



Full wwPDB EM Validation Report ⓘ

Mar 9, 2026 – 05:01 AM UTC

PDB ID : 8G0E / pdb_00008g0e
EMDB ID : EMD-29655
Title : Cryo-EM structure of TBAJ-876-bound Mycobacterium smegmatis ATP synthase rotational state 3
Authors : Courbon, G.M.; Rubinstein, J.L.
Deposited on : 2023-01-31
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

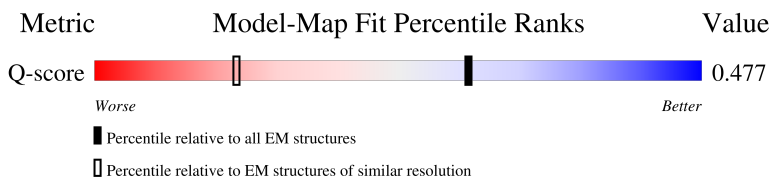
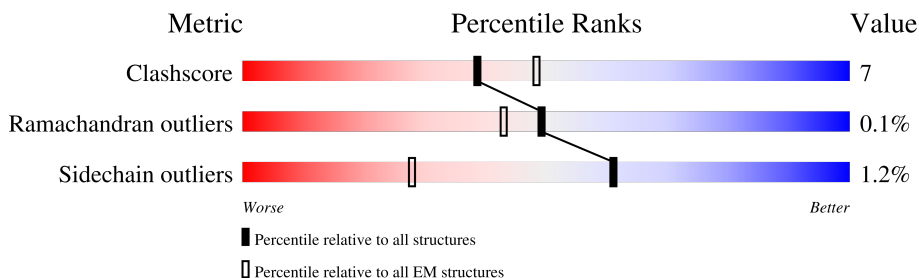
1 Overall quality at a glance

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.



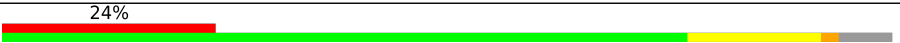
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	8728 (2.10 - 3.10)

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	548	 81% 12% • 6%
1	B	548	 78% 13% • 8%
1	C	548	 84% 9% 6%
2	D	475	 84% 12% ••

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Mol	Chain	Length	Quality of chain
2	E	475	
2	F	475	
3	G	307	
4	H	121	
5	b	170	
6	d	445	
7	1	86	
7	2	86	
7	3	86	
7	4	86	
7	5	86	
7	6	86	
7	7	86	
7	8	86	
7	9	86	
8	a	252	

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 35959 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 ATP synthase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	514	Total	C	N	O	S	0	0
			3837	2416	671	740	10		
1	B	506	Total	C	N	O	S	0	0
			3735	2363	660	702	10		
1	C	514	Total	C	N	O	S	0	0
			3781	2396	674	701	10		

- Molecule 2 is a protein called ATP synthase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	464	Total	C	N	O	S	0	0
			3508	2225	608	662	13		
2	E	464	Total	C	N	O	S	0	0
			3522	2230	609	670	13		
2	F	464	Total	C	N	O	S	0	0
			3512	2225	609	665	13		

- Molecule 3 is a protein called ATP synthase gamma chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	279	Total	C	N	O	S	0	0
			2103	1324	393	380	6		

- Molecule 4 is a protein called ATP synthase epsilon chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	117	Total	C	N	O		0	0
			835	528	160	147			

- Molecule 5 is a protein called ATP synthase subunit b.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	b	141	Total	C	N	O	S	0	0
			966	612	178	174	2		

- Molecule 6 is a protein called ATP synthase subunit b-delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	d	432	Total	C	N	O	S	0	0
			2893	1825	555	510	3		

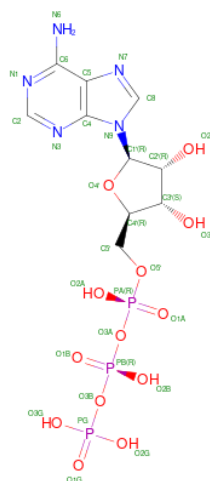
- Molecule 7 is a protein called ATP synthase subunit c.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	9	81	Total	C	N	O	S	0	0
			560	370	92	96	2		
7	1	81	Total	C	N	O	S	0	0
			556	368	91	95	2		
7	2	81	Total	C	N	O	S	0	0
			560	370	92	96	2		
7	3	81	Total	C	N	O	S	0	0
			554	367	89	96	2		
7	4	81	Total	C	N	O	S	0	0
			564	372	92	98	2		
7	5	81	Total	C	N	O	S	0	0
			560	370	92	96	2		
7	6	81	Total	C	N	O	S	0	0
			560	370	92	96	2		
7	7	81	Total	C	N	O	S	0	0
			556	368	91	95	2		
7	8	81	Total	C	N	O	S	0	0
			556	368	91	95	2		

- Molecule 8 is a protein called ATP synthase subunit a.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	a	234	Total	C	N	O	S	0	0
			1807	1222	286	293	6		

- Molecule 9 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

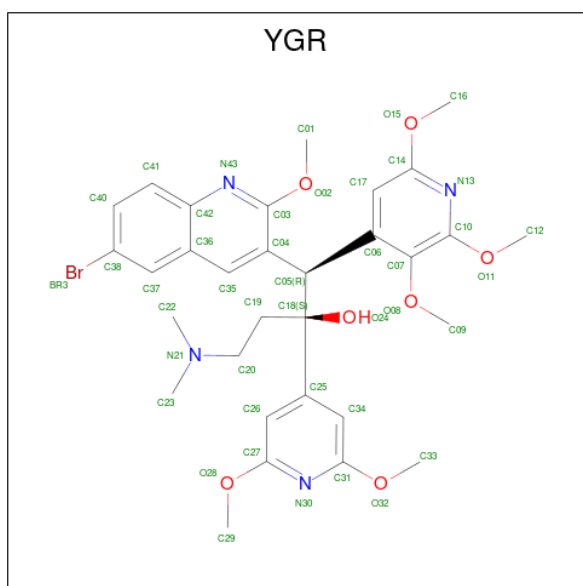
Mol	Chain	Residues	Atoms	AltConf
10	A	1	Total Mg 1 1	0
10	B	1	Total Mg 1 1	0
10	C	1	Total Mg 1 1	0
10	E	1	Total Mg 1 1	0

- Molecule 11 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			AltConf
11	F	1	Total	O	P	0
			5	4	1	

- Molecule 12 is (1R,2S)-1-(6-bromo-2-methoxyquinolin-3-yl)-2-(2,6-dimethoxypyridin-4-yl)-4-(dimethylamino)-1-(2,3,6-trimethoxypyridin-4-yl)butan-2-ol (CCD ID: YGR) (formula: $C_{31}H_{37}BrN_4O_7$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
12	9	1	Total	Br	C	N	O	0
			43	1	31	4	7	
12	3	1	Total	Br	C	N	O	0
			43	1	31	4	7	

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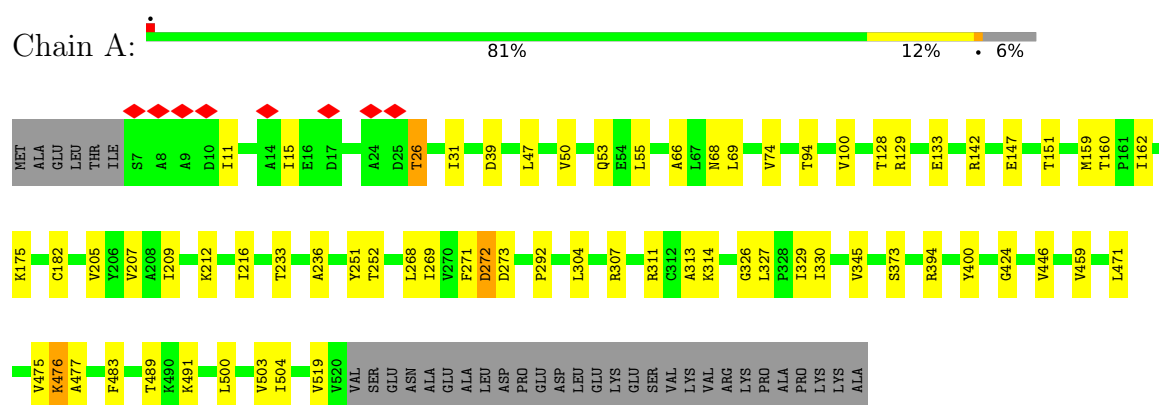
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Mol	Chain	Residues	Atoms					AltConf
12	4	1	Total 43	Br 1	C 31	N 4	O 7	0
12	5	1	Total 43	Br 1	C 31	N 4	O 7	0
12	6	1	Total 43	Br 1	C 31	N 4	O 7	0
12	7	1	Total 43	Br 1	C 31	N 4	O 7	0
12	8	1	Total 43	Br 1	C 31	N 4	O 7	0

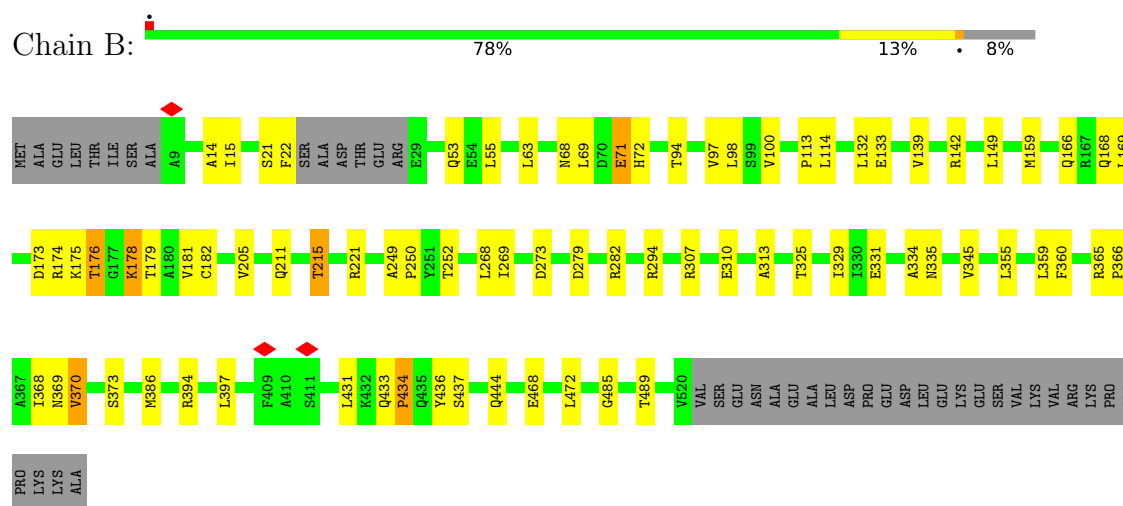
3 Residue-property plots

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.

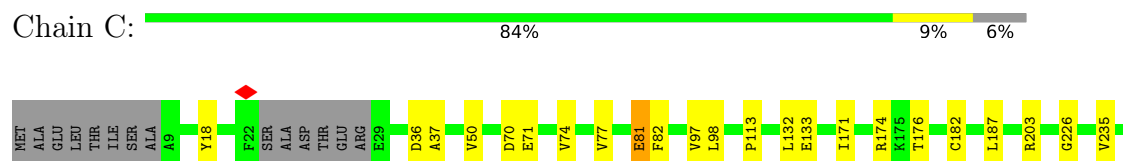
• Molecule 1: ATP synthase subunit alpha

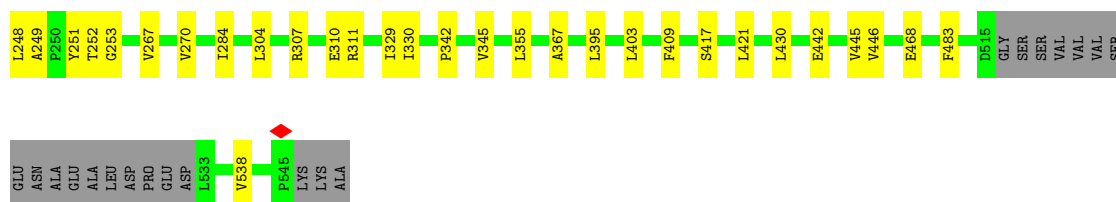


• Molecule 1: ATP synthase subunit alpha



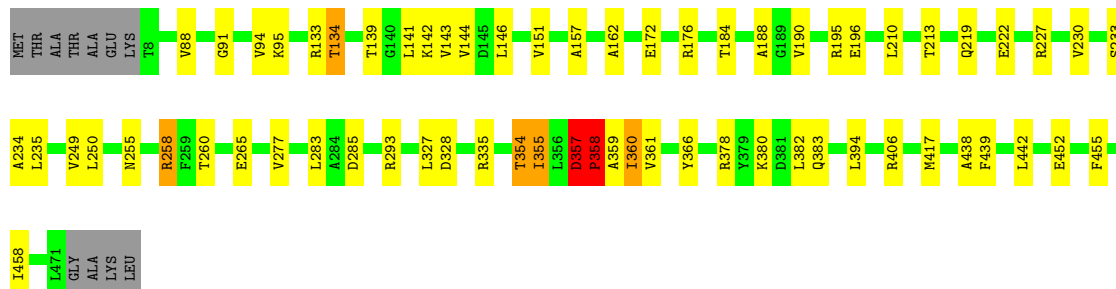
• Molecule 1: ATP synthase subunit alpha





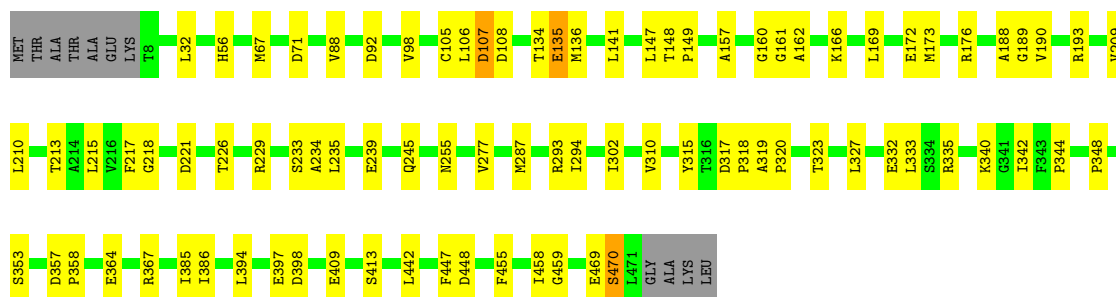
• Molecule 2: ATP synthase subunit beta

Chain D: 84% 12% ..



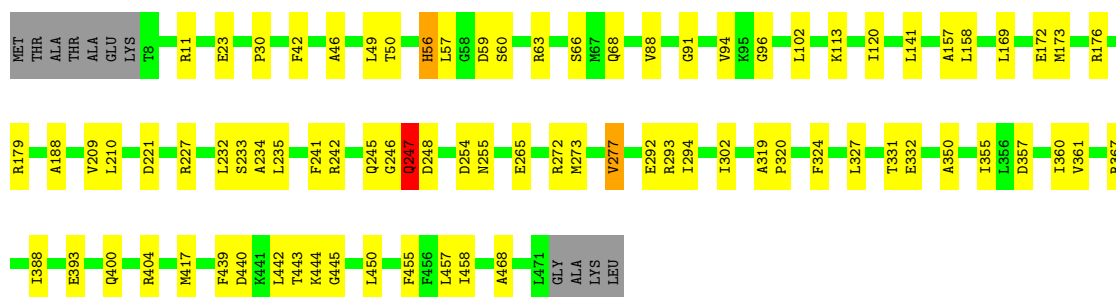
• Molecule 2: ATP synthase subunit beta

Chain E: 80% 17% ..



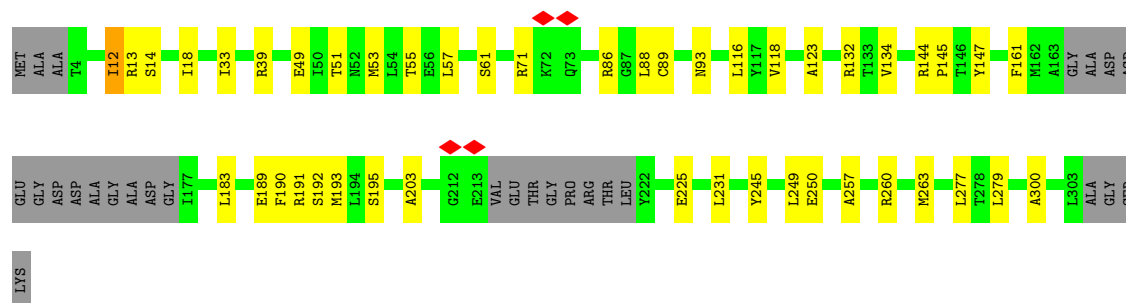
• Molecule 2: ATP synthase subunit beta

Chain F: 80% 17% .

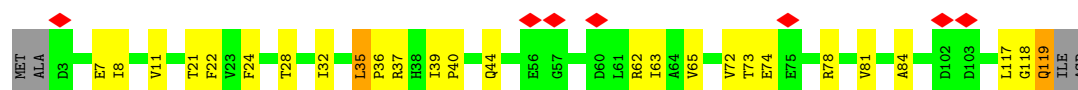
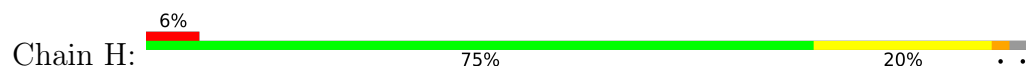


• Molecule 3: ATP synthase gamma chain

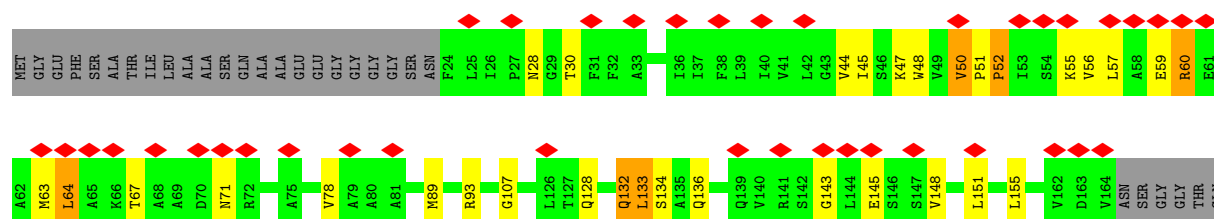
Chain G: 76% 14% 9%



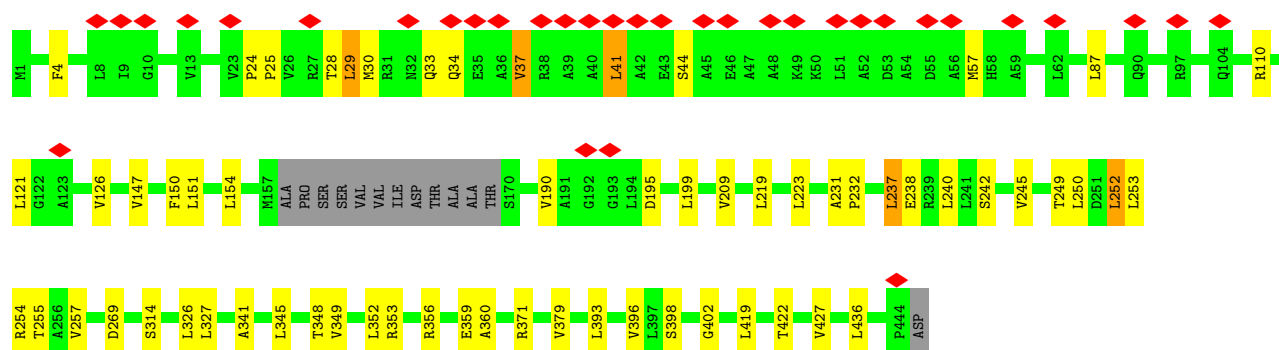
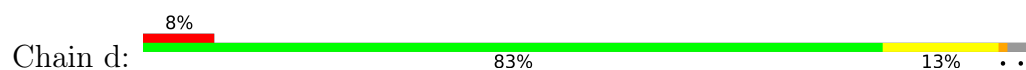
• Molecule 4: ATP synthase epsilon chain



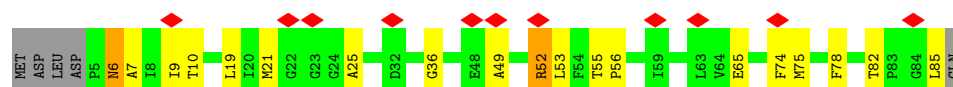
• Molecule 5: ATP synthase subunit b



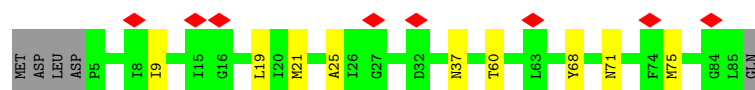
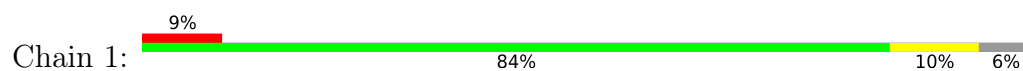
• Molecule 6: ATP synthase subunit b-delta



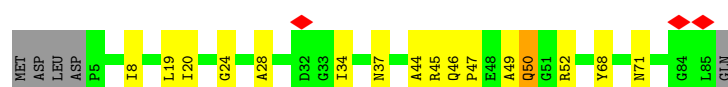
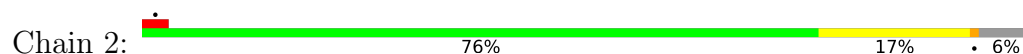
• Molecule 7: ATP synthase subunit c



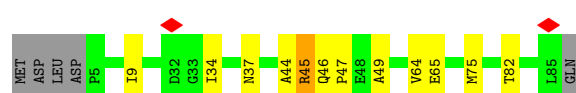
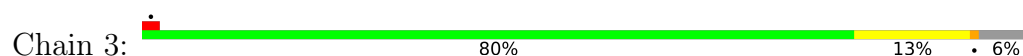
● Molecule 7: ATP synthase subunit c



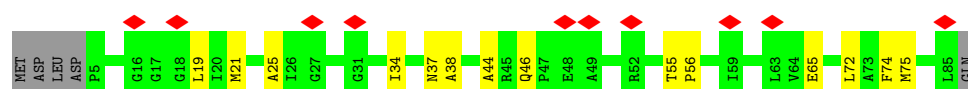
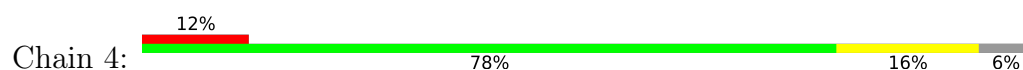
● Molecule 7: ATP synthase subunit c



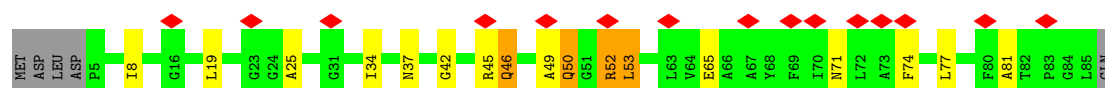
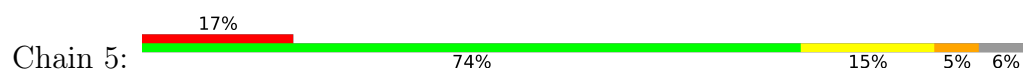
● Molecule 7: ATP synthase subunit c



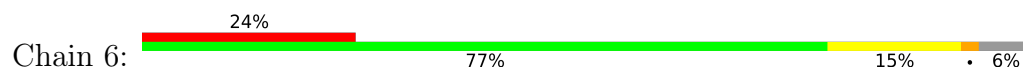
● Molecule 7: ATP synthase subunit c



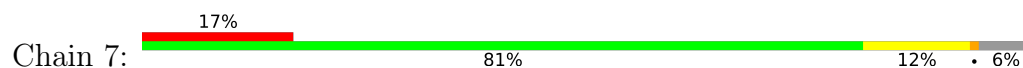
● Molecule 7: ATP synthase subunit c

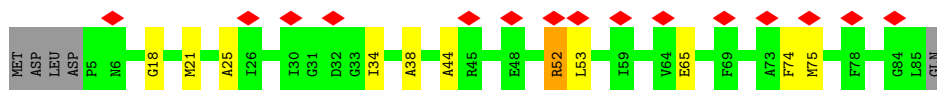


● Molecule 7: ATP synthase subunit c

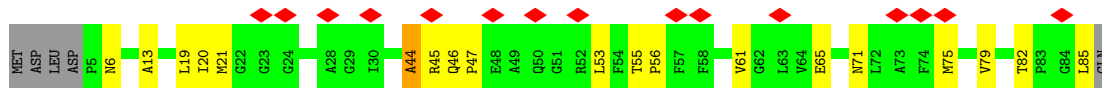
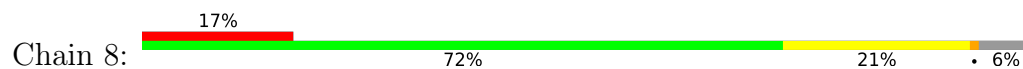


● Molecule 7: ATP synthase subunit c

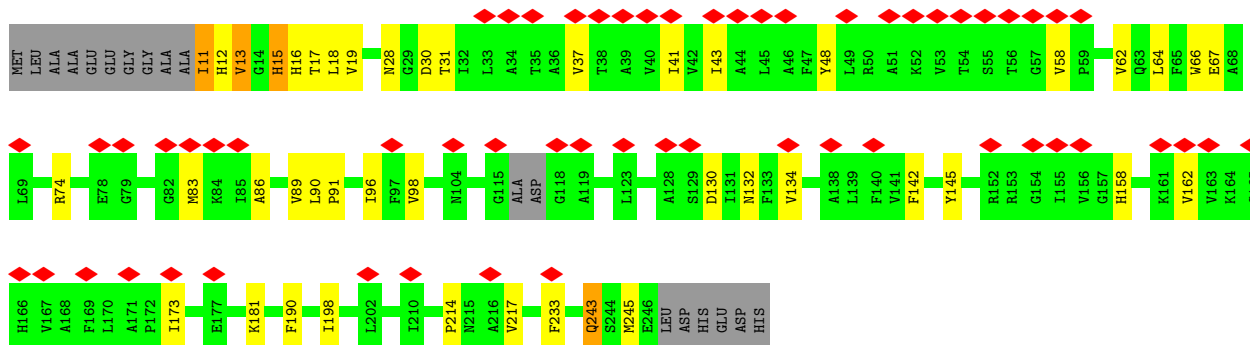
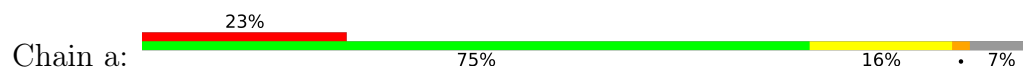




- Molecule 7: ATP synthase subunit c



- Molecule 8: ATP synthase subunit a



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	166169	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	44	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	2100	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	3.362	Depositor
Minimum map value	-0.699	Depositor
Average map value	0.013	Depositor
Map value standard deviation	0.126	Depositor
Recommended contour level	0.7	Depositor
Map size (Å)	329.59998, 329.59998, 329.59998	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.03, 1.03, 1.03	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: PO4, MG, YGR, ATP

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.41	6/3897 (0.2%)	0.49	2/5280 (0.0%)
1	B	0.80	28/3793 (0.7%)	0.64	4/5140 (0.1%)
1	C	0.35	4/3840 (0.1%)	0.48	1/5202 (0.0%)
2	D	0.48	9/3573 (0.3%)	0.56	3/4848 (0.1%)
2	E	0.48	6/3587 (0.2%)	0.53	4/4867 (0.1%)
2	F	0.44	1/3577 (0.0%)	0.52	2/4854 (0.0%)
3	G	0.44	1/2131 (0.0%)	0.60	4/2883 (0.1%)
4	H	0.31	1/846 (0.1%)	0.52	0/1154
5	b	0.65	0/975	0.84	7/1327 (0.5%)
6	d	0.45	0/2921	0.59	3/3993 (0.1%)
7	1	0.18	0/567	0.35	0/769
7	2	0.29	0/571	0.42	0/774
7	3	0.25	0/565	0.55	1/767 (0.1%)
7	4	0.18	0/575	0.35	0/779
7	5	0.65	2/571 (0.4%)	0.59	1/774 (0.1%)
7	6	0.47	2/571 (0.4%)	0.54	1/774 (0.1%)
7	7	0.52	0/567	0.49	0/769
7	8	0.69	3/567 (0.5%)	0.50	1/769 (0.1%)
7	9	0.59	1/571 (0.2%)	0.49	1/774 (0.1%)
8	a	0.36	0/1859	0.46	2/2541 (0.1%)
All	All	0.49	64/36124 (0.2%)	0.55	37/49038 (0.1%)

All (64) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	365	ARG	CA-C	-9.36	1.44	1.52
1	B	71	GLU	CA-C	-9.16	1.41	1.52
1	B	173	ASP	CA-C	-8.61	1.41	1.53
1	B	369	ASN	CA-C	-8.34	1.42	1.52
7	8	44	ALA	CA-C	-7.87	1.43	1.52
2	D	357	ASP	CA-C	-7.86	1.44	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	176	THR	CA-C	-7.26	1.42	1.52
1	B	433	GLN	CA-C	-7.17	1.44	1.52
1	A	476	LYS	CA-C	-7.15	1.43	1.52
1	B	179	THR	CA-C	-6.95	1.43	1.52
1	B	175	LYS	CA-C	-6.94	1.44	1.53
1	B	176	THR	CA-CB	-6.90	1.41	1.53
1	B	433	GLN	C-O	-6.78	1.16	1.24
1	C	81	GLU	CA-C	-6.77	1.44	1.52
1	B	215	THR	CA-C	-6.71	1.43	1.52
1	B	434	PRO	CA-CB	-6.67	1.44	1.53
1	A	272	ASP	CA-C	-6.66	1.45	1.53
2	D	134	THR	CA-C	-6.64	1.44	1.52
1	C	81	GLU	N-CA	-6.55	1.38	1.46
1	B	178	LYS	CA-C	-6.50	1.44	1.52
7	5	46	GLN	CA-C	-6.50	1.44	1.52
7	6	46	GLN	N-CA	-6.42	1.39	1.46
7	8	46	GLN	CA-C	-6.39	1.46	1.52
1	B	178	LYS	C-O	-6.36	1.16	1.24
1	B	211	GLN	C-O	-6.17	1.16	1.23
1	A	477	ALA	CA-C	-6.13	1.44	1.52
1	B	176	THR	N-CA	-6.09	1.38	1.46
7	8	44	ALA	N-CA	-6.07	1.39	1.46
2	E	398	ASP	CA-C	-6.04	1.44	1.52
1	A	272	ASP	C-O	-6.03	1.16	1.23
2	E	135	GLU	CA-C	-6.00	1.45	1.52
1	A	272	ASP	CA-CB	-5.99	1.43	1.52
1	B	175	LYS	CA-CB	-5.94	1.44	1.53
1	C	82	PHE	N-CA	-5.86	1.38	1.46
2	D	358	PRO	CA-CB	-5.83	1.45	1.53
1	B	370	VAL	CA-CB	-5.82	1.47	1.54
2	E	134	THR	CA-C	-5.81	1.47	1.53
1	B	175	LYS	N-CA	-5.77	1.38	1.46
7	9	52	ARG	CA-C	-5.73	1.44	1.52
1	C	82	PHE	CA-C	-5.64	1.44	1.52
1	B	365	ARG	C-O	-5.63	1.18	1.25
1	B	173	ASP	C-O	-5.59	1.16	1.23
2	E	315	TYR	CA-C	-5.43	1.45	1.52
7	5	52	ARG	CA-C	-5.41	1.45	1.52
7	6	46	GLN	CA-CB	-5.41	1.46	1.53
2	D	354	THR	C-O	-5.36	1.16	1.23
2	D	354	THR	CA-C	-5.35	1.45	1.52
1	A	475	VAL	C-O	-5.34	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	173	ASP	CA-CB	-5.32	1.45	1.53
1	B	369	ASN	CA-CB	-5.32	1.46	1.53
2	F	248	ASP	N-CA	-5.31	1.39	1.46
2	D	357	ASP	C-O	-5.30	1.17	1.24
3	G	12	ILE	CA-C	-5.29	1.44	1.52
2	D	355	ILE	C-O	-5.27	1.18	1.24
4	H	35	LEU	CA-CB	-5.25	1.47	1.54
1	B	436	TYR	CA-C	-5.25	1.45	1.53
1	B	174	ARG	CA-C	-5.24	1.46	1.52
2	E	315	TYR	C-O	-5.22	1.17	1.24
2	D	360	ILE	CA-C	-5.16	1.45	1.52
2	D	357	ASP	CA-CB	-5.08	1.46	1.54
2	E	136	MET	CA-C	-5.05	1.46	1.52
1	B	433	GLN	CA-CB	-5.04	1.45	1.53
1	B	359	LEU	C-N	-5.04	1.27	1.34
1	B	179	THR	C-O	-5.01	1.18	1.24

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	359	ALA	N-CA-C	11.54	123.86	111.28
1	B	178	LYS	N-CA-C	9.91	122.09	111.28
1	B	71	GLU	N-CA-C	-8.82	101.62	111.14
5	b	71	ASN	N-CA-C	-8.53	101.48	112.23
3	G	195	SER	N-CA-C	7.27	120.78	109.07
5	b	136	GLN	N-CA-C	-7.15	103.49	111.28
7	8	44	ALA	N-CA-C	-7.01	103.56	111.14
7	9	6	ASN	N-CA-C	-7.00	103.65	111.28
8	a	243	GLN	N-CA-C	6.75	118.29	111.07
8	a	15	HIS	N-CA-C	6.61	120.51	111.39
1	B	370	VAL	CB-CA-C	-6.45	103.43	112.14
2	E	106	LEU	N-CA-C	6.43	121.32	112.90
2	F	56	HIS	N-CA-C	-6.29	98.94	109.07
7	6	45	ARG	N-CA-C	6.18	120.02	112.23
6	d	30	MET	N-CA-C	-6.14	104.59	111.28
7	5	50	GLN	O-C-N	-6.11	115.08	122.11
2	F	247	GLN	N-CA-C	-6.08	99.81	109.23
1	A	26	THR	N-CA-CB	-5.75	102.03	110.49
5	b	60	ARG	N-CA-C	-5.74	104.93	111.07
6	d	360	ALA	N-CA-C	5.73	118.27	111.33
2	D	360	ILE	N-CA-C	-5.70	106.15	111.45
7	3	45	ARG	O-C-N	5.69	127.93	122.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	b	67	THR	N-CA-C	-5.66	105.19	111.36
6	d	41	LEU	N-CA-C	-5.63	105.14	111.28
1	C	367	ALA	N-CA-C	5.54	119.86	113.38
5	b	50	VAL	CA-C-N	-5.45	114.76	120.38
5	b	50	VAL	C-N-CA	-5.45	114.76	120.38
1	B	433	GLN	CB-CA-C	-5.44	101.56	109.45
1	A	272	ASP	CB-CA-C	-5.39	104.38	111.82
2	D	134	THR	N-CA-CB	5.39	117.89	109.97
2	E	315	TYR	N-CA-C	-5.39	106.48	113.16
2	E	397	GLU	N-CA-C	-5.38	106.88	113.50
3	G	13	ARG	N-CA-C	5.25	117.08	111.36
2	E	108	ASP	N-CA-CB	5.21	119.08	110.38
3	G	192	SER	O-C-N	-5.17	117.45	122.99
3	G	12	ILE	N-CA-C	-5.08	104.63	111.44
5	b	143	GLY	N-CA-C	-5.02	106.66	113.24

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3837	0	3834	45	0
1	B	3735	0	3730	52	0
1	C	3781	0	3787	40	0
2	D	3508	0	3495	44	0
2	E	3522	0	3503	57	0
2	F	3512	0	3497	55	0
3	G	2103	0	2129	31	0
4	H	835	0	806	22	0
5	b	966	0	936	25	0
6	d	2893	0	2767	57	0
7	1	556	0	571	15	0
7	2	560	0	577	20	0
7	3	554	0	566	18	0
7	4	564	0	581	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	5	560	0	577	18	0
7	6	560	0	577	14	0
7	7	556	0	571	13	0
7	8	556	0	571	18	0
7	9	560	0	577	23	0
8	a	1807	0	1857	38	0
9	A	31	0	12	1	0
9	B	31	0	12	0	0
9	C	31	0	12	0	0
9	E	31	0	12	1	0
10	A	1	0	0	0	0
10	B	1	0	0	0	0
10	C	1	0	0	0	0
10	E	1	0	0	0	0
11	F	5	0	0	0	0
12	3	43	0	0	1	0
12	4	43	0	0	1	0
12	5	43	0	0	1	0
12	6	43	0	0	1	0
12	7	43	0	0	1	0
12	8	43	0	0	1	0
12	9	43	0	0	1	0
All	All	35959	0	35557	520	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 7.

All (520) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:2:47:PRO:HG3	7:3:46:GLN:HE22	1.25	1.00
4:H:22:PHE:HD1	4:H:35:LEU:HG	1.28	0.99
1:C:235:VAL:HG11	1:C:252:THR:CG2	1.94	0.97
5:b:63:MET:HE1	6:d:41:LEU:HA	1.51	0.90
1:C:235:VAL:HG11	1:C:252:THR:HG23	1.55	0.89
3:G:33:ILE:HD13	3:G:193:MET:HE1	1.56	0.86
6:d:327:LEU:HD22	6:d:349:VAL:HG21	1.61	0.83
1:B:159:MET:CE	1:B:370:VAL:HG13	2.10	0.82
1:C:249:ALA:O	1:C:252:THR:HG22	1.79	0.82
1:C:113:PRO:HA	1:C:252:THR:HG21	1.64	0.79
2:D:195:ARG:NH1	2:D:196:GLU:OE2	2.17	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:169:LEU:HD21	2:E:344:PRO:HB3	1.67	0.76
4:H:22:PHE:CD1	4:H:35:LEU:HG	2.19	0.76
4:H:32:ILE:HD12	7:3:45:ARG:HH12	1.52	0.74
4:H:32:ILE:CD1	7:3:45:ARG:HH12	2.00	0.74
6:d:348:THR:O	6:d:352:LEU:HD12	1.90	0.72
7:8:6:ASN:OD1	7:8:85:LEU:HD11	1.91	0.71
2:D:157:ALA:HB2	2:D:327:LEU:HD13	1.73	0.71
7:9:82:THR:CG2	7:9:85:LEU:HD22	2.21	0.71
8:a:12:HIS:NE2	8:a:15:HIS:CG	2.59	0.71
1:B:159:MET:HE1	1:B:370:VAL:HG22	1.74	0.70
1:B:149:LEU:HD21	1:B:268:LEU:HD13	1.72	0.69
7:1:25:ALA:HB2	7:2:71:ASN:OD1	1.92	0.68
1:A:182:CYS:SG	1:A:329:ILE:HD11	2.34	0.68
7:4:44:ALA:HB2	7:5:53:LEU:HD11	1.76	0.67
1:B:21:SER:O	1:B:22:PHE:C	2.37	0.67
1:B:221:ARG:NE	2:D:133:ARG:HE	1.92	0.67
8:a:12:HIS:CD2	8:a:15:HIS:CG	2.83	0.67
7:2:47:PRO:HG3	7:3:46:GLN:NE2	2.07	0.66
1:B:113:PRO:HA	1:B:252:THR:HG21	1.78	0.66
7:2:46:GLN:HE22	7:2:52:ARG:HH22	1.42	0.65
7:7:44:ALA:HB2	7:8:53:LEU:HD21	1.78	0.65
2:E:67:MET:HE3	2:E:226:THR:HG22	1.78	0.65
1:C:182:CYS:SG	1:C:329:ILE:HD11	2.37	0.64
2:D:235:LEU:HD21	2:D:293:ARG:HB2	1.79	0.64
5:b:28:ASN:N	8:a:132:ASN:OD1	2.29	0.64
1:A:489:THR:HG22	1:A:491:LYS:HG2	1.80	0.64
2:F:319:ALA:HB3	2:F:320:PRO:HD3	1.80	0.63
1:A:159:MET:HE1	1:A:394:ARG:HA	1.80	0.63
1:B:15:ILE:HD11	6:d:219:LEU:HA	1.80	0.63
3:G:189:GLU:OE2	3:G:191:ARG:NH2	2.32	0.63
1:C:307:ARG:O	1:C:311:ARG:NH2	2.32	0.63
7:5:46:GLN:HG3	7:5:49:ALA:HB2	1.79	0.62
5:b:30:THR:OG1	8:a:28:ASN:N	2.32	0.62
8:a:134:VAL:HG21	8:a:190:PHE:HB2	1.81	0.62
2:D:277:VAL:HG12	2:D:277:VAL:O	1.99	0.62
8:a:86:ALA:HB2	8:a:245:MET:HE1	1.81	0.62
5:b:57:LEU:HD21	8:a:67:GLU:HG2	1.79	0.62
6:d:327:LEU:HD21	6:d:345:LEU:HB3	1.80	0.62
1:B:221:ARG:CZ	2:D:133:ARG:HE	2.12	0.62
1:B:279:ASP:OD1	1:B:282:ARG:NH1	2.33	0.62
5:b:145:GLU:O	5:b:148:VAL:HG12	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:ASP:OD1	1:C:71:GLU:N	2.32	0.61
2:E:302:ILE:HG23	2:E:302:ILE:O	2.00	0.61
6:d:254:ARG:O	6:d:255:THR:C	2.43	0.61
1:B:14:ALA:CB	6:d:240:LEU:HD12	2.30	0.60
2:F:11:ARG:NH1	2:F:23:GLU:OE1	2.33	0.60
2:F:172:GLU:OE2	2:F:176:ARG:NE	2.34	0.60
6:d:25:PRO:O	6:d:29:LEU:HD13	2.01	0.60
2:F:59:ASP:O	2:F:60:SER:OG	2.17	0.60
2:F:355:ILE:HG13	2:F:360:ILE:HG21	1.81	0.60
5:b:52:PRO:O	5:b:56:VAL:HG23	2.02	0.60
5:b:151:LEU:HD12	6:d:121:LEU:HD21	1.82	0.60
1:B:159:MET:HE1	1:B:370:VAL:CG2	2.32	0.60
5:b:151:LEU:CD1	6:d:121:LEU:HD21	2.31	0.59
7:2:44:ALA:O	7:3:46:GLN:NE2	2.35	0.59
2:D:438:ALA:O	2:D:442:LEU:HD23	2.02	0.59
7:5:65:GLU:OE1	12:5:601:YGR:N21	2.36	0.59
1:A:31:ILE:HD13	6:d:427:VAL:HG11	1.84	0.59
2:E:172:GLU:OE2	2:E:176:ARG:NE	2.34	0.59
2:E:319:ALA:HB3	2:E:320:PRO:HD3	1.84	0.59
2:D:142:LYS:NZ	2:D:458:ILE:O	2.31	0.59
1:A:26:THR:HG21	6:d:150:PHE:HE1	1.67	0.59
6:d:356:ARG:HB2	6:d:359:GLU:CB	2.33	0.59
3:G:49:GLU:OE1	4:H:78:ARG:NE	2.35	0.59
7:9:65:GLU:OE1	12:9:601:YGR:N21	2.36	0.58
2:F:404:ARG:HH22	2:F:444:LYS:C	2.10	0.58
7:2:47:PRO:CG	7:3:46:GLN:HE22	2.09	0.58
1:A:330:ILE:HD11	1:A:345:VAL:HG21	1.86	0.58
1:C:98:LEU:HD21	1:C:248:LEU:HD11	1.86	0.58
1:A:128:THR:HG22	1:A:129:ARG:N	2.17	0.58
6:d:237:LEU:HD23	6:d:238:GLU:N	2.18	0.58
1:B:178:LYS:HG2	1:B:355:LEU:HD12	1.85	0.58
3:G:147:TYR:CG	4:H:11:VAL:HG11	2.39	0.58
2:F:332:GLU:OE1	2:F:350:ALA:HB1	2.04	0.58
4:H:24:PHE:HE1	7:2:46:GLN:CD	2.11	0.58
7:9:78:PHE:CZ	7:8:79:VAL:HG11	2.39	0.58
1:B:182:CYS:SG	1:B:329:ILE:HD11	2.44	0.57
1:C:235:VAL:HG11	1:C:252:THR:HG21	1.81	0.57
7:9:6:ASN:O	7:9:7:ALA:C	2.47	0.57
2:D:88:VAL:HG11	2:D:233:SER:OG	2.04	0.57
2:E:294:ILE:HG22	2:E:294:ILE:O	2.04	0.57
3:G:57:LEU:O	3:G:61:SER:N	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:157:ALA:HB3	2:E:327:LEU:HD13	1.87	0.57
2:F:393:GLU:OE1	3:G:86:ARG:NH1	2.37	0.57
7:9:7:ALA:HA	7:1:9:ILE:HD11	1.85	0.57
1:C:395:LEU:HD13	2:F:457:LEU:HD11	1.86	0.56
2:E:193:ARG:NH1	9:E:600:ATP:O1G	2.34	0.56
1:A:39:ASP:OD1	2:F:272:ARG:NH1	2.39	0.56
6:d:356:ARG:HB2	6:d:359:GLU:HB3	1.87	0.56
2:E:287:MET:HE3	2:E:323:THR:HB	1.88	0.56
7:3:37:ASN:HB2	7:4:34:ILE:HG22	1.87	0.56
8:a:181:LYS:NZ	8:a:243:GLN:HE22	2.03	0.56
2:E:188:ALA:HB2	2:E:234:ALA:CB	2.35	0.56
7:3:75:MET:HE1	7:4:74:PHE:CE1	2.40	0.56
2:D:219:GLN:NE2	2:D:222:GLU:OE2	2.40	0.55
2:D:190:VAL:HG22	2:D:230:VAL:HG13	1.88	0.55
7:1:21:MET:HE3	7:2:71:ASN:ND2	2.22	0.55
1:A:15:ILE:HG13	6:d:121:LEU:HD22	1.89	0.55
2:D:328:ASP:O	2:D:354:THR:HG23	2.07	0.55
7:9:36:GLY:CA	7:1:60:THR:HG21	2.36	0.55
7:8:55:THR:HB	7:8:56:PRO:HD3	1.89	0.55
2:F:49:LEU:HD11	2:F:66:SER:OG	2.08	0.54
2:F:404:ARG:HH22	2:F:445:GLY:N	2.05	0.54
2:F:157:ALA:CB	2:F:327:LEU:HD13	2.37	0.54
7:6:37:ASN:OD1	7:7:38:ALA:HB2	2.07	0.54
1:B:294:ARG:NH1	2:E:317:ASP:OD2	2.34	0.54
8:a:12:HIS:CD2	8:a:15:HIS:CB	2.90	0.54
8:a:48:TYR:HD2	8:a:62:VAL:HG21	1.73	0.54
2:E:190:VAL:HG12	2:E:190:VAL:O	2.08	0.54
6:d:209:VAL:HG21	6:d:253:LEU:HD11	1.90	0.54
1:C:235:VAL:CG1	1:C:252:THR:CG2	2.80	0.54
2:E:342:ILE:HG23	2:E:413:SER:HB3	1.90	0.54
2:F:158:LEU:HD23	2:F:331:THR:HB	1.89	0.54
5:b:155:LEU:HD21	6:d:126:VAL:HG23	1.90	0.54
7:9:82:THR:HG23	7:9:85:LEU:HB2	1.90	0.53
1:C:77:VAL:HG13	1:C:284:ILE:HD12	1.90	0.53
2:E:67:MET:O	2:E:226:THR:HG23	2.09	0.53
6:d:34:GLN:CG	8:a:64:LEU:HD11	2.38	0.53
7:7:25:ALA:HB2	7:8:71:ASN:ND2	2.23	0.53
1:B:14:ALA:HB3	6:d:240:LEU:HD12	1.90	0.53
1:A:268:LEU:HD11	1:A:327:LEU:HD13	1.91	0.53
6:d:327:LEU:HD22	6:d:349:VAL:CG2	2.35	0.53
7:9:25:ALA:HB2	7:1:71:ASN:OD1	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:VAL:HG13	1:B:132:LEU:HD11	1.90	0.53
1:C:36:ASP:OD1	1:C:37:ALA:N	2.42	0.53
2:F:49:LEU:HD12	2:F:68:GLN:HB2	1.91	0.53
7:4:65:GLU:OE1	12:4:601:YGR:N21	2.41	0.53
1:C:430:LEU:HD12	1:C:468:GLU:OE1	2.09	0.52
7:2:37:ASN:HB2	7:3:34:ILE:HG22	1.90	0.52
1:B:334:ALA:HB1	3:G:279:LEU:HD21	1.92	0.52
5:b:133:LEU:HD21	6:d:110:ARG:HA	1.92	0.52
1:C:133:GLU:OE2	1:C:307:ARG:NH2	2.42	0.52
5:b:133:LEU:O	5:b:134:SER:C	2.51	0.52
7:3:75:MET:HE1	7:4:74:PHE:CZ	2.45	0.52
8:a:83:MET:HE1	8:a:90:LEU:HB2	1.91	0.52
1:B:176:THR:HG21	1:B:355:LEU:O	2.08	0.52
1:B:468:GLU:OE2	1:B:472:LEU:HD11	2.10	0.52
3:G:71:ARG:NH1	3:G:161:PHE:O	2.41	0.52
7:9:74:PHE:CD1	7:8:75:MET:HE1	2.44	0.52
8:a:86:ALA:HB1	8:a:89:VAL:CG1	2.40	0.52
1:C:18:TYR:CD2	6:d:326:LEU:HD12	2.45	0.52
2:F:57:LEU:HD11	2:F:63:ARG:CZ	2.39	0.52
2:F:91:GLY:O	2:F:94:VAL:HG22	2.09	0.52
1:B:142:ARG:NH1	1:B:310:GLU:O	2.37	0.52
7:5:46:GLN:CG	7:5:49:ALA:HB2	2.40	0.52
2:D:172:GLU:OE2	2:D:176:ARG:NE	2.43	0.52
2:E:386:ILE:HD11	2:E:394:LEU:HD11	1.90	0.51
1:C:18:TYR:CG	6:d:326:LEU:HD12	2.46	0.51
7:3:65:GLU:OE1	12:3:601:YGR:N21	2.43	0.51
7:2:46:GLN:CD	7:2:52:ARG:HH12	2.18	0.51
7:8:21:MET:HG3	7:8:75:MET:HE2	1.92	0.51
1:B:159:MET:CE	1:B:370:VAL:CG1	2.87	0.51
8:a:13:VAL:HG21	8:a:198:ILE:HD11	1.92	0.51
2:D:455:PHE:HA	2:D:458:ILE:HD11	1.93	0.51
1:A:128:THR:HG22	1:A:129:ARG:H	1.74	0.51
7:9:10:THR:HG23	7:1:9:ILE:CD1	2.41	0.51
7:7:21:MET:HG2	7:7:75:MET:HG3	1.93	0.51
7:4:21:MET:HE1	7:4:72:LEU:HA	1.93	0.51
8:a:66:TRP:CZ3	8:a:98:VAL:HG11	2.45	0.51
2:E:340:LYS:NZ	2:E:409:GLU:OE2	2.42	0.51
2:E:215:LEU:HD23	2:E:217:PHE:CZ	2.46	0.51
1:B:55:LEU:O	1:B:94:THR:OG1	2.21	0.50
1:C:235:VAL:CG1	1:C:252:THR:HG21	2.41	0.50
2:F:277:VAL:HG12	2:F:277:VAL:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:a:173:ILE:HD12	8:a:173:ILE:H	1.75	0.50
2:F:120:ILE:O	2:F:293:ARG:NH2	2.40	0.50
1:C:98:LEU:HD21	1:C:248:LEU:CD1	2.41	0.50
7:3:46:GLN:HG3	7:3:49:ALA:HB3	1.92	0.50
1:A:50:VAL:HG11	1:A:74:VAL:HG21	1.93	0.50
7:5:37:ASN:HB2	7:6:34:ILE:HG22	1.94	0.50
2:E:235:LEU:HD21	2:E:293:ARG:HB2	1.94	0.50
8:a:48:TYR:CD2	8:a:62:VAL:HG21	2.47	0.50
6:d:150:PHE:CZ	6:d:154:LEU:HD11	2.46	0.50
2:E:149:PRO:HD2	2:E:353:SER:OG	2.11	0.50
6:d:269:ASP:OD1	6:d:353:ARG:NH2	2.43	0.50
7:9:9:ILE:HG21	7:9:85:LEU:HA	1.94	0.50
6:d:422:THR:HG22	6:d:427:VAL:HG12	1.93	0.50
2:E:141:LEU:HD12	2:E:173:MET:HE1	1.94	0.49
7:3:37:ASN:OD1	7:4:38:ALA:HB2	2.12	0.49
7:9:10:THR:HG23	7:1:9:ILE:HD12	1.94	0.49
4:H:65:VAL:HG13	4:H:81:VAL:HG22	1.94	0.49
7:4:37:ASN:HB2	7:5:34:ILE:HG22	1.94	0.49
7:6:75:MET:HE1	7:7:74:PHE:CE1	2.47	0.49
2:D:188:ALA:HB2	2:D:234:ALA:CB	2.42	0.49
5:b:44:VAL:HG21	8:a:43:ILE:HD12	1.93	0.49
5:b:50:VAL:HG21	8:a:91:PRO:HA	1.94	0.49
1:A:326:GLY:C	1:A:327:LEU:HD12	2.37	0.49
2:F:158:LEU:HD22	2:F:169:LEU:HD23	1.95	0.49
2:F:388:ILE:HD12	3:G:18:ILE:HD11	1.94	0.49
6:d:195:ASP:O	6:d:199:LEU:N	2.37	0.49
3:G:89:CYS:SG	3:G:93:ASN:ND2	2.86	0.49
1:B:397:LEU:HD11	1:B:431:LEU:HD11	1.95	0.49
5:b:50:VAL:HB	5:b:51:PRO:HD3	1.95	0.49
7:4:72:LEU:HD11	7:5:74:PHE:CE2	2.47	0.49
6:d:341:ALA:O	6:d:345:LEU:HD23	2.12	0.49
2:D:91:GLY:O	2:D:94:VAL:HG22	2.13	0.49
2:E:469:GLU:O	2:E:470:SER:C	2.55	0.49
5:b:128:GLN:O	5:b:132:GLN:HG3	2.13	0.49
7:9:36:GLY:HA3	7:1:60:THR:HG21	1.95	0.49
1:B:133:GLU:OE2	1:B:307:ARG:NH2	2.46	0.48
5:b:107:GLY:HA3	6:d:87:LEU:HD13	1.95	0.48
8:a:86:ALA:HB1	8:a:89:VAL:HG12	1.95	0.48
1:A:160:THR:HG21	1:A:373:SER:OG	2.14	0.48
1:A:268:LEU:CD1	1:A:327:LEU:HD13	2.43	0.48
7:2:49:ALA:O	7:2:50:GLN:C	2.56	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:d:231:ALA:N	6:d:232:PRO:CD	2.77	0.48
6:d:349:VAL:HA	6:d:352:LEU:CD1	2.43	0.48
7:1:37:ASN:HB2	7:2:34:ILE:HG22	1.95	0.48
7:5:42:GLY:HA3	7:5:53:LEU:HD21	1.96	0.48
2:F:235:LEU:HD21	2:F:293:ARG:CB	2.43	0.48
7:2:28:ALA:HB2	7:2:68:TYR:HB2	1.95	0.48
7:4:21:MET:HE3	7:4:75:MET:CB	2.43	0.48
6:d:209:VAL:HG21	6:d:253:LEU:HD21	1.95	0.48
7:6:72:LEU:HD11	7:7:74:PHE:CE2	2.49	0.48
7:7:25:ALA:HB2	7:8:71:ASN:HD21	1.77	0.48
1:A:251:TYR:OH	1:A:304:LEU:HD12	2.14	0.48
2:D:378:ARG:HE	2:D:382:LEU:CD1	2.26	0.48
1:B:334:ALA:O	1:B:335:ASN:OD1	2.31	0.48
2:E:169:LEU:HD12	2:E:333:LEU:HD11	1.96	0.48
7:9:55:THR:HB	7:9:56:PRO:HD3	1.94	0.48
1:A:100:VAL:HG11	1:A:252:THR:HG23	1.95	0.48
2:E:357:ASP:OD1	2:E:358:PRO:HD2	2.13	0.48
1:B:149:LEU:CD2	1:B:268:LEU:HD13	2.41	0.48
2:D:227:ARG:NH1	2:D:265:GLU:OE1	2.43	0.48
2:F:355:ILE:CG1	2:F:360:ILE:HG21	2.43	0.48
7:2:8:ILE:HD12	7:2:8:ILE:H	1.79	0.47
1:B:176:THR:HG22	1:B:176:THR:O	2.14	0.47
7:8:65:GLU:OE1	12:8:601:YGR:N21	2.47	0.47
2:E:320:PRO:HA	2:E:323:THR:HG22	1.96	0.47
1:B:181:VAL:HG12	1:B:181:VAL:O	2.14	0.47
2:E:88:VAL:HG11	2:E:233:SER:OG	2.15	0.47
6:d:34:GLN:HE21	8:a:58:VAL:HG12	1.79	0.47
3:G:225:GLU:CD	7:2:45:ARG:HH22	2.23	0.47
6:d:24:PRO:O	6:d:28:THR:HG23	2.15	0.47
7:4:55:THR:HB	7:4:56:PRO:HD3	1.95	0.47
1:B:149:LEU:HD22	1:B:325:THR:HG21	1.95	0.47
1:C:442:GLU:O	1:C:445:VAL:HG22	2.14	0.47
2:D:162:ALA:O	2:D:335:ARG:NH1	2.48	0.47
2:D:355:ILE:HG23	2:D:360:ILE:HG21	1.97	0.47
2:F:393:GLU:HG3	3:G:88:LEU:HD11	1.96	0.47
6:d:190:VAL:HG21	6:d:252:LEU:CD1	2.43	0.47
6:d:349:VAL:HA	6:d:352:LEU:HD11	1.95	0.47
1:A:11:ILE:H	1:A:11:ILE:HD12	1.79	0.47
1:A:205:VAL:HG13	1:A:233:THR:HG23	1.97	0.47
1:C:252:THR:HG23	1:C:253:GLY:N	2.29	0.47
2:D:406:ARG:NH1	2:D:452:GLU:OE1	2.41	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:5:8:ILE:HD12	7:5:8:ILE:H	1.80	0.47
1:C:174:ARG:NH1	2:E:332:GLU:OE2	2.47	0.47
1:C:176:THR:HG22	1:C:355:LEU:HB3	1.96	0.47
2:D:172:GLU:OE2	2:D:176:ARG:NH1	2.48	0.47
6:d:356:ARG:HB2	6:d:359:GLU:HB2	1.97	0.47
6:d:379:VAL:CG1	6:d:419:LEU:HD11	2.44	0.47
7:6:37:ASN:HB2	7:7:34:ILE:HG22	1.95	0.47
1:C:132:LEU:HD12	1:C:248:LEU:HD11	1.97	0.46
2:E:442:LEU:HD23	2:E:447:PHE:HB2	1.96	0.46
1:B:53:GLN:HG3	2:E:71:ASP:OD1	2.14	0.46
2:D:95:LYS:HD3	2:D:184:THR:HG21	1.97	0.46
2:E:162:ALA:O	2:E:335:ARG:NH2	2.45	0.46
7:5:37:ASN:OD1	7:6:38:ALA:HB2	2.15	0.46
2:D:172:GLU:OE1	2:D:417:MET:N	2.48	0.46
2:D:255:ASN:O	2:D:258:ARG:HB2	2.16	0.46
7:8:45:ARG:NH2	7:8:45:ARG:HB3	2.30	0.46
1:A:151:THR:HG21	1:A:162:ILE:HD12	1.97	0.46
2:F:227:ARG:NH1	2:F:265:GLU:OE1	2.43	0.46
7:9:53:LEU:HD21	7:8:44:ALA:HA	1.98	0.46
1:A:272:ASP:HA	1:A:273:ASP:HA	1.74	0.46
3:G:118:VAL:CG1	3:G:123:ALA:HB3	2.45	0.46
3:G:245:TYR:CE2	3:G:249:LEU:HD11	2.51	0.46
2:D:133:ARG:HG2	2:D:133:ARG:NH2	2.31	0.46
2:D:133:ARG:HG2	2:D:133:ARG:HH21	1.81	0.46
3:G:39:ARG:NH2	3:G:145:PRO:O	2.43	0.46
7:5:25:ALA:HB2	7:6:71:ASN:OD1	2.16	0.46
1:A:142:ARG:CZ	1:A:313:ALA:HB2	2.45	0.46
8:a:12:HIS:O	8:a:13:VAL:C	2.59	0.46
8:a:142:PHE:HA	8:a:145:TYR:CE2	2.51	0.46
5:b:63:MET:O	5:b:64:LEU:C	2.56	0.46
6:d:34:GLN:HG3	8:a:64:LEU:HD11	1.97	0.46
1:B:434:PRO:HG2	1:B:437:SER:HB3	1.98	0.45
7:9:21:MET:HG2	7:9:75:MET:HG3	1.97	0.45
7:8:61:VAL:O	7:8:65:GLU:HG2	2.16	0.45
5:b:155:LEU:CD2	6:d:126:VAL:HG23	2.45	0.45
6:d:34:GLN:HG2	8:a:64:LEU:HD11	1.98	0.45
7:4:44:ALA:CB	7:5:53:LEU:HD11	2.45	0.45
2:E:210:LEU:HA	2:E:213:THR:HG22	1.97	0.45
2:F:157:ALA:HB3	2:F:327:LEU:HD13	1.98	0.45
7:6:8:ILE:HD12	7:6:8:ILE:H	1.80	0.45
2:F:367:ARG:NH1	2:F:440:ASP:OD2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:455:PHE:HA	2:F:458:ILE:HD13	1.98	0.45
3:G:57:LEU:HD21	3:G:231:LEU:HG	1.99	0.45
6:d:245:VAL:HB	6:d:249:THR:HG21	1.98	0.45
7:5:46:GLN:HG3	7:5:49:ALA:CB	2.45	0.45
1:A:471:LEU:HD11	1:A:504:ILE:HG12	1.99	0.45
2:E:148:THR:HG22	2:E:148:THR:O	2.17	0.45
7:5:49:ALA:O	7:5:50:GLN:C	2.60	0.45
1:A:446:VAL:HG21	1:A:483:PHE:HE1	1.81	0.45
2:E:209:VAL:HG12	2:E:209:VAL:O	2.16	0.45
2:E:310:VAL:HG23	2:E:310:VAL:O	2.16	0.45
4:H:32:ILE:HD11	7:3:45:ARG:HH12	1.80	0.45
7:2:46:GLN:HE22	7:2:52:ARG:NH2	2.11	0.45
2:D:151:VAL:HG21	2:D:328:ASP:HB3	1.97	0.45
1:C:77:VAL:CG1	1:C:284:ILE:HD12	2.46	0.45
2:D:382:LEU:HD23	2:D:394:LEU:HD13	1.99	0.45
2:E:105:CYS:SG	2:E:107:ASP:OD1	2.75	0.45
2:F:442:LEU:HD21	2:F:455:PHE:CE2	2.51	0.45
5:b:89:MET:O	5:b:93:ARG:HG2	2.17	0.45
1:A:292:PRO:HD3	2:F:273:MET:HE1	1.99	0.45
2:D:380:LYS:HA	2:D:383:GLN:HG3	1.99	0.45
2:F:50:THR:HG21	2:F:102:LEU:HD11	1.99	0.45
3:G:116:LEU:HD11	3:G:132:ARG:HD2	1.98	0.45
7:7:65:GLU:OE1	12:7:601:YGR:N21	2.50	0.45
8:a:15:HIS:CG	8:a:15:HIS:O	2.69	0.44
2:F:188:ALA:HB2	2:F:234:ALA:CB	2.47	0.44
3:G:88:LEU:HB3	3:G:263:MET:HE3	1.99	0.44
4:H:22:PHE:HD1	4:H:35:LEU:CG	2.15	0.44
7:2:20:ILE:HG23	7:2:71:ASN:HD22	1.82	0.44
2:E:277:VAL:HG12	2:E:318:PRO:HG2	1.99	0.44
7:4:25:ALA:HB2	7:5:71:ASN:OD1	2.16	0.44
1:A:175:LYS:NZ	9:A:600:ATP:O1G	2.42	0.44
2:D:210:LEU:HA	2:D:213:THR:HG22	1.99	0.44
3:G:183:LEU:CD1	3:G:203:ALA:HB3	2.47	0.44
1:A:400:TYR:HB2	1:A:424:GLY:HA3	1.99	0.44
1:C:171:ILE:HD11	1:C:342:PRO:HB3	1.99	0.44
2:E:32:LEU:HD12	2:E:56:HIS:CE1	2.52	0.44
1:C:403:LEU:HD23	1:C:421:LEU:HD12	2.00	0.44
7:6:46:GLN:HG2	7:6:49:ALA:HB3	1.99	0.44
1:A:207:VAL:HG12	1:A:209:ILE:HD11	1.98	0.44
2:D:188:ALA:HB2	2:D:234:ALA:HB3	1.98	0.44
2:E:189:GLY:O	2:E:217:PHE:O	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:VAL:HG21	2:E:98:VAL:HG11	1.98	0.44
6:d:147:VAL:O	6:d:151:LEU:HD12	2.18	0.44
8:a:30:ASP:OD1	8:a:31:THR:N	2.51	0.44
1:A:269:ILE:HG21	1:A:271:PHE:CZ	2.53	0.44
1:B:166:GLN:NE2	1:B:168:GLN:OE1	2.51	0.44
1:A:205:VAL:HG22	1:A:233:THR:HG22	2.00	0.43
1:B:360:PHE:CD1	1:B:360:PHE:C	2.96	0.43
2:F:235:LEU:HD21	2:F:293:ARG:HB2	2.00	0.43
4:H:37:ARG:O	7:3:47:PRO:HD2	2.17	0.43
7:9:82:THR:HG21	7:9:85:LEU:HD22	1.96	0.43
1:A:53:GLN:N	1:A:66:ALA:O	2.48	0.43
7:9:6:ASN:O	7:9:10:THR:HG22	2.18	0.43
8:a:11:ILE:CG2	8:a:13:VAL:HG23	2.48	0.43
8:a:130:ASP:OD1	8:a:132:ASN:N	2.50	0.43
1:A:133:GLU:OE1	1:A:307:ARG:NE	2.49	0.43
1:A:307:ARG:O	1:A:311:ARG:HD2	2.19	0.43
1:B:169:LEU:HD13	1:B:345:VAL:HG11	1.99	0.43
1:B:355:LEU:HD23	1:B:368:ILE:HD13	2.00	0.43
1:B:373:SER:O	1:B:394:ARG:NH2	2.50	0.43
1:C:50:VAL:HG11	1:C:74:VAL:HG21	2.01	0.43
2:E:448:ASP:OD1	2:E:448:ASP:N	2.50	0.43
2:E:458:ILE:HG23	2:E:459:GLY:N	2.33	0.43
8:a:12:HIS:CD2	8:a:15:HIS:CD2	3.06	0.43
1:B:68:ASN:C	1:B:69:LEU:HD12	2.43	0.43
3:G:53:MET:O	3:G:57:LEU:HD13	2.18	0.43
3:G:134:VAL:HG23	3:G:134:VAL:O	2.17	0.43
3:G:245:TYR:CZ	3:G:249:LEU:HD11	2.53	0.43
4:H:62:ARG:NH2	4:H:117:LEU:O	2.52	0.43
7:1:21:MET:HG2	7:1:75:MET:HG3	2.00	0.43
8:a:37:VAL:O	8:a:41:ILE:HG12	2.18	0.43
1:C:182:CYS:HB3	1:C:270:VAL:HG11	2.01	0.43
2:E:92:ASP:OD1	2:E:245:GLN:NE2	2.51	0.43
2:F:42:PHE:O	2:F:46:ALA:N	2.51	0.43
4:H:118:GLY:O	4:H:119:GLN:C	2.62	0.43
1:A:205:VAL:HB	1:A:269:ILE:HD13	2.01	0.43
1:B:159:MET:HE1	1:B:370:VAL:HG13	1.99	0.43
1:B:355:LEU:CD2	1:B:368:ILE:HD13	2.49	0.43
2:D:249:VAL:HG12	2:D:250:LEU:N	2.34	0.43
7:3:64:VAL:HG23	7:3:65:GLU:N	2.34	0.43
2:D:146:LEU:HD22	2:D:439:PHE:CD2	2.53	0.43
2:E:67:MET:HE1	2:E:229:ARG:HB2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:179:ARG:HD3	2:F:209:VAL:HG21	1.99	0.43
3:G:193:MET:CE	3:G:260:ARG:HD3	2.49	0.43
1:A:459:VAL:HG22	1:A:519:VAL:HG11	2.00	0.43
1:C:538:VAL:HG23	1:C:538:VAL:O	2.18	0.43
2:D:358:PRO:HG3	2:D:366:TYR:CD2	2.54	0.43
2:F:88:VAL:HG11	2:F:233:SER:OG	2.19	0.43
8:a:96:ILE:HG21	8:a:233:PHE:HD1	1.84	0.43
1:A:68:ASN:C	1:A:69:LEU:HD12	2.44	0.43
1:A:142:ARG:NH2	1:A:313:ALA:HB2	2.34	0.43
1:A:212:LYS:HA	2:F:292:GLU:OE2	2.19	0.43
1:C:97:VAL:HG12	1:C:98:LEU:N	2.34	0.43
1:C:249:ALA:HA	1:C:252:THR:HG22	2.00	0.43
7:4:72:LEU:HD11	7:5:74:PHE:HE2	1.84	0.43
1:B:249:ALA:HB3	1:B:250:PRO:HD3	2.00	0.42
2:E:277:VAL:HG22	2:E:277:VAL:O	2.19	0.42
3:G:12:ILE:CG1	3:G:277:LEU:HB3	2.49	0.42
7:6:44:ALA:HA	7:7:53:LEU:HD11	2.01	0.42
1:C:446:VAL:HG21	1:C:483:PHE:HE1	1.84	0.42
3:G:39:ARG:CZ	3:G:144:ARG:NH2	2.82	0.42
8:a:158:HIS:O	8:a:162:VAL:HG23	2.19	0.42
6:d:314:SER:HA	6:d:352:LEU:HD22	2.00	0.42
1:B:386:MET:HE1	1:B:444:GLN:HB2	2.01	0.42
2:D:139:THR:HG23	2:D:141:LEU:H	1.84	0.42
2:D:355:ILE:HD13	2:D:360:ILE:HD13	2.01	0.42
2:F:30:PRO:O	2:F:56:HIS:NE2	2.50	0.42
5:b:48:TRP:O	5:b:52:PRO:HG2	2.20	0.42
6:d:150:PHE:CE1	6:d:154:LEU:HD11	2.54	0.42
7:1:19:LEU:HG	7:2:19:LEU:HD13	2.01	0.42
1:B:159:MET:HE2	1:B:370:VAL:HG13	1.96	0.42
1:B:485:GLY:O	1:B:489:THR:HG22	2.19	0.42
1:C:310:GLU:HG2	2:F:221:ASP:HB3	2.02	0.42
2:F:357:ASP:O	2:F:361:VAL:HG22	2.18	0.42
4:H:73:THR:HG22	4:H:74:GLU:N	2.35	0.42
5:b:78:VAL:HG23	6:d:57:MET:HE1	2.01	0.42
7:6:65:GLU:OE1	12:6:601:YGR:N21	2.52	0.42
7:8:45:ARG:HB3	7:8:45:ARG:CZ	2.50	0.42
1:C:330:ILE:HD11	1:C:345:VAL:HG21	2.01	0.42
2:D:151:VAL:HG22	2:D:355:ILE:HG13	2.01	0.42
2:F:273:MET:HE3	3:G:300:ALA:CB	2.49	0.42
1:A:216:ILE:HD13	1:A:236:ALA:HB1	2.00	0.42
7:3:9:ILE:HG22	7:3:82:THR:HG21	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:254:ASP:HA	2:F:255:ASN:HA	1.83	0.42
3:G:12:ILE:HG13	3:G:277:LEU:HB3	2.00	0.42
5:b:63:MET:HE3	6:d:44:SER:HB2	2.01	0.42
7:9:19:LEU:HD13	7:8:19:LEU:HG	2.01	0.42
1:B:15:ILE:HD11	6:d:219:LEU:CA	2.50	0.42
2:D:357:ASP:O	2:D:361:VAL:HG22	2.19	0.42
6:d:219:LEU:O	6:d:223:LEU:HD13	2.20	0.42
6:d:398:SER:O	6:d:402:GLY:N	2.53	0.42
1:B:159:MET:HE1	1:B:370:VAL:CG1	2.49	0.42
2:E:364:GLU:OE2	2:E:367:ARG:NH1	2.53	0.42
3:G:12:ILE:O	3:G:12:ILE:HG22	2.20	0.42
4:H:40:PRO:HA	4:H:72:VAL:O	2.20	0.42
7:9:82:THR:HG23	7:9:82:THR:O	2.20	0.42
8:a:214:PRO:HA	8:a:217:VAL:HG12	2.02	0.42
1:A:128:THR:CG2	1:A:129:ARG:N	2.83	0.41
1:B:181:VAL:O	1:B:181:VAL:CG1	2.67	0.41
1:C:409:PHE:CE2	2:E:385:ILE:HD13	2.55	0.41
5:b:60:ARG:NH2	6:d:37:VAL:HG21	2.34	0.41
1:B:142:ARG:NH2	1:B:313:ALA:HB2	2.35	0.41
2:D:355:ILE:HD12	2:D:360:ILE:HG21	2.01	0.41
2:E:147:LEU:HD13	2:E:348:PRO:HB3	2.02	0.41
2:F:241:PHE:CD1	2:F:247:GLN:OE1	2.73	0.41
6:d:150:PHE:O	6:d:154:LEU:HD13	2.21	0.41
7:1:21:MET:HE1	7:1:68:TYR:CZ	2.54	0.41
7:4:19:LEU:HG	7:5:19:LEU:HD13	2.03	0.41
7:6:44:ALA:HB2	7:7:53:LEU:HD11	2.02	0.41
1:A:55:LEU:O	1:A:94:THR:OG1	2.34	0.41
4:H:8:ILE:CD1	4:H:63:ILE:HD13	2.49	0.41
8:a:12:HIS:CD2	8:a:15:HIS:HB2	2.55	0.41
1:A:47:LEU:O	1:A:50:VAL:HG12	2.20	0.41
1:B:97:VAL:O	1:B:98:LEU:C	2.62	0.41
1:C:417:SER:O	1:C:421:LEU:HD13	2.21	0.41
4:H:21:THR:O	4:H:35:LEU:HD23	2.20	0.41
7:4:46:GLN:O	7:4:46:GLN:HG2	2.19	0.41
7:6:47:PRO:HB3	7:7:52:ARG:NE	2.35	0.41
8:a:17:THR:HG23	8:a:18:LEU:HD13	2.02	0.41
1:A:147:GLU:OE2	1:A:314:LYS:NZ	2.41	0.41
1:A:500:LEU:HA	1:A:503:VAL:HG12	2.03	0.41
1:B:63:LEU:HD12	1:B:114:LEU:HD13	2.01	0.41
2:D:143:VAL:HG13	2:D:144:VAL:N	2.36	0.41
2:F:442:LEU:HD21	2:F:455:PHE:HE2	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:251:TYR:OH	1:C:304:LEU:HD12	2.19	0.41
2:E:442:LEU:HD21	2:E:455:PHE:CE1	2.55	0.41
7:9:49:ALA:HB1	7:9:52:ARG:NH1	2.36	0.41
7:7:18:GLY:HA2	7:8:20:ILE:HG13	2.03	0.41
2:E:235:LEU:HD13	2:E:294:ILE:HG12	2.02	0.41
2:F:439:PHE:O	2:F:443:THR:HG23	2.20	0.41
2:F:450:LEU:HD22	2:F:468:ALA:HB1	2.01	0.41
7:9:36:GLY:HA2	7:1:60:THR:HG21	2.02	0.41
7:1:25:ALA:HB1	7:2:24:GLY:HA2	2.02	0.41
7:4:21:MET:HE3	7:4:75:MET:HB3	2.02	0.41
7:5:77:LEU:HD12	7:5:81:ALA:HB3	2.02	0.41
8:a:74:ARG:HG2	8:a:74:ARG:HH21	1.86	0.41
2:E:217:PHE:O	2:E:218:GLY:C	2.63	0.41
2:E:239:GLU:OE1	2:E:293:ARG:NH1	2.46	0.41
4:H:7:GLU:OE2	4:H:78:ARG:NH2	2.43	0.41
4:H:28:THR:N	4:H:44:GLN:O	2.50	0.41
4:H:35:LEU:HB3	4:H:36:PRO:HD2	2.03	0.41
6:d:4:PHE:CD2	6:d:4:PHE:C	2.98	0.41
1:B:97:VAL:HG12	1:B:98:LEU:N	2.35	0.41
1:B:205:VAL:HB	1:B:269:ILE:CD1	2.50	0.41
2:E:161:GLY:O	2:E:166:LYS:NZ	2.52	0.41
2:E:310:VAL:HG12	2:E:320:PRO:CG	2.50	0.41
2:F:141:LEU:HD12	2:F:173:MET:CE	2.51	0.41
2:F:232:LEU:HD12	2:F:232:LEU:N	2.36	0.41
2:F:324:PHE:CE1	2:F:332:GLU:OE2	2.73	0.41
2:F:400:GLN:O	2:F:404:ARG:HG3	2.21	0.41
4:H:39:ILE:HG23	7:3:44:ALA:O	2.21	0.41
5:b:45:ILE:O	5:b:50:VAL:HG23	2.21	0.41
7:8:21:MET:HG2	7:8:75:MET:HG3	2.02	0.41
8:a:16:HIS:O	8:a:19:VAL:HG23	2.21	0.41
2:E:294:ILE:O	2:E:294:ILE:CG2	2.68	0.41
2:F:96:GLY:HA2	2:F:210:LEU:O	2.20	0.41
2:D:378:ARG:O	2:D:382:LEU:HD13	2.21	0.40
2:E:169:LEU:O	2:E:173:MET:HG2	2.21	0.40
2:F:294:ILE:O	2:F:294:ILE:HG22	2.20	0.40
7:9:52:ARG:NH1	7:8:47:PRO:HG3	2.36	0.40
7:1:21:MET:HB3	7:2:71:ASN:ND2	2.36	0.40
2:E:160:GLY:C	2:E:310:VAL:HG22	2.46	0.40
2:F:176:ARG:CZ	2:F:417:MET:HE1	2.51	0.40
4:H:81:VAL:HG11	4:H:84:ALA:HB2	2.03	0.40
7:6:32:ASP:OD1	7:6:64:VAL:HG11	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:260:THR:HG23	2:D:283:LEU:HD13	2.03	0.40
2:F:245:GLN:O	2:F:246:GLY:C	2.61	0.40
3:G:33:ILE:HD11	3:G:257:ALA:HA	2.04	0.40
3:G:51:THR:O	3:G:55:THR:HG23	2.21	0.40
6:d:393:LEU:HA	6:d:396:VAL:HG22	2.04	0.40
7:8:13:ALA:HB2	7:8:82:THR:HG23	2.03	0.40
1:A:205:VAL:HB	1:A:269:ILE:CD1	2.52	0.40
1:C:187:LEU:HD22	1:C:226:GLY:HA3	2.04	0.40
2:D:133:ARG:HD3	2:D:133:ARG:HA	1.87	0.40
6:d:254:ARG:O	6:d:257:VAL:N	2.55	0.40
1:C:203:ARG:O	1:C:267:VAL:HA	2.22	0.40
2:E:157:ALA:CB	2:E:327:LEU:HD13	2.50	0.40
2:E:221:ASP:OD1	2:E:221:ASP:O	2.39	0.40
2:F:242:ARG:HD3	2:F:302:ILE:HD12	2.03	0.40
3:G:190:PHE:H	3:G:250:GLU:HG2	1.85	0.40
5:b:47:LYS:HG3	5:b:48:TRP:CD1	2.56	0.40
6:d:436:LEU:HD23	6:d:436:LEU:C	2.46	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	512/548 (93%)	500 (98%)	12 (2%)	0	100	100
1	B	502/548 (92%)	488 (97%)	13 (3%)	1 (0%)	43	66
1	C	508/548 (93%)	494 (97%)	14 (3%)	0	100	100
2	D	462/475 (97%)	452 (98%)	10 (2%)	0	100	100
2	E	462/475 (97%)	444 (96%)	18 (4%)	0	100	100
2	F	462/475 (97%)	452 (98%)	9 (2%)	1 (0%)	43	66
3	G	273/307 (89%)	272 (100%)	1 (0%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	H	115/121 (95%)	114 (99%)	1 (1%)	0	100	100
5	b	139/170 (82%)	133 (96%)	6 (4%)	0	100	100
6	d	428/445 (96%)	419 (98%)	8 (2%)	1 (0%)	43	66
7	1	79/86 (92%)	76 (96%)	3 (4%)	0	100	100
7	2	79/86 (92%)	78 (99%)	1 (1%)	0	100	100
7	3	79/86 (92%)	78 (99%)	1 (1%)	0	100	100
7	4	79/86 (92%)	77 (98%)	2 (2%)	0	100	100
7	5	79/86 (92%)	77 (98%)	2 (2%)	0	100	100
7	6	79/86 (92%)	77 (98%)	2 (2%)	0	100	100
7	7	79/86 (92%)	76 (96%)	3 (4%)	0	100	100
7	8	79/86 (92%)	78 (99%)	1 (1%)	0	100	100
7	9	79/86 (92%)	76 (96%)	3 (4%)	0	100	100
8	a	230/252 (91%)	223 (97%)	7 (3%)	0	100	100
All	All	4804/5138 (94%)	4684 (98%)	117 (2%)	3 (0%)	49	70

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	72	HIS
6	d	237	LEU
2	F	277	VAL

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	399/441 (90%)	398 (100%)	1 (0%)	86	94
1	B	377/441 (86%)	372 (99%)	5 (1%)	61	82
1	C	379/441 (86%)	378 (100%)	1 (0%)	86	94
2	D	367/388 (95%)	362 (99%)	5 (1%)	59	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	371/388 (96%)	367 (99%)	4 (1%)	65	84
2	F	369/388 (95%)	367 (100%)	2 (0%)	81	92
3	G	205/238 (86%)	204 (100%)	1 (0%)	81	92
4	H	72/96 (75%)	71 (99%)	1 (1%)	59	81
5	b	79/124 (64%)	73 (92%)	6 (8%)	12	27
6	d	229/347 (66%)	222 (97%)	7 (3%)	35	63
7	1	49/56 (88%)	49 (100%)	0	100	100
7	2	50/56 (89%)	49 (98%)	1 (2%)	48	74
7	3	49/56 (88%)	49 (100%)	0	100	100
7	4	51/56 (91%)	51 (100%)	0	100	100
7	5	50/56 (89%)	47 (94%)	3 (6%)	17	37
7	6	50/56 (89%)	49 (98%)	1 (2%)	48	74
7	7	49/56 (88%)	48 (98%)	1 (2%)	48	74
7	8	49/56 (88%)	49 (100%)	0	100	100
7	9	50/56 (89%)	50 (100%)	0	100	100
8	a	185/200 (92%)	183 (99%)	2 (1%)	65	84
All	All	3479/3996 (87%)	3438 (99%)	41 (1%)	61	83

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	476	LYS
1	B	71	GLU
1	B	215	THR
1	B	273	ASP
1	B	331	GLU
1	B	366	PRO
1	C	81	GLU
2	D	134	THR
2	D	258	ARG
2	D	285	ASP
2	D	357	ASP
2	D	358	PRO
2	E	107	ASP
2	E	135	GLU
2	E	255	ASN

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Mol	Chain	Res	Type
2	E	470	SER
2	F	113	LYS
2	F	247	GLN
3	G	14	SER
4	H	119	GLN
5	b	52	PRO
5	b	55	LYS
5	b	59	GLU
5	b	64	LEU
5	b	132	GLN
5	b	133	LEU
6	d	29	LEU
6	d	33	GLN
6	d	37	VAL
6	d	242	SER
6	d	250	LEU
6	d	252	LEU
6	d	371	ARG
7	2	50	GLN
7	5	45	ARG
7	5	52	ARG
7	5	53	LEU
7	6	45	ARG
7	7	52	ARG
8	a	11	ILE
8	a	13	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	305	HIS
1	B	72	HIS
1	B	143	GLN
1	B	211	GLN
1	B	258	GLN
1	C	305	HIS
1	C	479	HIS
2	D	68	GLN
2	D	97	HIS
2	D	365	HIS
2	D	425	GLN
2	E	175	ASN

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Mol	Chain	Res	Type
2	E	208	ASN
2	E	306	GLN
2	E	326	HIS
2	F	97	HIS
2	F	175	ASN
2	F	326	HIS
5	b	132	GLN
5	b	136	GLN
6	d	34	GLN
6	d	220	ASN
6	d	273	HIS
7	9	6	ASN
7	2	71	ASN
7	3	46	GLN
8	a	229	GLN
8	a	243	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](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 4 are monoatomic - leaving 12 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
12	YGR	3	601	-	46,46,46	1.38	4 (8%)	60,66,66	1.35	9 (15%)
9	ATP	B	600	10	32,33,33	1.68	7 (21%)	48,52,52	1.84	13 (27%)
12	YGR	9	601	-	46,46,46	1.45	4 (8%)	60,66,66	1.43	9 (15%)
12	YGR	8	601	-	46,46,46	1.42	5 (10%)	60,66,66	1.34	7 (11%)
12	YGR	7	601	-	46,46,46	1.32	5 (10%)	60,66,66	1.31	8 (13%)
12	YGR	6	601	-	46,46,46	1.47	4 (8%)	60,66,66	1.43	10 (16%)
11	PO4	F	600	-	4,4,4	1.10	0	6,6,6	0.43	0
9	ATP	E	600	10	32,33,33	0.52	0	48,52,52	0.66	0
12	YGR	5	601	-	46,46,46	1.34	4 (8%)	60,66,66	1.32	8 (13%)
9	ATP	A	600	10	32,33,33	0.55	0	48,52,52	0.69	1 (2%)
9	ATP	C	600	10	32,33,33	0.49	0	48,52,52	0.70	1 (2%)
12	YGR	4	601	-	46,46,46	1.47	5 (10%)	60,66,66	1.36	8 (13%)

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
12	YGR	3	601	-	-	6/38/38/38	0/4/4/4
12	YGR	9	601	-	-	11/38/38/38	0/4/4/4
9	ATP	B	600	10	-	3/22/38/38	0/3/3/3
12	YGR	8	601	-	-	10/38/38/38	0/4/4/4
12	YGR	7	601	-	-	10/38/38/38	0/4/4/4
12	YGR	6	601	-	-	8/38/38/38	0/4/4/4
9	ATP	E	600	10	-	11/22/38/38	0/3/3/3
12	YGR	5	601	-	-	10/38/38/38	0/4/4/4
9	ATP	A	600	10	-	5/22/38/38	0/3/3/3
9	ATP	C	600	10	-	6/22/38/38	0/3/3/3
12	YGR	4	601	-	-	10/38/38/38	0/4/4/4

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	4	601	YGR	C18-C05	6.13	1.64	1.57
12	9	601	YGR	C18-C05	6.07	1.64	1.57
12	6	601	YGR	C18-C05	5.85	1.63	1.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	8	601	YGR	C18-C05	5.72	1.63	1.57
12	3	601	YGR	C18-C05	5.44	1.63	1.57
12	5	601	YGR	C18-C05	5.26	1.63	1.57
12	7	601	YGR	C18-C05	4.76	1.62	1.57
9	B	600	ATP	C5-N7	-4.47	1.30	1.39
9	B	600	ATP	C4-N9	-3.25	1.30	1.37
12	9	601	YGR	C18-C25	3.23	1.56	1.52
12	6	601	YGR	C18-C25	3.16	1.56	1.52
9	B	600	ATP	C8-N9	-3.06	1.32	1.37
12	4	601	YGR	C18-C25	2.89	1.56	1.52
9	B	600	ATP	C5-C4	2.77	1.44	1.39
12	3	601	YGR	C18-C25	2.73	1.56	1.52
9	B	600	ATP	PB-O3B	-2.69	1.56	1.59
12	8	601	YGR	C18-C25	2.60	1.56	1.52
12	6	601	YGR	C19-C18	2.49	1.57	1.54
12	7	601	YGR	C18-C25	2.40	1.55	1.52
12	9	601	YGR	O24-C18	-2.38	1.39	1.43
12	3	601	YGR	C10-N13	2.37	1.35	1.32
12	5	601	YGR	C18-C25	2.31	1.55	1.52
12	4	601	YGR	C10-N13	2.25	1.35	1.32
12	8	601	YGR	C10-N13	2.23	1.35	1.32
12	6	601	YGR	C10-N13	2.22	1.35	1.32
9	B	600	ATP	PG-O3G	-2.22	1.46	1.54
12	3	601	YGR	O24-C18	-2.21	1.39	1.43
12	7	601	YGR	C10-N13	2.18	1.35	1.32
12	4	601	YGR	O24-C18	-2.17	1.39	1.43
12	7	601	YGR	O24-C18	-2.13	1.39	1.43
12	7	601	YGR	BR3-C38	2.11	1.94	1.90
12	8	601	YGR	BR3-C38	2.11	1.94	1.90
12	5	601	YGR	C10-N13	2.11	1.35	1.32
12	9	601	YGR	C10-N13	2.11	1.35	1.32
12	5	601	YGR	O24-C18	-2.10	1.39	1.43
12	8	601	YGR	O24-C18	-2.10	1.39	1.43
9	B	600	ATP	C2'-C3'	-2.09	1.47	1.53
12	4	601	YGR	C19-C18	2.06	1.56	1.54

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	B	600	ATP	C5-C4-N3	-5.65	118.94	126.72
9	B	600	ATP	N3-C4-N9	4.97	135.62	127.17
12	4	601	YGR	C01-O02-C03	4.28	121.20	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	6	601	YGR	C01-O02-C03	4.13	121.06	117.20
9	B	600	ATP	N3-C2-N1	-3.62	123.11	128.58
9	B	600	ATP	C2-N3-C4	3.55	120.50	111.83
12	9	601	YGR	C25-C18-C05	3.52	115.63	110.54
12	8	601	YGR	C01-O02-C03	3.50	120.47	117.20
12	6	601	YGR	C03-C04-C05	3.49	124.42	119.77
12	7	601	YGR	C01-O02-C03	3.43	120.41	117.20
12	5	601	YGR	C25-C18-C05	3.33	115.36	110.54
12	5	601	YGR	C01-O02-C03	3.31	120.30	117.20
12	8	601	YGR	C25-C18-C05	3.22	115.20	110.54
12	3	601	YGR	C03-C04-C05	3.20	124.03	119.77
12	9	601	YGR	C01-O02-C03	3.18	120.18	117.20
12	9	601	YGR	C03-C04-C05	3.13	123.93	119.77
9	B	600	ATP	O3B-PB-O1B	-3.10	101.37	110.70
12	6	601	YGR	O02-C03-N43	-3.09	115.96	119.93
12	3	601	YGR	C25-C18-C05	3.03	114.93	110.54
12	3	601	YGR	O02-C03-N43	-3.00	116.08	119.93
12	4	601	YGR	C03-C04-C05	2.99	123.75	119.77
12	3	601	YGR	C35-C04-C03	-2.94	113.51	116.23
12	7	601	YGR	C35-C04-C03	-2.91	113.54	116.23
12	9	601	YGR	C35-C04-C03	-2.87	113.58	116.23
9	B	600	ATP	C4-N9-C8	2.87	108.75	105.74
12	5	601	YGR	C03-C04-C05	2.83	123.53	119.77
12	8	601	YGR	C35-C04-C03	-2.82	113.62	116.23
12	5	601	YGR	C35-C04-C03	-2.80	113.64	116.23
12	6	601	YGR	C35-C04-C03	-2.78	113.66	116.23
12	6	601	YGR	C25-C18-C05	2.78	114.57	110.54
12	8	601	YGR	C03-C04-C05	2.77	123.45	119.77
12	4	601	YGR	C35-C04-C03	-2.76	113.68	116.23
12	3	601	YGR	C01-O02-C03	2.75	119.77	117.20
12	4	601	YGR	C25-C18-C05	2.75	114.51	110.54
12	9	601	YGR	O02-C03-N43	-2.72	116.44	119.93
12	9	601	YGR	C10-N13-C14	-2.72	115.25	117.12
12	6	601	YGR	C10-N13-C14	-2.66	115.28	117.12
12	7	601	YGR	O02-C03-N43	-2.66	116.52	119.93
9	B	600	ATP	O2A-PA-O3A	2.60	114.31	107.27
12	8	601	YGR	O02-C03-N43	-2.60	116.59	119.93
12	7	601	YGR	C03-C04-C05	2.60	123.23	119.77
12	7	601	YGR	C10-N13-C14	-2.56	115.36	117.12
12	4	601	YGR	C10-N13-C14	-2.53	115.38	117.12
12	8	601	YGR	C10-N13-C14	-2.52	115.38	117.12
9	B	600	ATP	C4-C5-N7	-2.48	107.74	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	5	601	YGR	C10-N13-C14	-2.45	115.43	117.12
9	B	600	ATP	O2B-PB-O1B	2.43	123.75	112.44
12	4	601	YGR	O02-C03-N43	-2.42	116.82	119.93
12	4	601	YGR	O24-C18-C19	-2.42	104.83	108.26
12	5	601	YGR	O02-C03-N43	-2.42	116.83	119.93
12	6	601	YGR	O02-C03-C04	2.42	117.92	115.05
12	9	601	YGR	C06-C17-C14	2.41	120.76	118.61
12	3	601	YGR	C10-N13-C14	-2.40	115.46	117.12
12	7	601	YGR	C25-C18-C05	2.37	113.98	110.54
12	5	601	YGR	O24-C18-C19	-2.32	104.98	108.26
12	9	601	YGR	O24-C18-C19	-2.28	105.03	108.26
9	B	600	ATP	C5-N7-C8	2.27	107.01	103.45
12	7	601	YGR	O24-C18-C19	-2.24	105.09	108.26
12	3	601	YGR	O02-C03-C04	2.23	117.69	115.05
12	8	601	YGR	O24-C18-C19	-2.22	105.11	108.26
12	3	601	YGR	O24-C18-C19	-2.21	105.13	108.26
12	9	601	YGR	C17-C06-C07	-2.19	114.72	117.97
12	6	601	YGR	O24-C18-C19	-2.17	105.18	108.26
12	6	601	YGR	C06-C17-C14	2.17	120.55	118.61
12	4	601	YGR	C06-C17-C14	2.16	120.53	118.61
12	3	601	YGR	C06-C17-C14	2.14	120.52	118.61
9	B	600	ATP	N6-C6-N1	2.09	123.03	118.38
12	6	601	YGR	C17-C06-C07	-2.08	114.88	117.97
12	7	601	YGR	C06-C17-C14	2.08	120.46	118.61
9	C	600	ATP	O3'-C3'-C2'	-2.07	105.17	111.82
9	B	600	ATP	C2-N1-C6	2.05	122.10	118.73
12	5	601	YGR	C06-C17-C14	2.05	120.43	118.61
9	A	600	ATP	O3'-C3'-C2'	-2.04	105.29	111.82
9	B	600	ATP	N9-C8-N7	-2.03	111.05	113.94

There are no chirality outliers.

All (90) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	600	ATP	PB-O3B-PG-O2G
9	A	600	ATP	C5'-O5'-PA-O1A
9	B	600	ATP	PB-O3B-PG-O3G
9	C	600	ATP	O4'-C4'-C5'-O5'
9	E	600	ATP	PB-O3B-PG-O2G
9	E	600	ATP	C5'-O5'-PA-O2A
12	9	601	YGR	C04-C03-O02-C01
12	9	601	YGR	N43-C03-O02-C01

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Mol	Chain	Res	Type	Atoms
12	3	601	YGR	C07-C10-O11-C12
12	3	601	YGR	N13-C10-O11-C12
12	3	601	YGR	C26-C27-O28-C29
12	3	601	YGR	N30-C27-O28-C29
12	4	601	YGR	C04-C03-O02-C01
12	4	601	YGR	N43-C03-O02-C01
12	4	601	YGR	C07-C10-O11-C12
12	4	601	YGR	N13-C10-O11-C12
12	4	601	YGR	C34-C31-O32-C33
12	4	601	YGR	N30-C31-O32-C33
12	5	601	YGR	C04-C03-O02-C01
12	5	601	YGR	N43-C03-O02-C01
12	5	601	YGR	C07-C10-O11-C12
12	5	601	YGR	N13-C10-O11-C12
12	5	601	YGR	N13-C14-O15-C16
12	5	601	YGR	C26-C27-O28-C29
12	5	601	YGR	N30-C27-O28-C29
12	6	601	YGR	C04-C03-O02-C01
12	6	601	YGR	N43-C03-O02-C01
12	6	601	YGR	C07-C10-O11-C12
12	6	601	YGR	N13-C10-O11-C12
12	7	601	YGR	C07-C10-O11-C12
12	7	601	YGR	N13-C10-O11-C12
12	7	601	YGR	C17-C14-O15-C16
12	7	601	YGR	N13-C14-O15-C16
12	7	601	YGR	C26-C27-O28-C29
12	7	601	YGR	N30-C27-O28-C29
12	8	601	YGR	C04-C03-O02-C01
12	8	601	YGR	C07-C10-O11-C12
12	8	601	YGR	N13-C10-O11-C12
12	8	601	YGR	C17-C14-O15-C16
12	8	601	YGR	N13-C14-O15-C16
12	9	601	YGR	C26-C27-O28-C29
12	4	601	YGR	C26-C27-O28-C29
12	5	601	YGR	C17-C14-O15-C16
12	8	601	YGR	C26-C27-O28-C29
12	9	601	YGR	N30-C27-O28-C29
12	4	601	YGR	N30-C27-O28-C29
12	8	601	YGR	N30-C31-O32-C33
12	9	601	YGR	N13-C10-O11-C12
9	C	600	ATP	C3'-C4'-C5'-O5'
9	E	600	ATP	C3'-C4'-C5'-O5'

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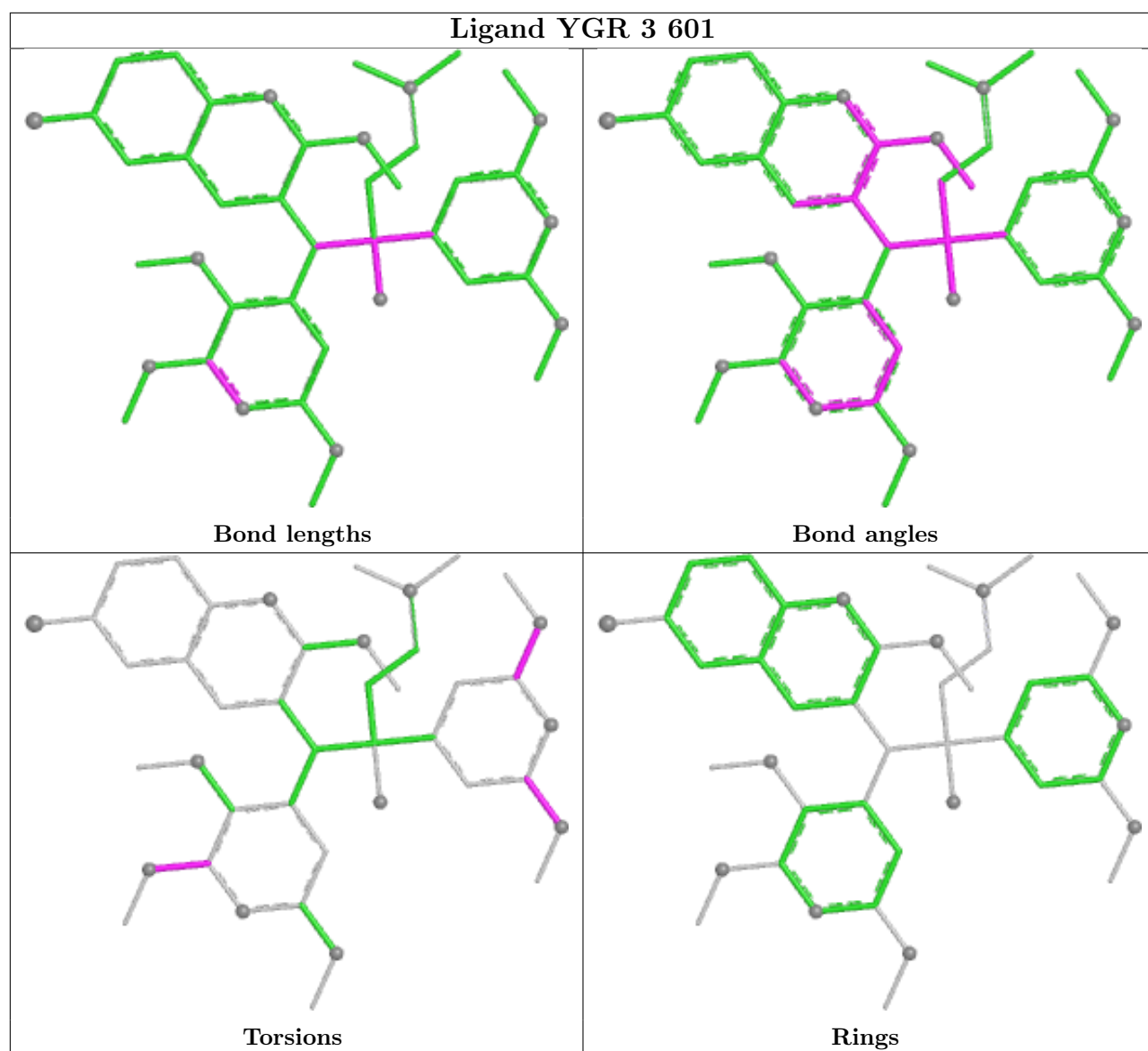
Mol	Chain	Res	Type	Atoms
12	8	601	YGR	C34-C31-O32-C33
12	9	601	YGR	N30-C31-O32-C33
12	8	601	YGR	N30-C27-O28-C29
12	7	601	YGR	C19-C20-N21-C23
12	8	601	YGR	N43-C03-O02-C01
9	E	600	ATP	O4'-C4'-C5'-O5'
12	9	601	YGR	N13-C14-O15-C16
12	3	601	YGR	N30-C31-O32-C33
12	5	601	YGR	N30-C31-O32-C33
12	6	601	YGR	N13-C14-O15-C16
12	6	601	YGR	N30-C31-O32-C33
12	9	601	YGR	C17-C14-O15-C16
12	9	601	YGR	C34-C31-O32-C33
12	3	601	YGR	C34-C31-O32-C33
12	5	601	YGR	C34-C31-O32-C33
12	6	601	YGR	C17-C14-O15-C16
12	6	601	YGR	C34-C31-O32-C33
12	9	601	YGR	C07-C10-O11-C12
12	7	601	YGR	C19-C20-N21-C22
12	7	601	YGR	C34-C31-O32-C33
12	7	601	YGR	N30-C31-O32-C33
12	4	601	YGR	C17-C14-O15-C16
12	4	601	YGR	N13-C14-O15-C16
9	A	600	ATP	PA-O3A-PB-O1B
9	E	600	ATP	PG-O3B-PB-O1B
9	E	600	ATP	C5'-O5'-PA-O1A
9	E	600	ATP	C5'-O5'-PA-O3A
9	E	600	ATP	PB-O3B-PG-O1G
9	C	600	ATP	PA-O3A-PB-O1B
9	C	600	ATP	PA-O3A-PB-O2B
9	B	600	ATP	PA-O3A-PB-O2B
9	B	600	ATP	PB-O3B-PG-O1G
9	A	600	ATP	PG-O3B-PB-O2B
9	A	600	ATP	PA-O3A-PB-O2B
9	C	600	ATP	PG-O3B-PB-O1B
9	C	600	ATP	PG-O3B-PB-O2B
9	E	600	ATP	PA-O3A-PB-O1B
9	E	600	ATP	PA-O3A-PB-O2B
12	9	601	YGR	C25-C18-C19-C20
9	E	600	ATP	PG-O3B-PB-O2B

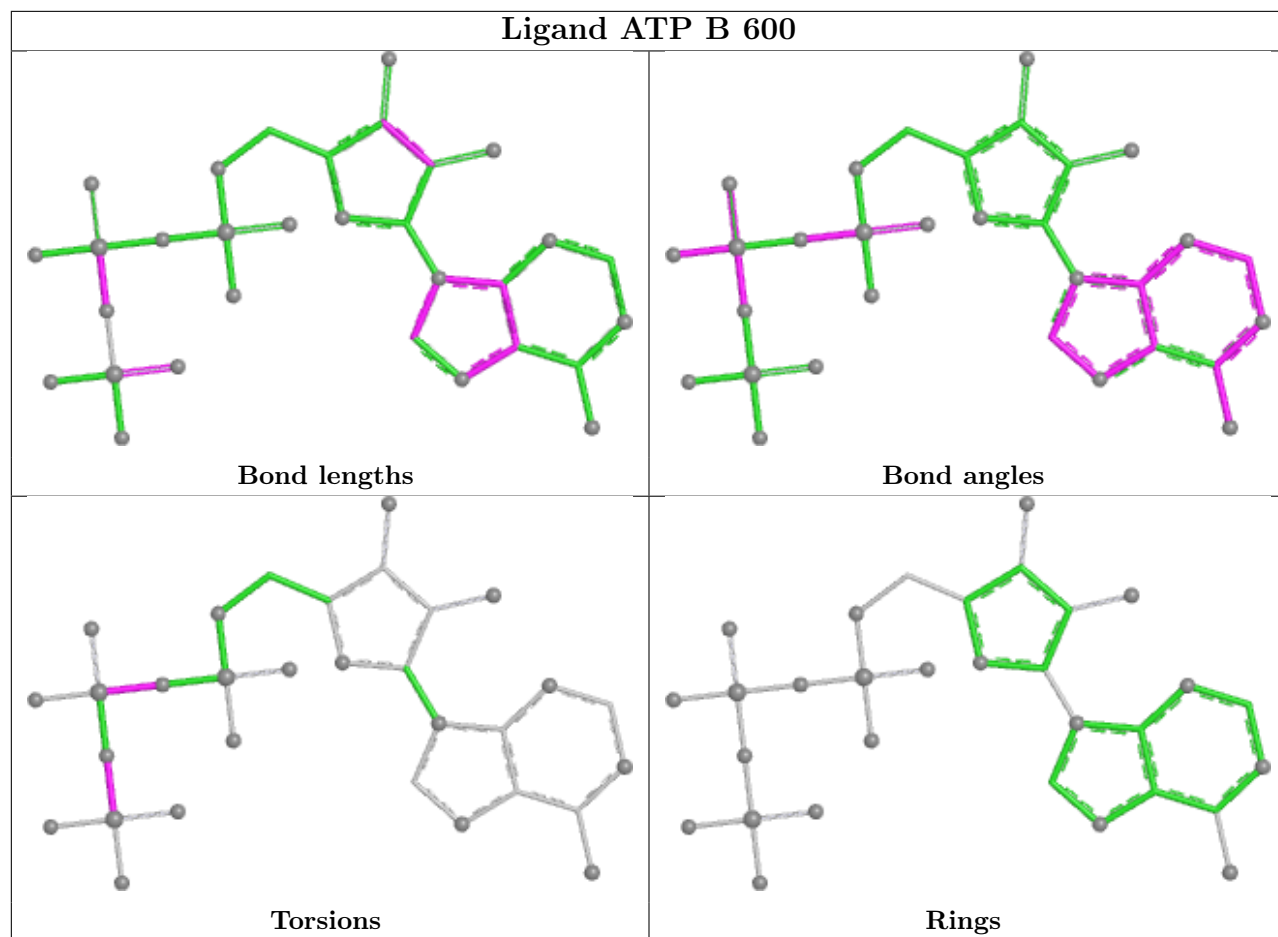
There are no ring outliers.

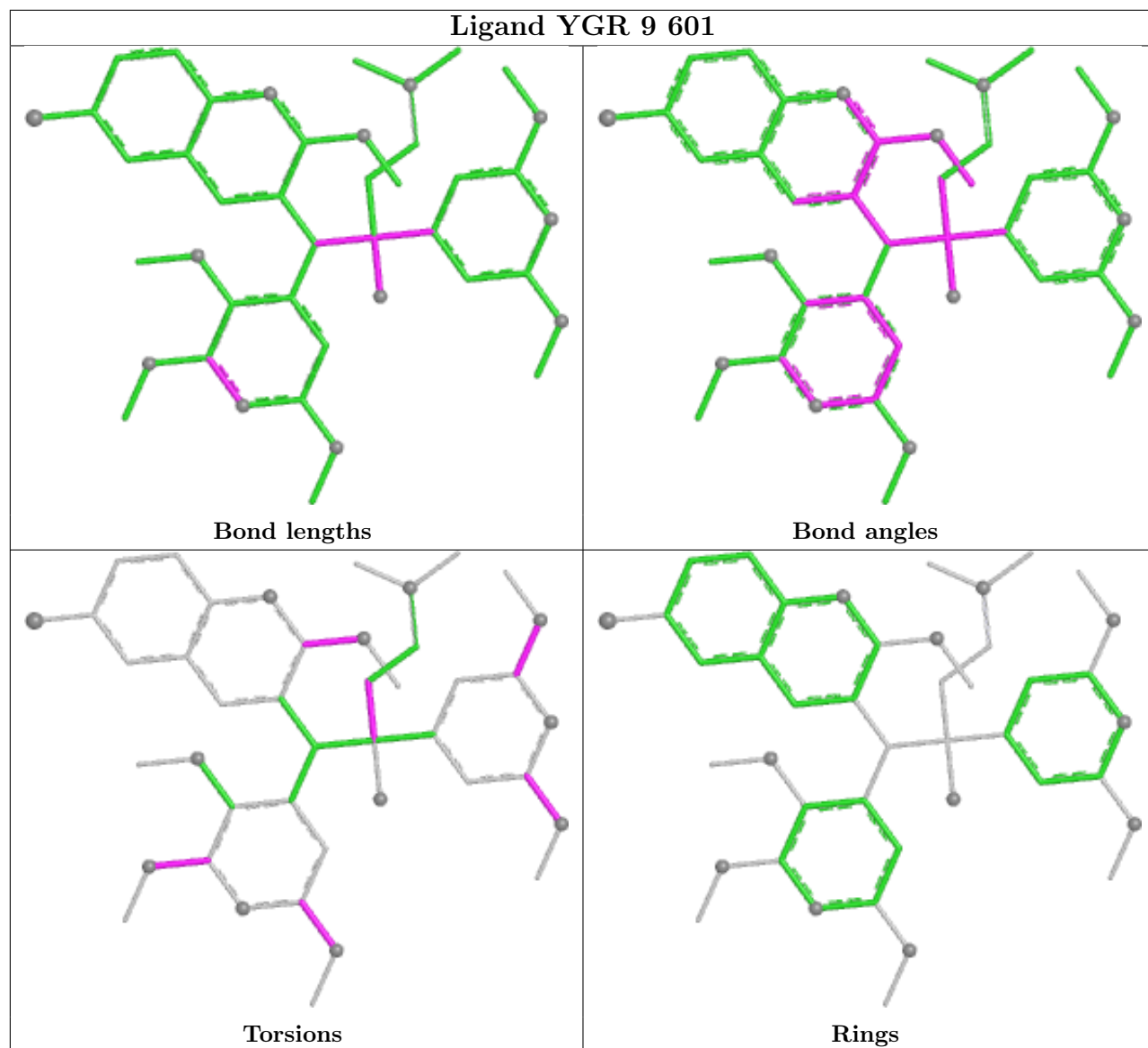
9 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
12	3	601	YGR	1	0
12	9	601	YGR	1	0
12	8	601	YGR	1	0
12	7	601	YGR	1	0
12	6	601	YGR	1	0
9	E	600	ATP	1	0
12	5	601	YGR	1	0
9	A	600	ATP	1	0
12	4	601	YGR	1	0

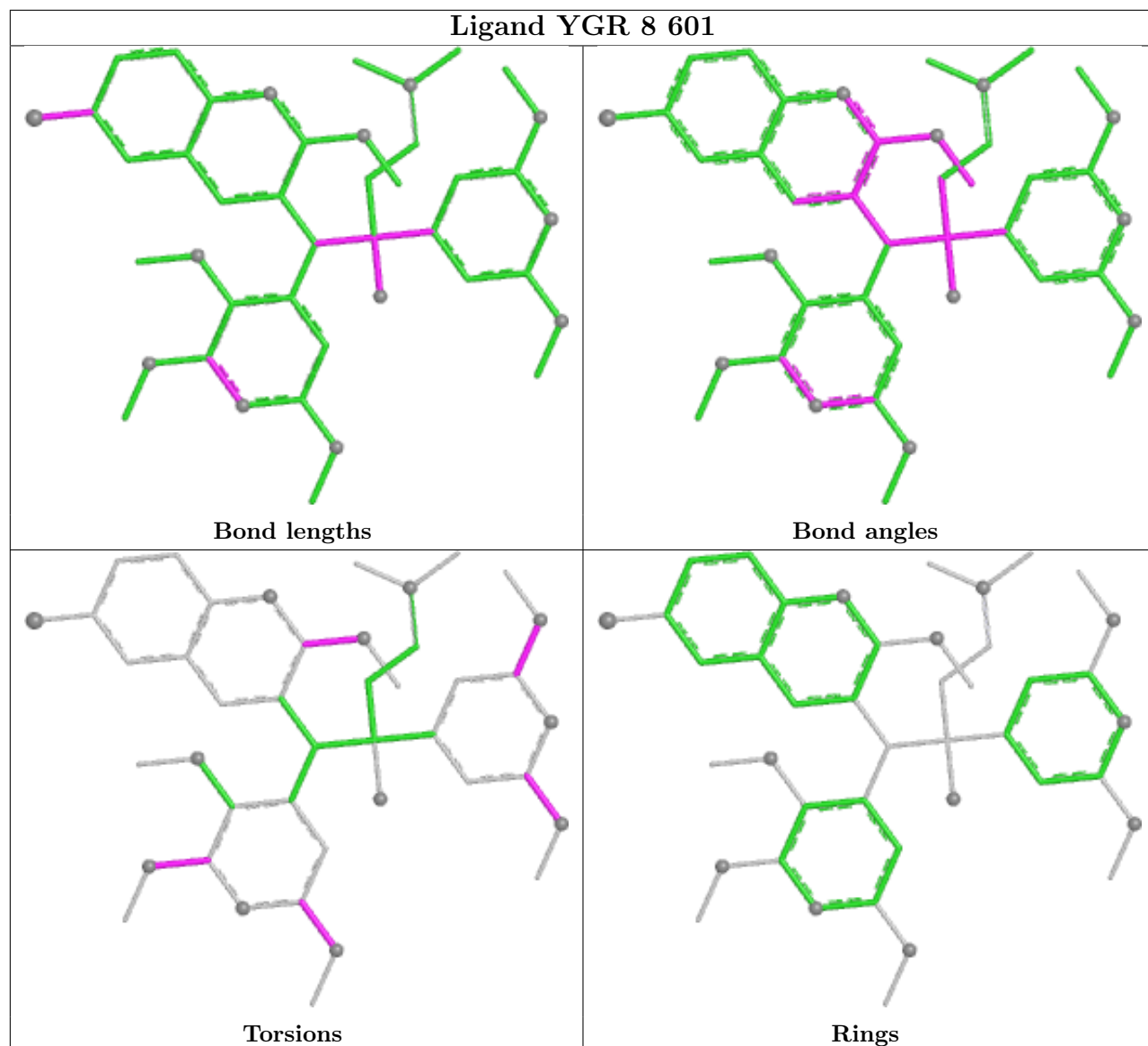
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

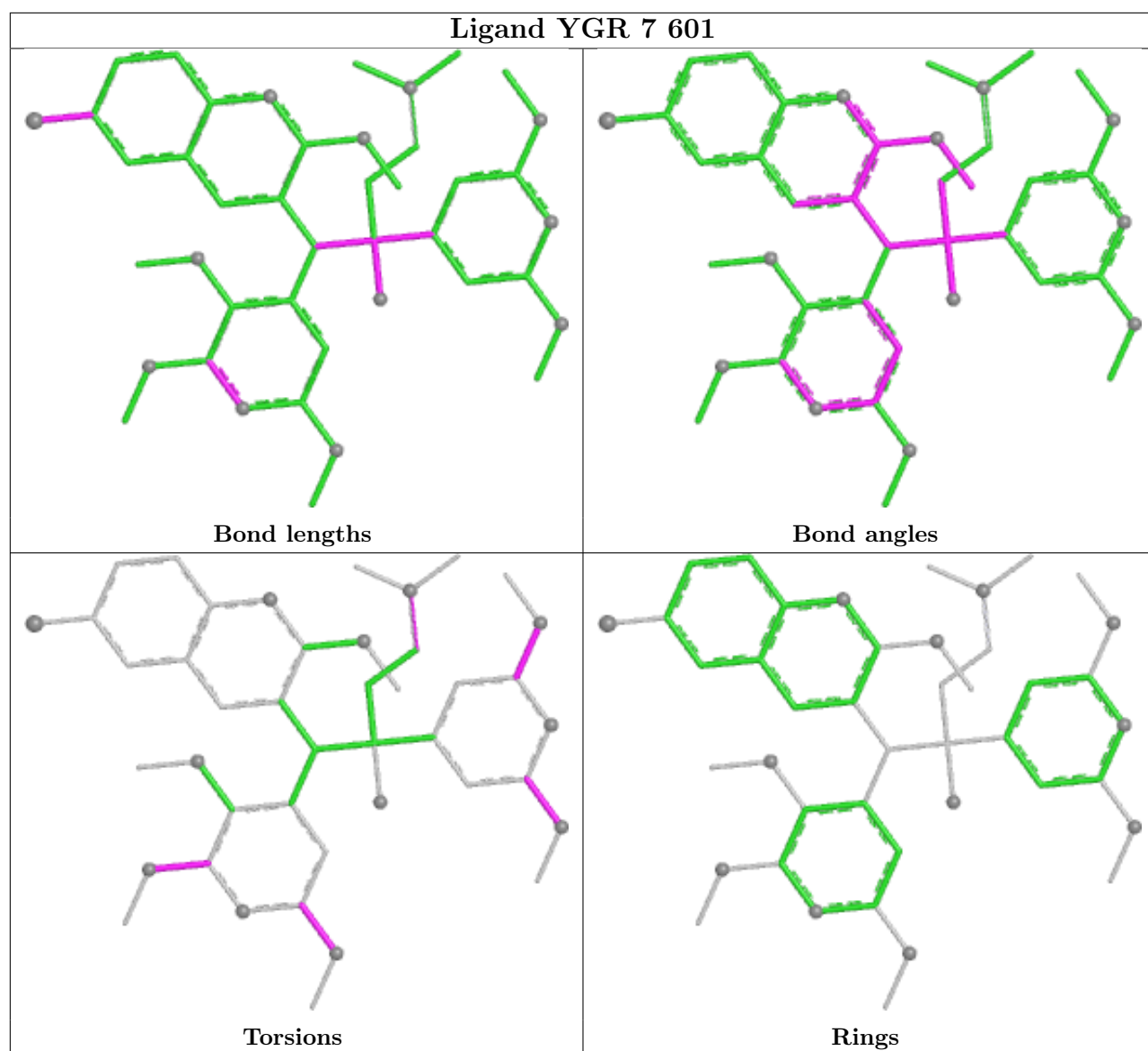


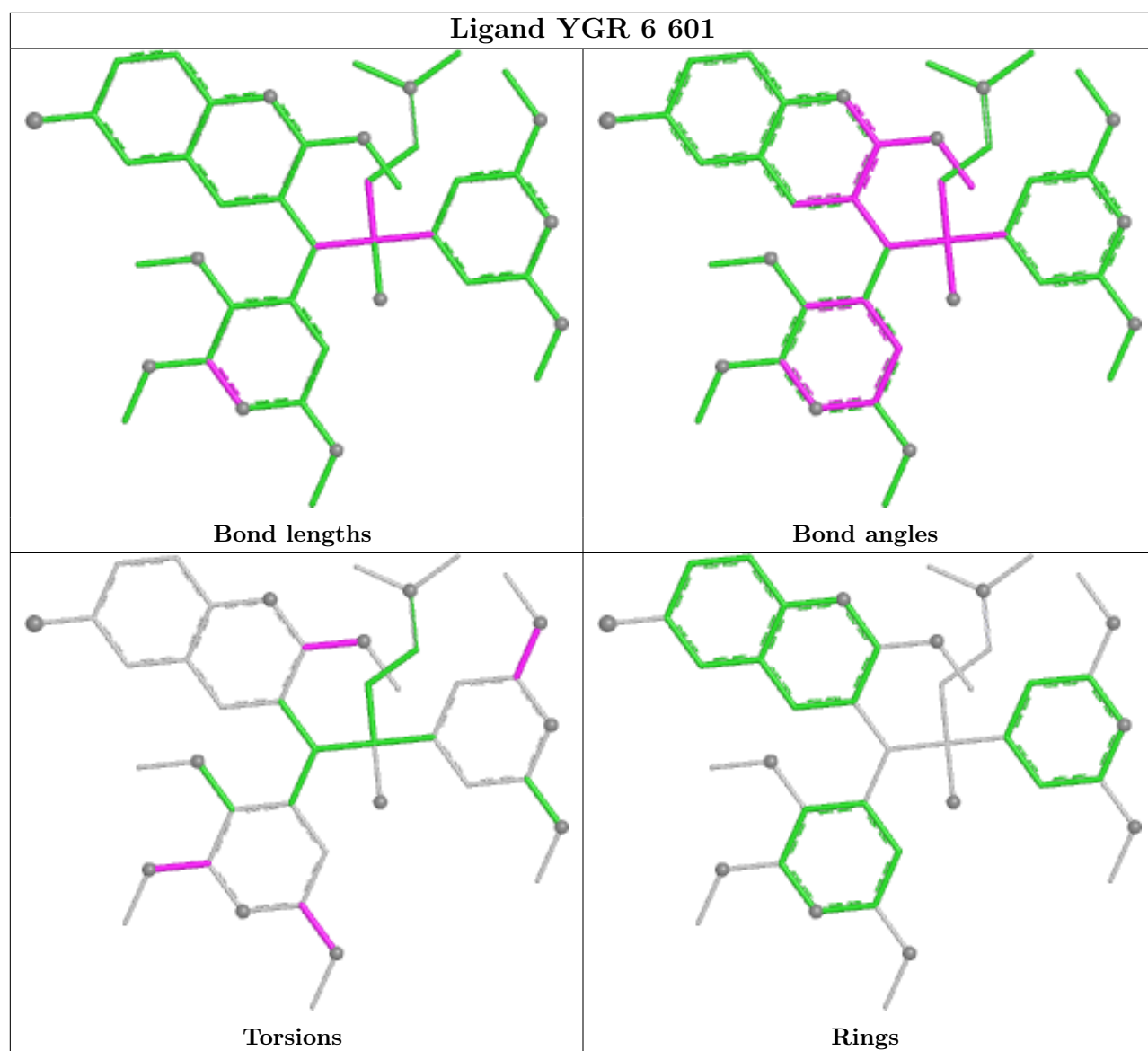


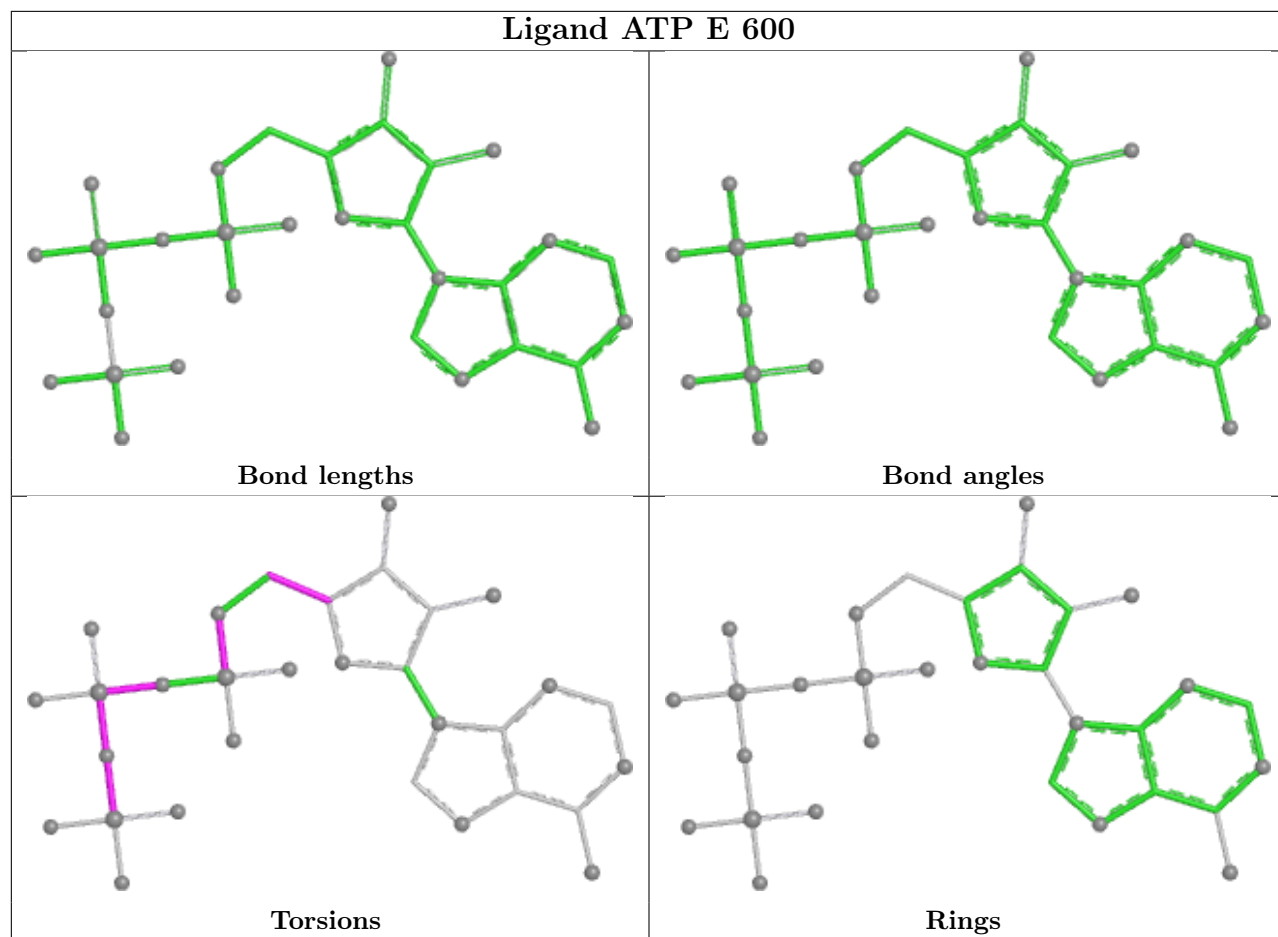


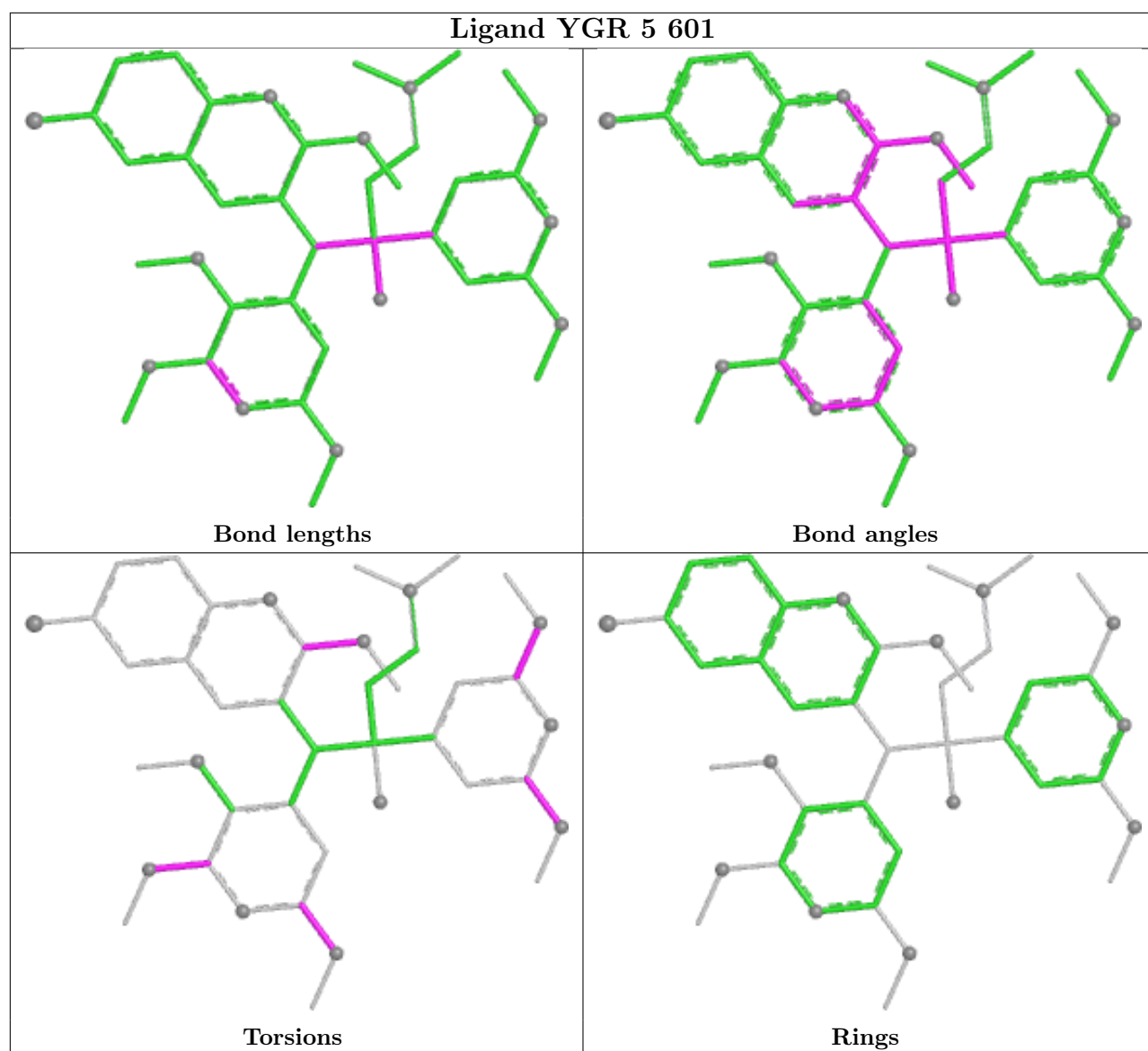
Ligand YGR 8 601

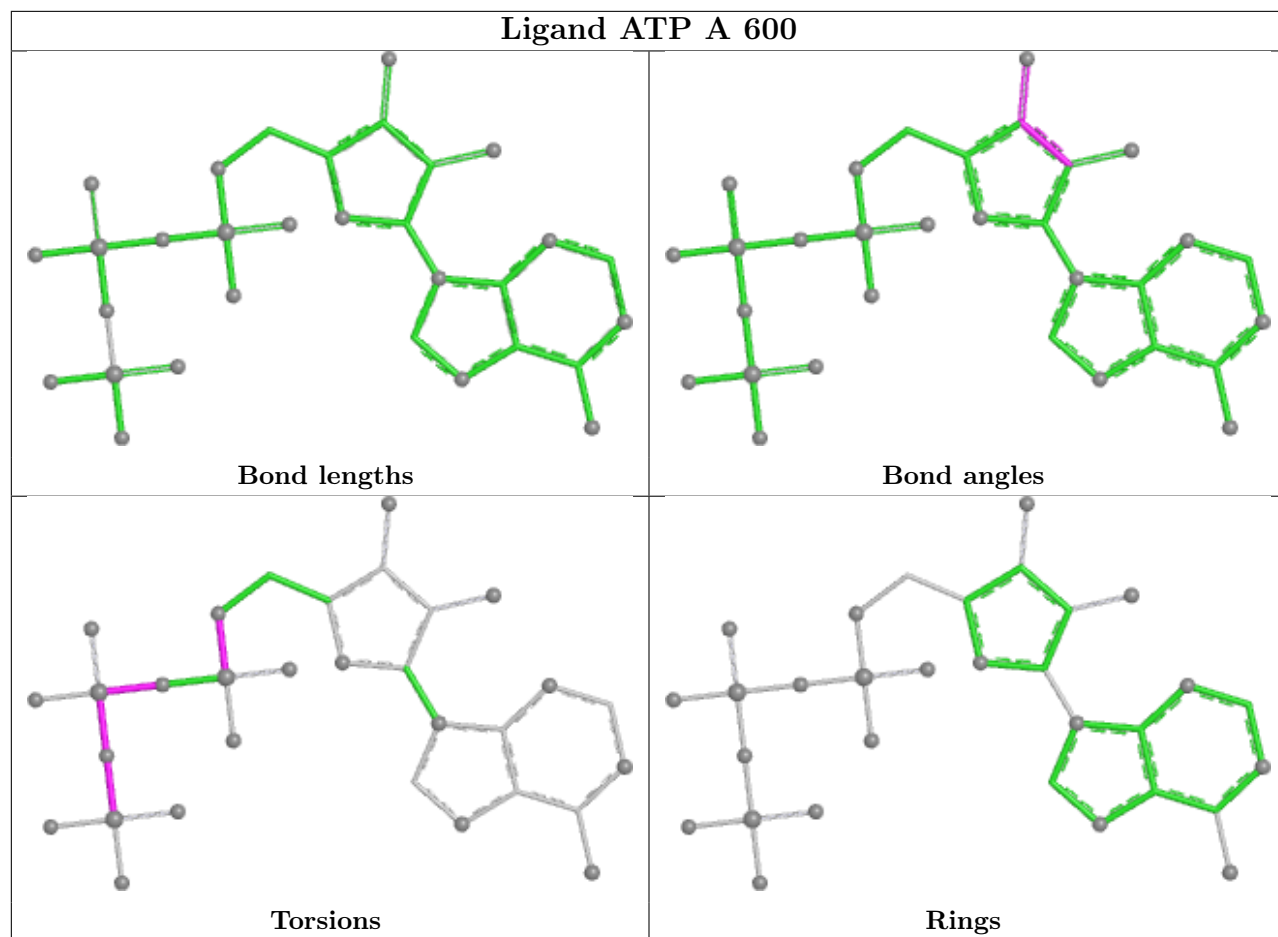


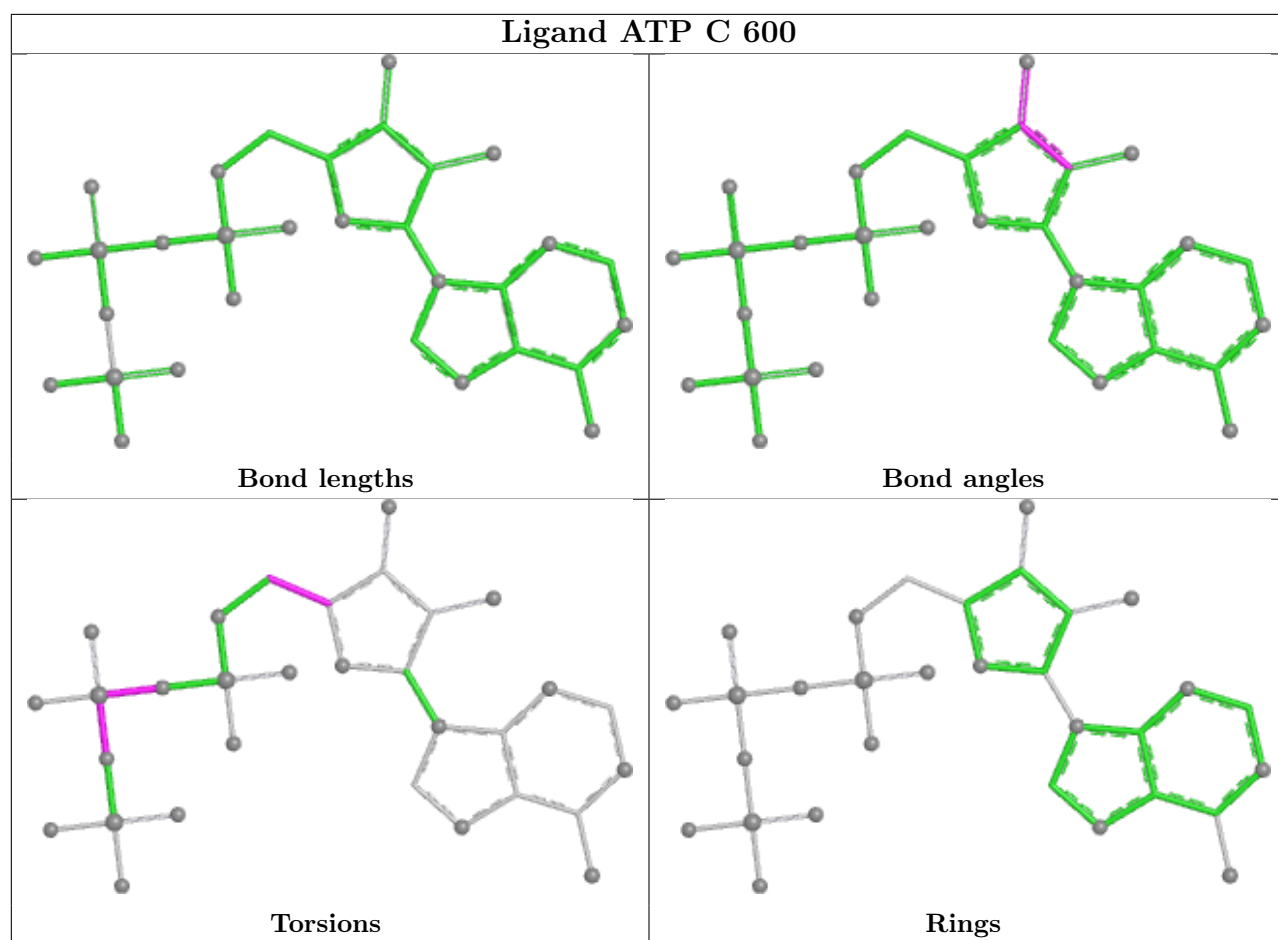


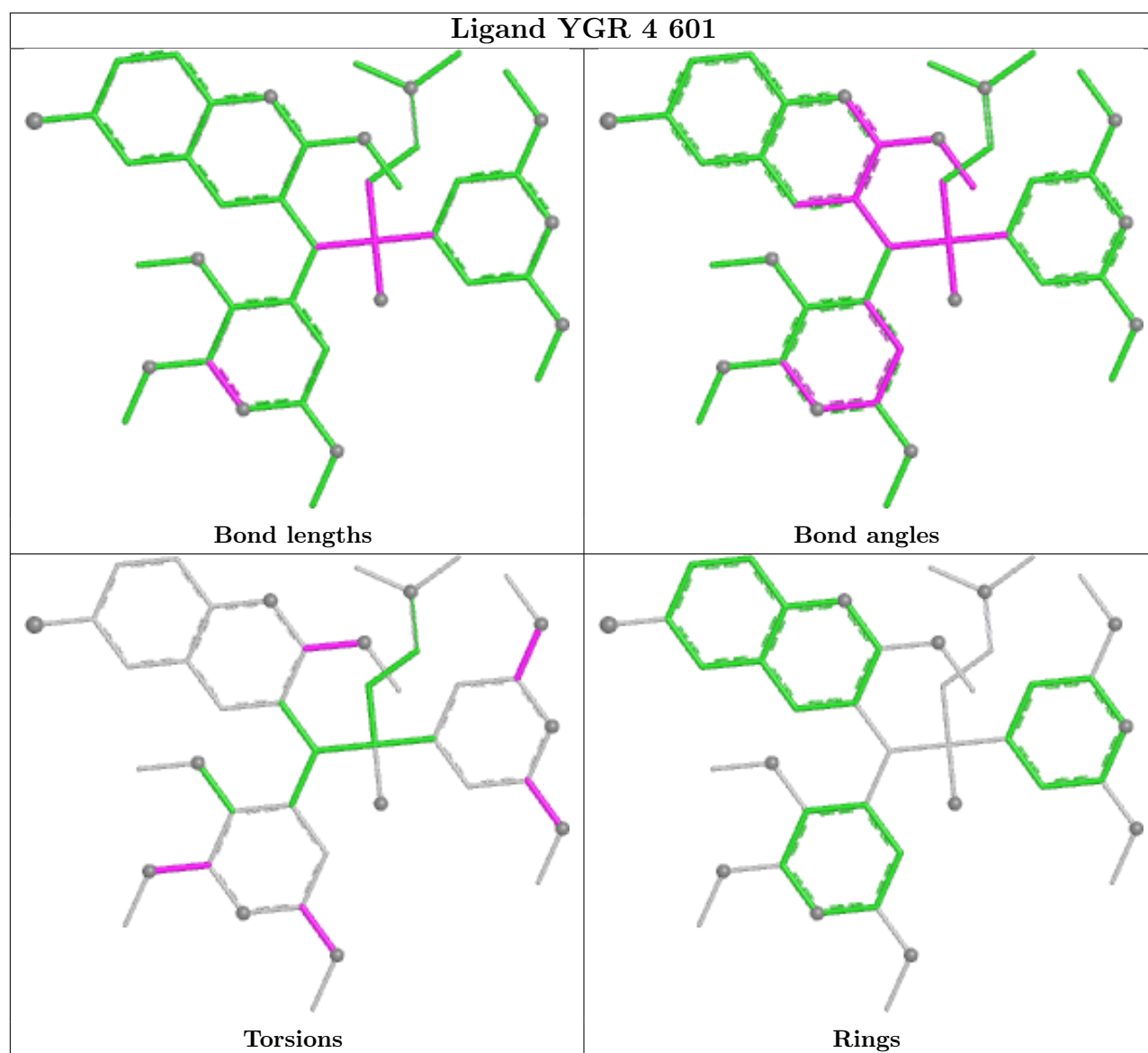












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

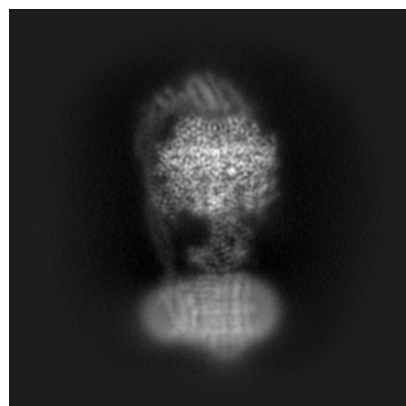
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29655. These allow visual inspection of the internal detail of the map and identification of artifacts.

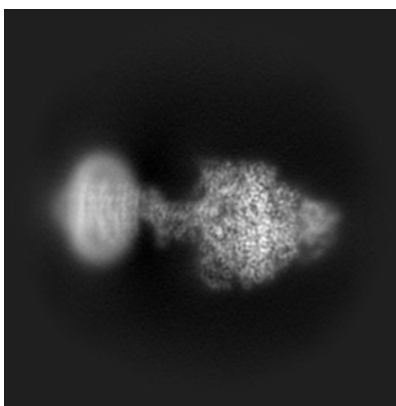
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

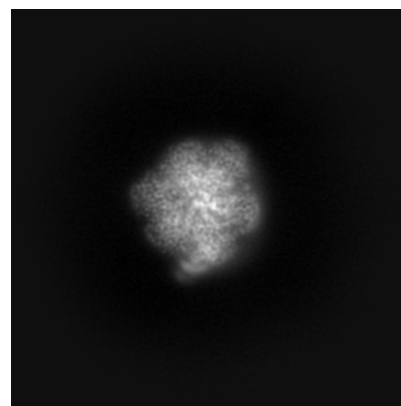
6.1.1 Primary map



X

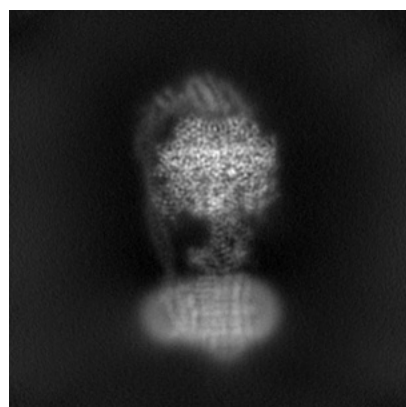


Y

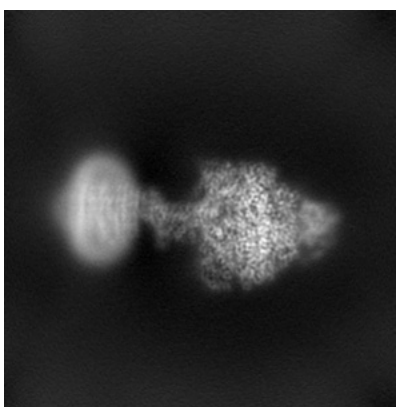


Z

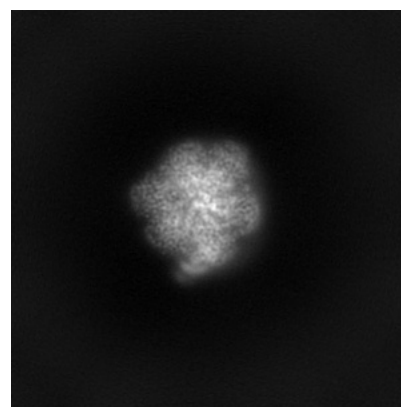
6.1.2 Raw map



X



Y

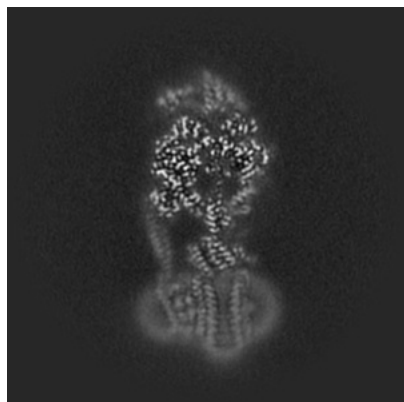


Z

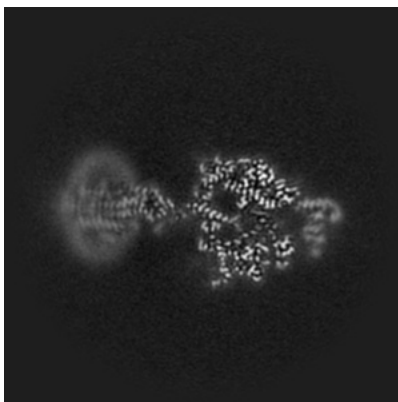
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

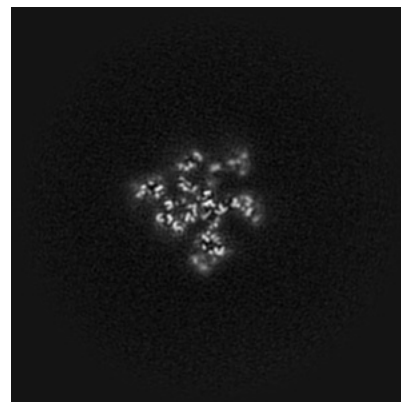
6.2.1 Primary map



X Index: 160

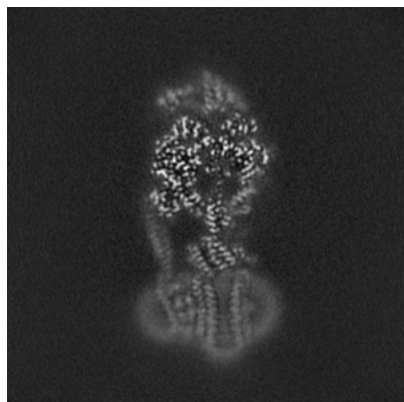


Y Index: 160

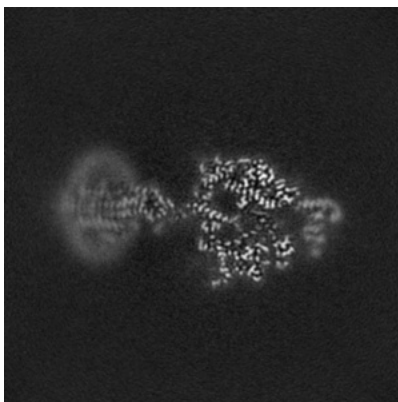


Z Index: 160

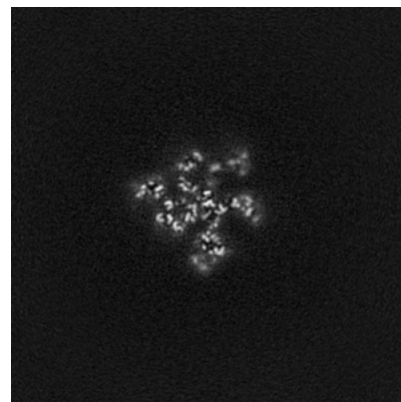
6.2.2 Raw map



X Index: 160



Y Index: 160

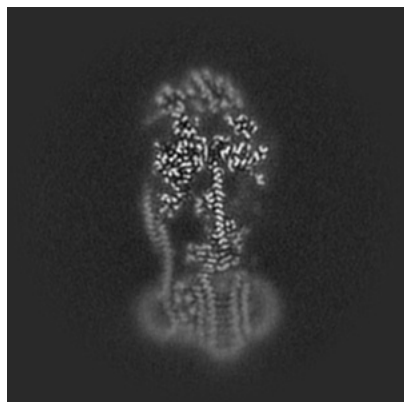


Z Index: 160

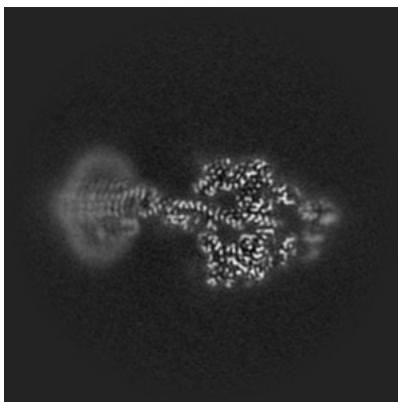
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

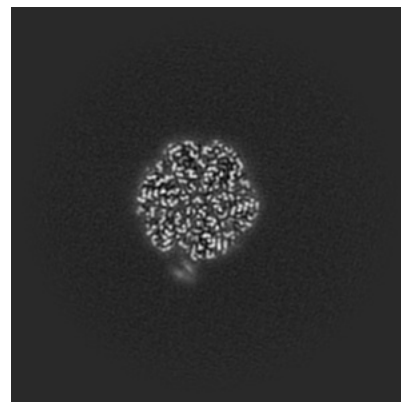
6.3.1 Primary map



X Index: 155

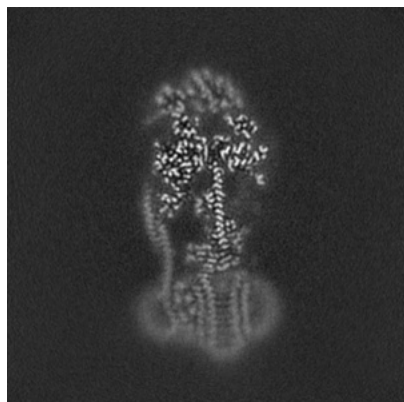


Y Index: 164

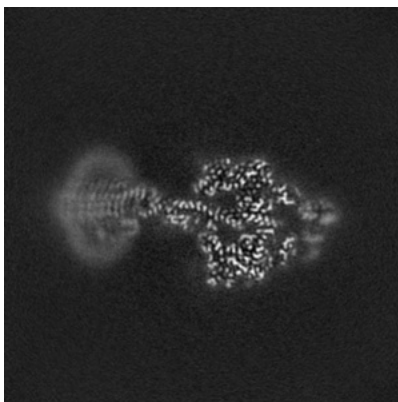


Z Index: 205

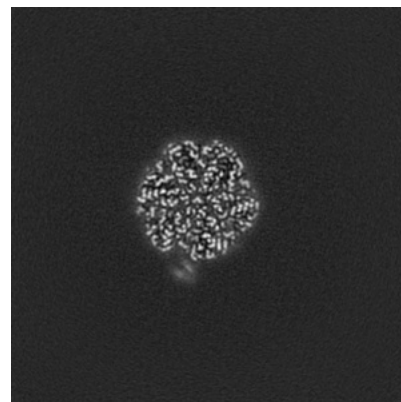
6.3.2 Raw map



X Index: 155



Y Index: 164

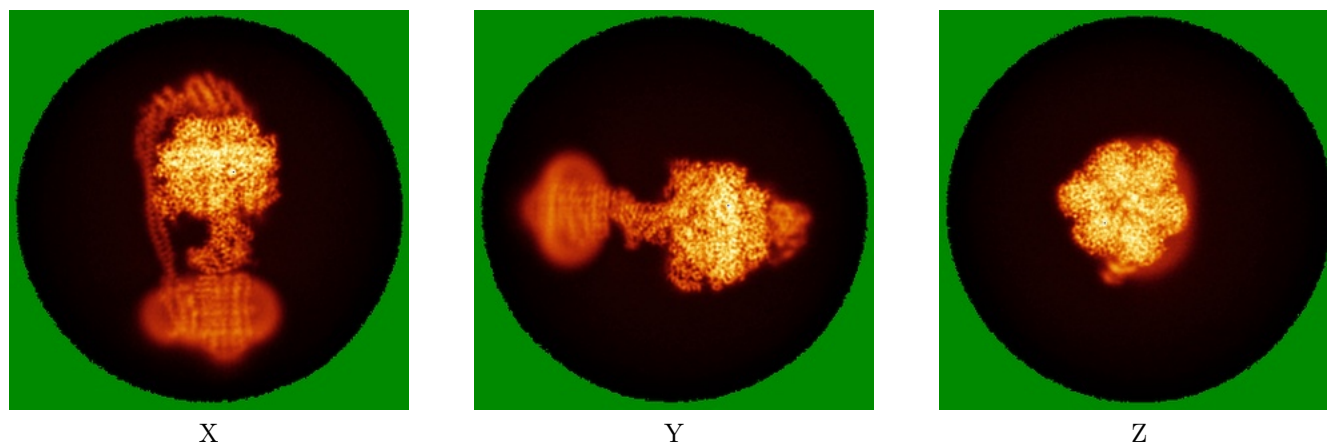


Z Index: 205

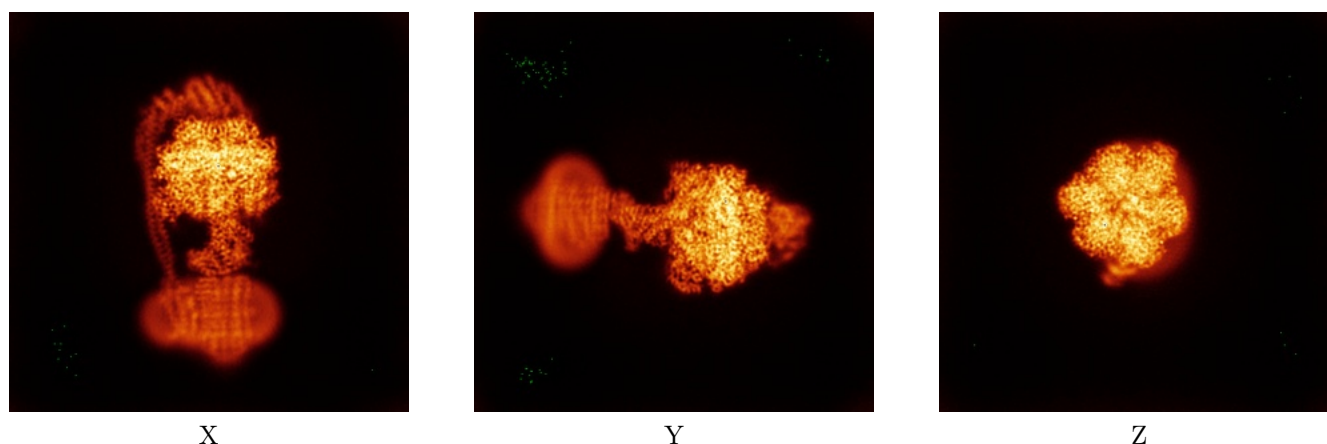
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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

6.5 Orthogonal surface views [i](#)

This section was not generated.

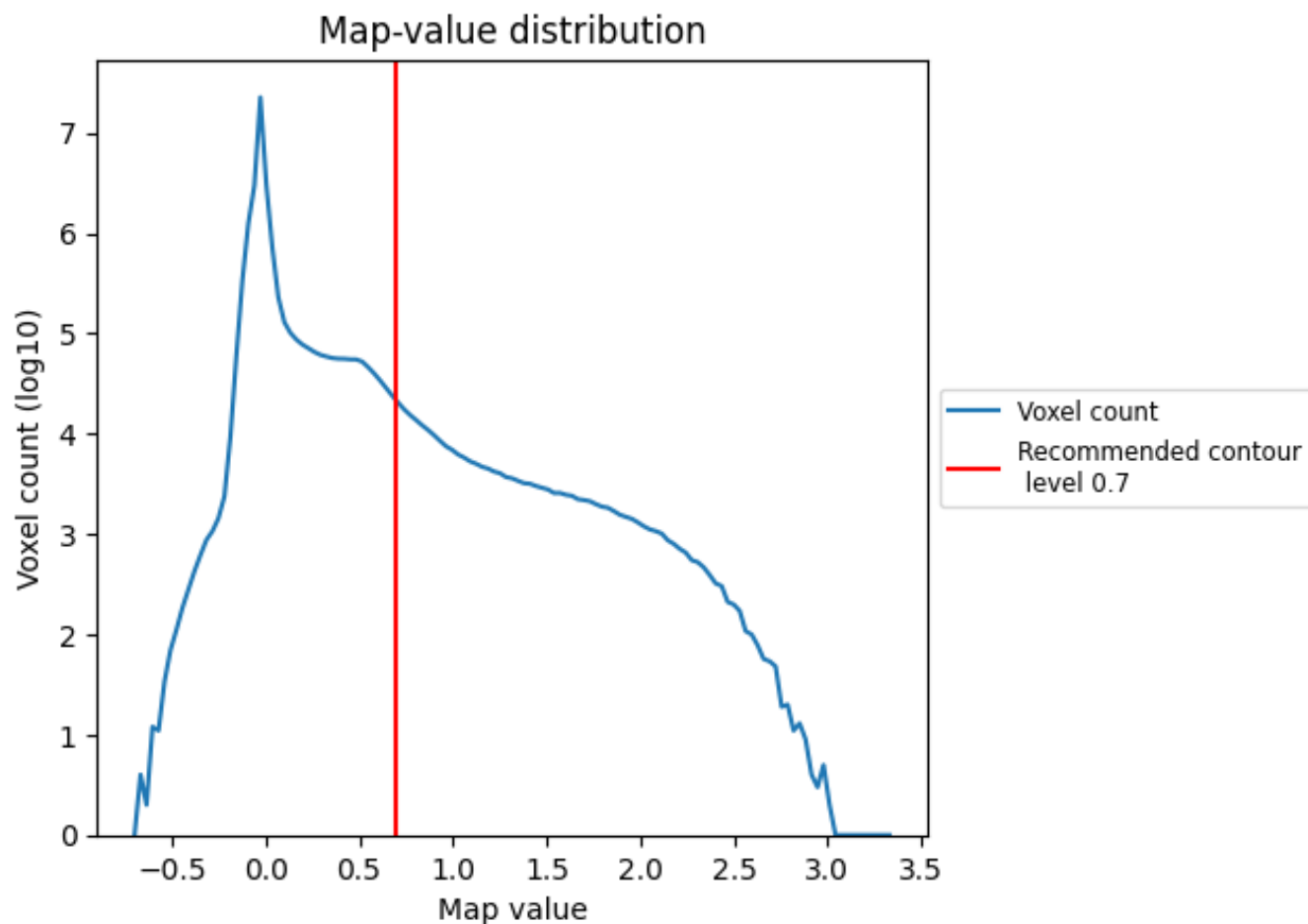
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

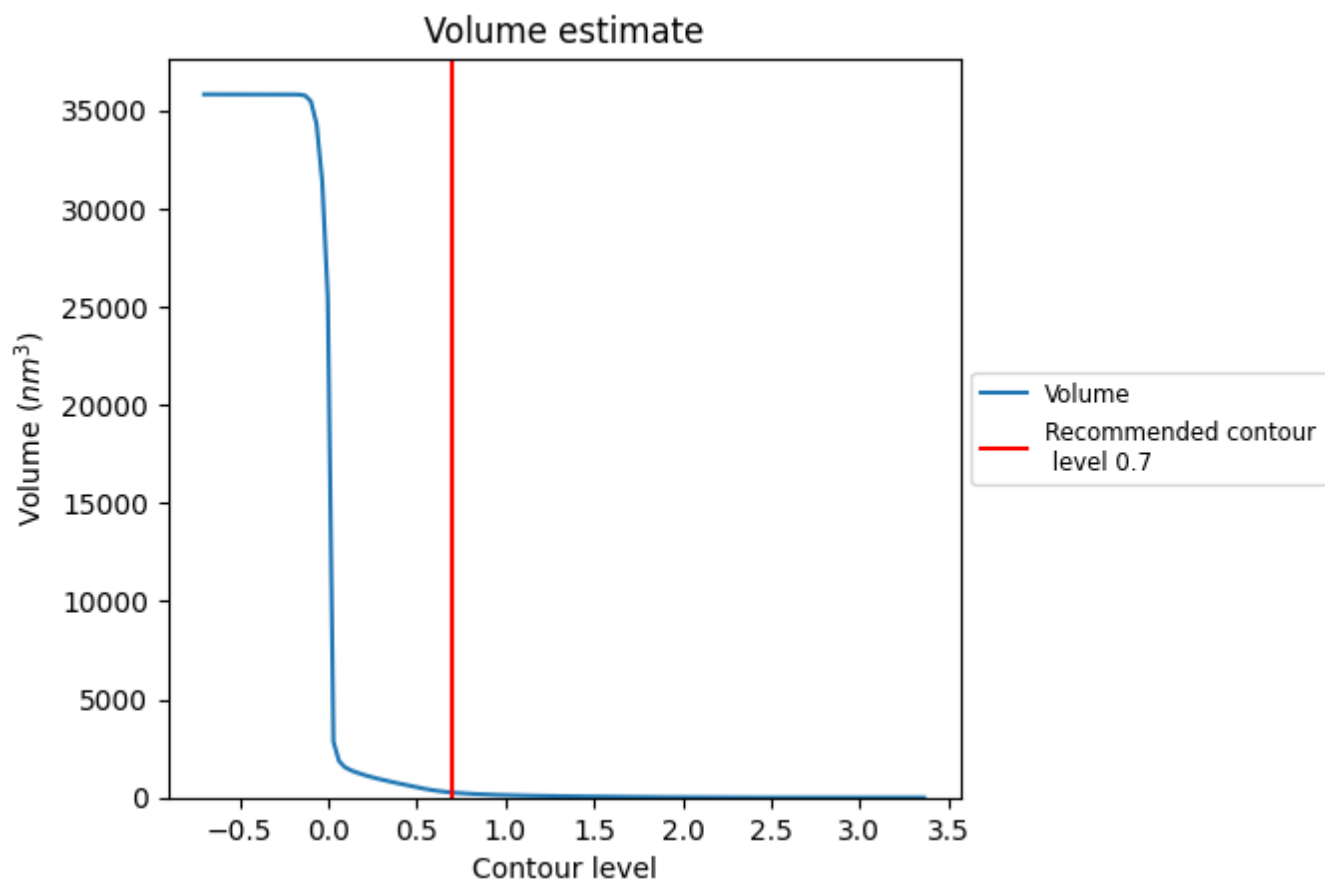
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

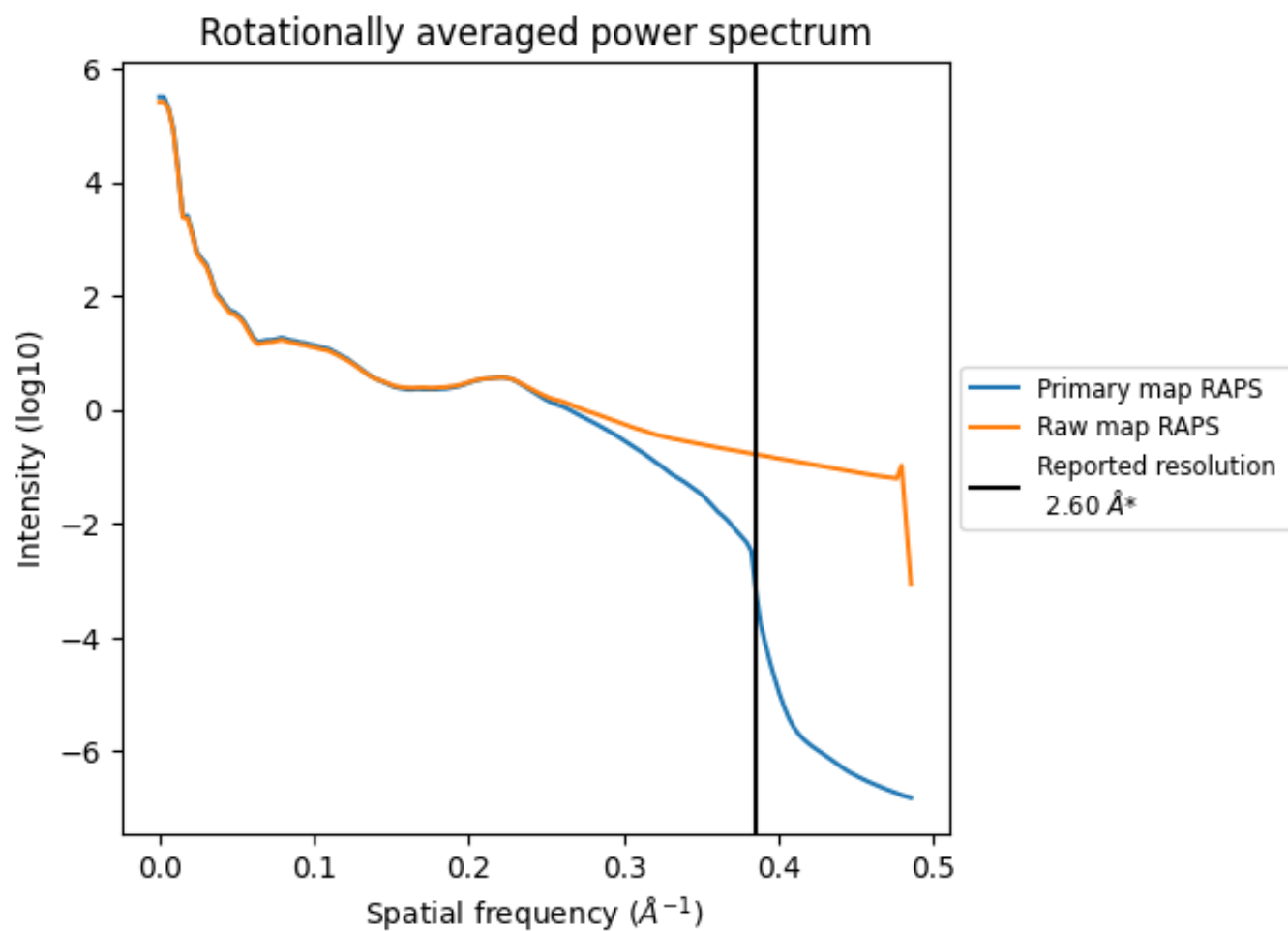
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 257 nm³; this corresponds to an approximate mass of 232 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 ⓘ

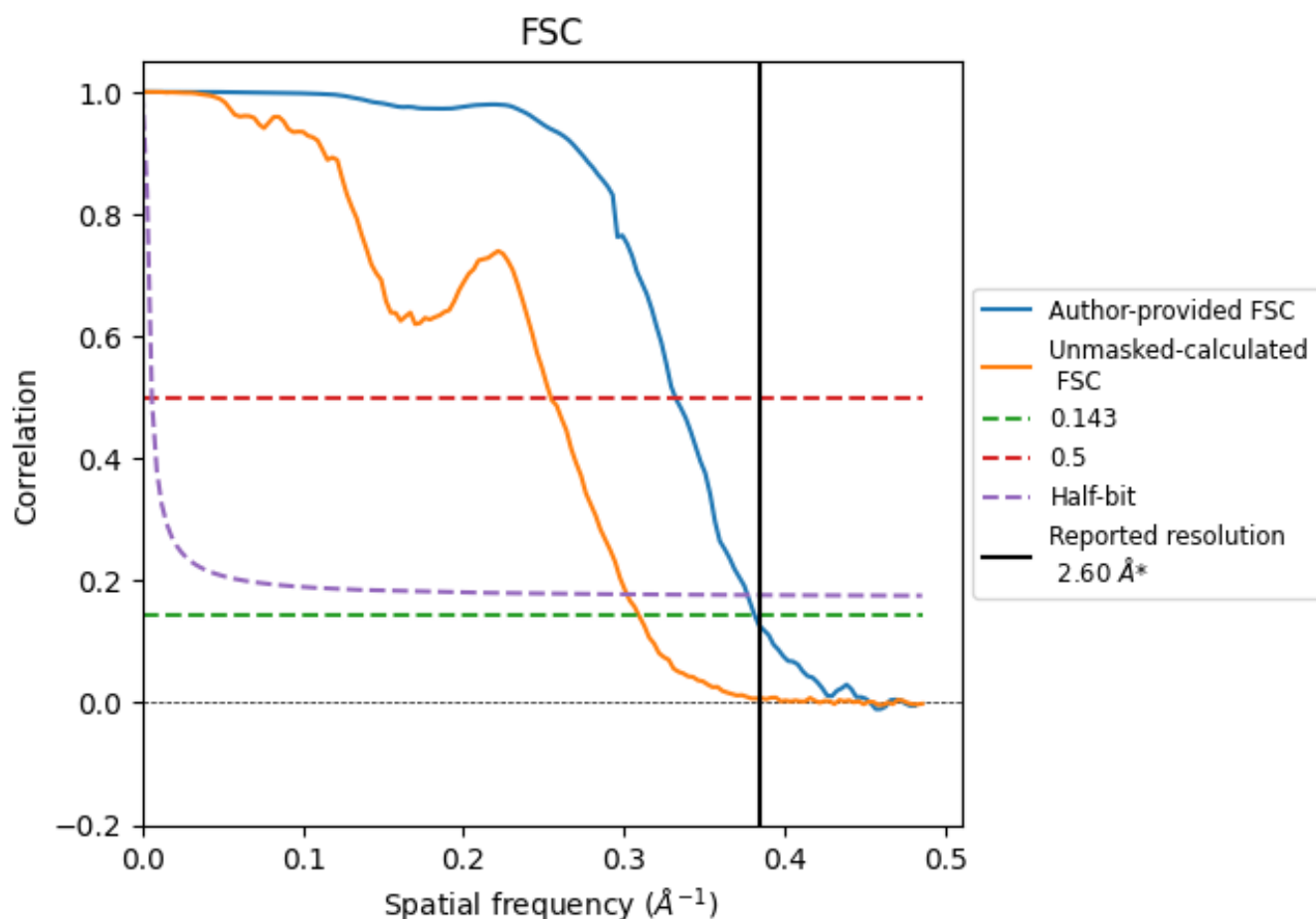


*Reported resolution corresponds to spatial frequency of $0.385 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of $0.385 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.60	-	-
Author-provided FSC curve	2.62	3.01	2.65
Unmasked-calculated*	3.24	3.93	3.32

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.24 differs from the reported value 2.6 by more than 10 %

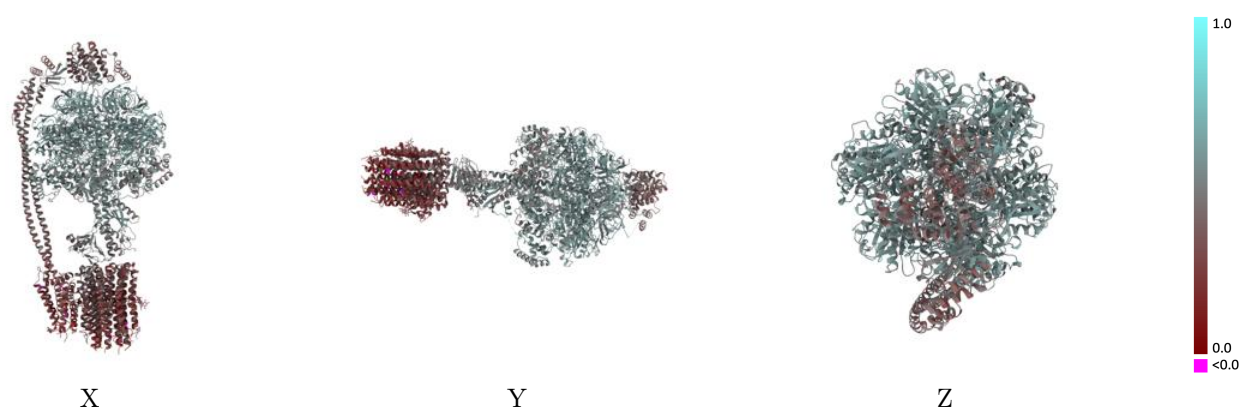
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29655 and PDB model 8G0E. Per-residue inclusion information can be found in section 3 on page 9.

9.1 Map-model overlay [i](#)

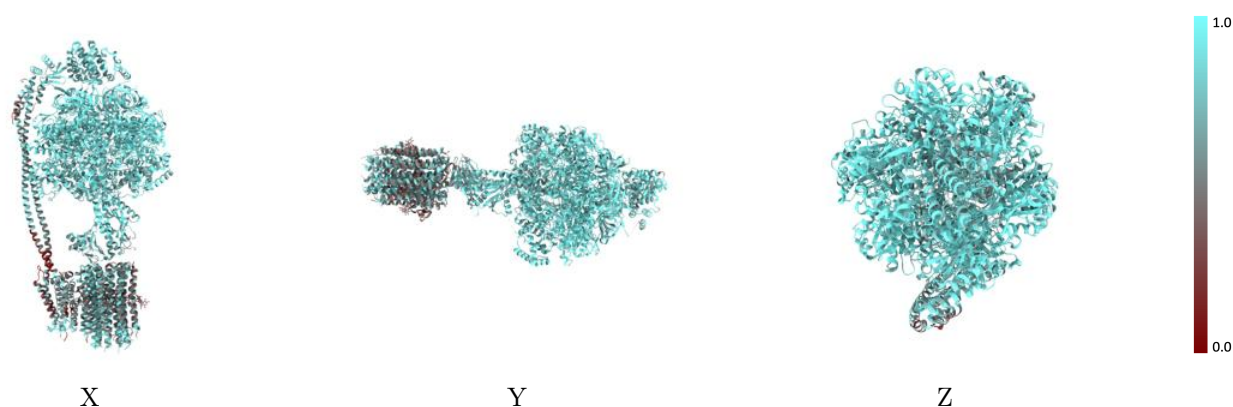
This section was not generated.

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



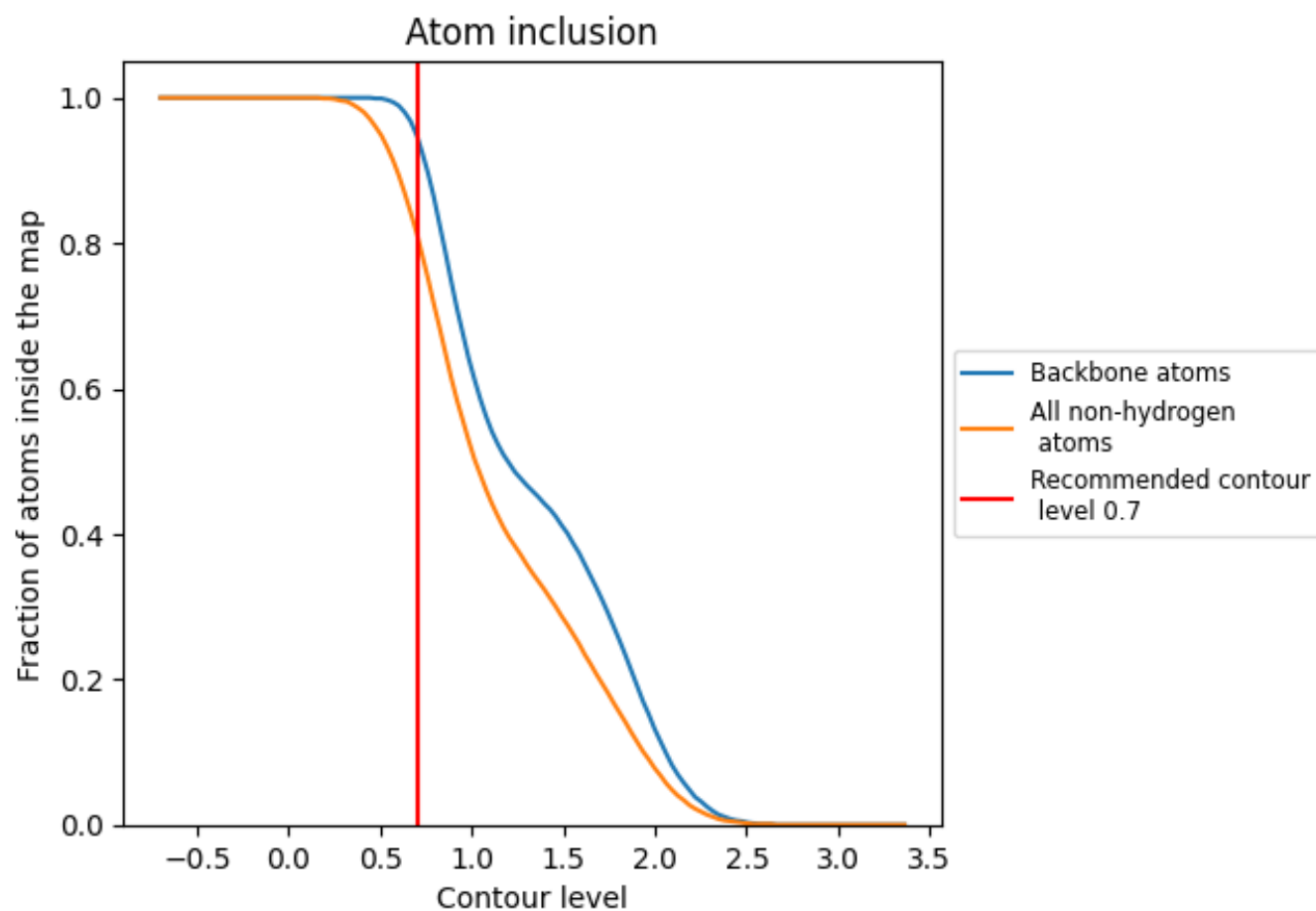
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



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 (0.7).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 81% 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 (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8130	0.4770
1	0.6620	0.3170
2	0.7080	0.3370
3	0.6650	0.3150
4	0.5900	0.2940
5	0.5790	0.2660
6	0.5720	0.2550
7	0.5940	0.2550
8	0.5770	0.2470
9	0.5790	0.2630
A	0.8920	0.5550
B	0.8880	0.5450
C	0.9120	0.5550
D	0.9050	0.5540
E	0.9090	0.5630
F	0.9110	0.5610
G	0.8600	0.5220
H	0.7550	0.4760
a	0.5390	0.2410
b	0.5750	0.3520
d	0.7390	0.3890

1.0

0.0

<0.0