



wwPDB EM Validation Summary Report ⓘ

Mar 9, 2026 – 01:58 PM UTC

PDB ID : 7RMP / pdb_00007rmp
EMDB ID : EMD-24577
Title : Structure of ACLY D1026A - substrates-asym
Authors : Wei, X.; Marmorstein, R.
Deposited on : 2021-07-28
Resolution : 2.70 Å(reported)

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

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

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<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

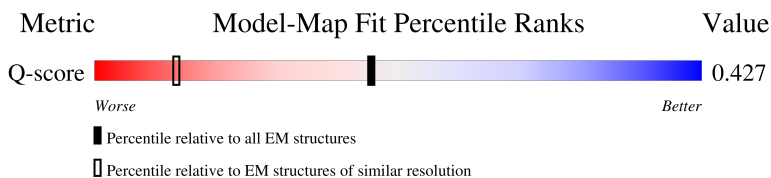
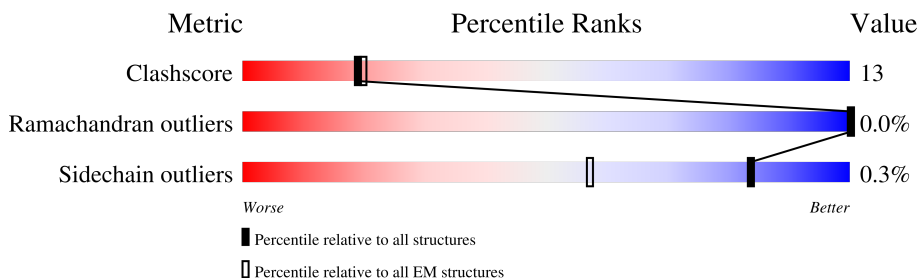
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.70 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	10327 (2.20 - 3.20)

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	1101	70% 23% 6%
1	B	1101	67% 27% 6%
1	C	1101	15% 65% 28% 6%
1	D	1101	68% 26% 6%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 32519 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.

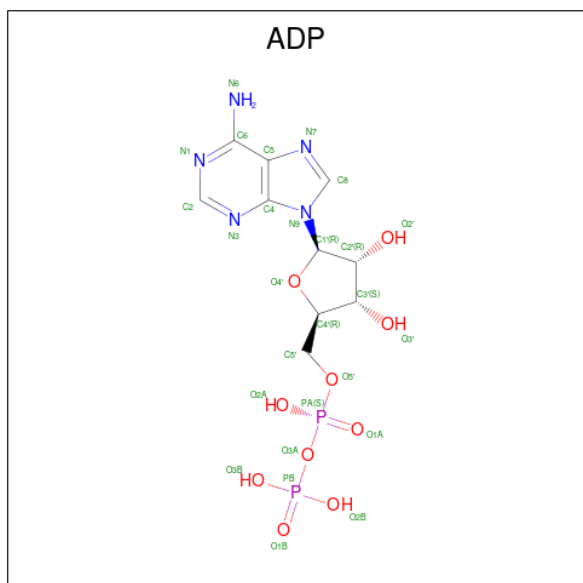
- Molecule 1 is a protein called ATP-citrate synthase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1032	Total	C	N	O	S	2	0
			7989	5114	1354	1474	47		
1	B	1032	Total	C	N	O	S	2	0
			7989	5114	1354	1474	47		
1	C	1032	Total	C	N	O	S	2	0
			7989	5114	1354	1474	47		
1	D	1032	Total	C	N	O	S	2	0
			7989	5114	1354	1474	47		

There are 4 discrepancies between the modelled and reference sequences:

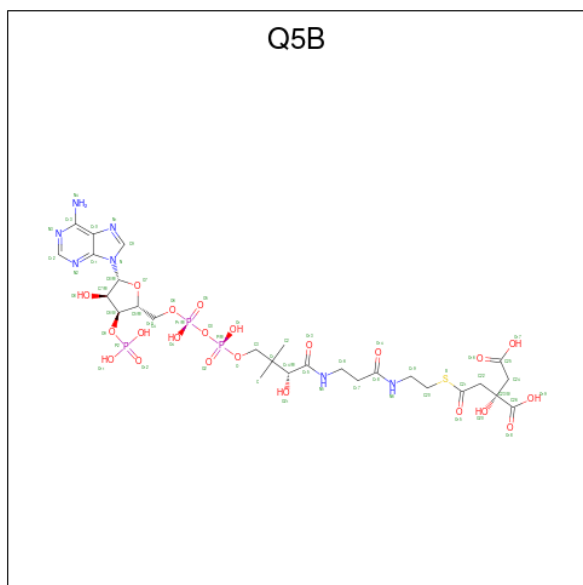
Chain	Residue	Modelled	Actual	Comment	Reference
A	1026	ALA	ASP	engineered mutation	UNP P53396
B	1026	ALA	ASP	engineered mutation	UNP P53396
C	1026	ALA	ASP	engineered mutation	UNP P53396
D	1026	ALA	ASP	engineered mutation	UNP P53396

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



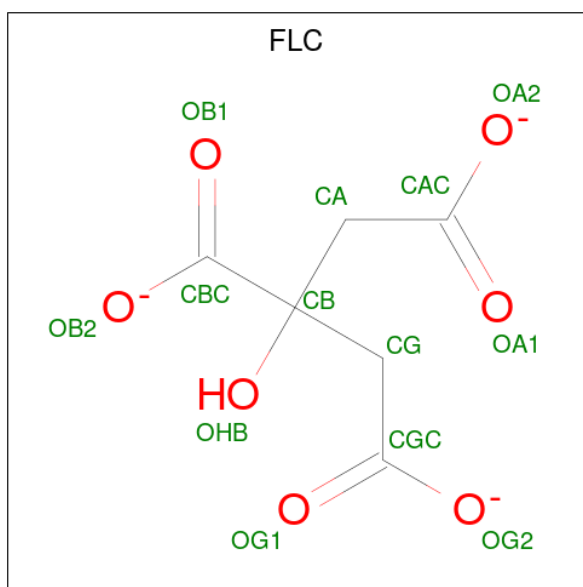
Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	B	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	D	1	Total	C	N	O	P	0
			27	10	5	10	2	

- Molecule 3 is (3S)-citryl-Coenzyme A (CCD ID: Q5B) (formula: $C_{27}H_{42}N_7O_{22}P_3S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf	
3	A	1	Total 60	C 27	N 7	O 22	P 3	S 1	0
3	B	1	Total 60	C 27	N 7	O 22	P 3	S 1	0
3	D	1	Total 60	C 27	N 7	O 22	P 3	S 1	0

- Molecule 4 is CITRATE ANION (CCD ID: FLC) (formula: $C_6H_5O_7$).

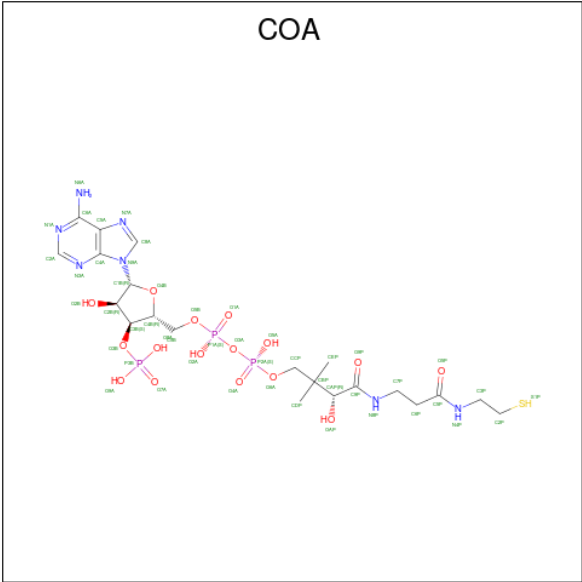


Mol	Chain	Residues	Atoms			AltConf
4	A	1	Total	C	O	0
			13	6	7	
4	B	1	Total	C	O	0
			13	6	7	
4	C	1	Total	C	O	0
			13	6	7	
4	D	1	Total	C	O	0
			13	6	7	

- Molecule 5 is UNKNOWN LIGAND (CCD ID: UNL) (formula:).

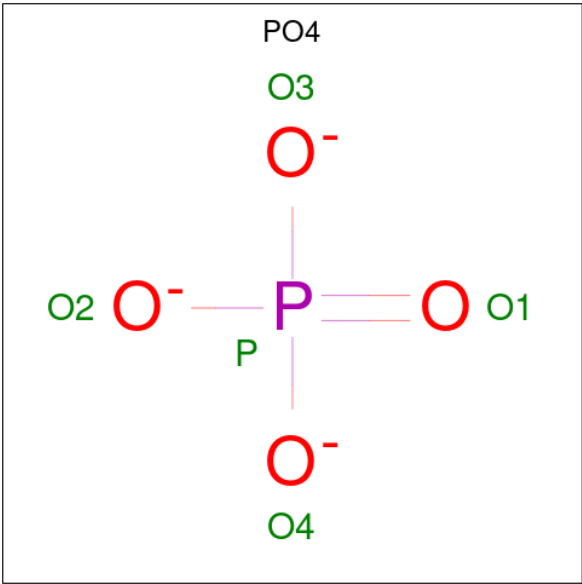
Mol	Chain	Residues	Atoms		AltConf
5	A	1	Total	C	0
			1	1	
5	C	1	Total	C	0
			1	1	

- Molecule 6 is COENZYME A (CCD ID: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						AltConf
6	A	1	Total	C	N	O	P	S	0
			48	21	7	16	3	1	
6	B	1	Total	C	N	O	P	S	0
			48	21	7	16	3	1	
6	C	1	Total	C	N	O	P	S	0
			48	21	7	16	3	1	
6	D	1	Total	C	N	O	P	S	0
			48	21	7	16	3	1	

- Molecule 7 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			AltConf
7	A	1	Total	O	P	0
			5	4	1	
7	B	1	Total	O	P	0
			5	4	1	
7	D	1	Total	O	P	0
			5	4	1	

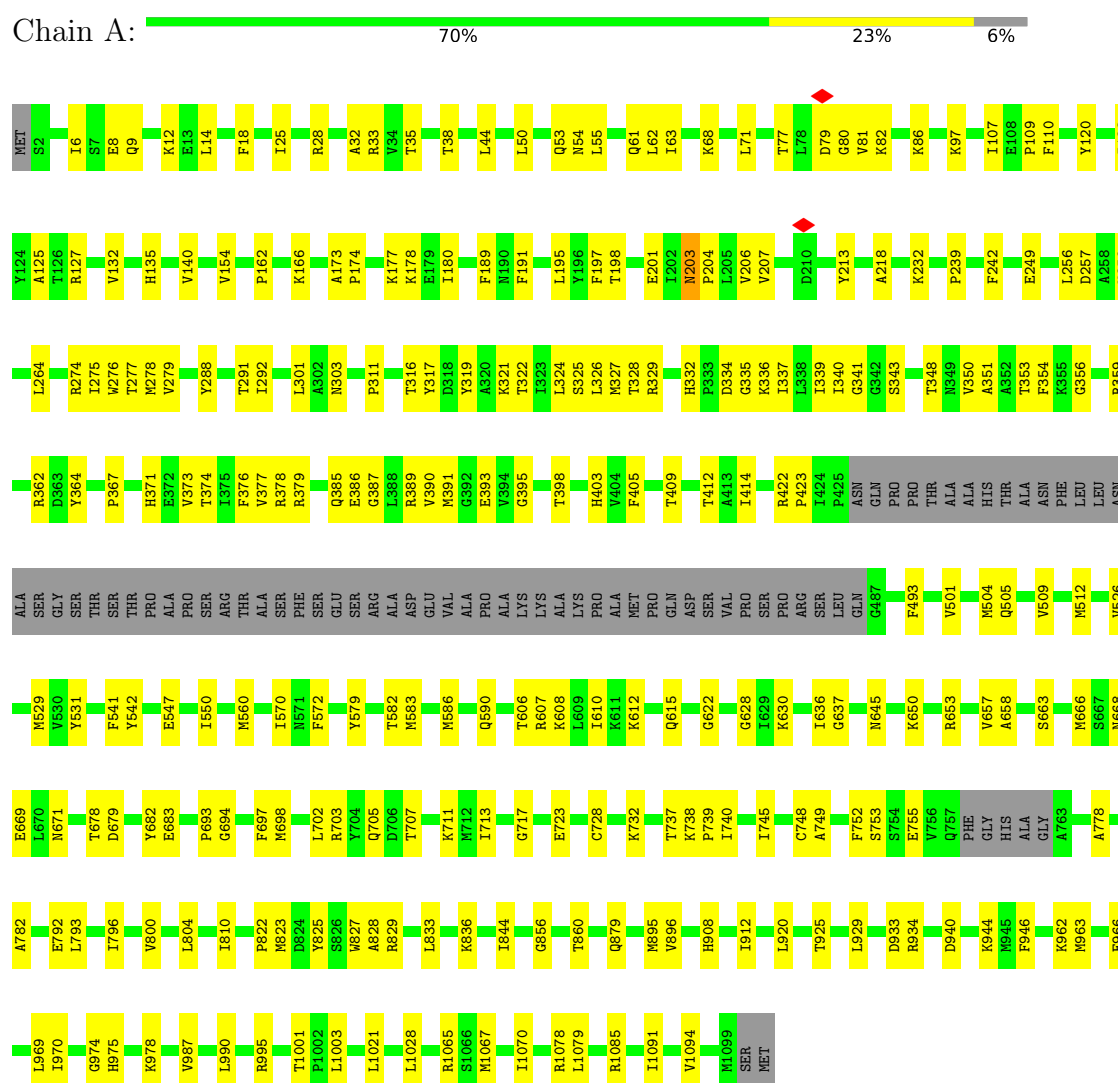
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		AltConf
8	A	10	Total	O	0
			10	10	
8	B	9	Total	O	0
			9	9	
8	C	13	Total	O	0
			13	13	
8	D	9	Total	O	0
			9	9	

3 Residue-property plots

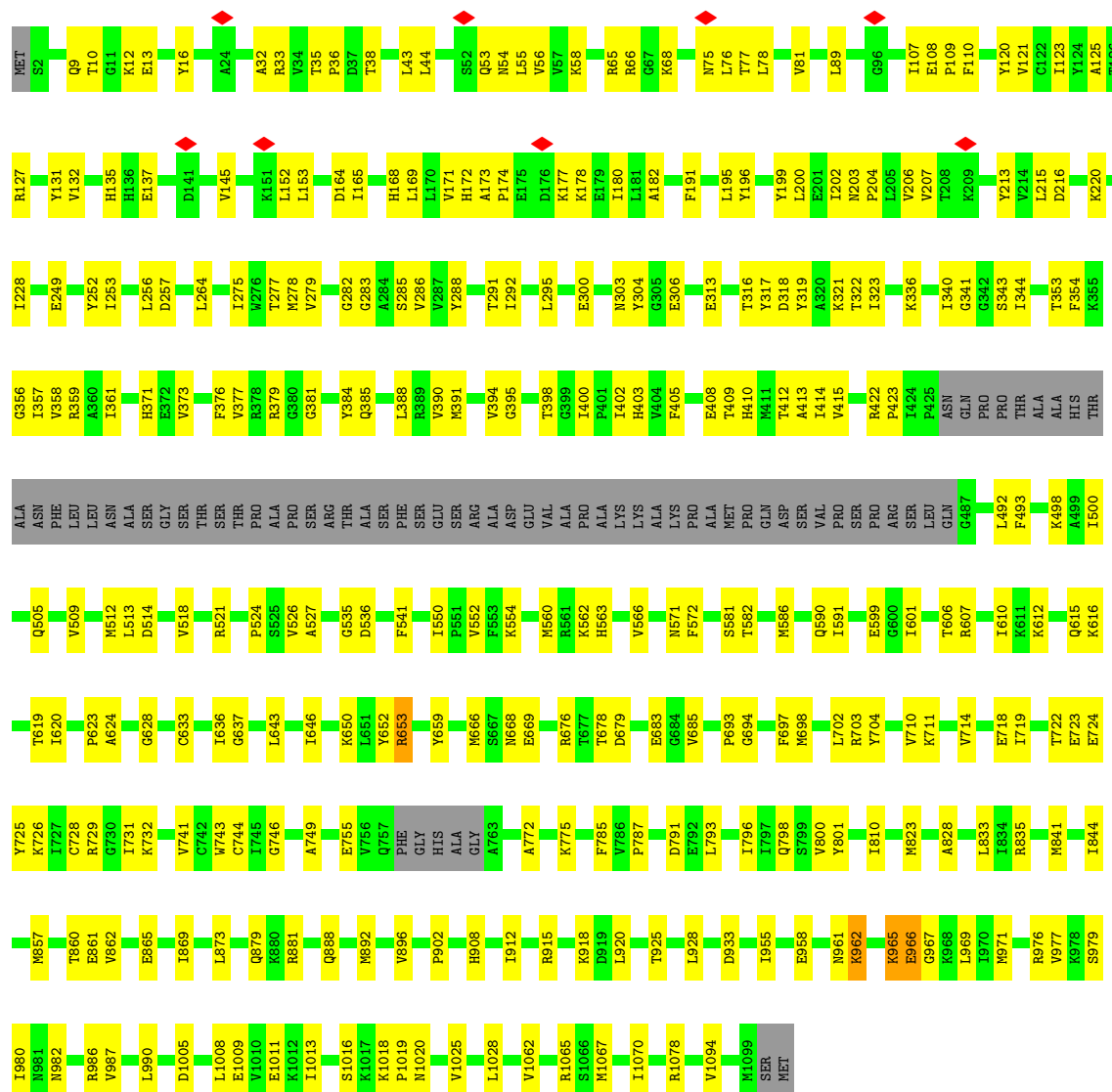
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

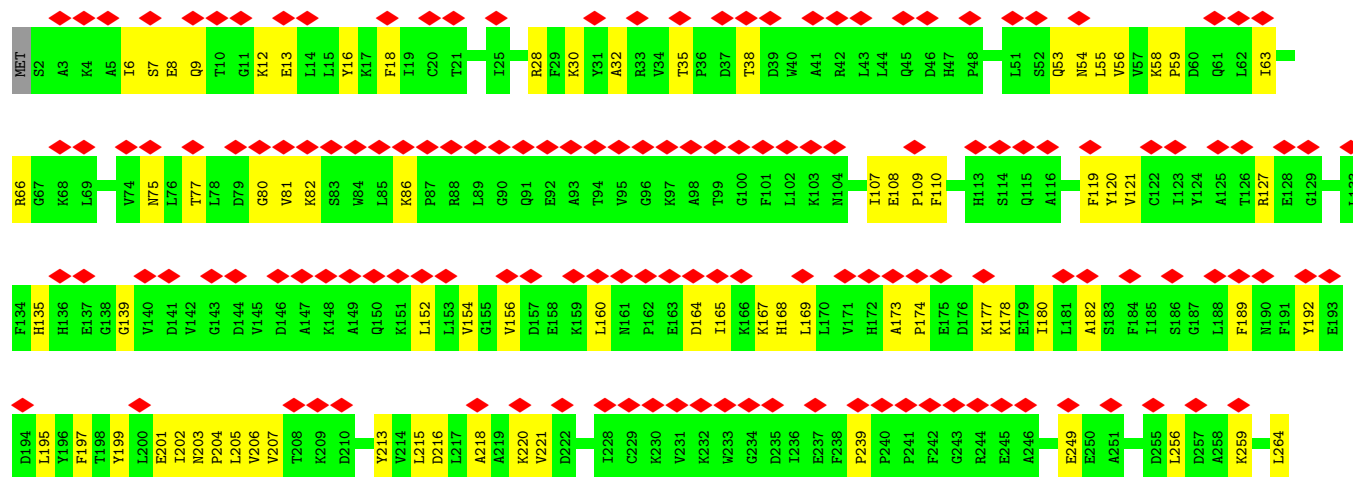


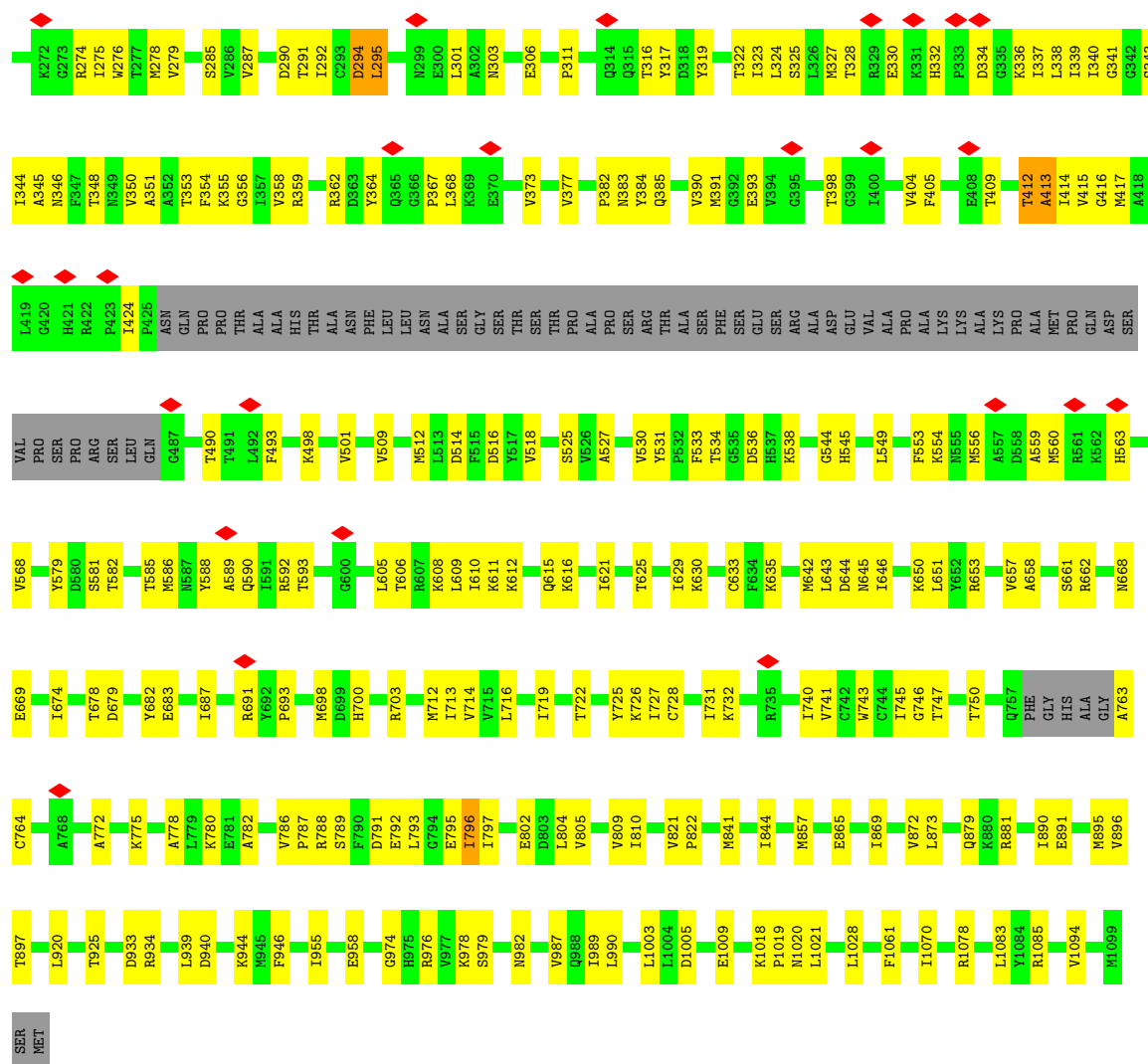
• Molecule 1: ATP-citrate synthase





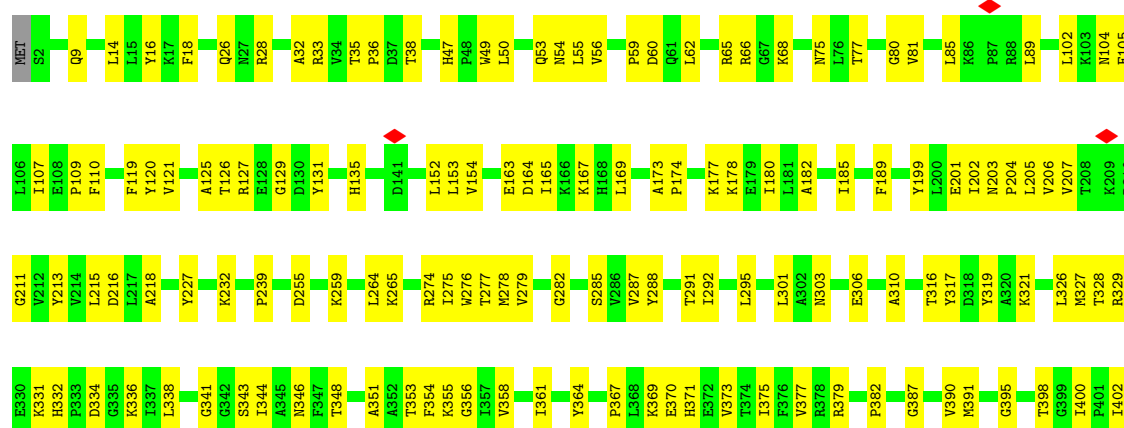
• Molecule 1: ATP-citrate synthase





Molecule 1: ATP-citrate synthase

Chain D: 68% 26% 6%



K1018	P1019	N1020	L1021	N1024	L1028	G1064	R1065	S1066	M1067	R1078	L1079	R1085	H1086	V1094	L1095	M1099	SER	MET																												
L831	R835	I844	M857	I869	L873	Q879	R880	R881	E891	N895	V896	T897	A898	H908	I912	R915	T925	L929	D933	R934	K964	L969	I970	N971	Y977	K978	S979	I980	N981	Y987	Q988	I989	L990	K991	R995	Q996	H997	K1017								
E718	I719	G720	E724	I727	C728	I731	K732	E733	G734	R735	L736	T737	K738	P739	I740	V741	C742	C748	A749	T750	M751	F752	E755	V756	Q757	PHE	GLY	HIS	ALA	GLY	A763	C764	Q767	A768	S769	E770	K775	L779	V784	D791	Y801	L804	V809	I810	M823	
P602	E603	A604	L605	T606	R607	K608	L609	A613	D614	Q615	V618	T619	I620	G628	C633	I636	G637	M642	R653	V657	R662	S663	Q664	M666	S667	N668	E669	R676	T677	T678	D679	E683	P693	F697	M698	R703	T707	K711	M712	I713	V714					
PRO	ALA	MET	PRO	GLN	ASP	SER	VAL	PRO	SER	PRO	ARG	SER	LEU	G487	I500	Q505	V509	M512	L513	R521	V526	F541	I548	L549	I550	M556	M560	V566	THR	I570	N571	F572	R576	Y579	T582	M583	F584	T585	M586	N587	Y588	I594				
H410	M411	T412	A413	I414	V415	G416	M417	L419	I424	F425	ASN	GLN	PRO	PRO	THR	ALA	ALA	HIS	THR	ALA	ASN	PHE	LEU	LEU	ASN	ALA	SER	GLY	SER	THR	THR	PRO	ALA	PRO	SER	ARG	ALA	ASP	GLU	VAL	ALA	PRO	ALA	LYS	ALA	LYS

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	100062	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 ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	3.039	Depositor
Minimum map value	-1.208	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.151	Depositor
Recommended contour level	0.15	Depositor
Map size (Å)	182.59999, 182.59999, 182.59999	wwPDB
Map dimensions	220, 220, 220	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	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: COA, Q5B, FLC, PO4, UNL, ADP

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.16	0/8167	0.34	0/11054
1	B	0.21	0/8167	0.37	0/11054
1	C	0.15	0/8167	0.35	0/11054
1	D	0.17	0/8167	0.34	0/11054
All	All	0.18	0/32668	0.35	0/44216

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7989	0	8040	183	0
1	B	7989	0	8040	217	0
1	C	7989	0	8040	236	0
1	D	7989	0	8040	220	0
2	A	27	0	12	1	0
2	B	27	0	12	3	0
2	D	27	0	12	2	0
3	A	60	0	0	0	0
3	B	60	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	60	0	0	2	0
4	A	13	0	5	1	0
4	B	13	0	5	1	0
4	C	13	0	5	1	0
4	D	13	0	5	1	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	A	48	0	32	3	0
6	B	48	0	32	3	0
6	C	48	0	32	9	0
6	D	48	0	32	8	0
7	A	5	0	0	0	0
7	B	5	0	0	0	0
7	D	5	0	0	0	0
8	A	10	0	0	1	0
8	B	9	0	0	0	0
8	C	13	0	0	1	0
8	D	9	0	0	0	0
All	All	32519	0	32344	829	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 13.

The worst 5 of 829 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:279:VAL:HG12	1:D:341:GLY:H	1.35	0.90
1:B:500:ILE:HG12	1:B:566:VAL:HG11	1.56	0.88
1:B:512:MET:HE2	1:B:637:GLY:H	1.44	0.83
1:B:925:THR:HG23	1:D:925:THR:HG23	1.62	0.82
1:A:350:VAL:O	1:A:354:PHE:HB2	1.79	0.82

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	1028/1101 (93%)	1000 (97%)	28 (3%)	0	100	100
1	B	1028/1101 (93%)	991 (96%)	37 (4%)	0	100	100
1	C	1028/1101 (93%)	990 (96%)	37 (4%)	1 (0%)	48	73
1	D	1028/1101 (93%)	1005 (98%)	23 (2%)	0	100	100
All	All	4112/4404 (93%)	3986 (97%)	125 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	413	ALA

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	852/908 (94%)	851 (100%)	1 (0%)	88	96
1	B	852/908 (94%)	848 (100%)	4 (0%)	81	92
1	C	852/908 (94%)	848 (100%)	4 (0%)	81	92
1	D	852/908 (94%)	851 (100%)	1 (0%)	88	96
All	All	3408/3632 (94%)	3398 (100%)	10 (0%)	84	94

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	412	THR
1	C	796	ILE
1	D	615	GLN
1	B	965	LYS
1	B	966	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22

such sidechains are listed below:

Mol	Chain	Res	Type
1	C	671	ASN
1	C	879	GLN
1	C	850	GLN
1	D	54	ASN
1	B	672	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 2 are unknown - leaving 17 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	COA	D	2101	-	47,50,50	1.13	5 (10%)	69,75,75	1.56	10 (14%)
2	ADP	B	1201	-	28,29,29	1.41	4 (14%)	43,45,45	1.86	10 (23%)
2	ADP	D	2102	-	28,29,29	1.42	4 (14%)	43,45,45	1.84	9 (20%)
3	Q5B	A	1202	-	60,62,62	2.87	22 (36%)	87,93,93	1.73	19 (21%)
4	FLC	A	1203	-	12,12,12	1.14	0	17,17,17	1.35	2 (11%)
3	Q5B	D	2103	-	60,62,62	2.87	22 (36%)	87,93,93	1.82	20 (22%)
3	Q5B	B	1202	-	60,62,62	2.93	20 (33%)	87,93,93	1.88	23 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FLC	C	2102	-	12,12,12	1.08	0	17,17,17	1.33	1 (5%)
6	COA	C	2101	-	47,50,50	1.14	5 (10%)	69,75,75	1.56	10 (14%)
4	FLC	B	1203	-	12,12,12	1.09	0	17,17,17	1.42	1 (5%)
7	PO4	D	2105	-	4,4,4	1.01	0	6,6,6	0.45	0
2	ADP	A	1201	-	28,29,29	1.42	4 (14%)	43,45,45	1.84	9 (20%)
7	PO4	B	1205	-	4,4,4	0.98	0	6,6,6	0.47	0
6	COA	B	1204	-	47,50,50	1.12	5 (10%)	69,75,75	1.57	11 (15%)
6	COA	A	1205	-	47,50,50	1.16	4 (8%)	69,75,75	1.50	10 (14%)
7	PO4	A	1206	-	4,4,4	1.01	0	6,6,6	0.50	0
4	FLC	D	2104	-	12,12,12	1.12	0	17,17,17	1.36	2 (11%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '–' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	COA	D	2101	-	-	17/48/64/64	0/3/3/3
2	ADP	B	1201	-	-	4/16/32/32	0/3/3/3
2	ADP	D	2102	-	-	2/16/32/32	0/3/3/3
3	Q5B	A	1202	-	-	22/66/83/83	0/3/3/3
4	FLC	A	1203	-	-	8/16/16/16	-
3	Q5B	D	2103	-	-	22/66/83/83	0/3/3/3
3	Q5B	B	1202	-	-	32/66/83/83	0/3/3/3
4	FLC	C	2102	-	-	6/16/16/16	-
6	COA	C	2101	-	-	15/48/64/64	0/3/3/3
4	FLC	B	1203	-	-	3/16/16/16	-
2	ADP	A	1201	-	-	8/16/32/32	0/3/3/3
6	COA	B	1204	-	-	13/48/64/64	0/3/3/3
6	COA	A	1205	-	-	12/48/64/64	0/3/3/3
4	FLC	D	2104	-	-	7/16/16/16	-

The worst 5 of 95 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1202	Q5B	P1-O3	8.65	1.68	1.59
3	D	2103	Q5B	P1-O3	8.52	1.68	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1202	Q5B	P1-O3	8.26	1.68	1.59
3	D	2103	Q5B	C6-C5	-8.14	1.31	1.52
3	A	1202	Q5B	C6-C5	-8.13	1.31	1.52

The worst 5 of 137 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1201	ADP	C5-C4-N3	-5.95	118.52	126.72
2	D	2102	ADP	C5-C4-N3	-5.95	118.52	126.72
3	B	1202	Q5B	C22-C21-S	5.94	121.05	113.56
6	B	1204	COA	C5A-C4A-N3A	-5.85	118.66	126.72
6	D	2101	COA	C5A-C4A-N3A	-5.81	118.71	126.72

There are no chirality outliers.

5 of 171 torsion outliers are listed below:

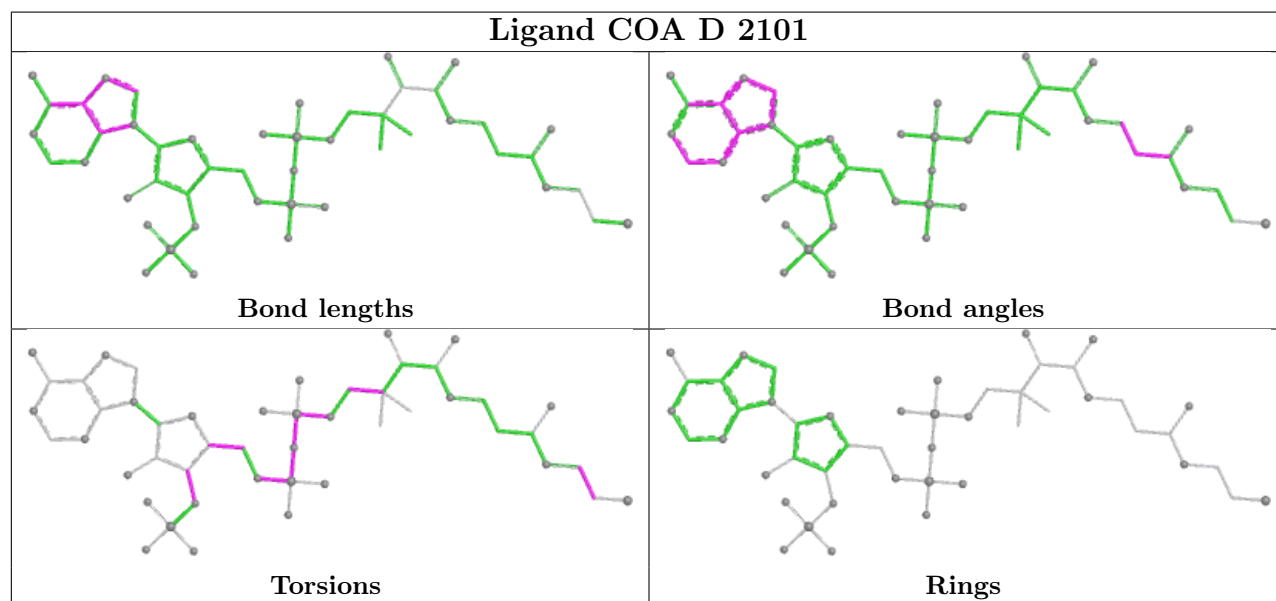
Mol	Chain	Res	Type	Atoms
2	A	1201	ADP	C5'-O5'-PA-O1A
2	A	1201	ADP	C5'-O5'-PA-O2A
2	A	1201	ADP	C5'-O5'-PA-O3A
2	B	1201	ADP	C5'-O5'-PA-O1A
2	B	1201	ADP	C5'-O5'-PA-O2A

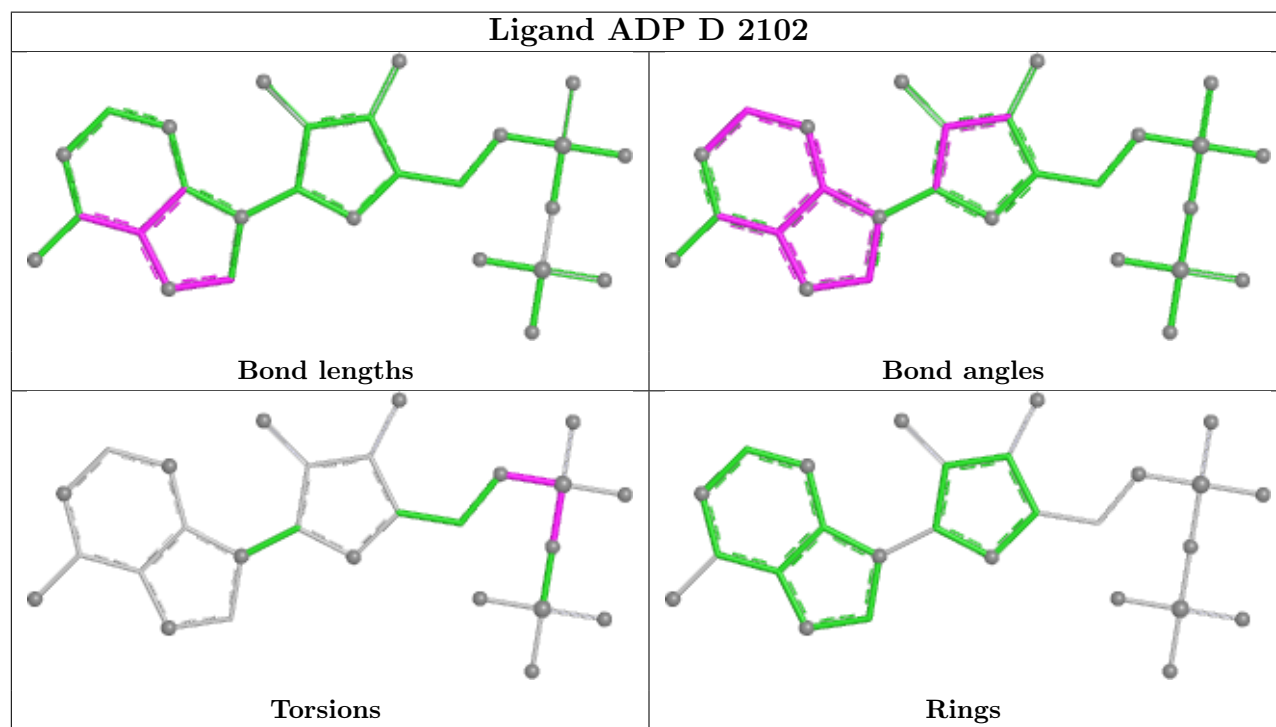
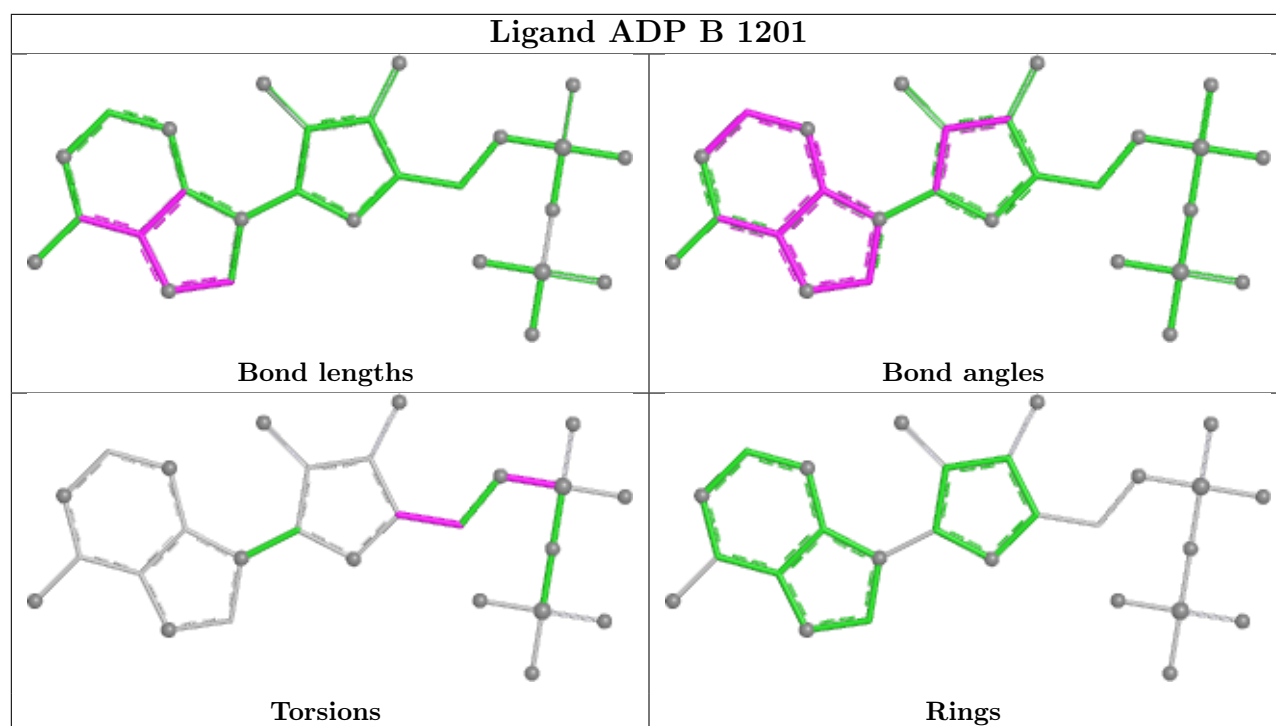
There are no ring outliers.

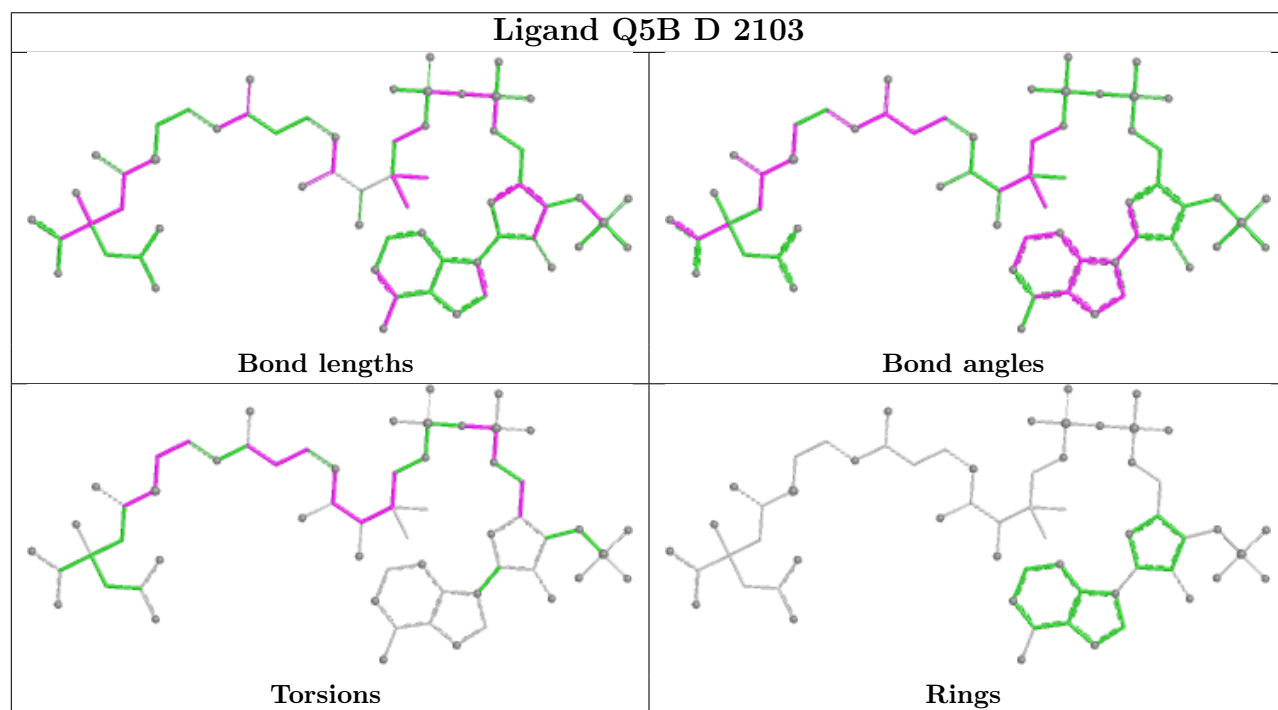
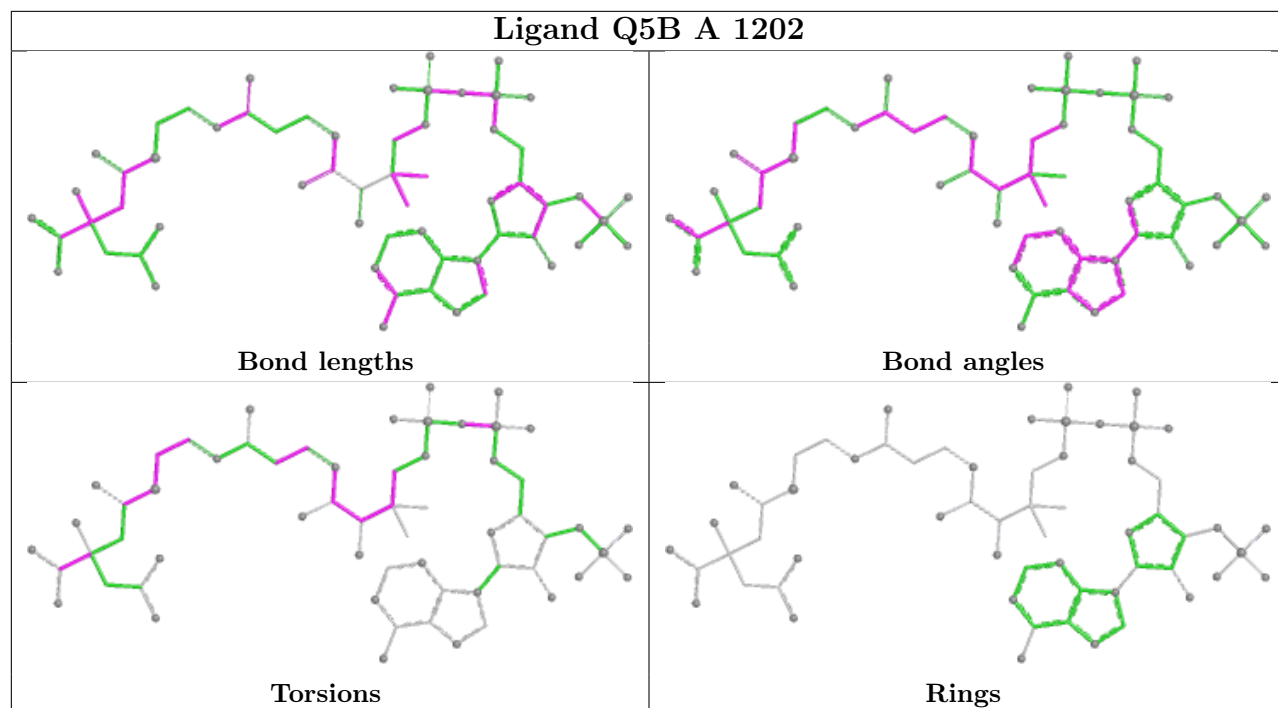
13 monomers are involved in 37 short contacts:

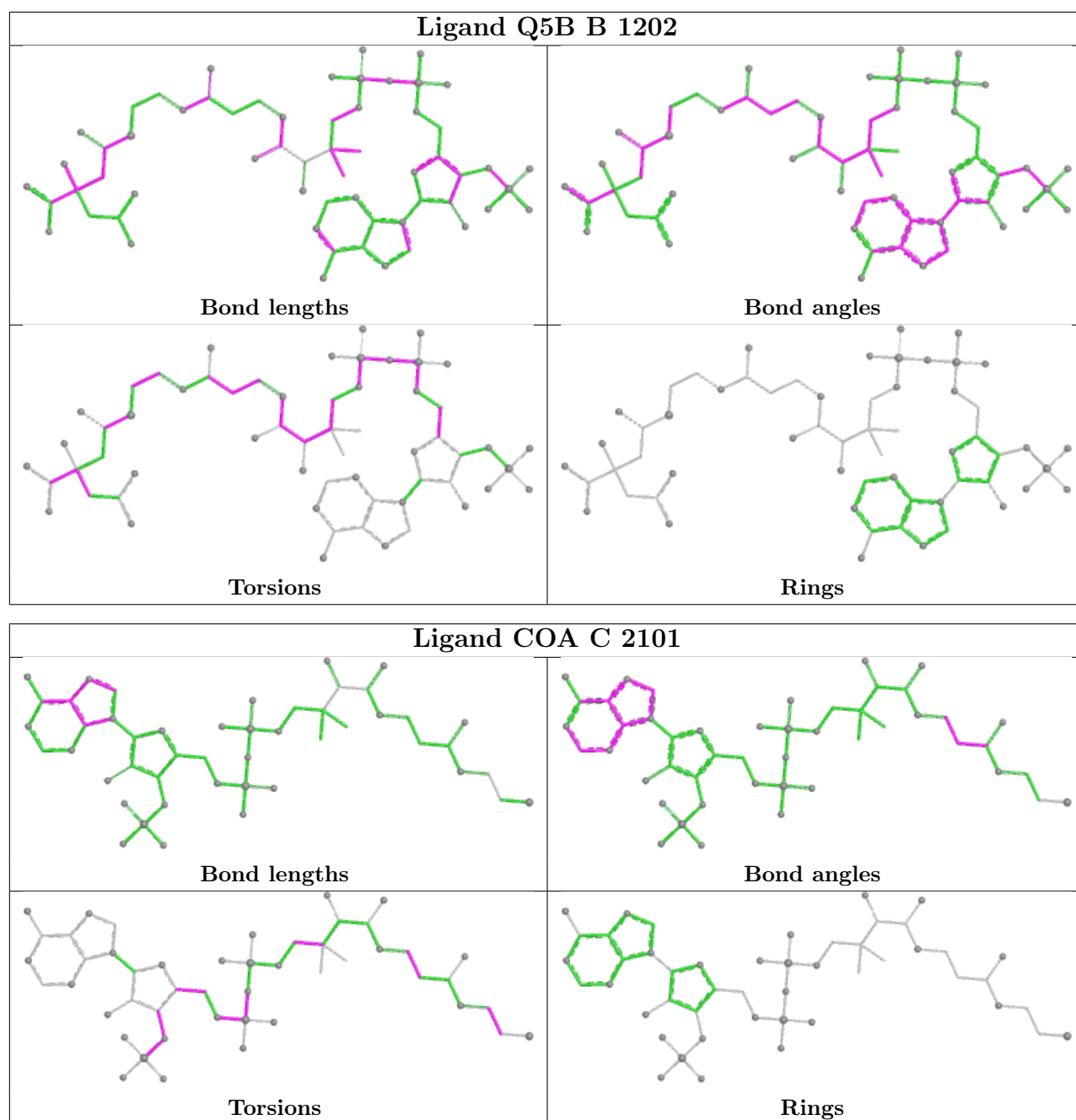
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	2101	COA	8	0
2	B	1201	ADP	3	0
2	D	2102	ADP	2	0
4	A	1203	FLC	1	0
3	D	2103	Q5B	2	0
3	B	1202	Q5B	3	0
4	C	2102	FLC	1	0
6	C	2101	COA	9	0
4	B	1203	FLC	1	0
2	A	1201	ADP	1	0
6	B	1204	COA	3	0
6	A	1205	COA	3	0
4	D	2104	FLC	1	0

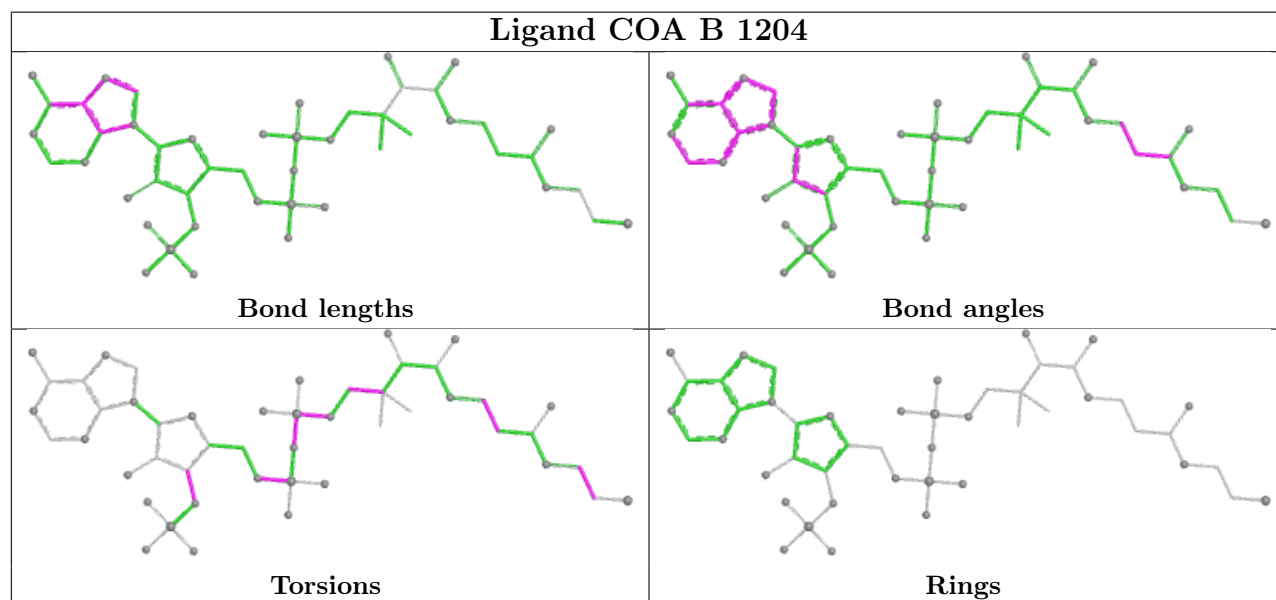
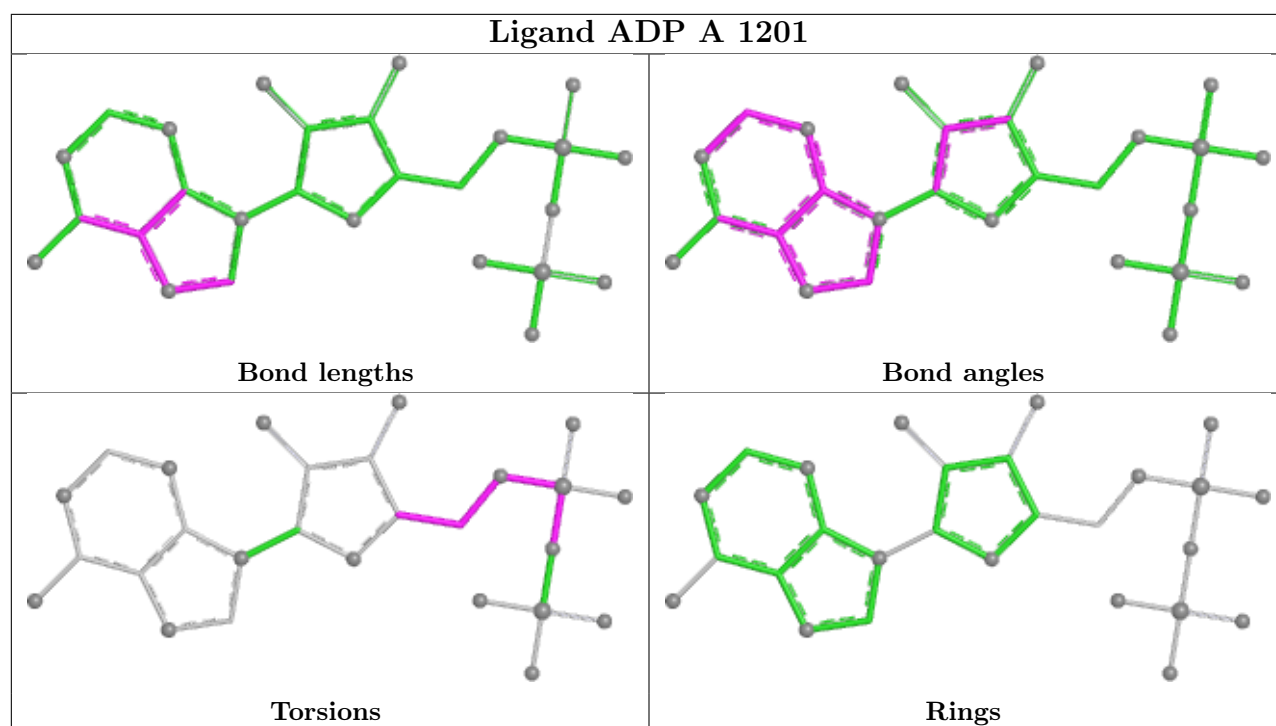
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

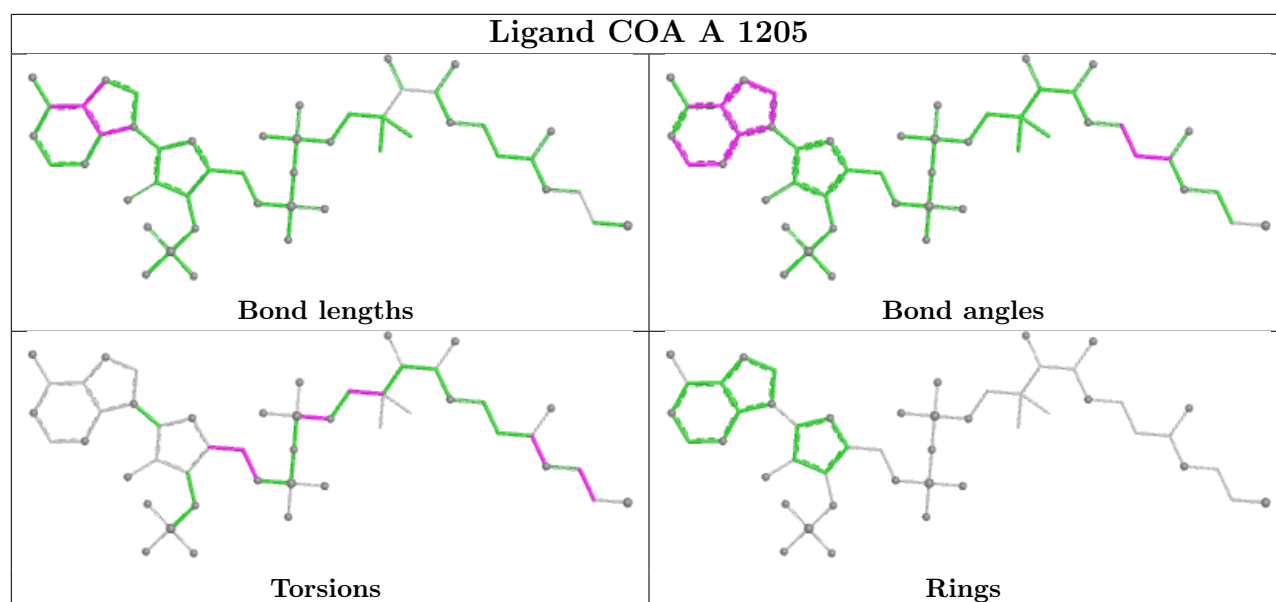












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

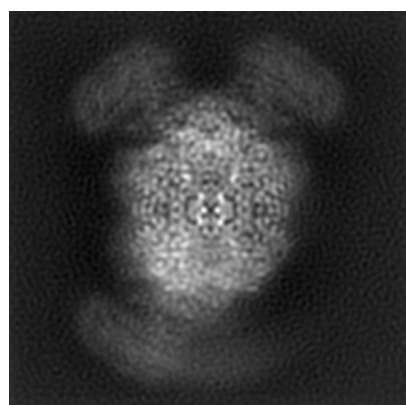
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24577. These allow visual inspection of the internal detail of the map and identification of artifacts.

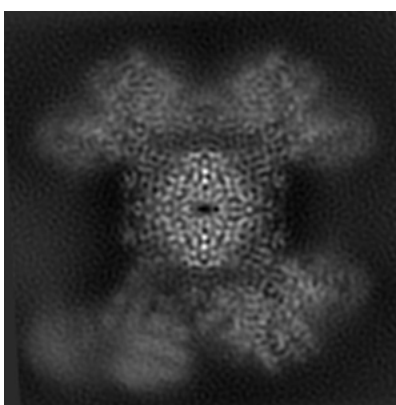
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

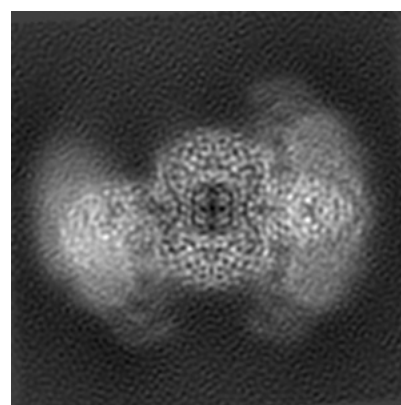
6.1.1 Primary map



X



Y

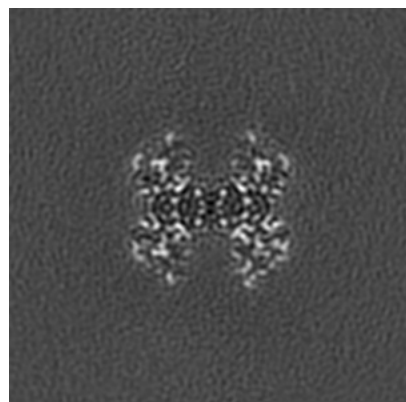


Z

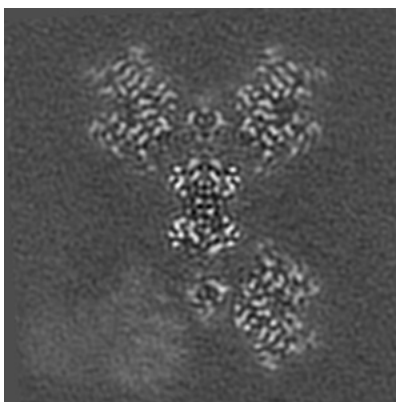
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

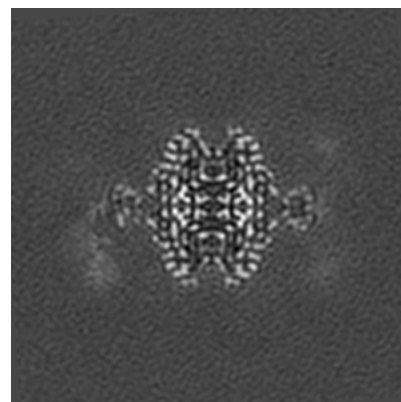
6.2.1 Primary map



X Index: 110



Y Index: 110

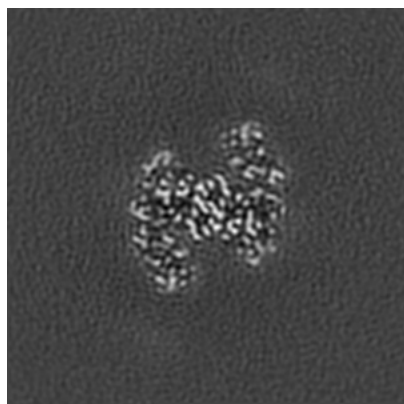


Z Index: 110

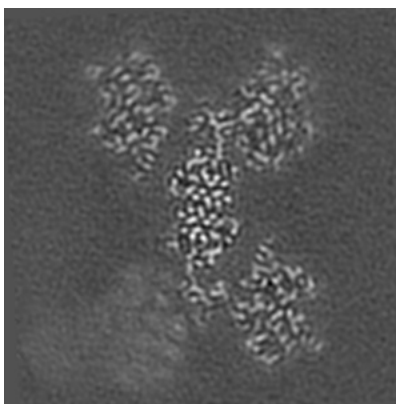
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

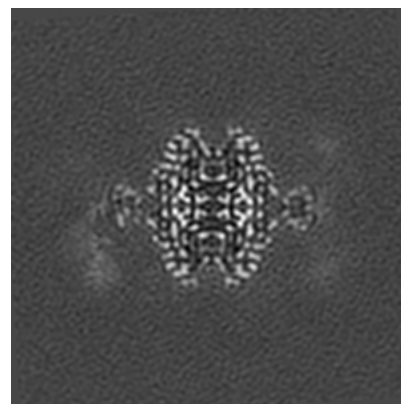
6.3.1 Primary map



X Index: 123



Y Index: 105

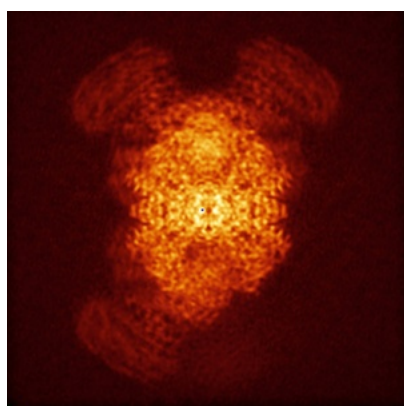


Z Index: 110

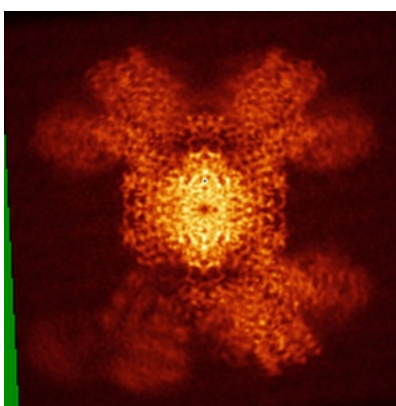
The images above show the largest variance slices of the map in three orthogonal directions.

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

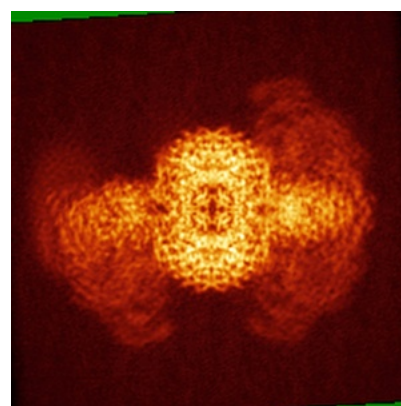
6.4.1 Primary map



X



Y

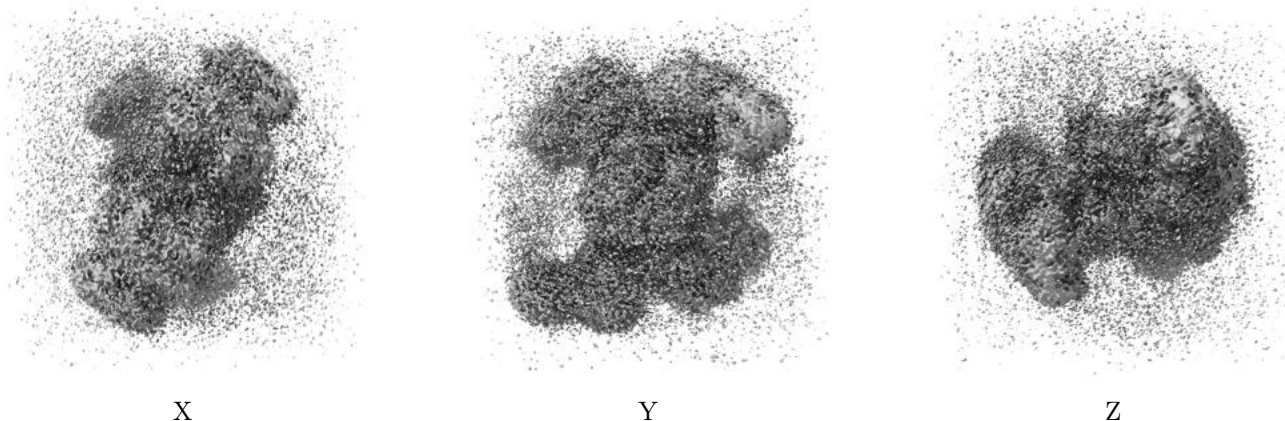


Z

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

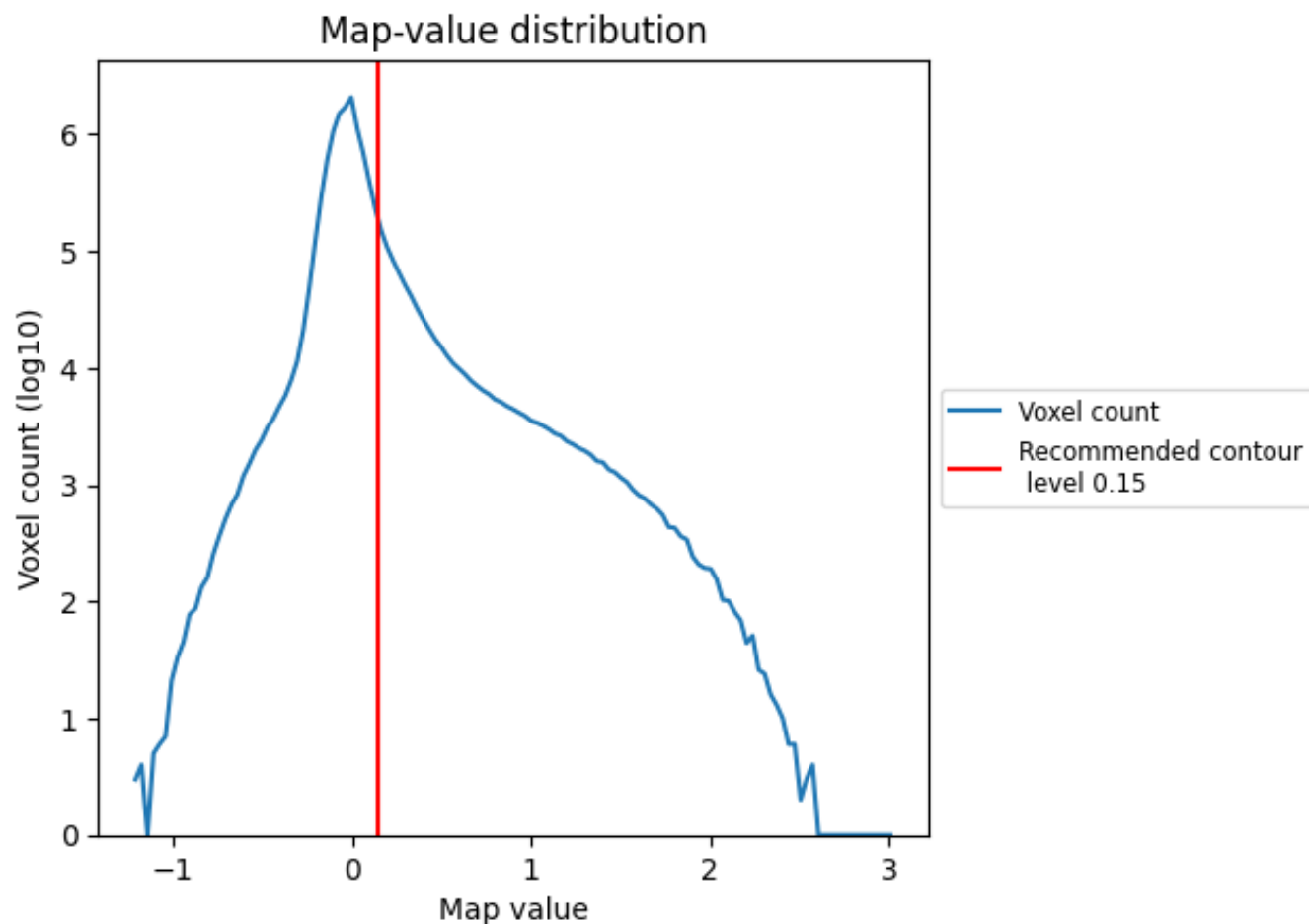
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

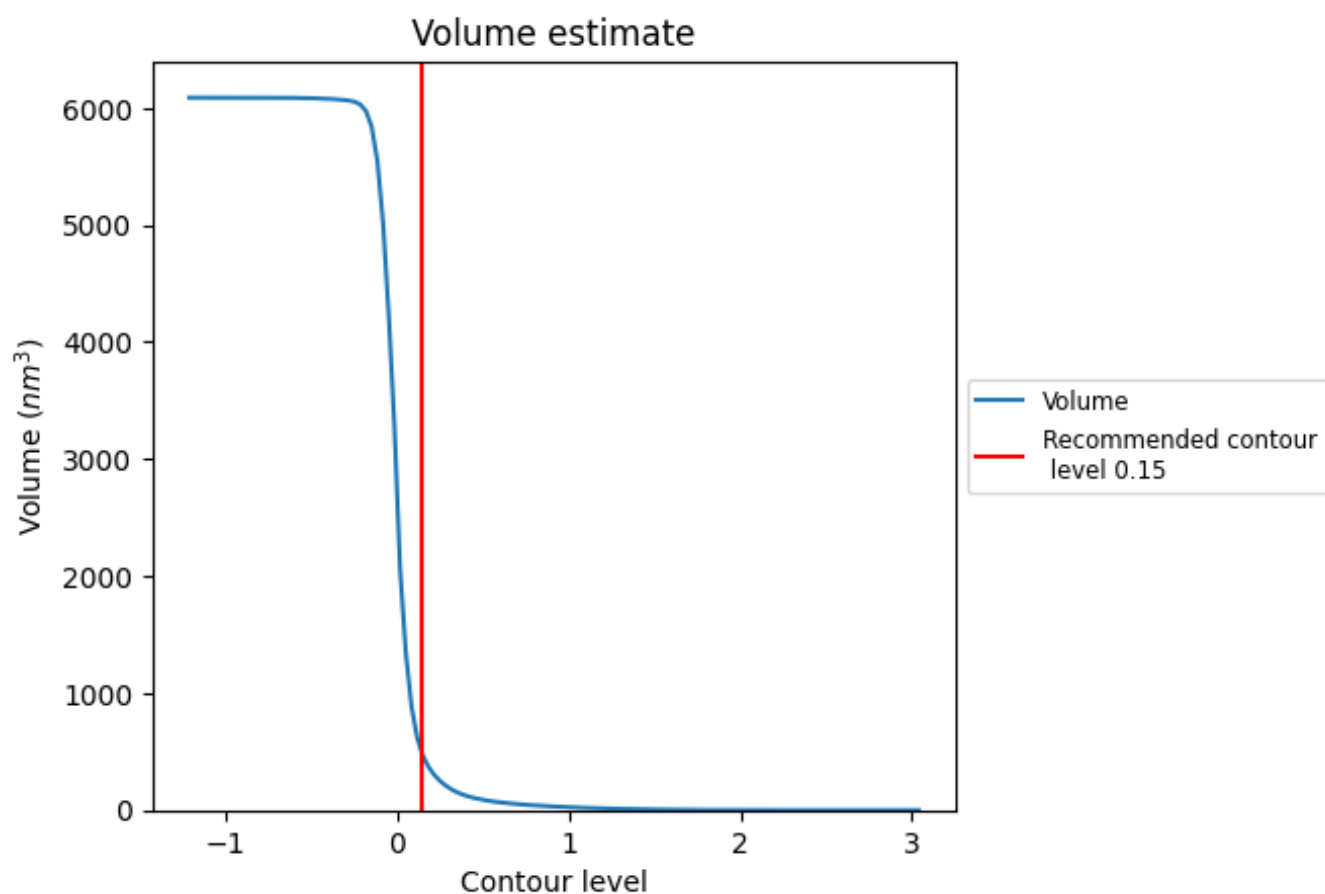
This section contains the results of statistical analysis of the map.

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.

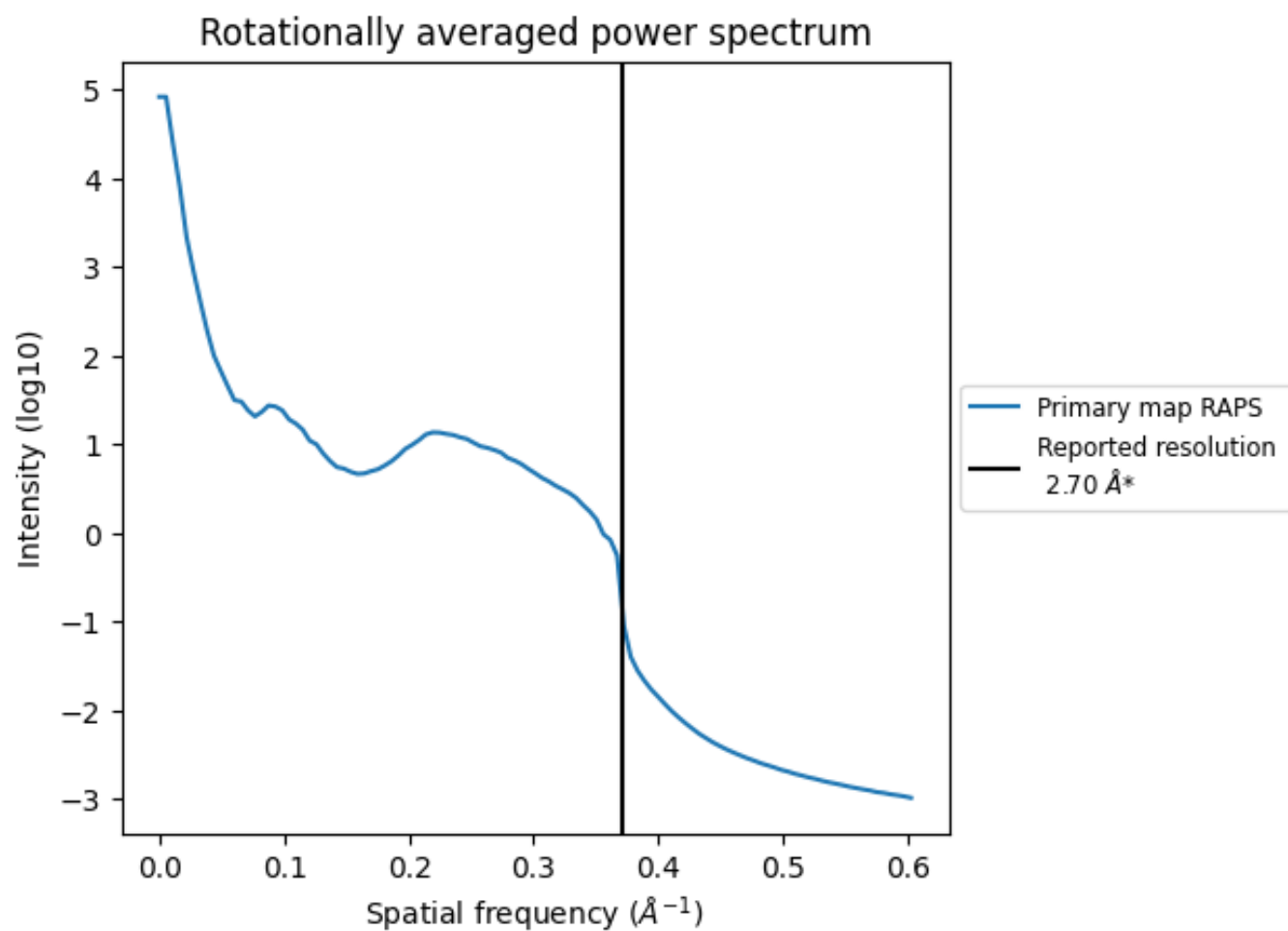
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 474 nm^3 ; this corresponds to an approximate mass of 428 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 0.370 $\text{\AA}^{-1}$

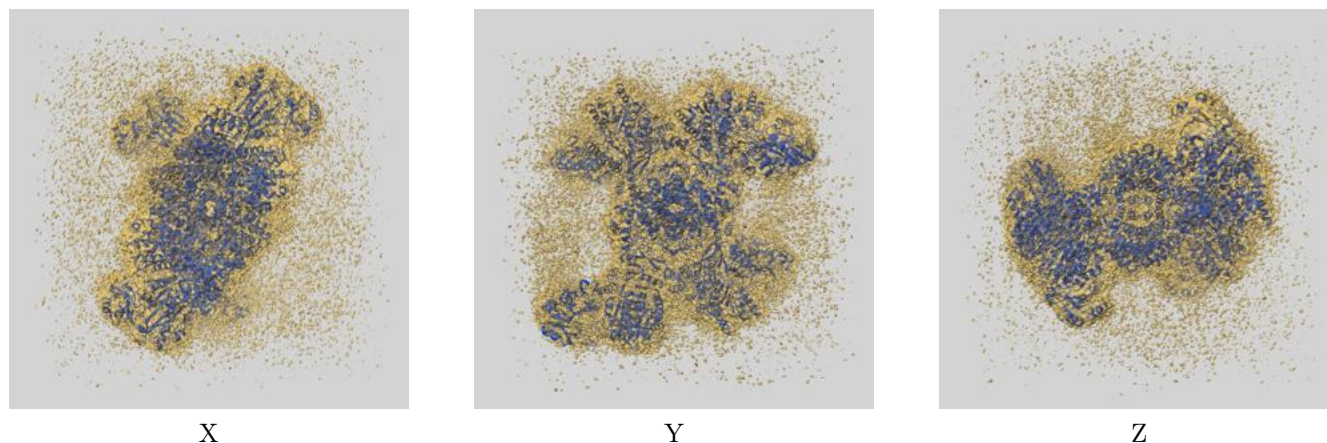
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

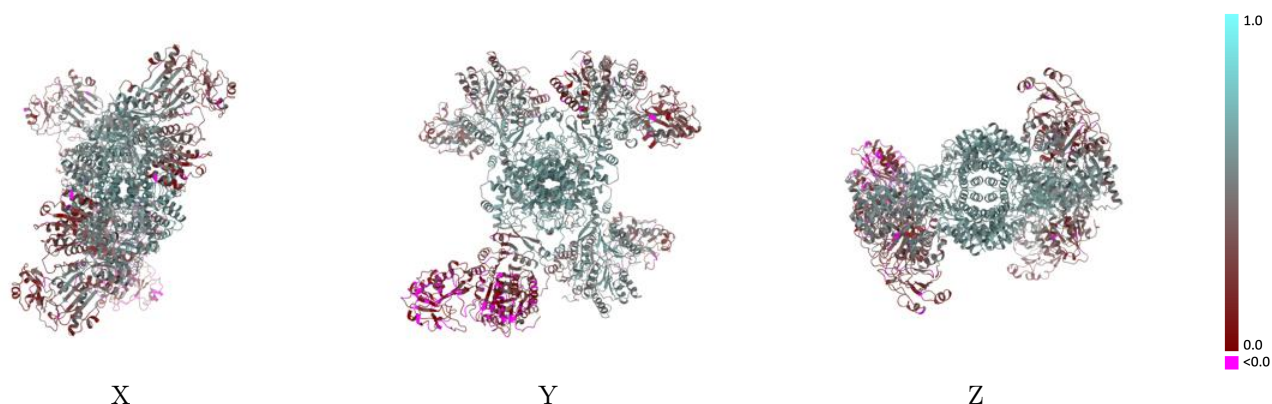
This section contains information regarding the fit between EMDB map EMD-24577 and PDB model 7RMP. Per-residue inclusion information can be found in [section 3](#) on [page 8](#).

9.1 Map-model overlay [i](#)



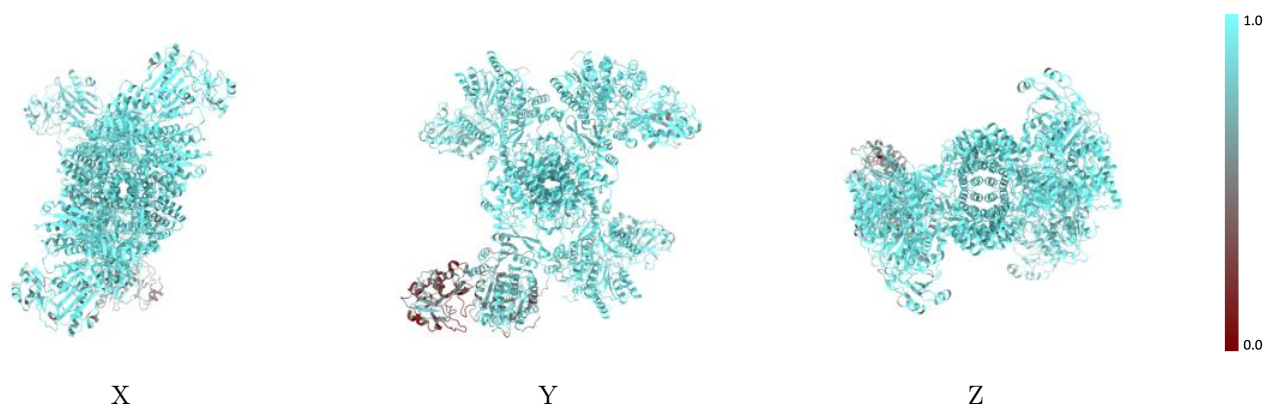
The images above show the 3D surface view of the map at the recommended contour level 0.15 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



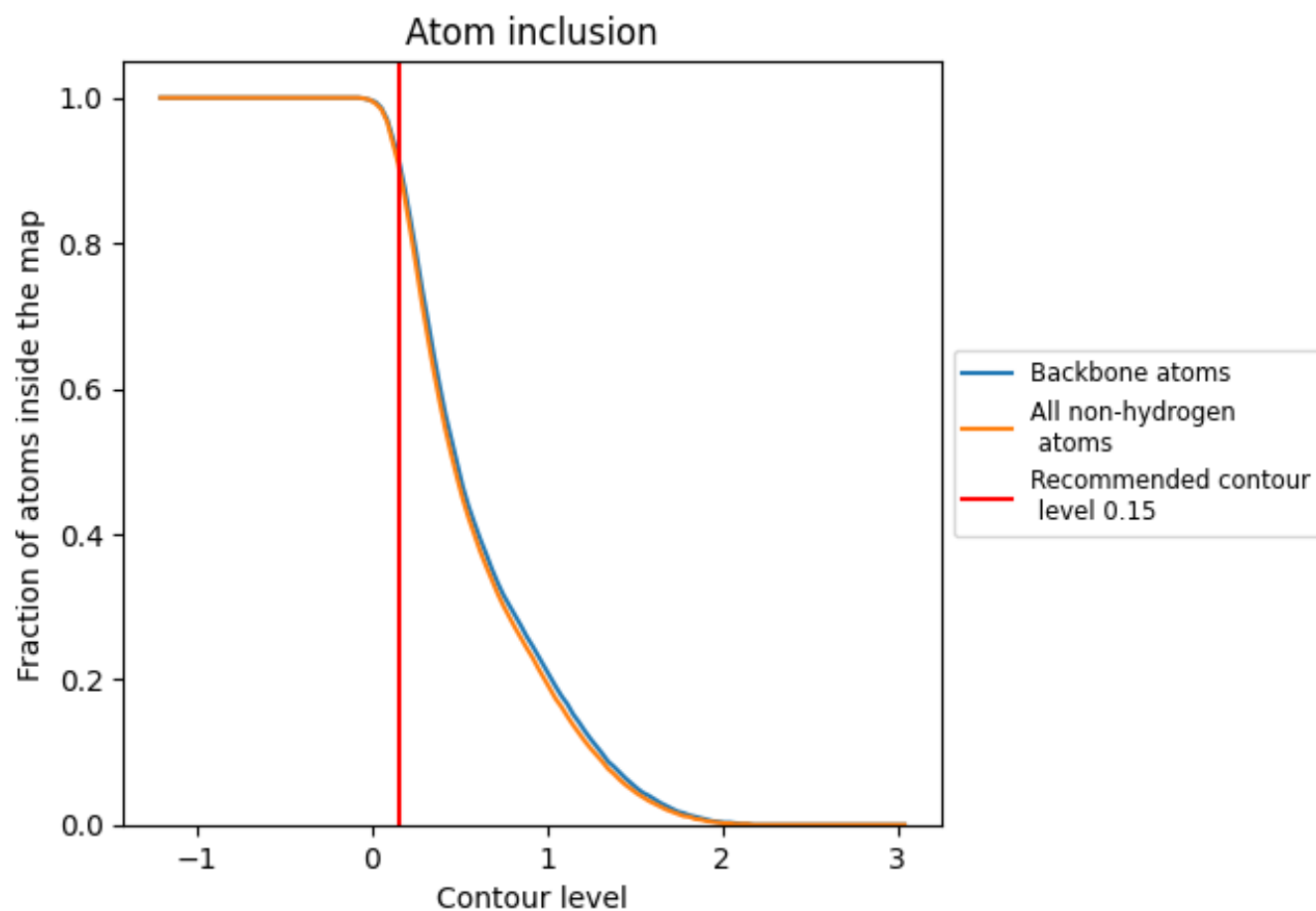
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.15).

9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 91% 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 (0.15) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9080	0.4270
A	0.9610	0.4810
B	0.9530	0.4620
C	0.7580	0.2750
D	0.9620	0.4860

