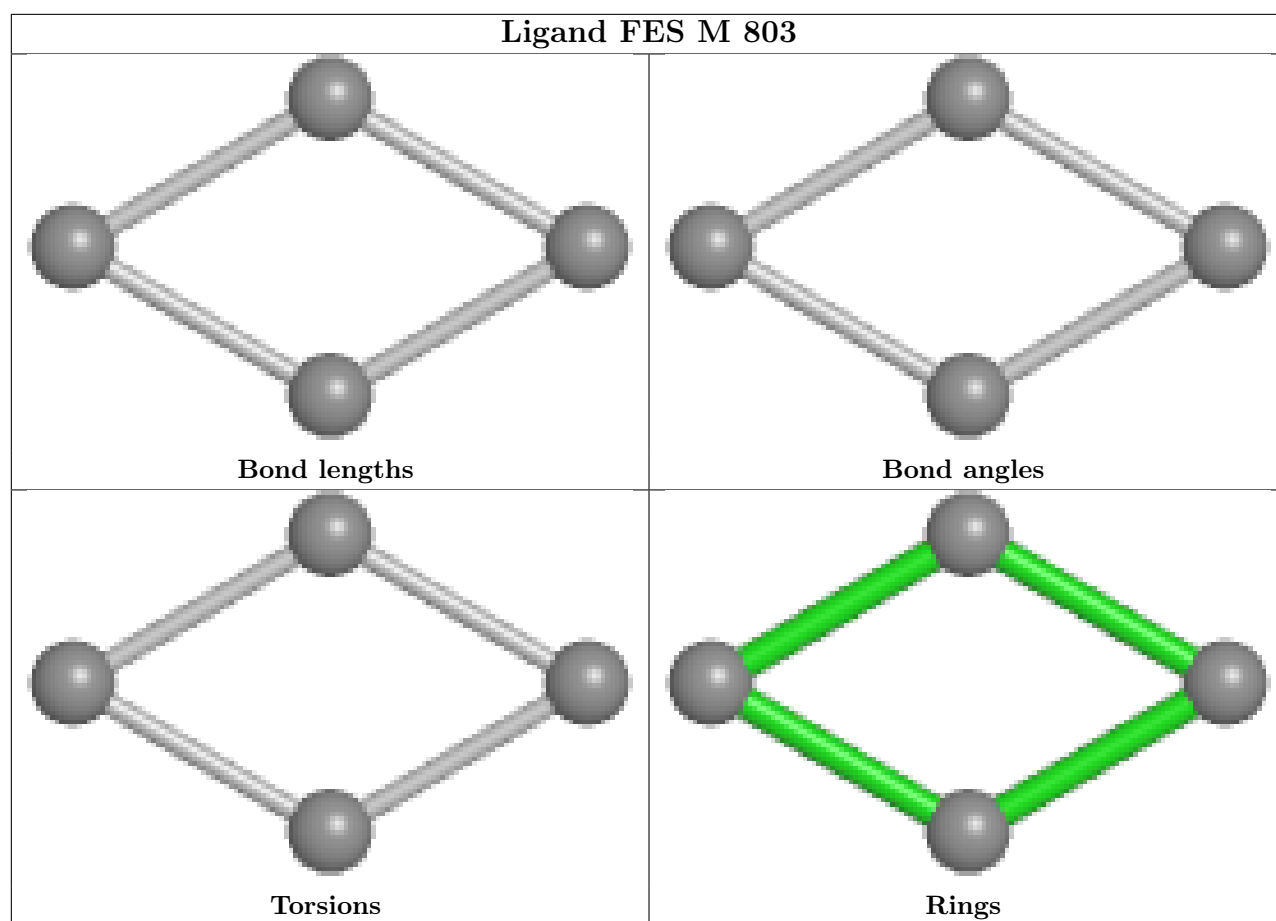


Ligand NDP J 401

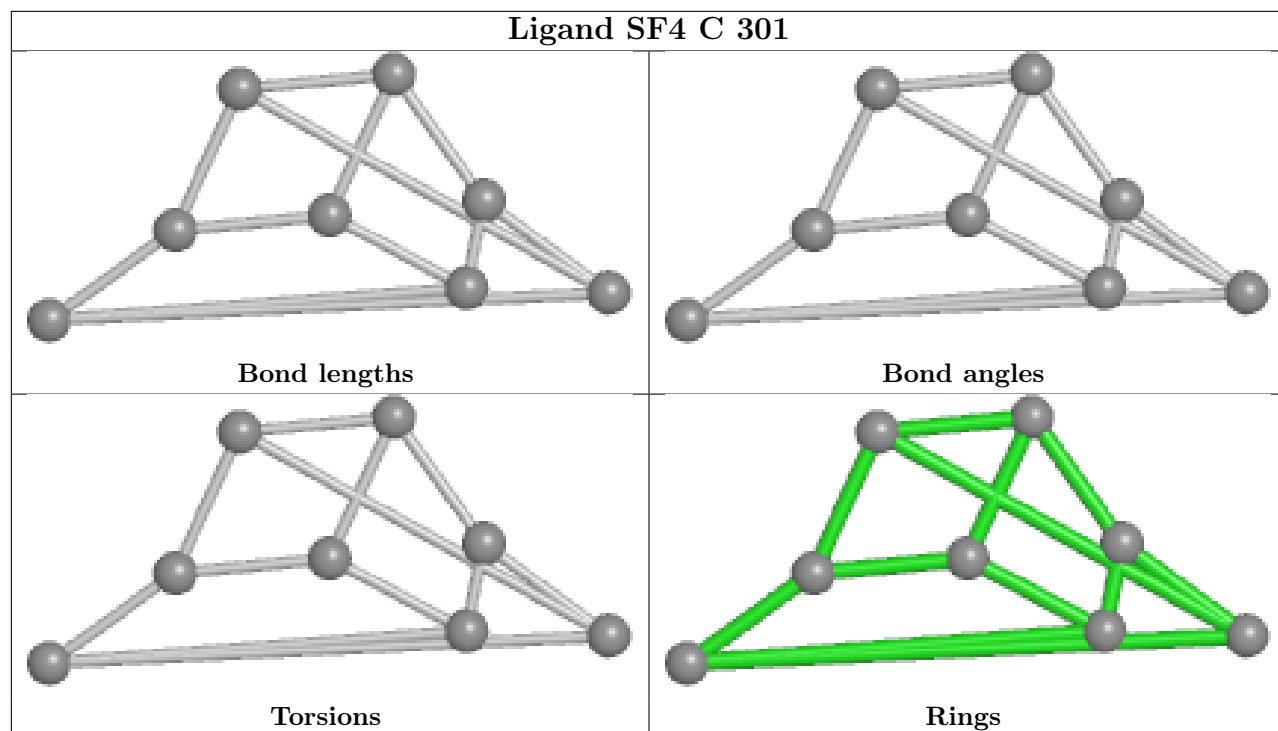




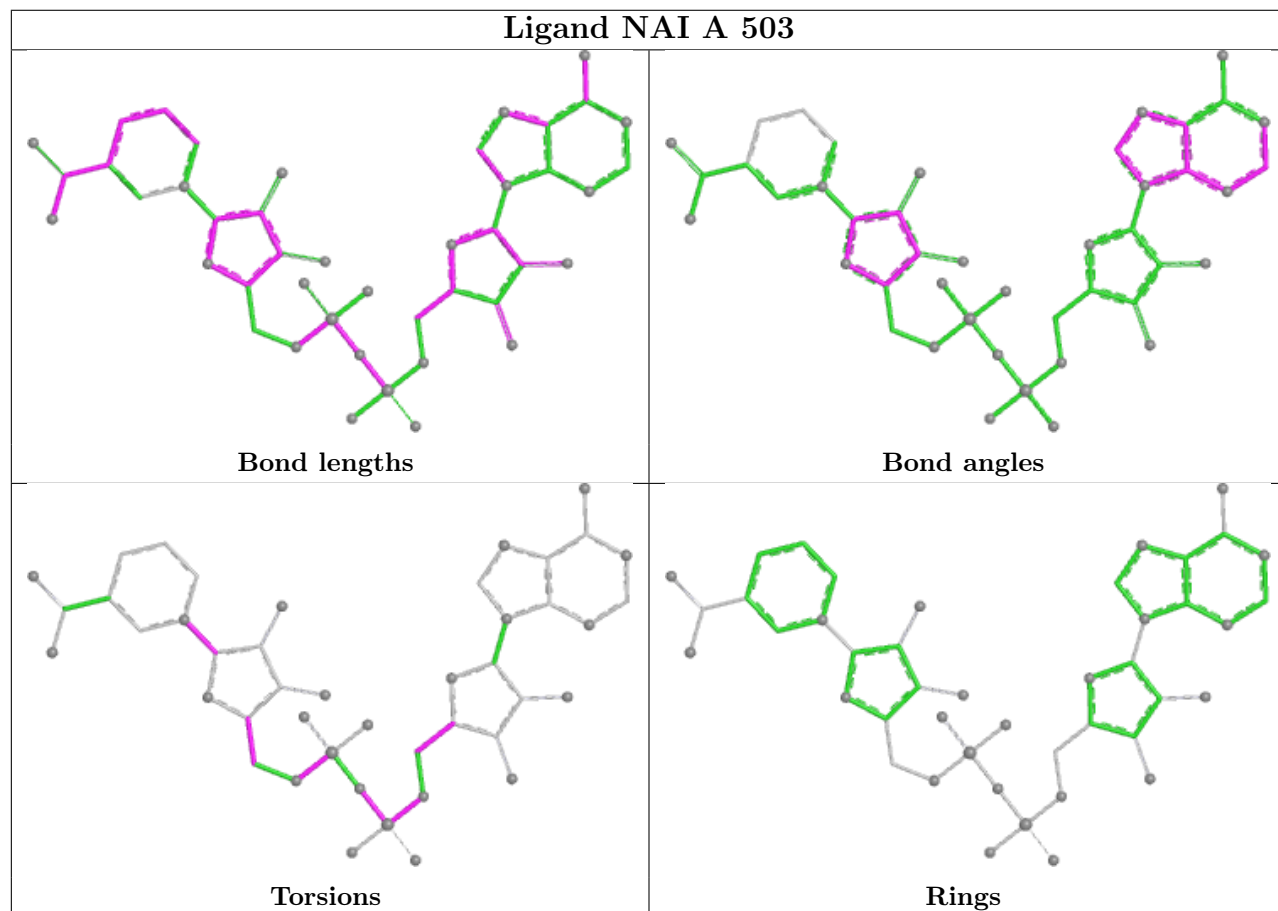


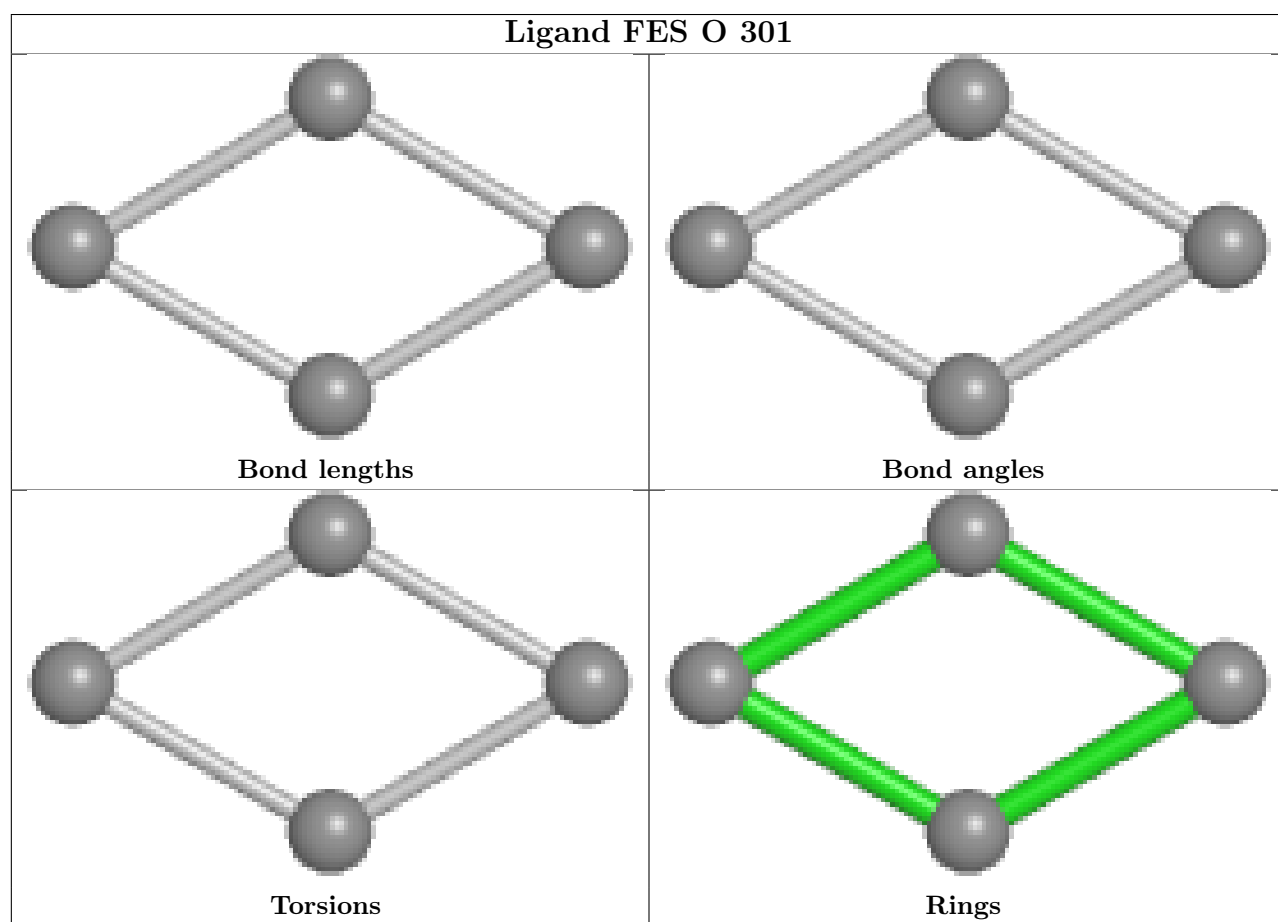
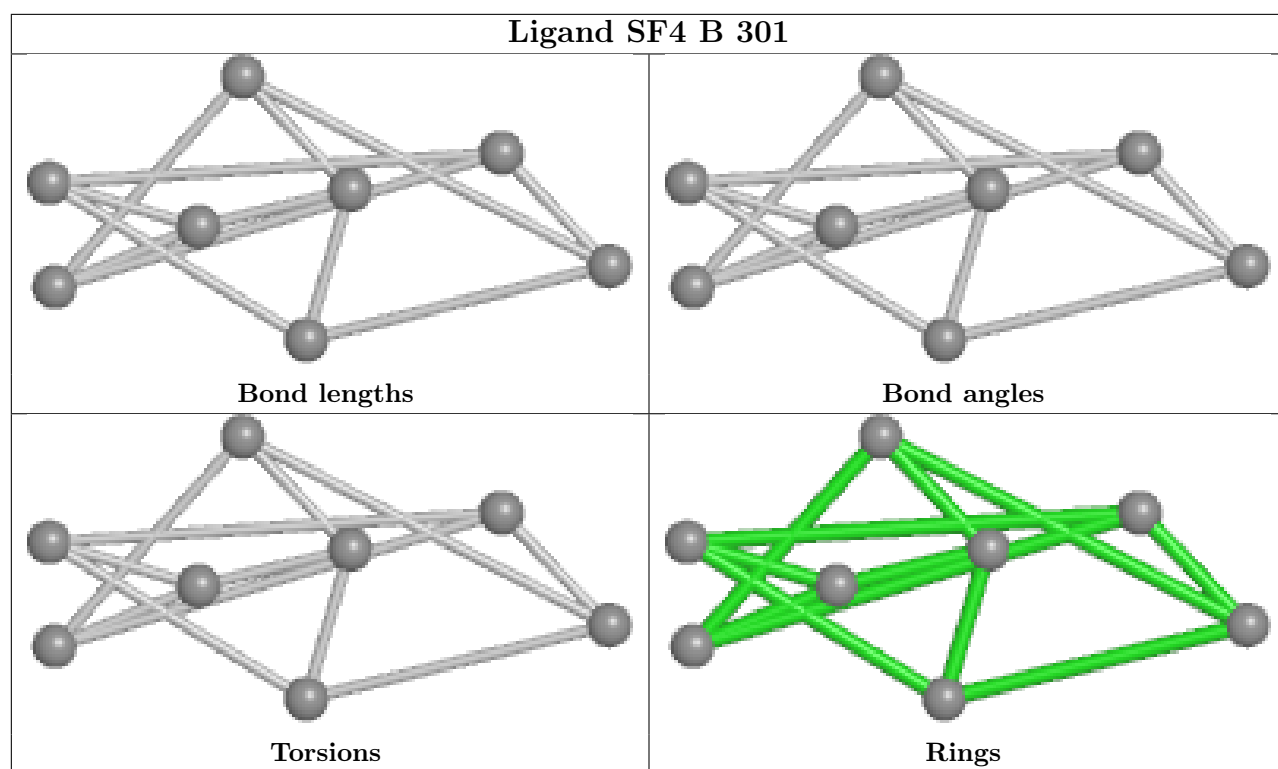


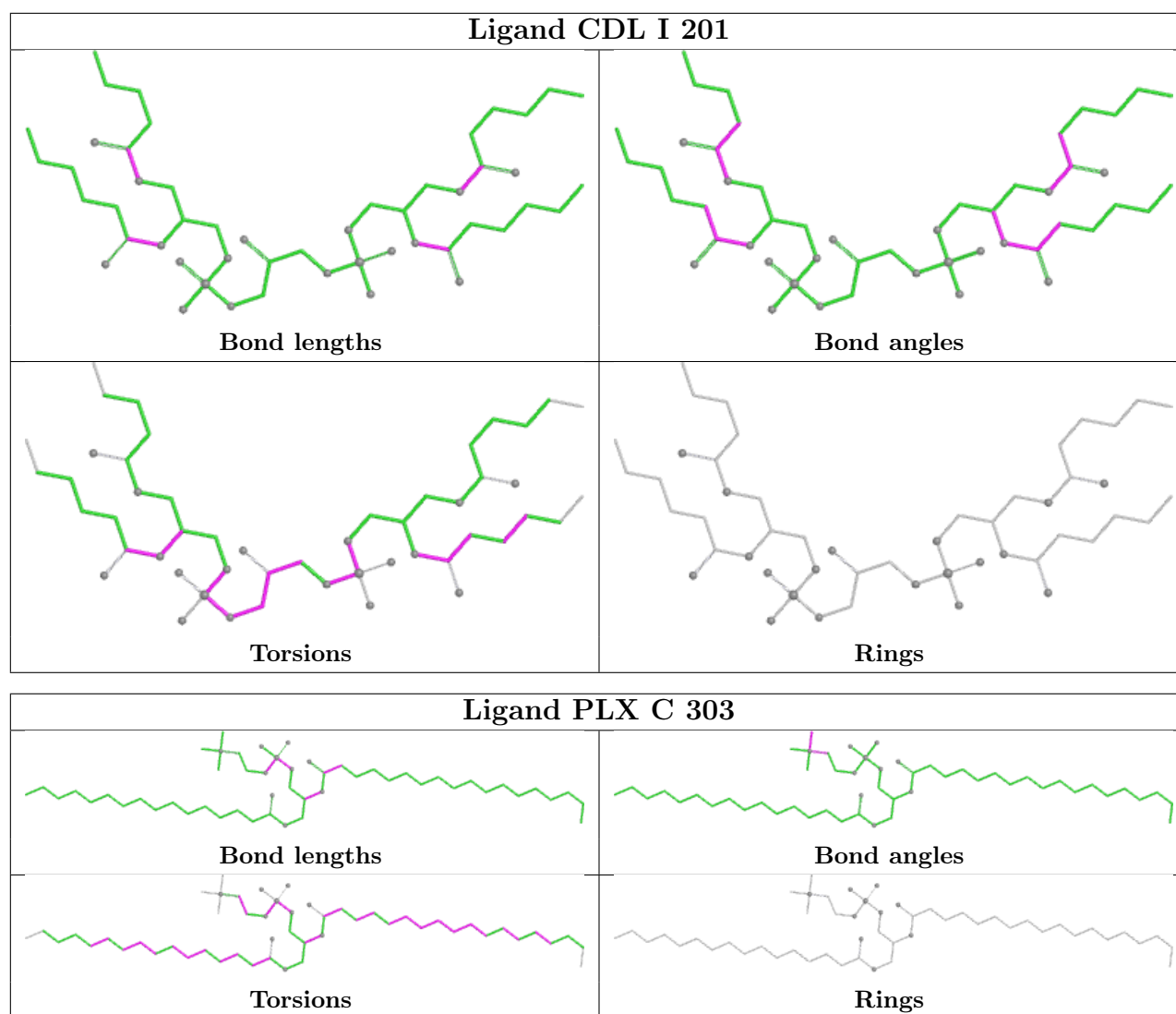
Ligand SF4 C 301



Ligand NAI A 503







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.

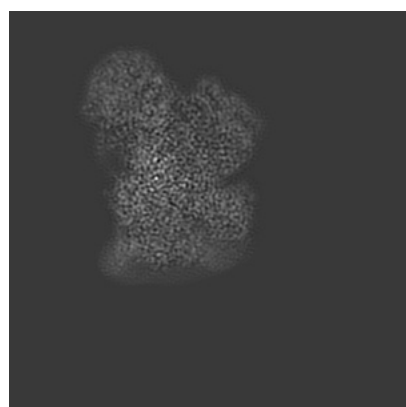
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32154. These allow visual inspection of the internal detail of the map and identification of artifacts.

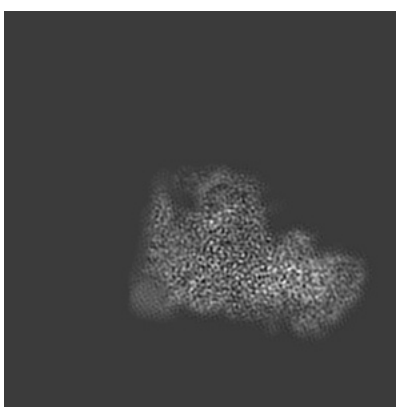
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

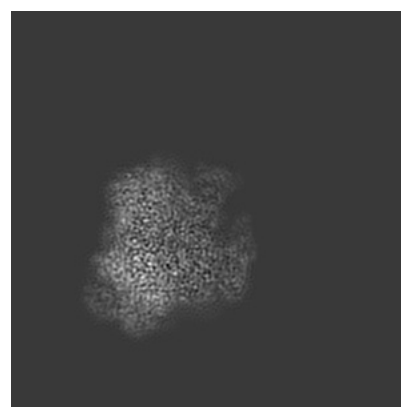
6.1.1 Primary map



X



Y



Z

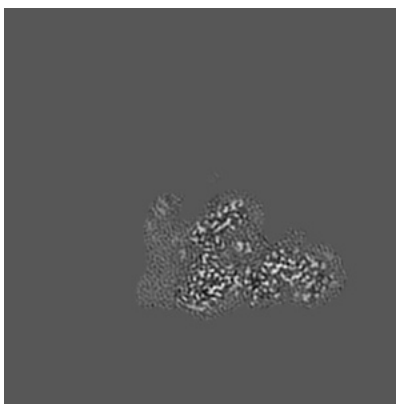
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

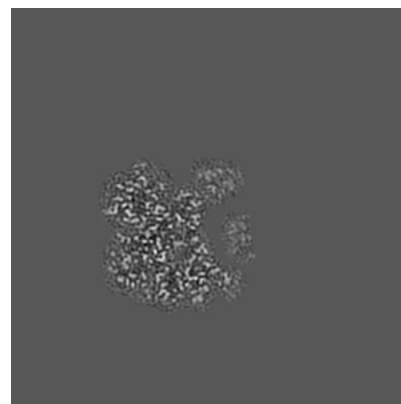
6.2.1 Primary map



X Index: 256



Y Index: 256

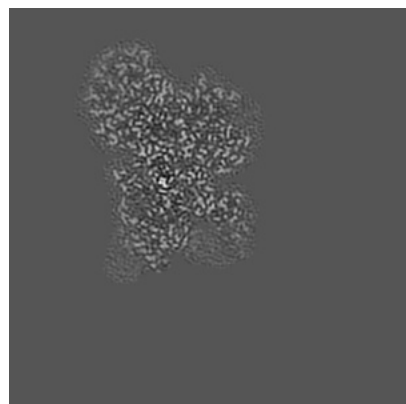


Z Index: 256

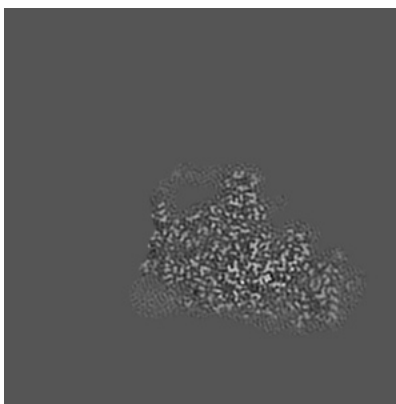
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

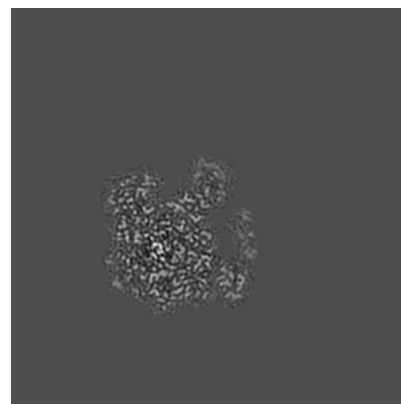
6.3.1 Primary map



X Index: 174



Y Index: 187

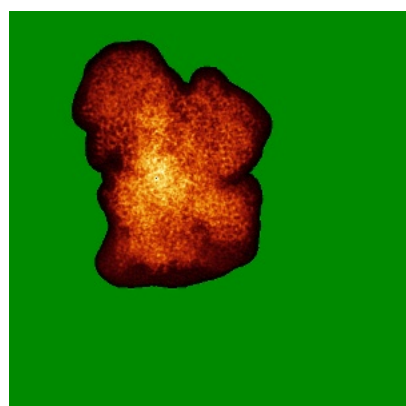


Z Index: 273

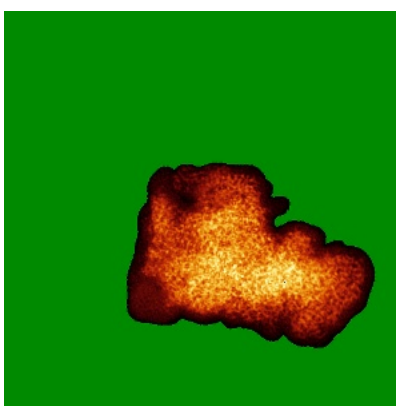
The images above show the largest variance slices of the map in three orthogonal directions.

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

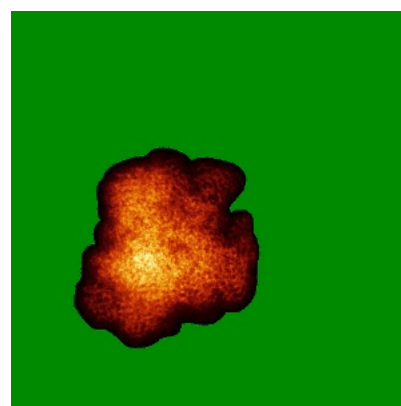
6.4.1 Primary map



X



Y

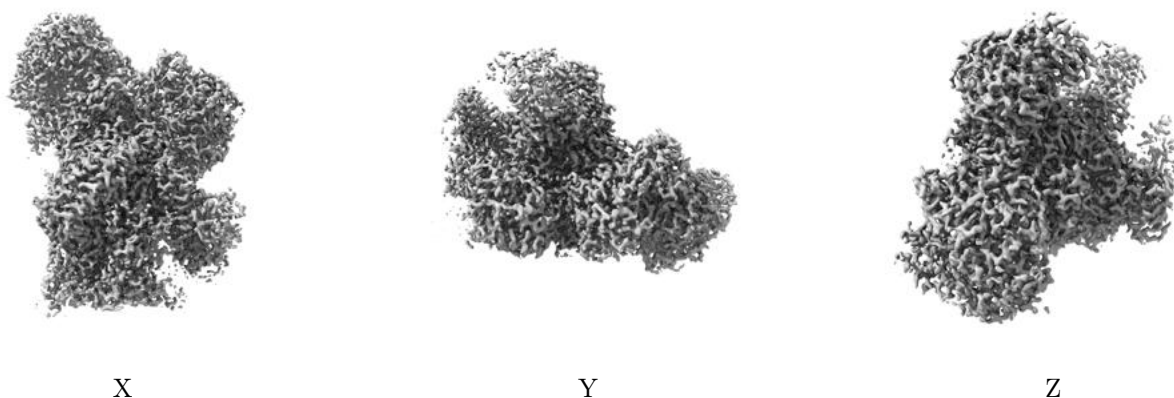


Z

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 0.0165. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

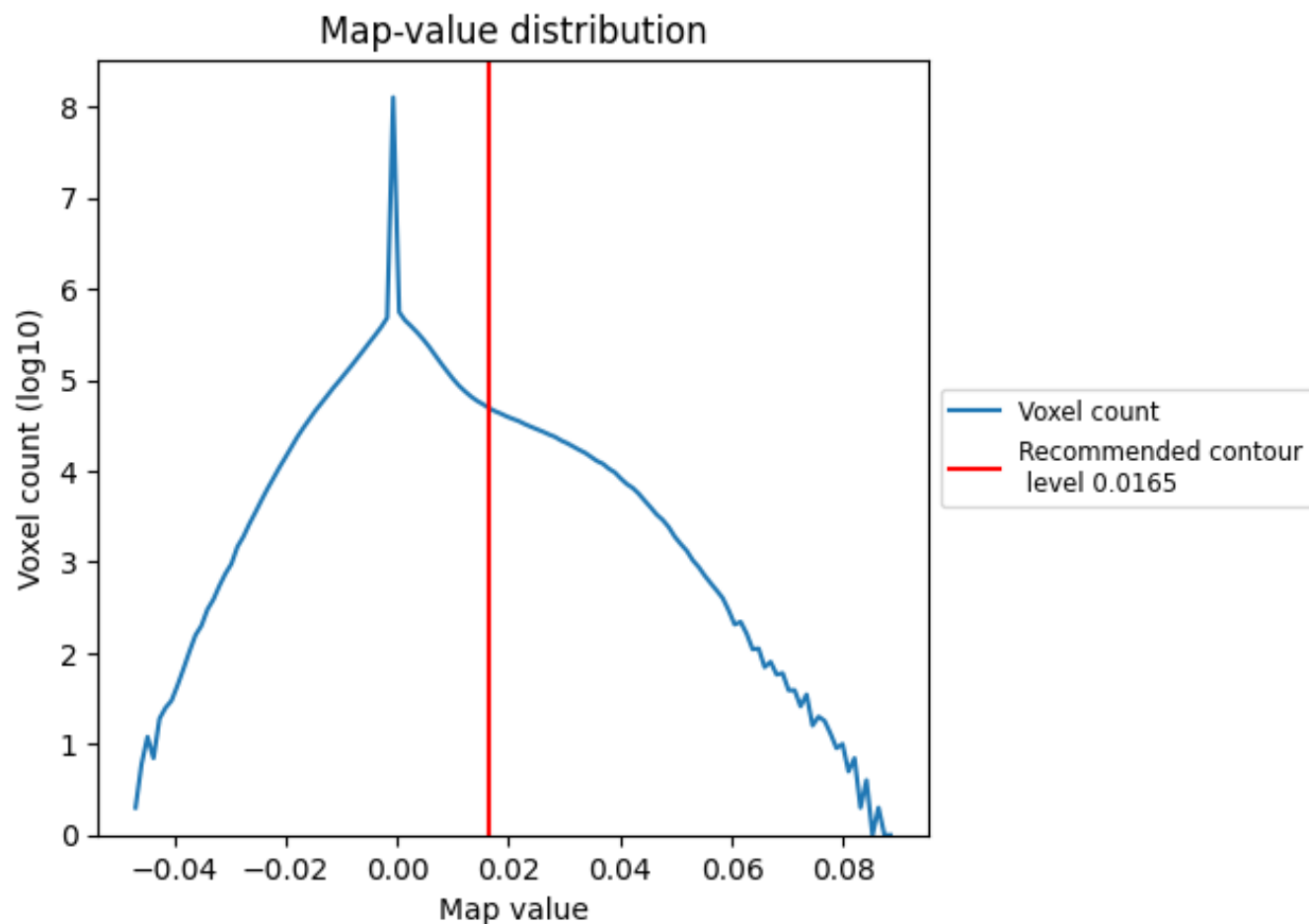
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

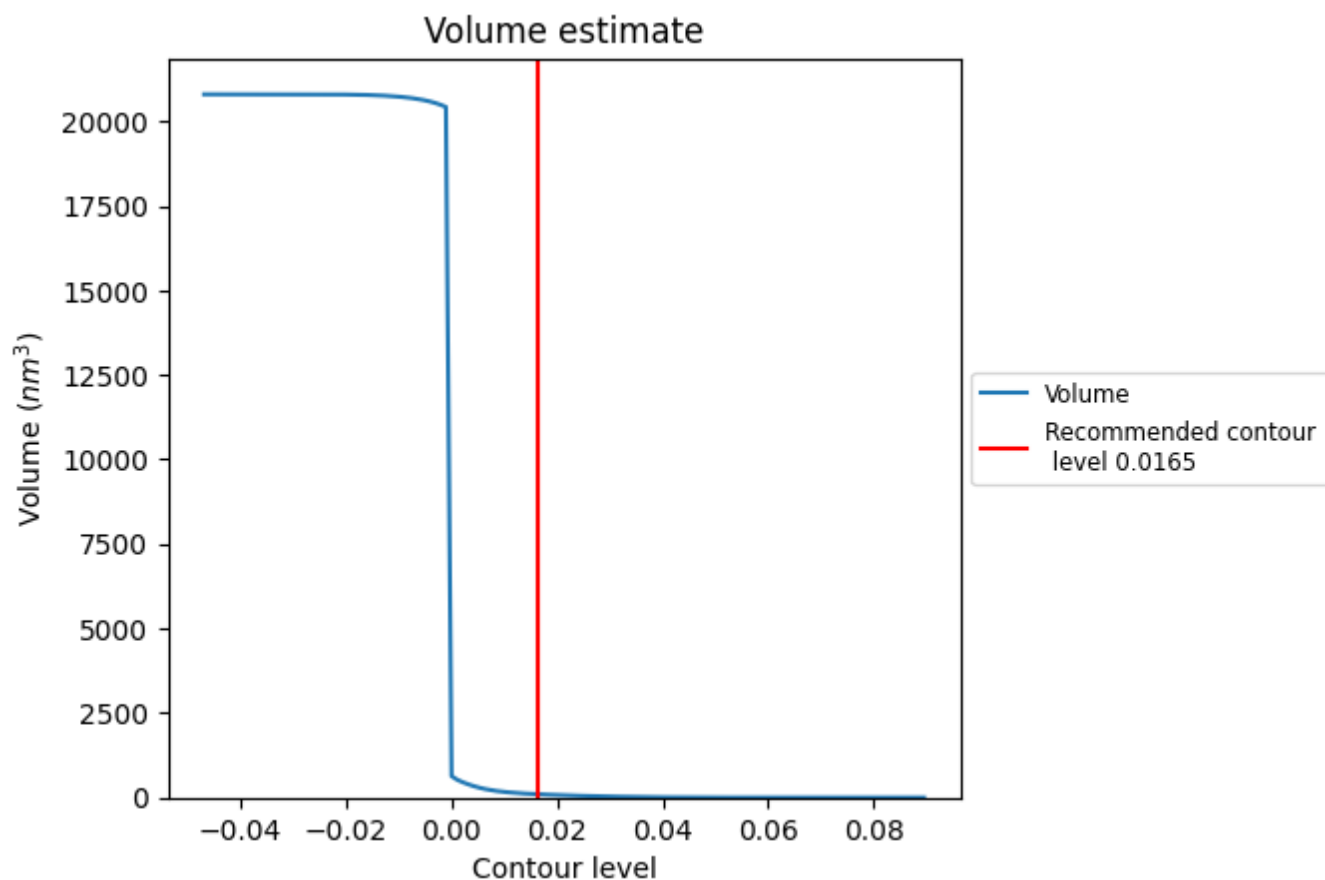
This section contains the results of statistical analysis of the map.

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.

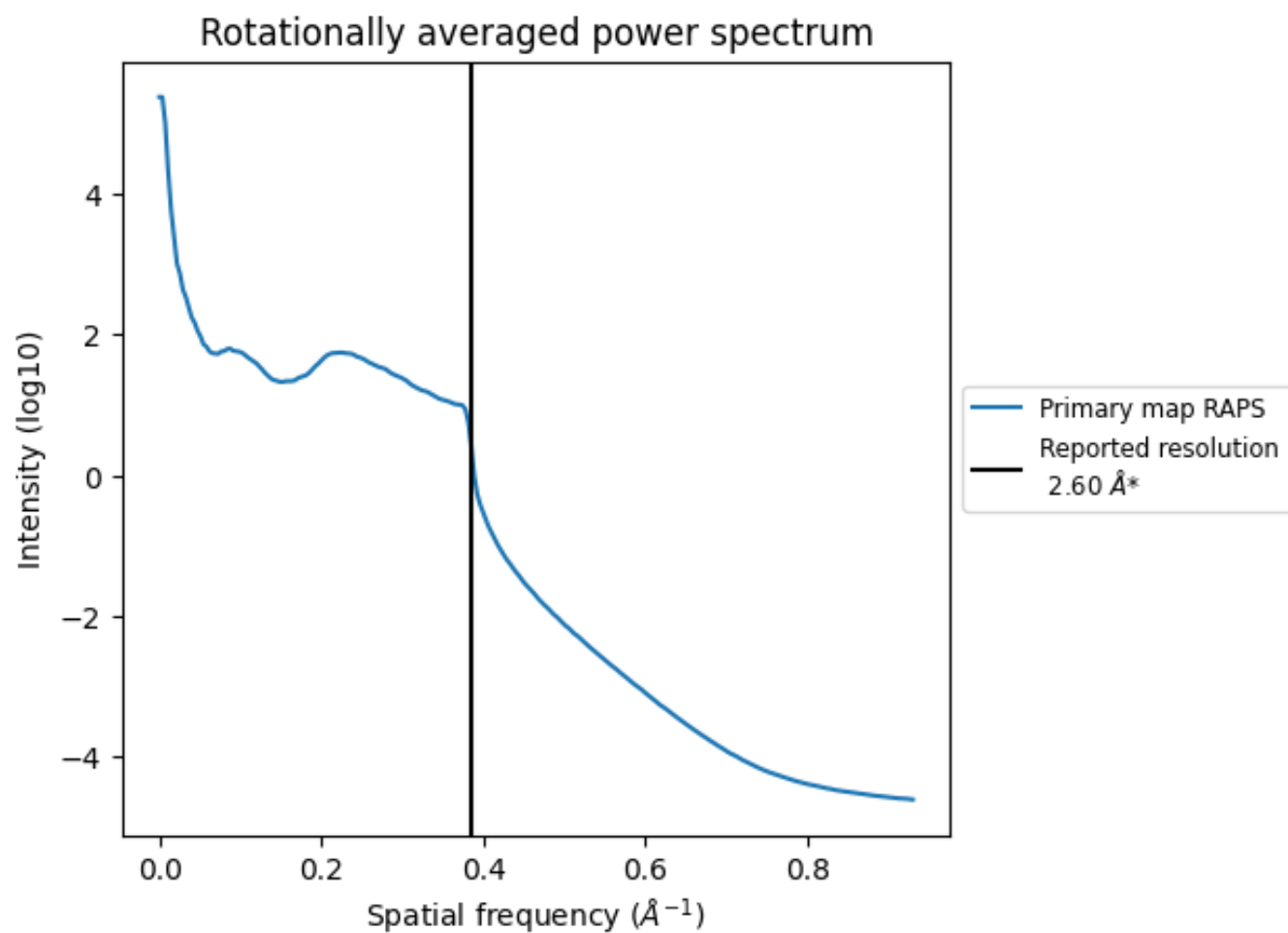
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 97 nm^3 ; this corresponds to an approximate mass of 88 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 0.385 Å⁻¹

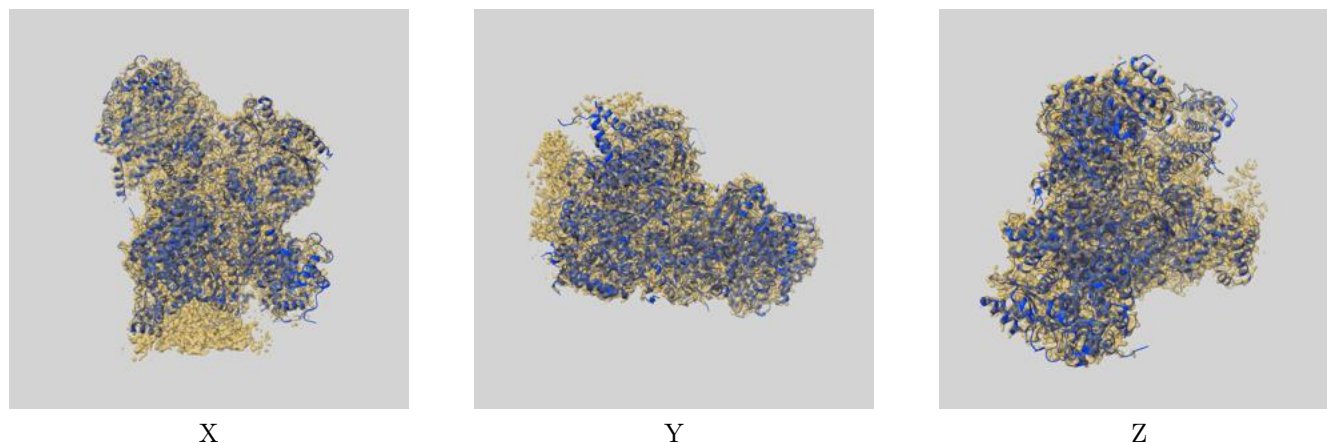
8 Fourier-Shell correlation ⓘ

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

9 Map-model fit [i](#)

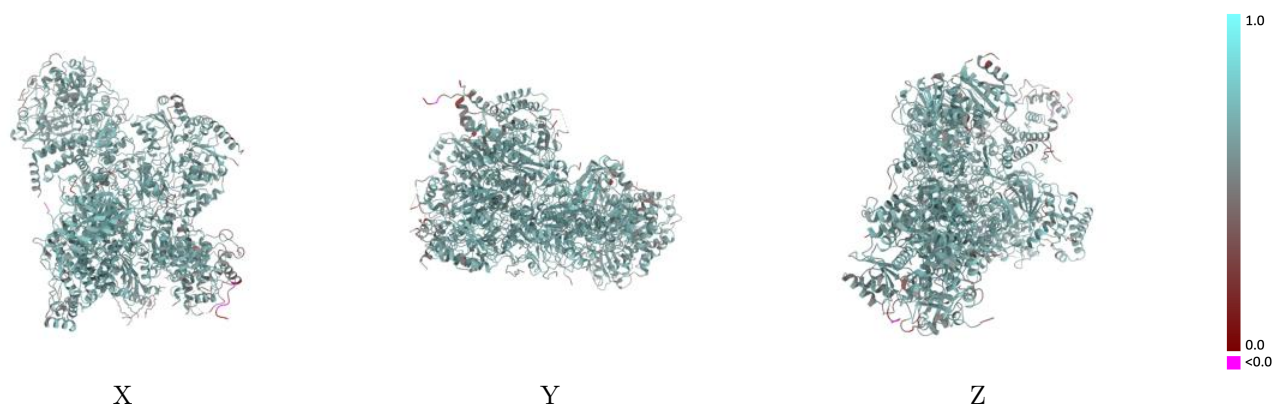
This section contains information regarding the fit between EMDB map EMD-32154 and PDB model 7VWJ. Per-residue inclusion information can be found in section [3](#) on page [14](#).

9.1 Map-model overlay [i](#)



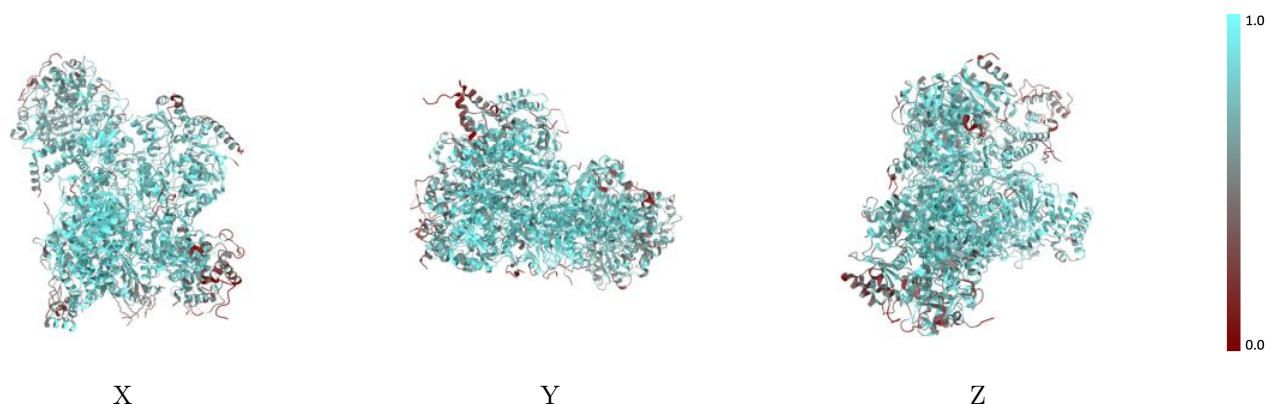
The images above show the 3D surface view of the map at the recommended contour level 0.0165 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



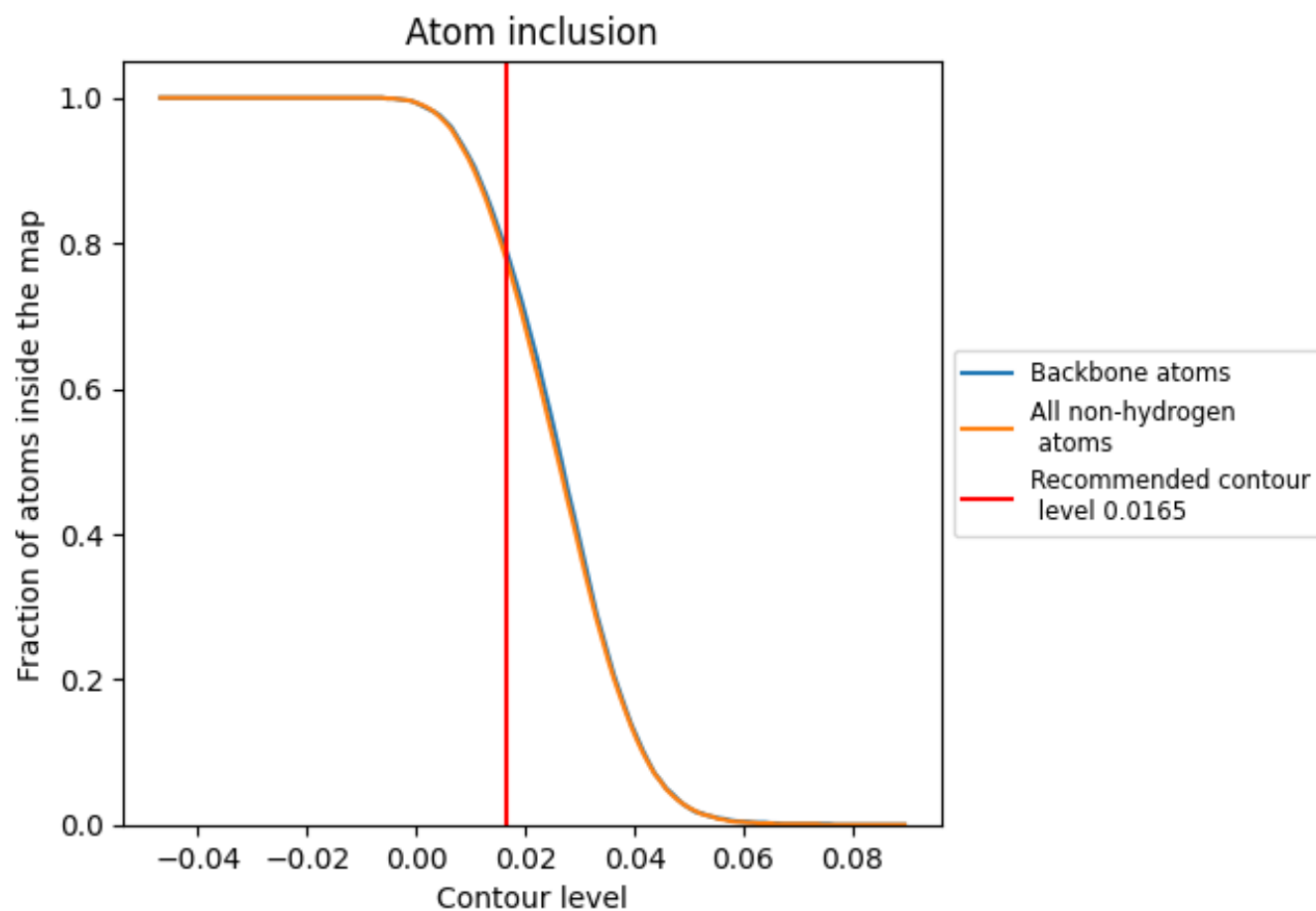
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0165).

9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 78% 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 (0.0165) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7780	 0.6310
A	 0.7440	 0.6220
B	 0.9070	 0.6670
C	 0.8330	 0.6450
E	 0.7820	 0.6290
F	 0.6760	 0.5890
G	 0.3710	 0.4750
H	 0.7640	 0.6190
I	 0.6420	 0.5880
J	 0.7580	 0.6190
K	 0.5230	 0.5520
L	 0.8170	 0.6410
M	 0.8390	 0.6500
N	 0.7020	 0.6240
O	 0.6350	 0.5950
P	 0.9060	 0.6670
Q	 0.9070	 0.6710
T	 0.7380	 0.6210
W	 0.5930	 0.5710

