



Full wwPDB EM Validation Report ⓘ

Nov 20, 2024 – 12:17 PM JST

PDB ID : 9IZ1
EMDB ID : EMD-61008
Title : dmCTPS tetramer with dATP dUTP dGTP and DON
Authors : Guo, C.J.; Liu, J.L.
Deposited on : 2024-07-31
Resolution : 2.73 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.dev113
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

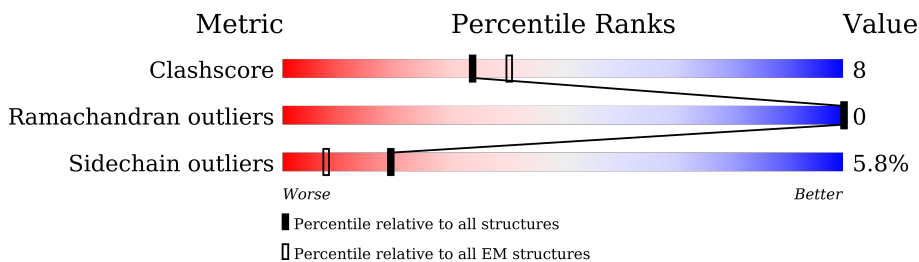
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.73 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	556	
1	B	556	
1	C	556	
1	D	556	

2 Entry composition [i](#)

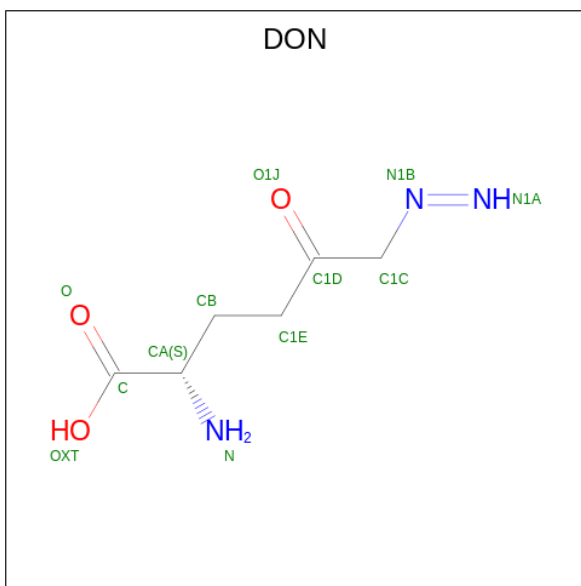
There are 6 unique types of molecules in this entry. The entry contains 17948 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 CTP synthase.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	556	Total 4384	C 2779	N 770	O 813	S 22	0	0
1	B	556	Total 4384	C 2779	N 770	O 813	S 22	0	0
1	C	556	Total 4384	C 2779	N 770	O 813	S 22	0	0
1	D	556	Total 4384	C 2779	N 770	O 813	S 22	0	0

- Molecule 2 is 6-DIAZENYL-5-OXO-L-NORLEUCINE (three-letter code: DON) (formula: C₆H₁₁N₃O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
2	A	1	Total 10	C 6	N 1	O 3	0
2	B	1	Total 10	C 6	N 1	O 3	0

Continued on next page...

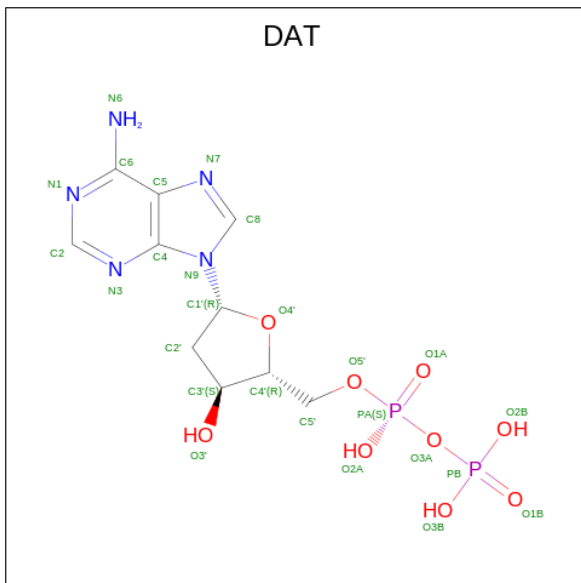
Continued from previous page...

Mol	Chain	Residues	Atoms				AltConf
2	C	1	Total	C	N	O	0
			10	6	1	3	
2	D	1	Total	C	N	O	0
			10	6	1	3	

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	A	4	Total	Mg	0
			4	4	
3	B	4	Total	Mg	0
			4	4	
3	C	5	Total	Mg	0
			5	5	
3	D	3	Total	Mg	0
			3	3	

- Molecule 4 is 2'-DEOXYADENOSINE-5'-DIPHOSPHATE (three-letter code: DAT) (formula: C₁₀H₁₅N₅O₉P₂) (labeled as "Ligand of Interest" by depositor).



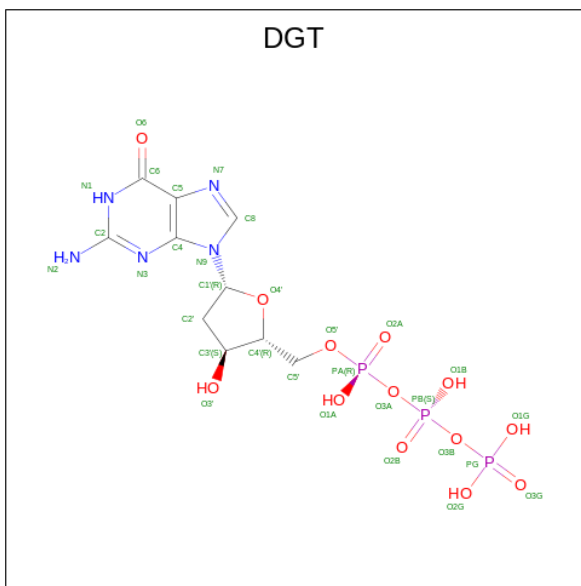
Mol	Chain	Residues	Atoms					AltConf
4	A	1	Total	C	N	O	P	0
			26	10	5	9	2	
4	B	1	Total	C	N	O	P	0
			26	10	5	9	2	

Continued on next page...

Continued from previous page...

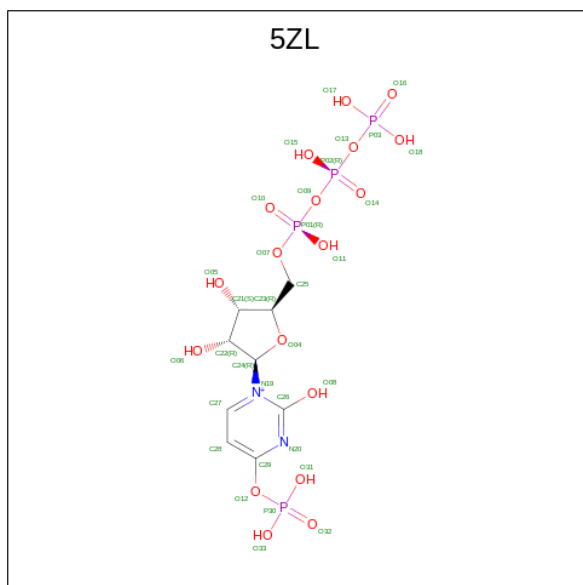
Mol	Chain	Residues	Atoms					AltConf
4	C	1	Total	C	N	O	P	0
			26	10	5	9	2	
4	D	1	Total	C	N	O	P	0
			26	10	5	9	2	

- Molecule 5 is 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (three-letter code: DGT) (formula: C₁₀H₁₆N₅O₁₃P₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is [[[2 {R},3 {S},4 {R},5 {R}]-3,4-bis(oxidanyl)-5-(2-oxidanyl-4-phosphonoxy-pyrimidin-1-yl)oxolan-2-yl]methoxy-oxidanyl-phosphoryl] phosphono hydrogen phosphate (three-letter code: 5ZL) (formula: C₉H₁₇N₂O₁₈P₄) (labeled as "Ligand of Interest" by depositor).

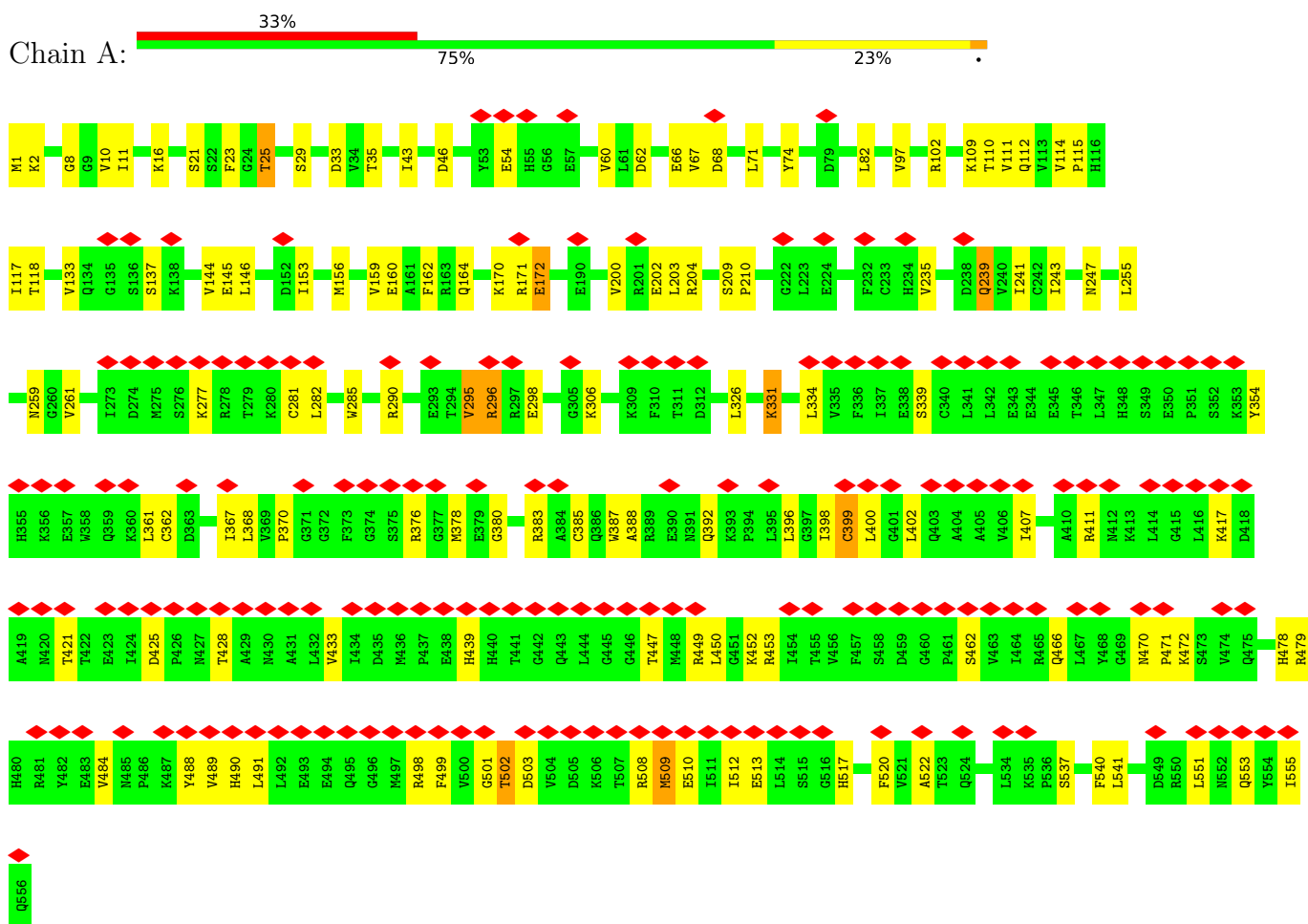


Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
6	A	1	Total	C	N	O	P	0
			32	9	2	17	4	
6	B	1	Total	C	N	O	P	0
			32	9	2	17	4	
6	C	1	Total	C	N	O	P	0
			32	9	2	17	4	
6	D	1	Total	C	N	O	P	0
			32	9	2	17	4	

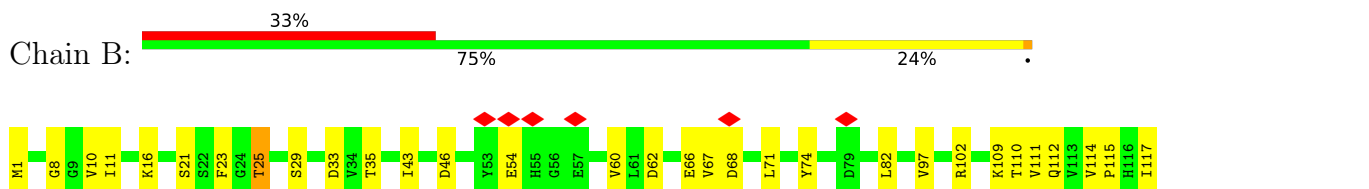
3 Residue-property plots

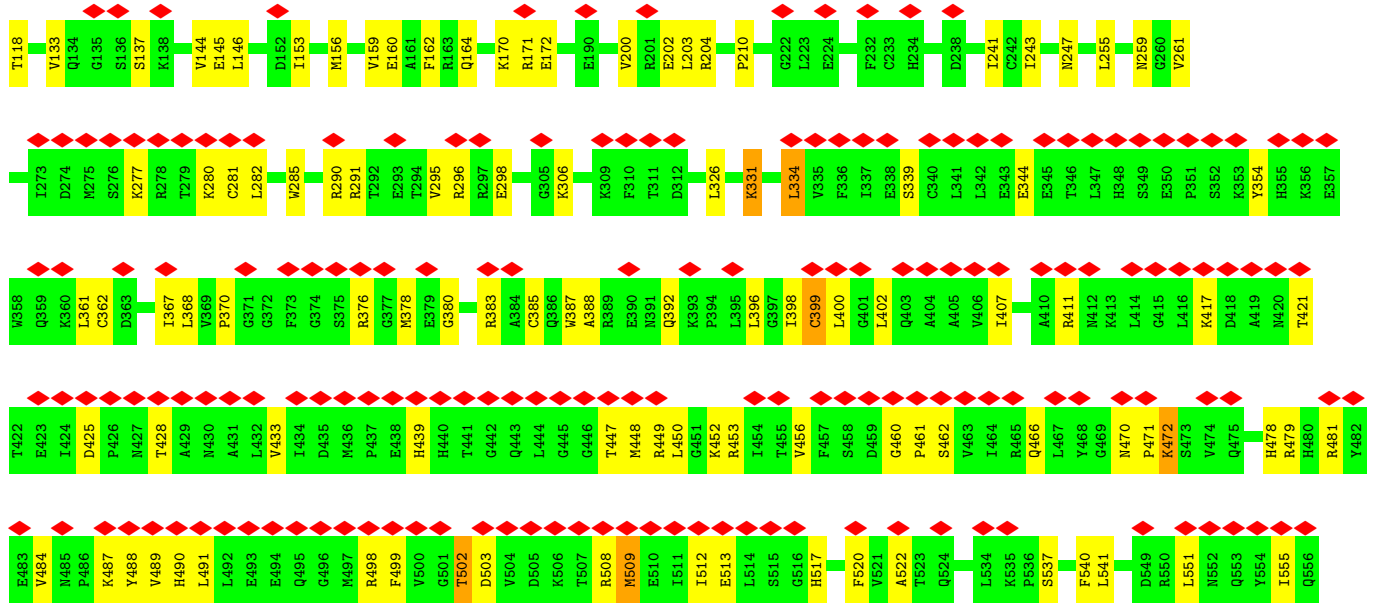
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

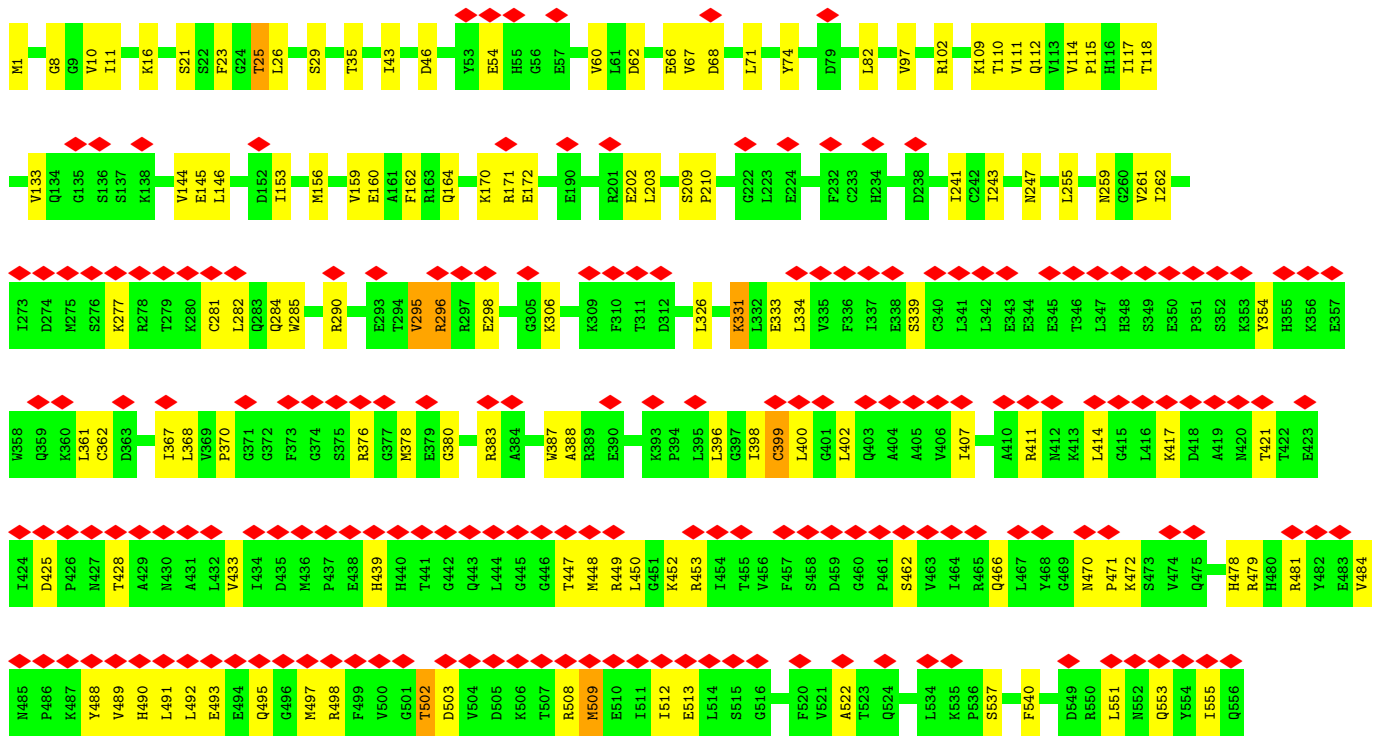
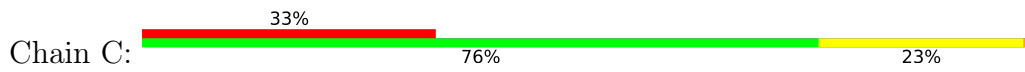


- Molecule 1: CTP synthase

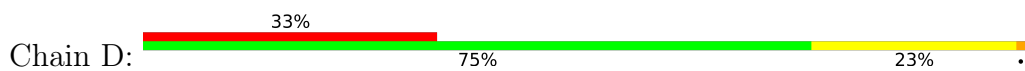




• Molecule 1: CTP synthase



• Molecule 1: CTP synthase



M1	G8	G9	V10	I11	K16	F23	G24	T25	L26	S29	D33	V34	T35	I43	D46	Y53	E54	H55	G56	E57	V60	L61	D62	E66	V67	D68	L71	Y74	D79	L82	V87	R102	K109	T110	V111	Q112	V113	V114	P115	H116	I117	V261	T118															
V133	Q134	G135	S136	S137	K138	V144	E145	L146	D152	I153	M156	V159	A161	F162	R163	Q164	K170	R171	E172	E190	P191	K192	R201	E202	L203	P210	G222	L223	E224	F232	C233	H234	D238	I241	C242	I243	N247	L255	N259	G260	I262																	
I273	D274	M275	S276	K277	R278	T279	K280	C281	L282	Q283	Q284	W285	R290	E293	T294	V295	R296	R297	E298	G305	K306	K309	F310	T311	D312	L326	K331	L332	E333	L334	V335	F336	I337	E338	S339	C340	L341	L342	E343	E344	E345	T346	L347	H348	S349	E350	P351	S352	K353	Y354	H355	K356						
E357	W358	Q359	L360	C362	D363	I367	L368	V369	P370	G371	G372	F373	G374	S375	R376	G377	M378	E379	G380	K381	I382	R383	A384	W387	A388	R389	E390	K393	P394	L395	L396	G397	I398	C399	L400	G401	L402	Q403	A404	A405	V406	I407	A410	R411	N412	K413	L414	G415	L416	K417	D418	A419	N420	T421				
T422	E423	I424	D425	P426	M427	T428	A429	M430	A431	L432	V433	I434	D435	M436	P437	E438	H439	H440	T441	G442	Q443	L444	G445	G446	T447	M448	R449	L450	G451	K452	R453	I454	T455	V456	F457	S458	D459	G460	P461	S462	V463	I464	R465	Q466	L467	Y468	G469	N470	P471	K472	S473	V474	Q475	H478	R479	H480	R481	Y482
E483	V484	M485	P486	K487	Y488	V489	H490	L491	L492	E493	E494	Q495	G496	M497	R498	F499	V500	G501	T502	D503	V504	D505	K506	T507	R508	M509	E510	I511	L512	E513	L514	S515	G516	F520	V521	A522	T523	Q524	L534	K535	P536	S537	F540	D549	R550	L551	N552	Q553	Y554	I555	Q556							

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	639301	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 ($e^-/\text{\AA}^2$)	59.5	Depositor
Minimum defocus (nm)	5000	Depositor
Maximum defocus (nm)	20000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.238	Depositor
Minimum map value	-0.101	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.008	Depositor
Recommended contour level	0.07	Depositor
Map size (\AA)	262.4, 262.4, 262.4	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.025, 1.025, 1.025	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: MG, DGT, 5ZL, DON, DAT

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	0.25	0/4473	0.49	0/6055
1	B	0.25	0/4473	0.49	0/6055
1	C	0.25	0/4473	0.49	0/6055
1	D	0.25	0/4473	0.49	0/6055
All	All	0.25	0/17892	0.49	0/24220

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	4384	0	4430	77	0
1	B	4384	0	4430	79	0
1	C	4384	0	4430	72	0
1	D	4384	0	4430	74	0
2	A	10	0	7	0	0
2	B	10	0	7	0	0
2	C	10	0	7	0	0
2	D	10	0	7	0	0
3	A	4	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	0	0	0
3	C	5	0	0	0	0
3	D	3	0	0	0	0
4	A	26	0	12	0	0
4	B	26	0	12	0	0
4	C	26	0	12	0	0
4	D	26	0	12	0	0
5	A	31	0	12	0	0
5	B	31	0	12	0	0
5	C	31	0	12	0	0
5	D	31	0	12	0	0
6	A	32	0	0	0	0
6	B	32	0	0	0	0
6	C	32	0	0	0	0
6	D	32	0	0	0	0
All	All	17948	0	17844	286	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (286) 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:C:203:LEU:HD23	1:C:210:PRO:HG3	1.75	0.68
1:A:203:LEU:HD23	1:A:210:PRO:HG3	1.76	0.67
1:B:203:LEU:HD23	1:B:210:PRO:HG3	1.75	0.67
1:D:203:LEU:HD23	1:D:210:PRO:HG3	1.75	0.67
1:B:54:GLU:HG2	1:D:110:THR:HG23	1.77	0.66
1:A:2:LYS:NZ	1:A:172:GLU:OE1	2.29	0.66
1:A:54:GLU:HG2	1:C:110:THR:HG23	1.77	0.65
1:A:114:VAL:O	1:A:118:THR:OG1	2.13	0.65
1:C:114:VAL:O	1:C:118:THR:OG1	2.13	0.65
1:D:114:VAL:O	1:D:118:THR:OG1	2.13	0.64
1:D:368:LEU:HD12	1:D:396:LEU:HB3	1.78	0.64
1:B:114:VAL:O	1:B:118:THR:OG1	2.13	0.64
1:C:368:LEU:HD12	1:C:396:LEU:HB3	1.79	0.64
1:B:110:THR:HG23	1:D:54:GLU:HG2	1.78	0.64
1:A:110:THR:HG23	1:C:54:GLU:HG2	1.78	0.63
1:B:450:LEU:O	1:B:453:ARG:NH2	2.33	0.61
1:A:425:ASP:O	1:A:428:THR:OG1	2.18	0.61
1:B:407:ILE:HD11	1:B:421:THR:HG23	1.82	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:425:ASP:O	1:C:428:THR:OG1	2.18	0.60
1:A:450:LEU:O	1:A:453:ARG:NH2	2.33	0.60
1:A:160:GLU:OE1	1:A:164:GLN:NE2	2.34	0.60
1:D:367:ILE:HD12	1:D:388:ALA:HB2	1.83	0.60
1:A:407:ILE:HD11	1:A:421:THR:HG23	1.82	0.60
1:D:407:ILE:HD11	1:D:421:THR:HG23	1.83	0.60
1:C:160:GLU:OE1	1:C:164:GLN:NE2	2.35	0.60
1:B:425:ASP:O	1:B:428:THR:OG1	2.18	0.60
1:C:433:VAL:HA	1:C:484:VAL:HA	1.84	0.60
1:B:368:LEU:HD12	1:B:396:LEU:HB3	1.83	0.60
1:B:160:GLU:OE1	1:B:164:GLN:NE2	2.34	0.59
1:D:425:ASP:O	1:D:428:THR:OG1	2.18	0.59
1:C:367:ILE:HD12	1:C:388:ALA:HB2	1.84	0.59
1:C:407:ILE:HD11	1:C:421:THR:HG23	1.83	0.59
1:D:160:GLU:OE1	1:D:164:GLN:NE2	2.35	0.59
1:B:498:ARG:HH21	1:B:513:GLU:HB3	1.68	0.59
1:C:10:VAL:HG23	1:C:11:ILE:HG23	1.85	0.59
1:A:368:LEU:HD12	1:A:396:LEU:HB3	1.83	0.58
1:A:502:THR:HA	1:A:509:MET:HA	1.85	0.58
1:D:362:CYS:HA	1:D:387:TRP:CH2	2.39	0.58
1:A:498:ARG:HH21	1:A:513:GLU:HB3	1.68	0.58
1:C:362:CYS:HA	1:C:387:TRP:CH2	2.39	0.58
1:D:433:VAL:HA	1:D:484:VAL:HA	1.85	0.58
1:A:433:VAL:HA	1:A:484:VAL:HA	1.86	0.58
1:C:502:THR:OG1	1:C:503:ASP:N	2.36	0.58
1:C:488:TYR:HA	1:C:491:LEU:HD12	1.86	0.58
1:B:367:ILE:HD12	1:B:388:ALA:HB2	1.86	0.58
1:D:10:VAL:HG23	1:D:11:ILE:HG23	1.85	0.57
1:A:10:VAL:HG23	1:A:11:ILE:HG23	1.86	0.57
1:D:354:TYR:OH	1:D:383:ARG:NH2	2.38	0.57
1:D:488:TYR:HA	1:D:491:LEU:HD12	1.86	0.57
1:A:362:CYS:HA	1:A:387:TRP:CH2	2.40	0.57
1:A:367:ILE:HD12	1:A:388:ALA:HB2	1.86	0.57
1:B:362:CYS:HA	1:B:387:TRP:CH2	2.40	0.57
1:C:354:TYR:OH	1:C:383:ARG:NH2	2.38	0.57
1:A:488:TYR:HA	1:A:491:LEU:HD12	1.87	0.57
1:A:74:TYR:OH	1:A:145:GLU:OE1	2.24	0.56
1:A:502:THR:OG1	1:A:503:ASP:N	2.38	0.56
1:B:354:TYR:OH	1:B:383:ARG:NH2	2.39	0.56
1:C:498:ARG:HH21	1:C:513:GLU:HB3	1.70	0.56
1:A:354:TYR:OH	1:A:383:ARG:NH2	2.39	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:VAL:HG23	1:B:11:ILE:HG23	1.86	0.56
1:D:498:ARG:HH21	1:D:513:GLU:HB3	1.70	0.56
1:B:74:TYR:OH	1:B:145:GLU:OE1	2.23	0.56
1:D:462:SER:OG	1:D:513:GLU:OE1	2.24	0.56
1:C:74:TYR:OH	1:C:145:GLU:OE1	2.24	0.56
1:D:502:THR:HA	1:D:509:MET:HA	1.88	0.56
1:D:74:TYR:OH	1:D:145:GLU:OE1	2.24	0.56
1:C:502:THR:HA	1:C:509:MET:HA	1.88	0.55
1:D:502:THR:OG1	1:D:503:ASP:N	2.36	0.55
1:A:235:VAL:HG12	1:A:239:GLN:HB3	1.89	0.55
1:C:462:SER:OG	1:C:513:GLU:OE1	2.24	0.55
1:B:102:ARG:NH2	1:D:46:ASP:OD1	2.39	0.55
1:B:282:LEU:HD23	1:B:285:TRP:CD1	2.42	0.54
1:B:462:SER:OG	1:B:513:GLU:OE1	2.26	0.54
1:C:282:LEU:HD23	1:C:285:TRP:CD1	2.42	0.54
1:A:282:LEU:HD23	1:A:285:TRP:CD1	2.42	0.54
1:B:46:ASP:OD1	1:D:102:ARG:NH2	2.40	0.54
1:D:450:LEU:O	1:D:453:ARG:NH2	2.40	0.54
1:A:102:ARG:NH2	1:C:46:ASP:OD1	2.39	0.54
1:C:450:LEU:O	1:C:453:ARG:NH2	2.40	0.54
1:D:421:THR:HG22	1:D:433:VAL:HB	1.90	0.54
1:B:488:TYR:HA	1:B:491:LEU:HD12	1.90	0.53
1:D:282:LEU:HD23	1:D:285:TRP:CD1	2.42	0.53
1:B:421:THR:HG22	1:B:433:VAL:HB	1.91	0.53
1:B:502:THR:OG1	1:B:503:ASP:N	2.38	0.53
1:A:46:ASP:OD1	1:C:102:ARG:NH2	2.40	0.53
1:B:453:ARG:HB2	1:B:478:HIS:CE1	2.44	0.53
1:A:453:ARG:HB2	1:A:478:HIS:CE1	2.44	0.53
1:C:421:THR:HG22	1:C:433:VAL:HB	1.90	0.53
1:B:502:THR:HA	1:B:509:MET:HA	1.91	0.53
1:C:449:ARG:HH21	1:C:453:ARG:HG3	1.73	0.53
1:A:421:THR:HG22	1:A:433:VAL:HB	1.91	0.53
1:D:449:ARG:HH21	1:D:453:ARG:HG3	1.74	0.53
1:B:449:ARG:HH21	1:B:453:ARG:HG3	1.74	0.52
1:A:462:SER:OG	1:A:513:GLU:OE1	2.26	0.52
1:A:498:ARG:N	1:A:513:GLU:O	2.31	0.52
1:B:295:VAL:HG23	1:B:331:LYS:HG3	1.92	0.52
1:A:449:ARG:HH21	1:A:453:ARG:HG3	1.74	0.52
1:B:498:ARG:N	1:B:513:GLU:O	2.30	0.51
1:B:433:VAL:HA	1:B:484:VAL:HA	1.91	0.51
1:C:82:LEU:HD23	1:C:133:VAL:HG21	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:453:ARG:HB2	1:D:478:HIS:CE1	2.46	0.51
1:D:439:HIS:CE1	1:D:449:ARG:HG3	2.46	0.51
1:A:82:LEU:HD23	1:A:133:VAL:HG21	1.93	0.51
1:C:537:SER:HB3	1:C:540:PHE:HD2	1.75	0.51
1:A:411:ARG:HG3	1:A:417:LYS:HA	1.93	0.50
1:B:112:GLN:HB2	1:B:115:PRO:HD2	1.93	0.50
1:C:453:ARG:HB2	1:C:478:HIS:CE1	2.46	0.50
1:D:112:GLN:HB2	1:D:115:PRO:HD2	1.93	0.50
1:D:537:SER:HB3	1:D:540:PHE:HD2	1.75	0.50
1:C:112:GLN:HB2	1:C:115:PRO:HD2	1.93	0.50
1:C:295:VAL:HG23	1:C:331:LYS:HG3	1.93	0.50
1:A:112:GLN:HB2	1:A:115:PRO:HD2	1.93	0.50
1:B:439:HIS:CE1	1:B:449:ARG:HG3	2.47	0.50
1:B:62:ASP:OD1	1:B:452:LYS:HD3	2.12	0.49
1:C:439:HIS:CE1	1:C:449:ARG:HG3	2.46	0.49
1:A:62:ASP:OD1	1:A:452:LYS:HD3	2.12	0.49
1:A:439:HIS:CE1	1:A:449:ARG:HG3	2.47	0.49
1:A:537:SER:HB3	1:A:540:PHE:HD2	1.77	0.49
1:A:243:ILE:HG23	1:A:255:LEU:HD23	1.94	0.49
1:B:411:ARG:HG3	1:B:417:LYS:HA	1.93	0.49
1:D:82:LEU:HD23	1:D:133:VAL:HG21	1.93	0.49
1:B:537:SER:HB3	1:B:540:PHE:HD2	1.77	0.49
1:C:433:VAL:HG22	1:C:484:VAL:HG12	1.94	0.49
1:B:82:LEU:HD23	1:B:133:VAL:HG21	1.93	0.49
1:D:295:VAL:HG23	1:D:331:LYS:HG3	1.95	0.49
1:A:489:VAL:HG21	1:A:509:MET:SD	2.53	0.48
1:A:512:ILE:HG12	1:A:522:ALA:HB3	1.94	0.48
1:B:512:ILE:HG12	1:B:522:ALA:HB3	1.94	0.48
1:D:62:ASP:OD1	1:D:452:LYS:HD3	2.12	0.48
1:D:433:VAL:HG22	1:D:484:VAL:HG12	1.95	0.48
1:C:62:ASP:OD1	1:C:452:LYS:HD3	2.13	0.48
1:D:367:ILE:HD11	1:D:387:TRP:CZ3	2.48	0.48
1:C:479:ARG:O	1:C:479:ARG:HG3	2.14	0.48
1:A:479:ARG:O	1:A:479:ARG:HG3	2.14	0.48
1:D:241:ILE:HG23	1:D:259:ASN:HD22	1.79	0.48
1:B:241:ILE:HG23	1:B:259:ASN:HD22	1.79	0.48
1:C:367:ILE:HD11	1:C:387:TRP:CZ3	2.49	0.48
1:D:306:LYS:O	1:D:376:ARG:NH2	2.47	0.47
1:D:448:MET:HE2	1:D:481:ARG:HH21	1.78	0.47
1:C:306:LYS:O	1:C:376:ARG:NH2	2.48	0.47
1:C:411:ARG:HG3	1:C:417:LYS:HA	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:LYS:O	1:A:376:ARG:NH2	2.48	0.47
1:B:306:LYS:O	1:B:376:ARG:NH2	2.48	0.47
1:A:295:VAL:HG23	1:A:331:LYS:HG3	1.96	0.47
1:B:489:VAL:HG11	1:B:509:MET:HE2	1.96	0.47
1:D:479:ARG:O	1:D:479:ARG:HG3	2.14	0.47
1:C:489:VAL:HG21	1:C:509:MET:SD	2.55	0.47
1:B:479:ARG:O	1:B:479:ARG:HG3	2.14	0.46
1:C:448:MET:HE2	1:C:481:ARG:HH21	1.81	0.46
1:B:367:ILE:HD11	1:B:387:TRP:CZ3	2.51	0.46
1:B:43:ILE:HD12	1:D:111:VAL:HG12	1.98	0.46
1:A:367:ILE:HD11	1:A:387:TRP:CZ3	2.51	0.46
1:D:489:VAL:HG21	1:D:509:MET:SD	2.55	0.46
1:D:411:ARG:HG3	1:D:417:LYS:HA	1.96	0.46
1:B:29:SER:HB3	1:B:281:CYS:HB2	1.97	0.46
1:A:43:ILE:HD12	1:C:111:VAL:HG12	1.98	0.46
1:A:241:ILE:HG23	1:A:259:ASN:HD22	1.80	0.46
1:D:296:ARG:HA	1:D:296:ARG:HD3	1.75	0.46
1:A:21:SER:O	1:A:25:THR:OG1	2.32	0.45
1:B:111:VAL:HG12	1:D:43:ILE:HD12	1.99	0.45
1:C:29:SER:HB3	1:C:281:CYS:HB2	1.99	0.45
1:A:399:CYS:HB3	1:A:400:LEU:H	1.68	0.45
1:C:241:ILE:HG23	1:C:259:ASN:HD22	1.80	0.45
1:B:153:ILE:HG21	1:D:114:VAL:HG23	1.99	0.45
1:B:114:VAL:HG23	1:D:153:ILE:HG21	1.99	0.45
1:A:114:VAL:HG23	1:C:153:ILE:HG21	1.99	0.45
1:A:153:ILE:HG21	1:C:114:VAL:HG23	1.99	0.45
1:A:111:VAL:HG12	1:C:43:ILE:HD12	1.99	0.45
1:B:392:GLN:HA	1:B:517:HIS:CD2	2.52	0.45
1:C:399:CYS:HB3	1:C:400:LEU:H	1.68	0.45
1:A:200:VAL:O	1:A:204:ARG:HG3	2.18	0.44
1:A:433:VAL:HG22	1:A:484:VAL:HG12	1.99	0.44
1:D:472:LYS:HE3	1:D:472:LYS:HB3	1.83	0.44
1:A:29:SER:HB3	1:A:281:CYS:HB2	2.00	0.44
1:A:453:ARG:HA	1:A:453:ARG:HD3	1.81	0.44
1:B:472:LYS:HE3	1:B:472:LYS:HB3	1.84	0.44
1:C:21:SER:O	1:C:25:THR:OG1	2.34	0.44
1:A:392:GLN:HA	1:A:517:HIS:CD2	2.52	0.44
1:C:97:VAL:HG21	1:C:117:ILE:HD13	1.99	0.44
1:D:29:SER:HB3	1:D:281:CYS:HB2	1.99	0.44
1:D:512:ILE:HG12	1:D:522:ALA:HB3	2.00	0.44
1:C:512:ILE:HG12	1:C:522:ALA:HB3	2.00	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:VAL:O	1:B:204:ARG:HG3	2.18	0.44
1:A:97:VAL:HG21	1:A:117:ILE:HD13	1.99	0.43
1:D:23:PHE:HE1	1:D:261:VAL:HG11	1.83	0.43
1:D:60:VAL:HG22	1:D:66:GLU:HG3	2.00	0.43
1:D:97:VAL:HG21	1:D:117:ILE:HD13	1.98	0.43
1:C:402:LEU:HD11	1:C:512:ILE:HG23	2.01	0.43
1:D:402:LEU:HD11	1:D:512:ILE:HG23	2.00	0.43
1:C:23:PHE:HE1	1:C:261:VAL:HG11	1.84	0.43
1:B:60:VAL:HG22	1:B:66:GLU:HG3	2.00	0.43
1:B:97:VAL:HG21	1:B:117:ILE:HD13	1.99	0.43
1:B:23:PHE:HE1	1:B:261:VAL:HG11	1.84	0.43
1:A:144:VAL:HG11	1:A:162:PHE:CZ	2.54	0.43
1:A:370:PRO:HA	1:A:398:ILE:HB	2.00	0.43
1:B:67:VAL:HG21	1:B:71:LEU:HD22	2.00	0.43
1:B:370:PRO:HA	1:B:398:ILE:HB	2.00	0.43
1:B:448:MET:HE2	1:B:481:ARG:HH21	1.84	0.43
1:B:509:MET:SD	1:B:509:MET:N	2.90	0.43
1:C:60:VAL:HG22	1:C:66:GLU:HG3	2.00	0.43
1:C:296:ARG:HA	1:C:296:ARG:HD3	1.75	0.43
1:A:67:VAL:HG21	1:A:71:LEU:HD22	2.00	0.43
1:A:498:ARG:HE	1:A:498:ARG:HB2	1.62	0.43
1:B:67:VAL:HB	1:B:71:LEU:HD13	2.01	0.43
1:B:449:ARG:NH1	1:B:508:ARG:HG2	2.33	0.43
1:C:361:LEU:HG	1:C:387:TRP:CZ3	2.54	0.43
1:A:23:PHE:HE1	1:A:261:VAL:HG11	1.84	0.42
1:A:489:VAL:HG13	1:A:499:PHE:CD2	2.54	0.42
1:C:243:ILE:HG23	1:C:255:LEU:HD23	2.01	0.42
1:D:361:LEU:HG	1:D:387:TRP:CZ3	2.54	0.42
1:A:385:CYS:SG	1:A:520:PHE:HZ	2.42	0.42
1:B:385:CYS:SG	1:B:520:PHE:HZ	2.42	0.42
1:D:67:VAL:HG21	1:D:71:LEU:HD22	2.00	0.42
1:A:553:GLN:OE1	1:A:553:GLN:N	2.52	0.42
1:C:67:VAL:HG21	1:C:71:LEU:HD22	2.00	0.42
1:D:67:VAL:HB	1:D:71:LEU:HD13	2.01	0.42
1:A:60:VAL:HG22	1:A:66:GLU:HG3	2.01	0.42
1:C:209:SER:HA	1:C:210:PRO:HD3	1.93	0.42
1:C:370:PRO:HA	1:C:398:ILE:HB	2.01	0.42
1:C:551:LEU:O	1:C:555:ILE:HG12	2.19	0.42
1:A:296:ARG:HA	1:A:296:ARG:HD3	1.74	0.42
1:A:551:LEU:O	1:A:555:ILE:HG12	2.19	0.42
1:B:8:GLY:CA	1:B:16:LYS:HD3	2.50	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:GLY:CA	1:D:16:LYS:HD3	2.50	0.42
1:A:339:SER:O	1:A:380:GLY:HA3	2.19	0.42
1:C:449:ARG:NH1	1:C:508:ARG:HG2	2.35	0.42
1:C:553:GLN:OE1	1:C:553:GLN:N	2.52	0.42
1:D:192:LYS:HA	1:D:192:LYS:HD2	1.78	0.42
1:C:144:VAL:HG11	1:C:162:PHE:CZ	2.55	0.42
1:A:209:SER:HA	1:A:210:PRO:HD3	1.94	0.42
1:A:449:ARG:NH1	1:A:508:ARG:HG2	2.35	0.42
1:B:144:VAL:HG11	1:B:162:PHE:CZ	2.54	0.42
1:B:489:VAL:HG13	1:B:499:PHE:CD2	2.54	0.42
1:C:67:VAL:HB	1:C:71:LEU:HD13	2.01	0.42
1:D:243:ILE:HG23	1:D:255:LEU:HD23	2.00	0.42
1:A:67:VAL:HB	1:A:71:LEU:HD13	2.01	0.41
1:B:243:ILE:HG23	1:B:255:LEU:HD23	2.00	0.41
1:B:551:LEU:O	1:B:555:ILE:HG12	2.19	0.41
1:C:339:SER:O	1:C:380:GLY:HA3	2.20	0.41
1:D:144:VAL:HG11	1:D:162:PHE:CZ	2.54	0.41
1:D:280:LYS:HE2	1:D:280:LYS:HB3	1.88	0.41
1:B:334:LEU:HD23	1:B:334:LEU:HA	1.84	0.41
1:B:402:LEU:HD11	1:B:512:ILE:HG23	2.02	0.41
1:D:370:PRO:HA	1:D:398:ILE:HB	2.01	0.41
1:D:382:ILE:HD13	1:D:411:ARG:HH12	1.85	0.41
1:B:344:GLU:H	1:B:344:GLU:HG2	1.70	0.41
1:B:487:LYS:HA	1:B:487:LYS:HD3	1.86	0.41
1:D:551:LEU:O	1:D:555:ILE:HG12	2.19	0.41
1:A:8:GLY:CA	1:A:16:LYS:HD3	2.50	0.41
1:B:280:LYS:HE2	1:B:280:LYS:HB3	1.90	0.41
1:A:361:LEU:HG	1:A:387:TRP:CZ3	2.55	0.41
1:B:339:SER:O	1:B:380:GLY:HA3	2.20	0.41
1:C:298:GLU:OE1	1:C:333:GLU:HB2	2.21	0.41
1:C:470:ASN:N	1:C:471:PRO:HD3	2.36	0.41
1:A:470:ASN:N	1:A:471:PRO:HD3	2.36	0.41
1:C:8:GLY:CA	1:C:16:LYS:HD3	2.50	0.41
1:D:298:GLU:OE1	1:D:333:GLU:HB2	2.21	0.41
1:D:493:GLU:HA	1:D:497:MET:O	2.20	0.41
1:A:402:LEU:HD11	1:A:512:ILE:HG23	2.02	0.41
1:B:33:ASP:OD2	1:B:137:SER:HB2	2.21	0.41
1:B:291:ARG:O	1:B:295:VAL:HG13	2.20	0.41
1:B:361:LEU:HG	1:B:387:TRP:CZ3	2.55	0.41
1:B:453:ARG:HA	1:B:453:ARG:HD3	1.81	0.41
1:C:156:MET:HA	1:C:159:VAL:HG22	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:414:LEU:HD23	1:C:492:LEU:HD23	2.03	0.41
1:D:339:SER:O	1:D:380:GLY:HA3	2.20	0.41
1:D:414:LEU:HD23	1:D:492:LEU:HD23	2.03	0.41
1:B:456:VAL:N	1:B:502:THR:O	2.31	0.41
1:D:156:MET:HA	1:D:159:VAL:HG22	2.03	0.41
1:D:449:ARG:NH1	1:D:508:ARG:HG2	2.35	0.41
1:B:399:CYS:HB3	1:B:400:LEU:H	1.68	0.40
1:B:460:GLY:HA2	1:B:461:PRO:HD3	1.96	0.40
1:C:493:GLU:HA	1:C:497:MET:O	2.20	0.40
1:D:456:VAL:N	1:D:502:THR:O	2.32	0.40
1:B:21:SER:O	1:B:25:THR:OG1	2.33	0.40
1:B:156:MET:HA	1:B:159:VAL:HG22	2.03	0.40
1:D:470:ASN:N	1:D:471:PRO:HD3	2.36	0.40
1:A:33:ASP:OD2	1:A:137:SER:HB2	2.21	0.40
1:B:470:ASN:N	1:B:471:PRO:HD3	2.36	0.40
1:C:479:ARG:HG2	1:C:479:ARG:HH11	1.87	0.40
1:D:26:LEU:HD13	1:D:262:ILE:HG13	2.03	0.40
1:D:33:ASP:OD2	1:D:137:SER:HB2	2.22	0.40
1:A:156:MET:HA	1:A:159:VAL:HG22	2.04	0.40
1:A:501:GLY:O	1:A:510:GLU:N	2.30	0.40
1:C:26:LEU:HD13	1:C:262:ILE:HG13	2.04	0.40
1:D:399:CYS:HB3	1:D:400:LEU:H	1.68	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	554/556 (100%)	540 (98%)	14 (2%)	0	100	100
1	B	554/556 (100%)	539 (97%)	15 (3%)	0	100	100
1	C	554/556 (100%)	540 (98%)	14 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	554/556 (100%)	539 (97%)	15 (3%)	0	100	100
All	All	2216/2224 (100%)	2158 (97%)	58 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	486/486 (100%)	457 (94%)	29 (6%)	16	28
1	B	486/486 (100%)	459 (94%)	27 (6%)	17	32
1	C	486/486 (100%)	458 (94%)	28 (6%)	17	30
1	D	486/486 (100%)	457 (94%)	29 (6%)	16	28
All	All	1944/1944 (100%)	1831 (94%)	113 (6%)	19	30

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	25	THR
1	A	35	THR
1	A	68	ASP
1	A	109	LYS
1	A	146	LEU
1	A	170	LYS
1	A	171	ARG
1	A	172	GLU
1	A	202	GLU
1	A	239	GLN
1	A	247	ASN
1	A	277	LYS
1	A	290	ARG
1	A	295	VAL
1	A	296	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	298	GLU
1	A	326	LEU
1	A	331	LYS
1	A	334	LEU
1	A	378	MET
1	A	399	CYS
1	A	447	THR
1	A	466	GLN
1	A	472	LYS
1	A	490	HIS
1	A	502	THR
1	A	509	MET
1	A	541	LEU
1	B	1	MET
1	B	25	THR
1	B	35	THR
1	B	68	ASP
1	B	109	LYS
1	B	146	LEU
1	B	170	LYS
1	B	171	ARG
1	B	172	GLU
1	B	202	GLU
1	B	247	ASN
1	B	277	LYS
1	B	290	ARG
1	B	296	ARG
1	B	298	GLU
1	B	326	LEU
1	B	331	LYS
1	B	334	LEU
1	B	378	MET
1	B	399	CYS
1	B	447	THR
1	B	466	GLN
1	B	472	LYS
1	B	490	HIS
1	B	502	THR
1	B	509	MET
1	B	541	LEU
1	C	1	MET
1	C	25	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	35	THR
1	C	68	ASP
1	C	109	LYS
1	C	146	LEU
1	C	170	LYS
1	C	171	ARG
1	C	172	GLU
1	C	202	GLU
1	C	247	ASN
1	C	277	LYS
1	C	284	GLN
1	C	290	ARG
1	C	295	VAL
1	C	296	ARG
1	C	326	LEU
1	C	331	LYS
1	C	334	LEU
1	C	378	MET
1	C	399	CYS
1	C	447	THR
1	C	466	GLN
1	C	472	LYS
1	C	490	HIS
1	C	495	GLN
1	C	502	THR
1	C	509	MET
1	D	1	MET
1	D	25	THR
1	D	35	THR
1	D	68	ASP
1	D	109	LYS
1	D	146	LEU
1	D	170	LYS
1	D	171	ARG
1	D	172	GLU
1	D	202	GLU
1	D	247	ASN
1	D	277	LYS
1	D	284	GLN
1	D	290	ARG
1	D	295	VAL
1	D	296	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	326	LEU
1	D	331	LYS
1	D	334	LEU
1	D	378	MET
1	D	399	CYS
1	D	436	MET
1	D	447	THR
1	D	466	GLN
1	D	472	LYS
1	D	490	HIS
1	D	495	GLN
1	D	502	THR
1	D	509	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 32 ligands modelled in this entry, 16 are monoatomic - leaving 16 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	DGT	B	607	3	26,33,33	3.29	14 (53%)	32,52,52	1.53	8 (25%)
2	DON	C	601	1	8,9,11	0.79	0	10,11,13	1.40	2 (20%)
2	DON	B	602	1	8,9,11	0.80	0	10,11,13	1.41	1 (10%)
4	DAT	B	606	3	24,28,28	0.87	1 (4%)	28,43,43	1.32	4 (14%)
4	DAT	D	605	3	24,28,28	0.87	1 (4%)	28,43,43	1.32	4 (14%)
6	5ZL	A	607	3	29,33,34	2.83	11 (37%)	36,52,54	1.10	3 (8%)
6	5ZL	B	608	3	29,33,34	2.83	11 (37%)	36,52,54	1.10	3 (8%)
5	DGT	D	606	3	26,33,33	3.28	14 (53%)	32,52,52	1.53	8 (25%)
4	DAT	A	605	3	24,28,28	0.87	1 (4%)	28,43,43	1.32	4 (14%)
6	5ZL	C	608	3	29,33,34	2.83	11 (37%)	36,52,54	1.10	3 (8%)
6	5ZL	D	607	3	29,33,34	2.84	11 (37%)	36,52,54	1.10	3 (8%)
2	DON	D	601	1	8,9,11	0.79	0	10,11,13	1.41	2 (20%)
5	DGT	C	607	3	26,33,33	3.28	14 (53%)	32,52,52	1.53	8 (25%)
2	DON	A	601	1	8,9,11	0.79	0	10,11,13	1.41	2 (20%)
4	DAT	C	606	3	24,28,28	0.87	1 (4%)	28,43,43	1.32	4 (14%)
5	DGT	A	606	3	26,33,33	3.28	14 (53%)	32,52,52	1.54	8 (25%)

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. ^{1,2} means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	DGT	B	607	3	-	8/18/34/34	0/3/3/3
2	DON	C	601	1	-	2/9/9/12	-
2	DON	B	602	1	-	2/9/9/12	-
4	DAT	B	606	3	-	4/12/28/28	0/3/3/3
4	DAT	D	605	3	-	4/12/28/28	0/3/3/3
6	5ZL	A	607	3	-	15/24/39/43	0/2/2/2
6	5ZL	B	608	3	-	15/24/39/43	0/2/2/2
5	DGT	D	606	3	-	8/18/34/34	0/3/3/3
4	DAT	A	605	3	-	4/12/28/28	0/3/3/3
6	5ZL	C	608	3	-	15/24/39/43	0/2/2/2
6	5ZL	D	607	3	-	15/24/39/43	0/2/2/2
2	DON	D	601	1	-	2/9/9/12	-
5	DGT	C	607	3	-	8/18/34/34	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DON	A	601	1	-	2/9/9/12	-
4	DAT	C	606	3	-	4/12/28/28	0/3/3/3
5	DGT	A	606	3	-	8/18/34/34	0/3/3/3

All (104) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	606	DGT	C2'-C3'	-7.52	1.33	1.52
6	C	608	5ZL	O04-C23	7.52	1.61	1.45
6	D	607	5ZL	O04-C23	7.52	1.61	1.45
6	B	608	5ZL	O04-C23	7.51	1.61	1.45
5	C	607	DGT	C2'-C3'	-7.51	1.33	1.52
5	B	607	DGT	C2'-C3'	-7.51	1.33	1.52
5	A	606	DGT	C2'-C3'	-7.50	1.33	1.52
6	A	607	5ZL	O04-C23	7.49	1.61	1.45
6	A	607	5ZL	C21-C23	-6.37	1.35	1.53
6	B	608	5ZL	C21-C23	-6.36	1.35	1.53
6	C	608	5ZL	C21-C23	-6.36	1.35	1.53
6	D	607	5ZL	C21-C23	-6.36	1.35	1.53
5	D	606	DGT	C2-N3	5.52	1.46	1.33
5	A	606	DGT	C2-N3	5.52	1.46	1.33
5	B	607	DGT	C2-N3	5.52	1.46	1.33
5	C	607	DGT	C2-N3	5.48	1.46	1.33
6	C	608	5ZL	P30-O12	5.40	1.67	1.59
6	A	607	5ZL	P30-O12	5.39	1.67	1.59
6	B	608	5ZL	P30-O12	5.39	1.67	1.59
6	D	607	5ZL	P30-O12	5.38	1.67	1.59
5	C	607	DGT	O4'-C4'	5.28	1.56	1.45
5	D	606	DGT	O4'-C4'	5.26	1.56	1.45
5	B	607	DGT	O4'-C4'	5.25	1.56	1.45
5	A	606	DGT	O4'-C4'	5.25	1.56	1.45
5	D	606	DGT	C4-N3	5.20	1.50	1.37
5	A	606	DGT	C4-N3	5.20	1.50	1.37
5	B	607	DGT	C4-N3	5.20	1.50	1.37
5	C	607	DGT	C4-N3	5.19	1.50	1.37
5	A	606	DGT	O4'-C1'	-5.19	1.30	1.42
5	B	607	DGT	O4'-C1'	-5.17	1.30	1.42
5	C	607	DGT	O4'-C1'	-5.17	1.30	1.42
5	D	606	DGT	O4'-C1'	-5.15	1.30	1.42
6	C	608	5ZL	O04-C24	-5.03	1.31	1.42
6	D	607	5ZL	O04-C24	-5.03	1.31	1.42
6	A	607	5ZL	O04-C24	-5.01	1.31	1.42

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	608	5ZL	O04-C24	-5.01	1.31	1.42
5	C	607	DGT	C2-N2	4.79	1.45	1.34
5	D	606	DGT	C2-N2	4.76	1.45	1.34
5	A	606	DGT	C2-N2	4.76	1.45	1.34
5	B	607	DGT	C2-N2	4.76	1.45	1.34
6	A	607	5ZL	O08-C26	4.40	1.44	1.25
6	C	608	5ZL	O08-C26	4.40	1.44	1.25
6	D	607	5ZL	O08-C26	4.38	1.44	1.25
6	B	608	5ZL	O08-C26	4.37	1.44	1.25
5	B	607	DGT	C2'-C1'	4.11	1.63	1.52
5	A	606	DGT	C2'-C1'	4.09	1.63	1.52
5	C	607	DGT	C2'-C1'	4.07	1.63	1.52
5	D	606	DGT	C2'-C1'	4.07	1.63	1.52
6	D	607	5ZL	C26-N19	-3.91	1.32	1.38
5	A	606	DGT	C5'-C4'	-3.90	1.39	1.51
5	D	606	DGT	C5'-C4'	-3.89	1.39	1.51
5	B	607	DGT	C5'-C4'	-3.87	1.39	1.51
6	C	608	5ZL	C26-N19	-3.87	1.33	1.38
5	C	607	DGT	C5'-C4'	-3.87	1.39	1.51
6	A	607	5ZL	C26-N19	-3.84	1.33	1.38
6	B	608	5ZL	C26-N19	-3.80	1.33	1.38
5	B	607	DGT	C6-N1	3.71	1.43	1.37
5	A	606	DGT	C6-N1	3.68	1.43	1.37
5	C	607	DGT	C6-N1	3.68	1.43	1.37
5	D	606	DGT	C6-N1	3.67	1.43	1.37
6	A	607	5ZL	O05-C21	3.59	1.51	1.43
6	C	608	5ZL	O05-C21	3.58	1.50	1.43
6	D	607	5ZL	O05-C21	3.58	1.50	1.43
6	B	608	5ZL	O05-C21	3.55	1.50	1.43
6	D	607	5ZL	P01-O07	3.38	1.73	1.59
6	A	607	5ZL	P01-O07	3.38	1.73	1.59
6	B	608	5ZL	P01-O07	3.36	1.72	1.59
6	C	608	5ZL	P01-O07	3.35	1.72	1.59
5	B	607	DGT	C5-C6	3.23	1.54	1.47
5	D	606	DGT	C5-C6	3.23	1.54	1.47
5	C	607	DGT	C5-C6	3.19	1.53	1.47
5	A	606	DGT	C5-C6	3.17	1.53	1.47
5	B	607	DGT	C2-N1	2.68	1.44	1.37
5	C	607	DGT	C2-N1	2.68	1.44	1.37
5	A	606	DGT	C2-N1	2.68	1.44	1.37
5	D	606	DGT	C2-N1	2.65	1.44	1.37
5	B	607	DGT	O3'-C3'	2.64	1.48	1.43

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	606	DGT	O3'-C3'	2.64	1.48	1.43
5	A	606	DGT	O3'-C3'	2.62	1.48	1.43
5	C	607	DGT	O3'-C3'	2.60	1.48	1.43
6	D	607	5ZL	O12-C29	2.44	1.43	1.40
6	A	607	5ZL	O12-C29	2.41	1.43	1.40
6	B	608	5ZL	O12-C29	2.39	1.43	1.40
5	A	606	DGT	C5-C4	-2.37	1.37	1.43
6	C	608	5ZL	O12-C29	2.37	1.43	1.40
5	B	607	DGT	C5-C4	-2.37	1.37	1.43
5	C	607	DGT	C5-C4	-2.37	1.37	1.43
5	D	606	DGT	C5-C4	-2.36	1.37	1.43
4	A	605	DAT	C5-C4	2.28	1.47	1.40
4	B	606	DAT	C5-C4	2.28	1.47	1.40
4	D	605	DAT	C5-C4	2.28	1.47	1.40
4	C	606	DAT	C5-C4	2.26	1.46	1.40
6	A	607	5ZL	C22-C24	2.23	1.58	1.52
6	B	608	5ZL	C22-C24	2.20	1.58	1.52
5	A	606	DGT	O6-C6	-2.19	1.18	1.23
5	C	607	DGT	O6-C6	-2.19	1.18	1.23
5	B	607	DGT	O6-C6	-2.19	1.18	1.23
6	C	608	5ZL	C22-C24	2.18	1.58	1.52
6	D	607	5ZL	C22-C24	2.18	1.58	1.52
5	D	606	DGT	O6-C6	-2.18	1.18	1.23
6	A	607	5ZL	C27-N19	-2.11	1.33	1.35
6	C	608	5ZL	C27-N19	-2.06	1.33	1.35
6	D	607	5ZL	C27-N19	-2.06	1.33	1.35
6	B	608	5ZL	C27-N19	-2.05	1.33	1.35

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	606	DGT	C5-C6-N1	3.57	120.26	113.95
5	D	606	DGT	C5-C6-N1	3.55	120.23	113.95
5	C	607	DGT	C5-C6-N1	3.55	120.22	113.95
5	B	607	DGT	C5-C6-N1	3.54	120.21	113.95
4	B	606	DAT	N3-C2-N1	-3.23	123.62	128.68
4	C	606	DAT	N3-C2-N1	-3.23	123.62	128.68
4	A	605	DAT	N3-C2-N1	-3.23	123.64	128.68
4	D	605	DAT	N3-C2-N1	-3.22	123.65	128.68
5	A	606	DGT	C2-N1-C6	-3.21	119.19	125.10
5	C	607	DGT	C2-N1-C6	-3.19	119.22	125.10
5	B	607	DGT	C2-N1-C6	-3.19	119.22	125.10

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	606	DGT	C2-N1-C6	-3.16	119.28	125.10
2	A	601	DON	OXT-C-O	-3.05	117.17	124.09
2	D	601	DON	OXT-C-O	-3.04	117.18	124.09
2	B	602	DON	OXT-C-O	-3.04	117.18	124.09
2	C	601	DON	OXT-C-O	-3.02	117.23	124.09
5	A	606	DGT	PB-O3B-PG	-2.85	123.05	132.83
5	B	607	DGT	PB-O3B-PG	-2.85	123.05	132.83
5	C	607	DGT	PB-O3B-PG	-2.85	123.06	132.83
5	D	606	DGT	PB-O3B-PG	-2.83	123.11	132.83
4	B	606	DAT	C4-C5-N7	-2.73	106.56	109.40
4	D	605	DAT	C4-C5-N7	-2.72	106.57	109.40
4	A	605	DAT	C4-C5-N7	-2.71	106.57	109.40
4	C	606	DAT	C4-C5-N7	-2.67	106.62	109.40
5	A	606	DGT	C8-N7-C5	2.65	108.04	102.99
5	B	607	DGT	C8-N7-C5	2.63	108.00	102.99
5	D	606	DGT	C8-N7-C5	2.63	108.00	102.99
5	C	607	DGT	C8-N7-C5	2.62	107.98	102.99
4	D	605	DAT	PA-O3A-PB	-2.57	123.99	132.83
4	B	606	DAT	PA-O3A-PB	-2.57	124.00	132.83
4	C	606	DAT	PA-O3A-PB	-2.57	124.01	132.83
4	A	605	DAT	PA-O3A-PB	-2.56	124.03	132.83
6	D	607	5ZL	C28-C29-N20	-2.33	118.17	123.31
6	B	608	5ZL	C28-C29-N20	-2.33	118.18	123.31
6	A	607	5ZL	C28-C29-N20	-2.32	118.19	123.31
6	C	608	5ZL	C28-C29-N20	-2.32	118.20	123.31
5	A	606	DGT	O6-C6-C5	-2.32	119.84	124.37
5	B	607	DGT	O6-C6-C5	-2.30	119.87	124.37
5	D	606	DGT	O6-C6-C5	-2.29	119.90	124.37
5	C	607	DGT	O6-C6-C5	-2.25	119.97	124.37
5	D	606	DGT	C3'-C2'-C1'	2.25	108.17	102.54
5	C	607	DGT	C3'-C2'-C1'	2.24	108.14	102.54
6	C	608	5ZL	P02-O13-P03	-2.24	125.15	132.83
6	D	607	5ZL	P02-O13-P03	-2.24	125.15	132.83
6	B	608	5ZL	P02-O13-P03	-2.23	125.16	132.83
6	A	607	5ZL	P02-O13-P03	-2.23	125.17	132.83
5	A	606	DGT	C3'-C2'-C1'	2.22	108.11	102.54
5	B	607	DGT	C3'-C2'-C1'	2.22	108.11	102.54
5	A	606	DGT	PA-O3A-PB	-2.18	125.36	132.83
5	D	606	DGT	PA-O3A-PB	-2.17	125.37	132.83
5	B	607	DGT	PA-O3A-PB	-2.17	125.37	132.83
5	C	607	DGT	PA-O3A-PB	-2.16	125.41	132.83
5	A	606	DGT	C2'-C3'-C4'	2.14	107.22	102.76

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	607	DGT	C2'-C3'-C4'	2.14	107.22	102.76
5	D	606	DGT	C2'-C3'-C4'	2.12	107.18	102.76
5	C	607	DGT	C2'-C3'-C4'	2.12	107.18	102.76
6	A	607	5ZL	P02-O09-P01	-2.10	125.63	132.83
6	C	608	5ZL	P02-O09-P01	-2.10	125.63	132.83
6	D	607	5ZL	P02-O09-P01	-2.09	125.64	132.83
6	B	608	5ZL	P02-O09-P01	-2.09	125.65	132.83
4	A	605	DAT	O3B-PB-O2B	2.08	115.58	107.64
4	B	606	DAT	O3B-PB-O2B	2.08	115.58	107.64
4	C	606	DAT	O3B-PB-O2B	2.08	115.58	107.64
4	D	605	DAT	O3B-PB-O2B	2.07	115.54	107.64
2	D	601	DON	CB-C1E-C1D	-2.04	112.48	114.57
2	C	601	DON	CB-C1E-C1D	-2.01	112.51	114.57
2	A	601	DON	CB-C1E-C1D	-2.01	112.51	114.57

There are no chirality outliers.

All (116) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	605	DAT	C5'-O5'-PA-O1A
4	B	606	DAT	C5'-O5'-PA-O1A
4	C	606	DAT	C5'-O5'-PA-O1A
4	D	605	DAT	C5'-O5'-PA-O1A
5	A	606	DGT	C5'-O5'-PA-O3A
5	A	606	DGT	C5'-O5'-PA-O2A
5	A	606	DGT	O4'-C4'-C5'-O5'
5	B	607	DGT	C5'-O5'-PA-O3A
5	B	607	DGT	C5'-O5'-PA-O2A
5	B	607	DGT	O4'-C4'-C5'-O5'
5	C	607	DGT	C5'-O5'-PA-O3A
5	C	607	DGT	C5'-O5'-PA-O2A
5	C	607	DGT	O4'-C4'-C5'-O5'
5	D	606	DGT	C5'-O5'-PA-O3A
5	D	606	DGT	C5'-O5'-PA-O2A
5	D	606	DGT	O4'-C4'-C5'-O5'
6	A	607	5ZL	O04-C23-C25-O07
6	A	607	5ZL	C25-O07-P01-O09
6	A	607	5ZL	P02-O13-P03-O17
6	B	608	5ZL	O04-C23-C25-O07
6	B	608	5ZL	C25-O07-P01-O09
6	B	608	5ZL	P02-O13-P03-O17
6	C	608	5ZL	O04-C23-C25-O07

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
6	C	608	5ZL	C25-O07-P01-O09
6	C	608	5ZL	P02-O13-P03-O17
6	D	607	5ZL	O04-C23-C25-O07
6	D	607	5ZL	C25-O07-P01-O09
6	D	607	5ZL	P02-O13-P03-O17
6	A	607	5ZL	C21-C23-C25-O07
6	B	608	5ZL	C21-C23-C25-O07
6	C	608	5ZL	C21-C23-C25-O07
6	D	607	5ZL	C21-C23-C25-O07
5	A	606	DGT	C3'-C4'-C5'-O5'
5	B	607	DGT	C3'-C4'-C5'-O5'
5	C	607	DGT	C3'-C4'-C5'-O5'
5	D	606	DGT	C3'-C4'-C5'-O5'
2	A	601	DON	O1J-C1D-C1E-CB
2	B	602	DON	O1J-C1D-C1E-CB
2	A	601	DON	C1C-C1D-C1E-CB
2	B	602	DON	C1C-C1D-C1E-CB
2	C	601	DON	O1J-C1D-C1E-CB
2	C	601	DON	C1C-C1D-C1E-CB
2	D	601	DON	O1J-C1D-C1E-CB
2	D	601	DON	C1C-C1D-C1E-CB
6	A	607	5ZL	P01-O09-P02-O14
6	B	608	5ZL	P01-O09-P02-O14
6	C	608	5ZL	P01-O09-P02-O14
6	D	607	5ZL	P01-O09-P02-O14
5	A	606	DGT	PA-O3A-PB-O3B
5	B	607	DGT	PA-O3A-PB-O3B
5	C	607	DGT	PA-O3A-PB-O3B
5	D	606	DGT	PA-O3A-PB-O3B
6	A	607	5ZL	N20-C29-O12-P30
6	B	608	5ZL	N20-C29-O12-P30
6	C	608	5ZL	N20-C29-O12-P30
6	D	607	5ZL	N20-C29-O12-P30
6	A	607	5ZL	C25-O07-P01-O10
6	A	607	5ZL	C25-O07-P01-O11
6	B	608	5ZL	C25-O07-P01-O10
6	B	608	5ZL	C25-O07-P01-O11
6	C	608	5ZL	C25-O07-P01-O10
6	C	608	5ZL	C25-O07-P01-O11
6	D	607	5ZL	C25-O07-P01-O10
6	D	607	5ZL	C25-O07-P01-O11
4	A	605	DAT	O4'-C4'-C5'-O5'

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
4	B	606	DAT	O4'-C4'-C5'-O5'
4	C	606	DAT	O4'-C4'-C5'-O5'
4	D	605	DAT	O4'-C4'-C5'-O5'
5	A	606	DGT	PG-O3B-PB-O1B
5	B	607	DGT	PG-O3B-PB-O1B
5	C	607	DGT	PG-O3B-PB-O1B
5	D	606	DGT	PG-O3B-PB-O1B
6	A	607	5ZL	P02-O09-P01-O10
6	A	607	5ZL	P03-O13-P02-O14
6	B	608	5ZL	P02-O09-P01-O10
6	B	608	5ZL	P03-O13-P02-O14
6	C	608	5ZL	P02-O09-P01-O10
6	C	608	5ZL	P03-O13-P02-O14
6	D	607	5ZL	P02-O09-P01-O10
6	D	607	5ZL	P03-O13-P02-O14
4	A	605	DAT	C4'-C5'-O5'-PA
4	B	606	DAT	C4'-C5'-O5'-PA
4	C	606	DAT	C4'-C5'-O5'-PA
4	D	605	DAT	C4'-C5'-O5'-PA
6	A	607	5ZL	C28-C29-O12-P30
6	B	608	5ZL	C28-C29-O12-P30
6	C	608	5ZL	C28-C29-O12-P30
6	D	607	5ZL	C28-C29-O12-P30
6	A	607	5ZL	P02-O13-P03-O18
6	B	608	5ZL	P02-O13-P03-O18
6	C	608	5ZL	P02-O13-P03-O18
6	D	607	5ZL	P02-O13-P03-O18
4	A	605	DAT	C5'-O5'-PA-O3A
4	B	606	DAT	C5'-O5'-PA-O3A
4	C	606	DAT	C5'-O5'-PA-O3A
4	D	605	DAT	C5'-O5'-PA-O3A
5	A	606	DGT	PA-O3A-PB-O1B
5	B	607	DGT	PA-O3A-PB-O1B
5	C	607	DGT	PA-O3A-PB-O1B
5	D	606	DGT	PA-O3A-PB-O1B
6	A	607	5ZL	P02-O09-P01-O11
6	A	607	5ZL	P01-O09-P02-O15
6	B	608	5ZL	P02-O09-P01-O11
6	B	608	5ZL	P01-O09-P02-O15
6	C	608	5ZL	P02-O09-P01-O11
6	C	608	5ZL	P01-O09-P02-O15
6	D	607	5ZL	P02-O09-P01-O11

Continued on next page...

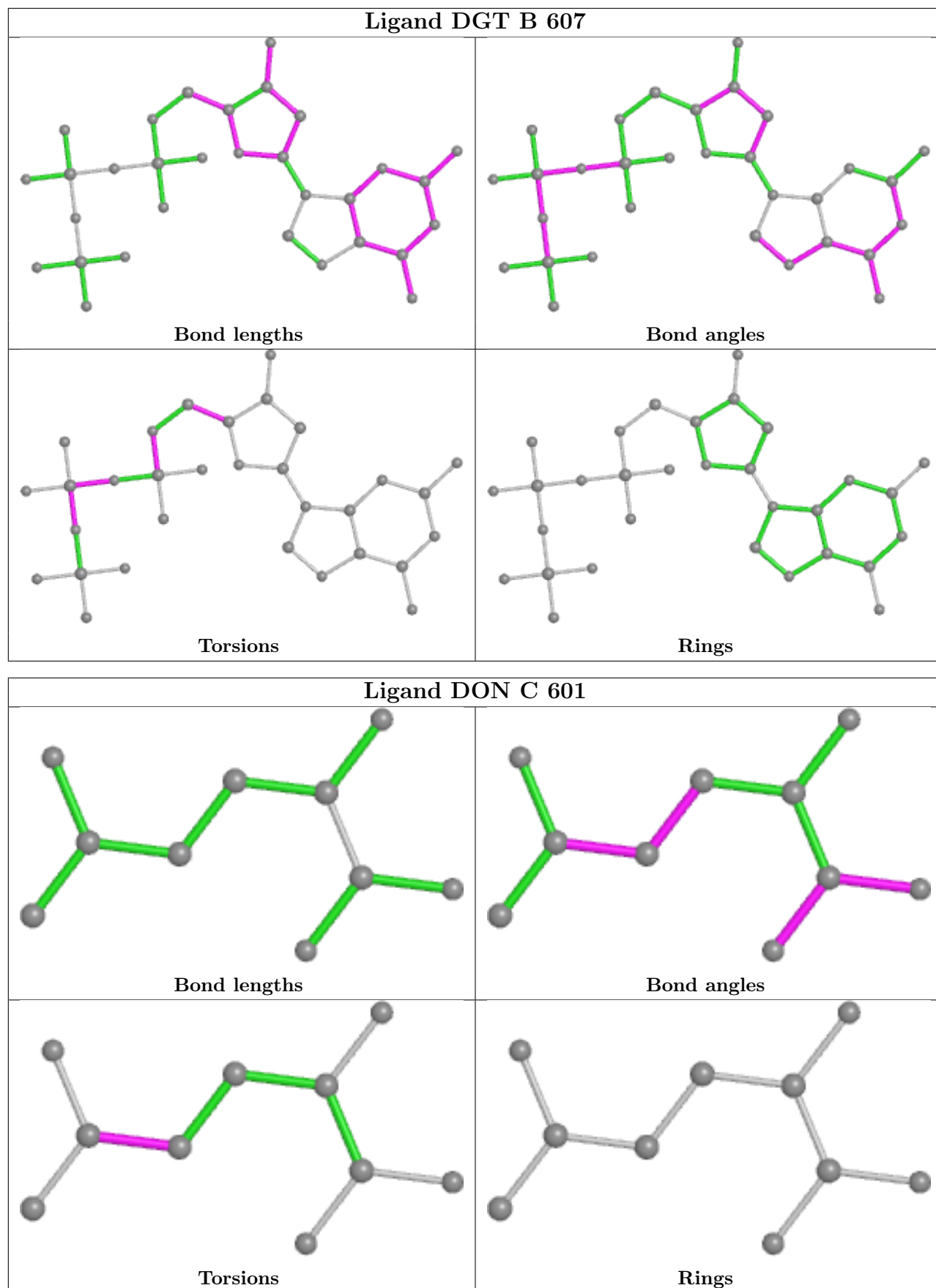
Continued from previous page...

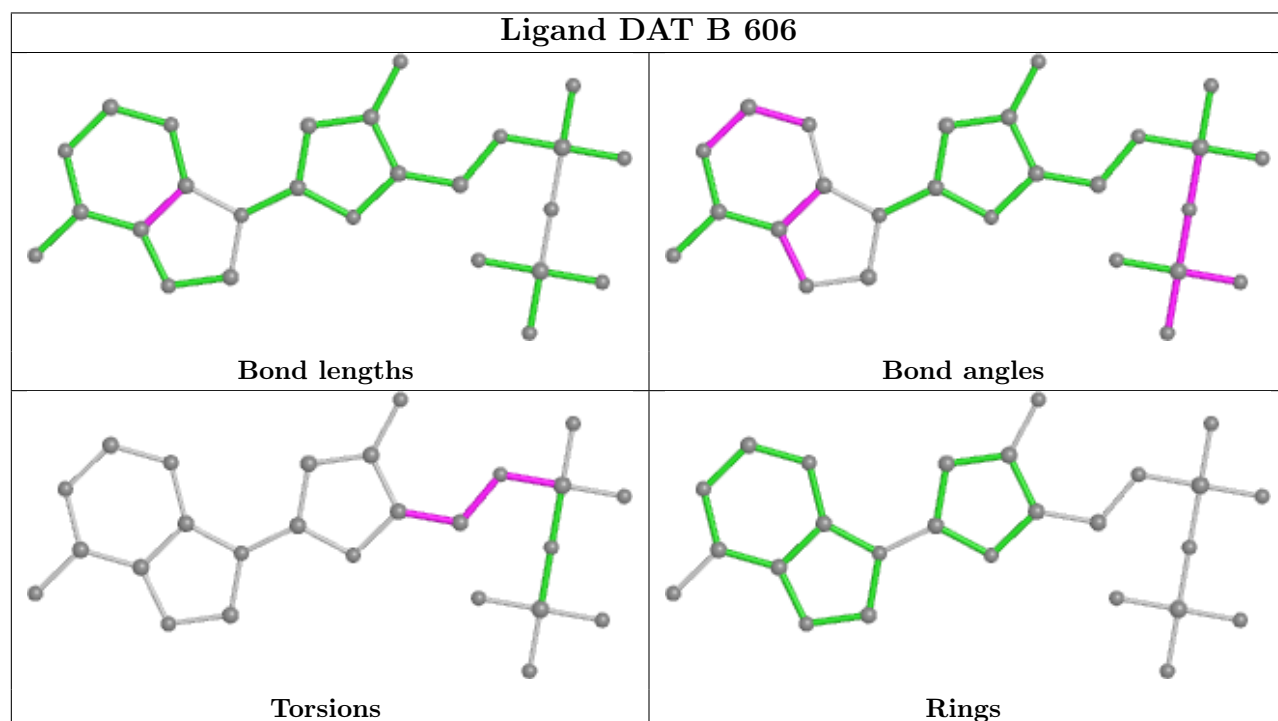
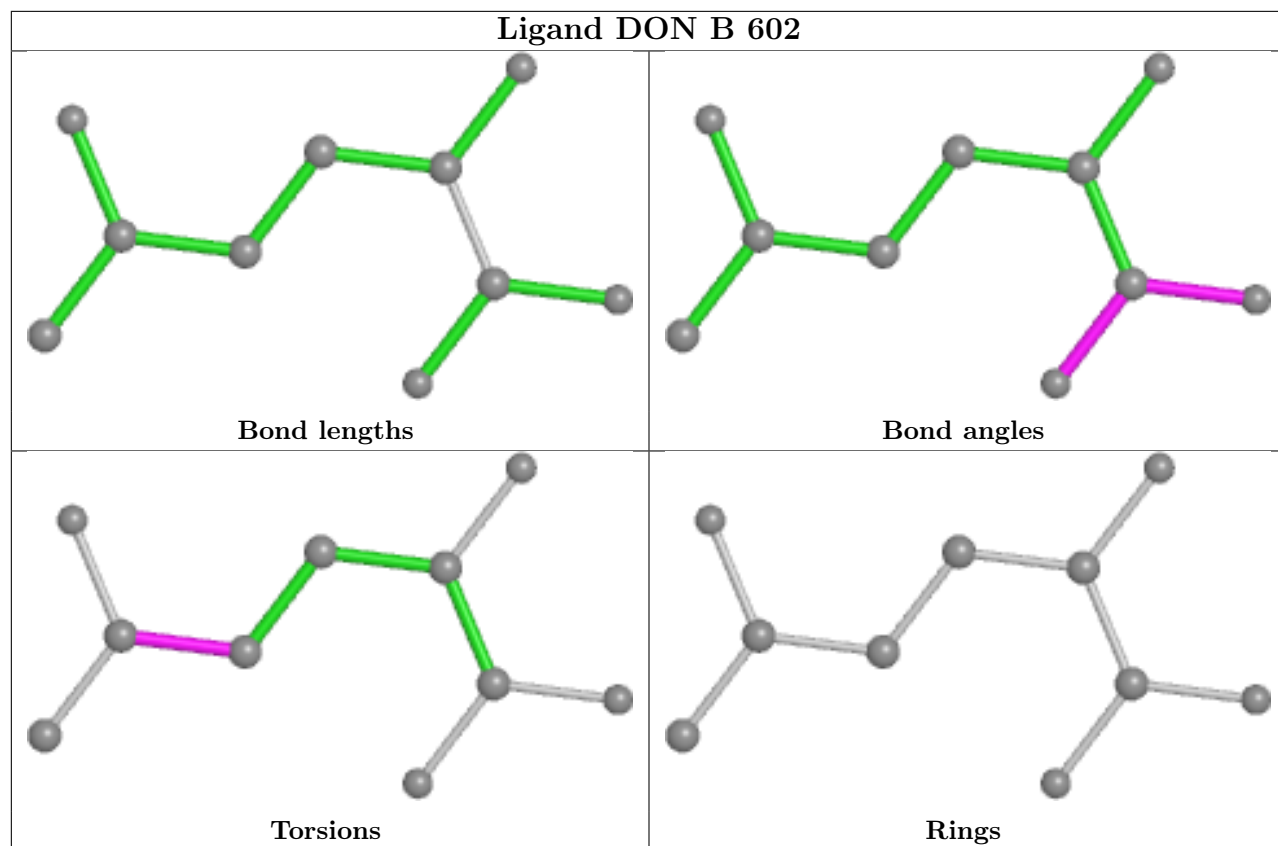
Mol	Chain	Res	Type	Atoms
6	D	607	5ZL	P01-O09-P02-O15
5	A	606	DGT	C5'-O5'-PA-O1A
5	B	607	DGT	C5'-O5'-PA-O1A
5	C	607	DGT	C5'-O5'-PA-O1A
5	D	606	DGT	C5'-O5'-PA-O1A
6	A	607	5ZL	P02-O13-P03-O16
6	B	608	5ZL	P02-O13-P03-O16
6	C	608	5ZL	P02-O13-P03-O16
6	D	607	5ZL	P02-O13-P03-O16

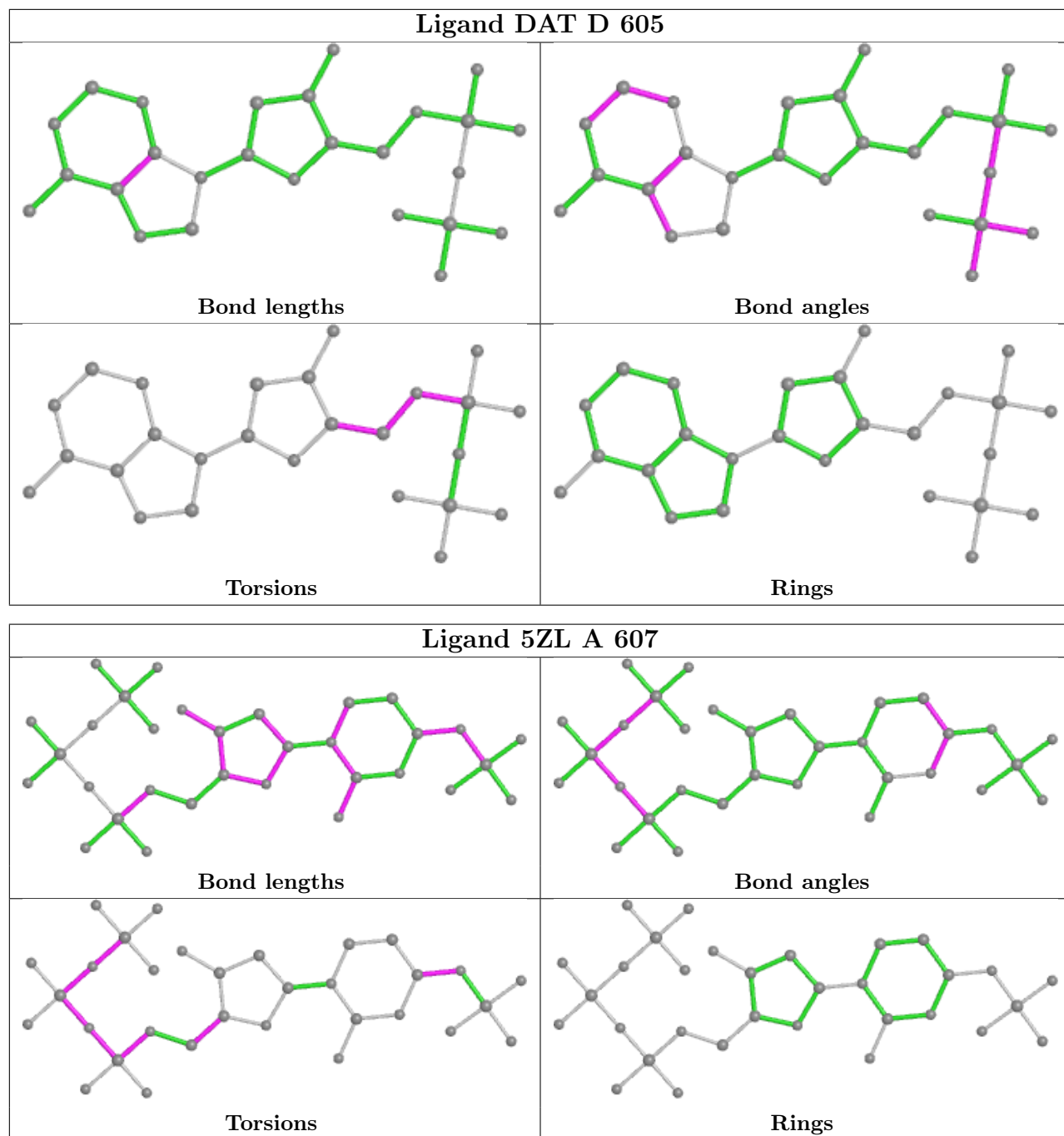
There are no ring outliers.

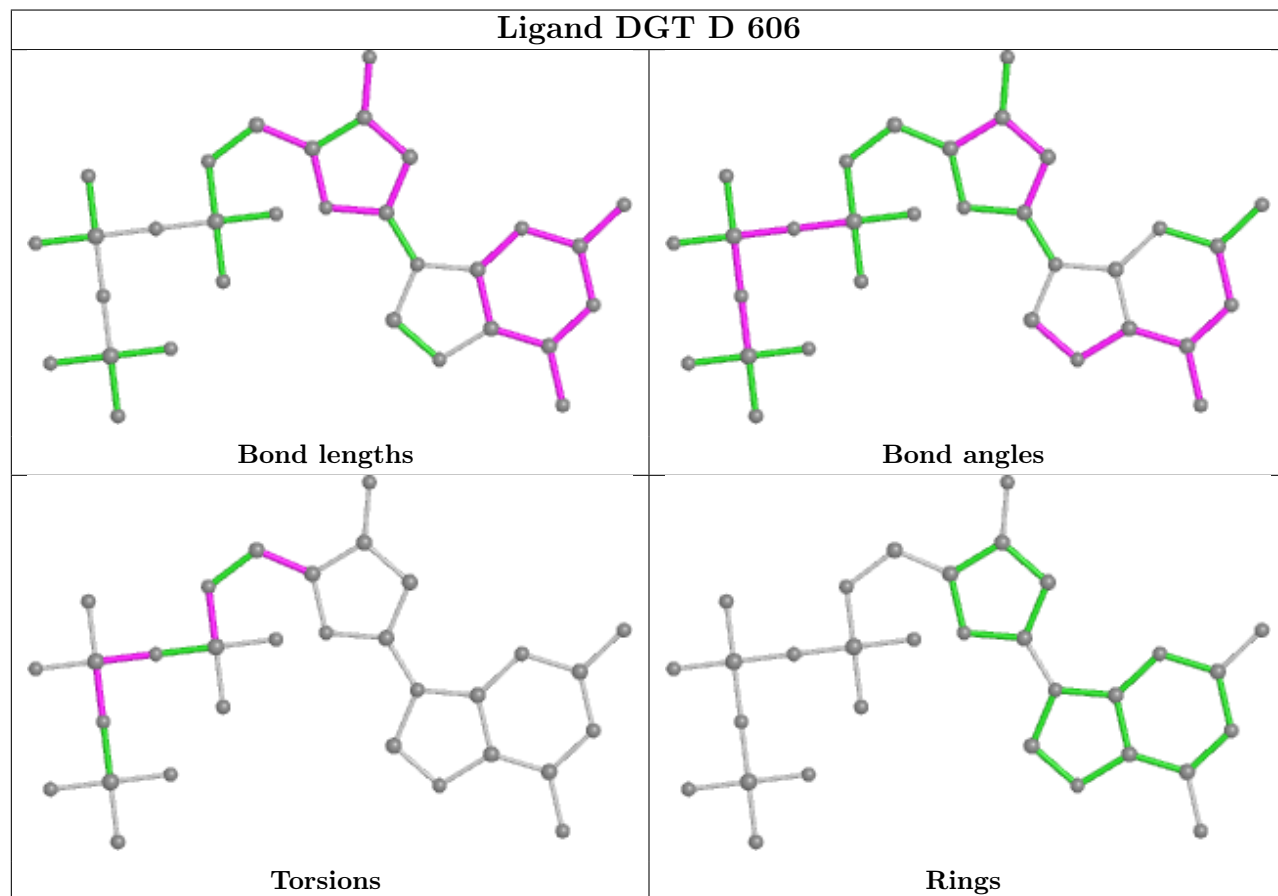
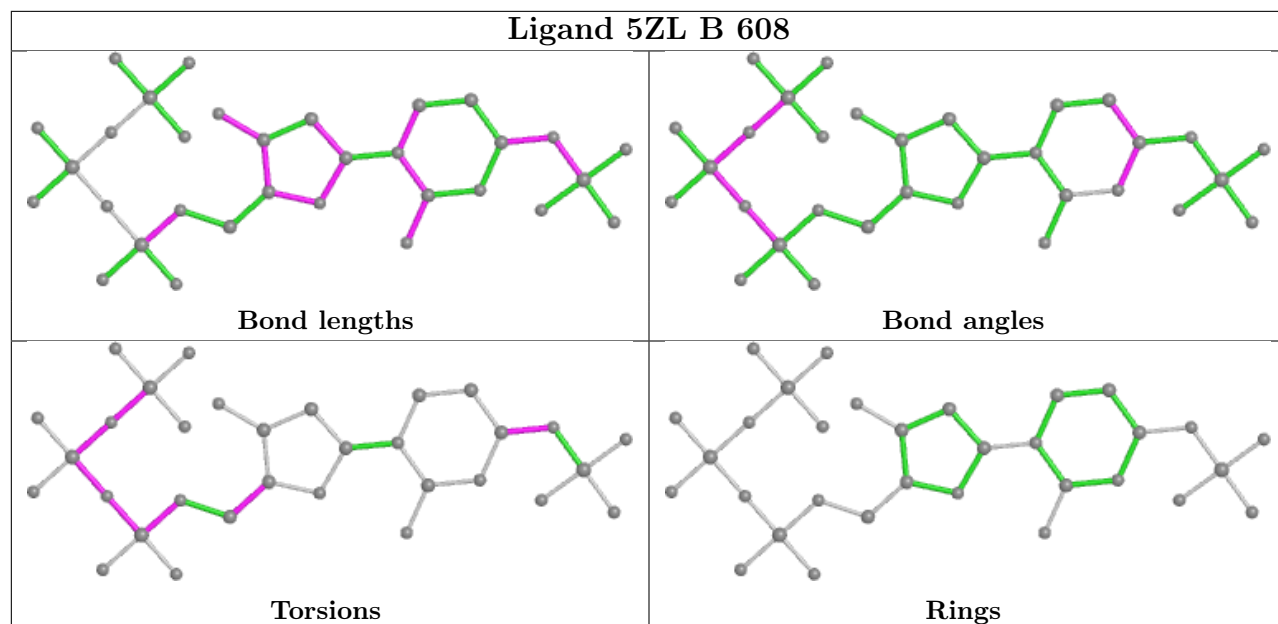
No monomer is involved in short contacts.

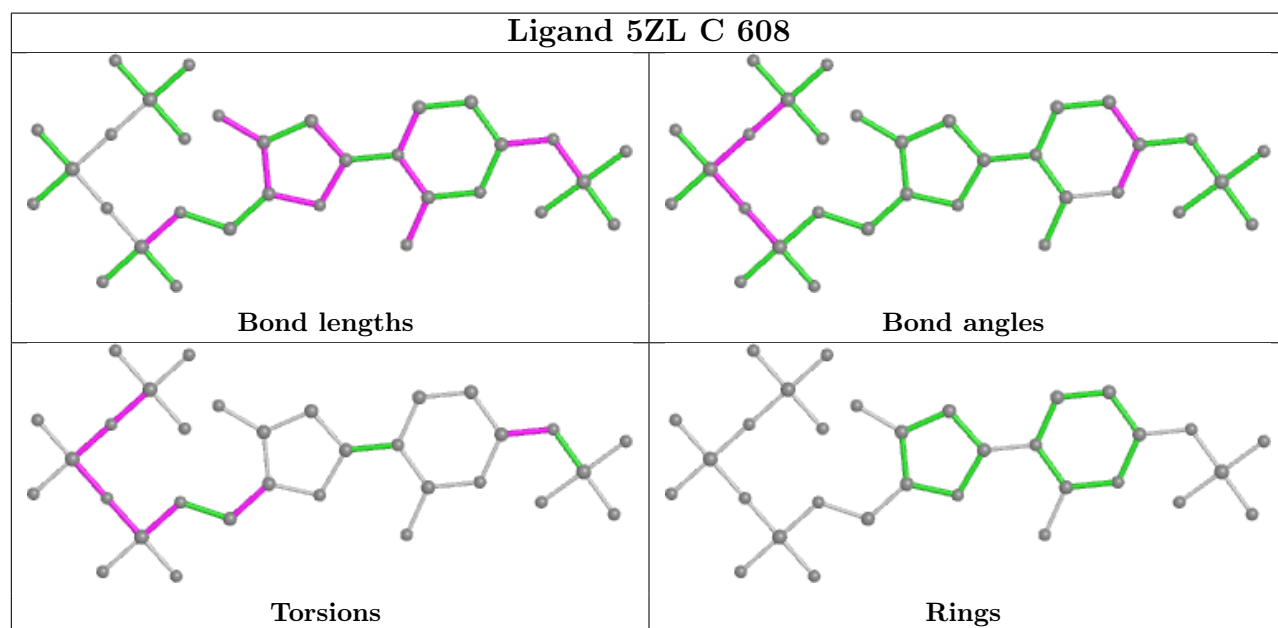
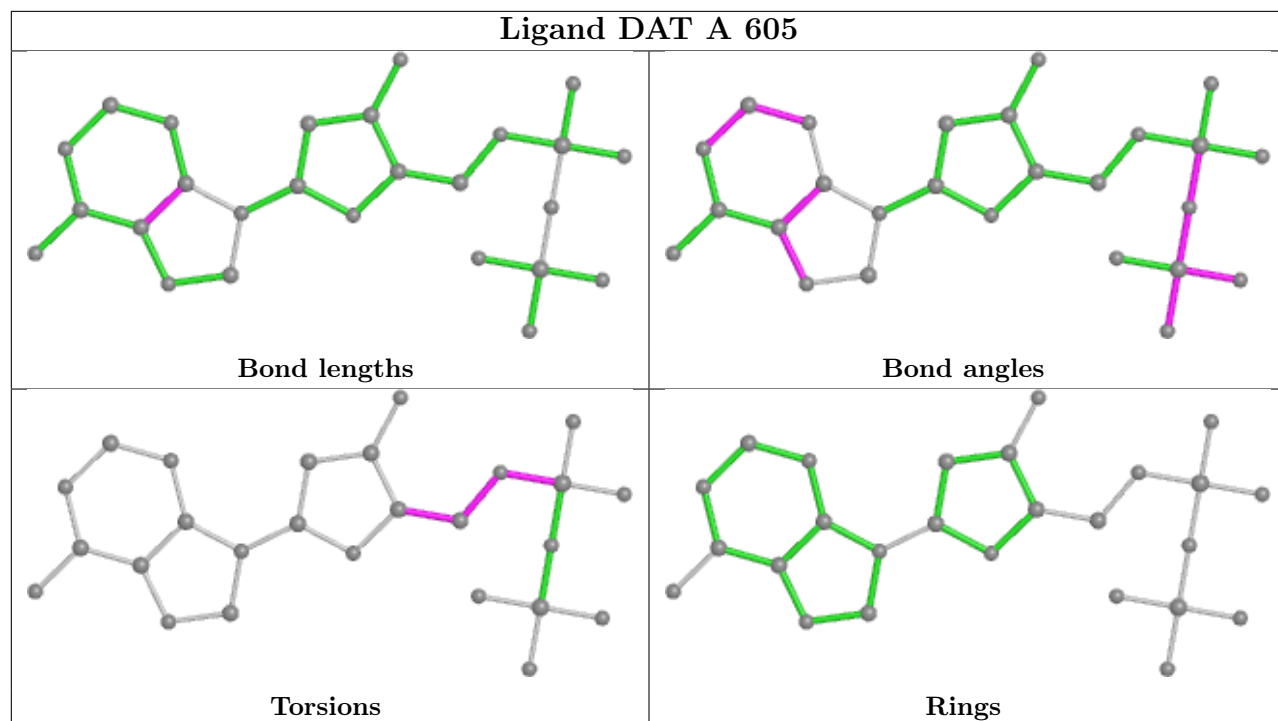
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

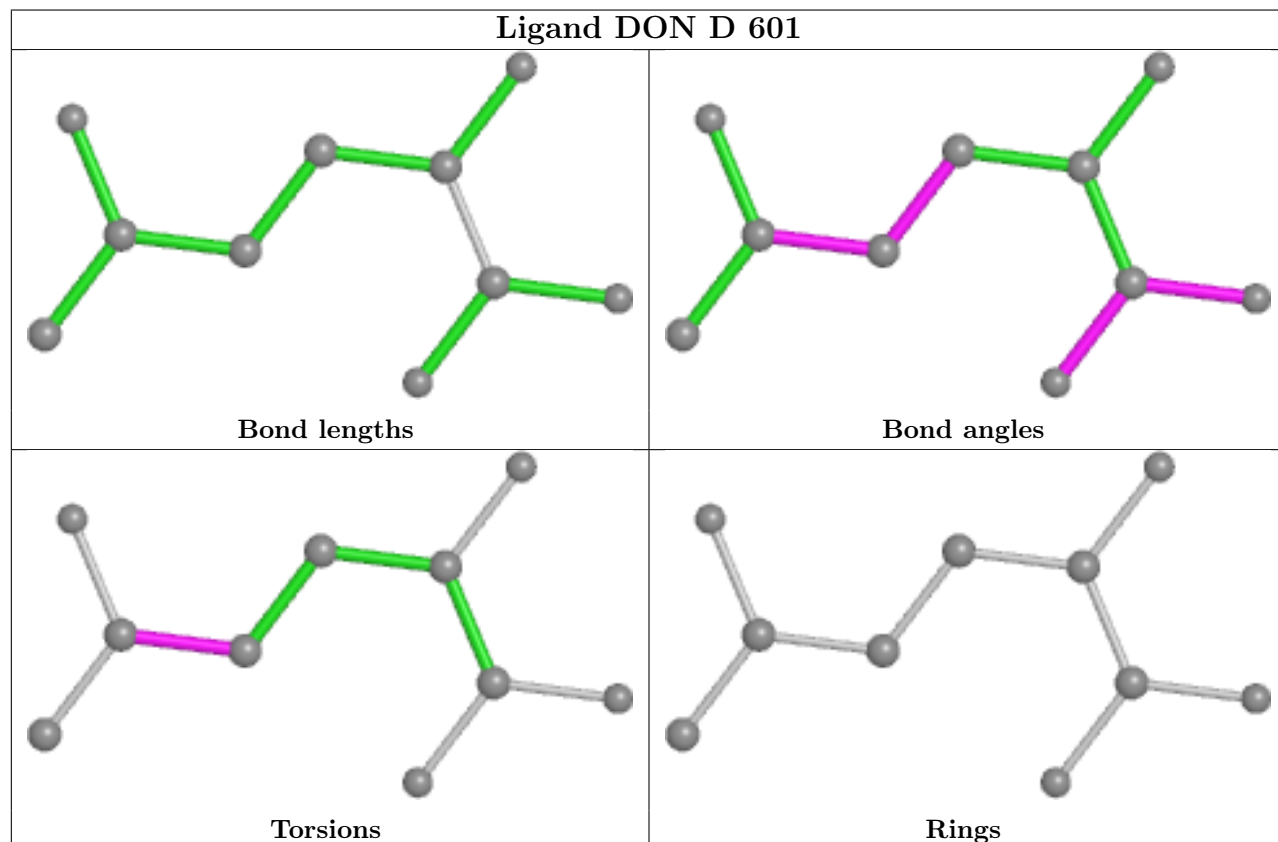
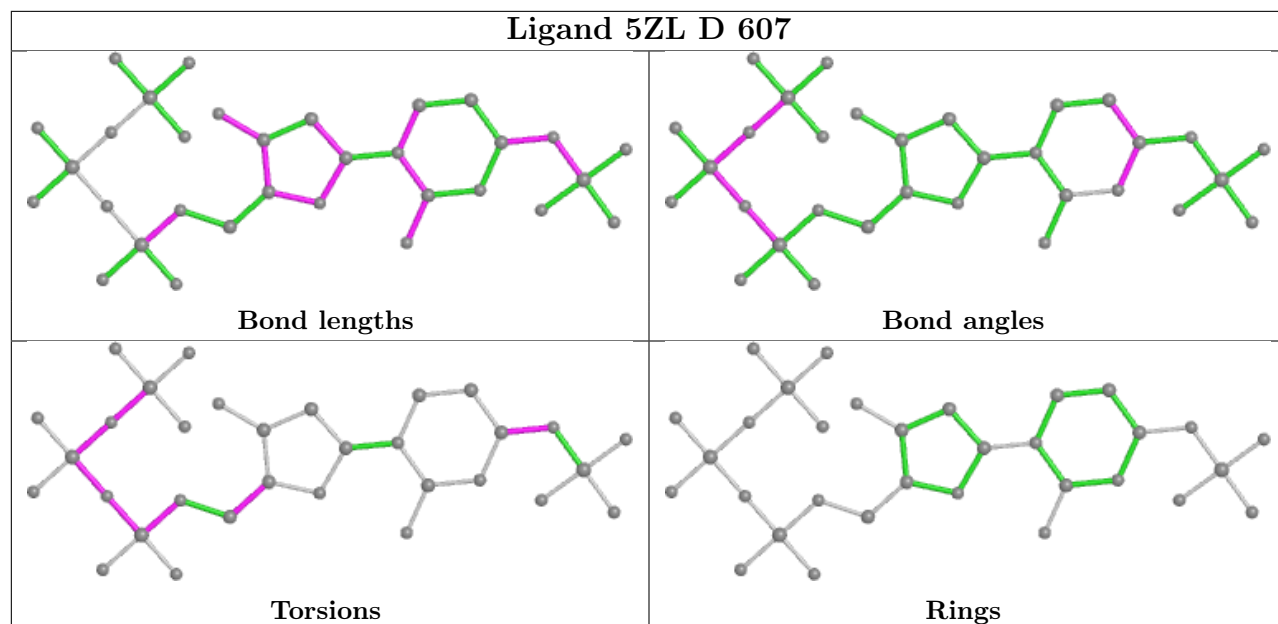


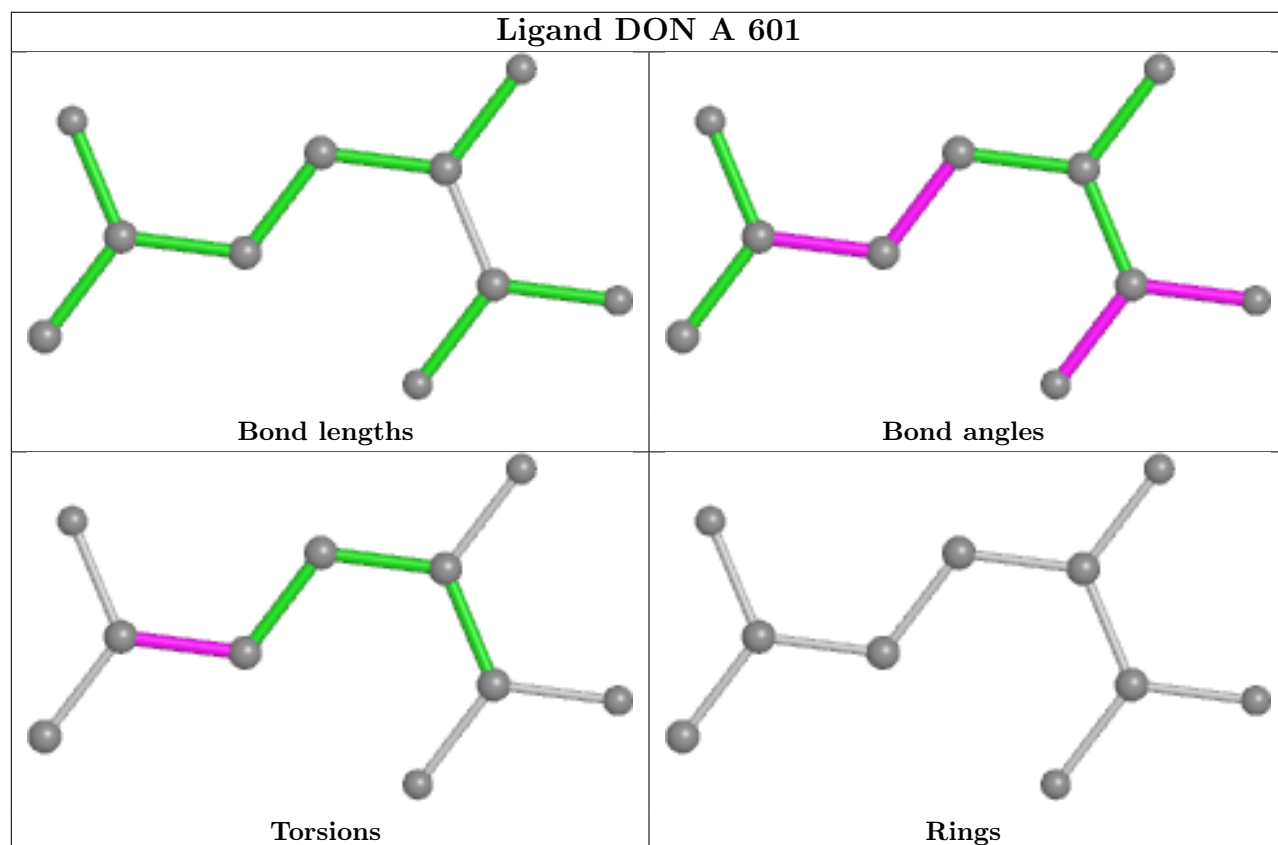
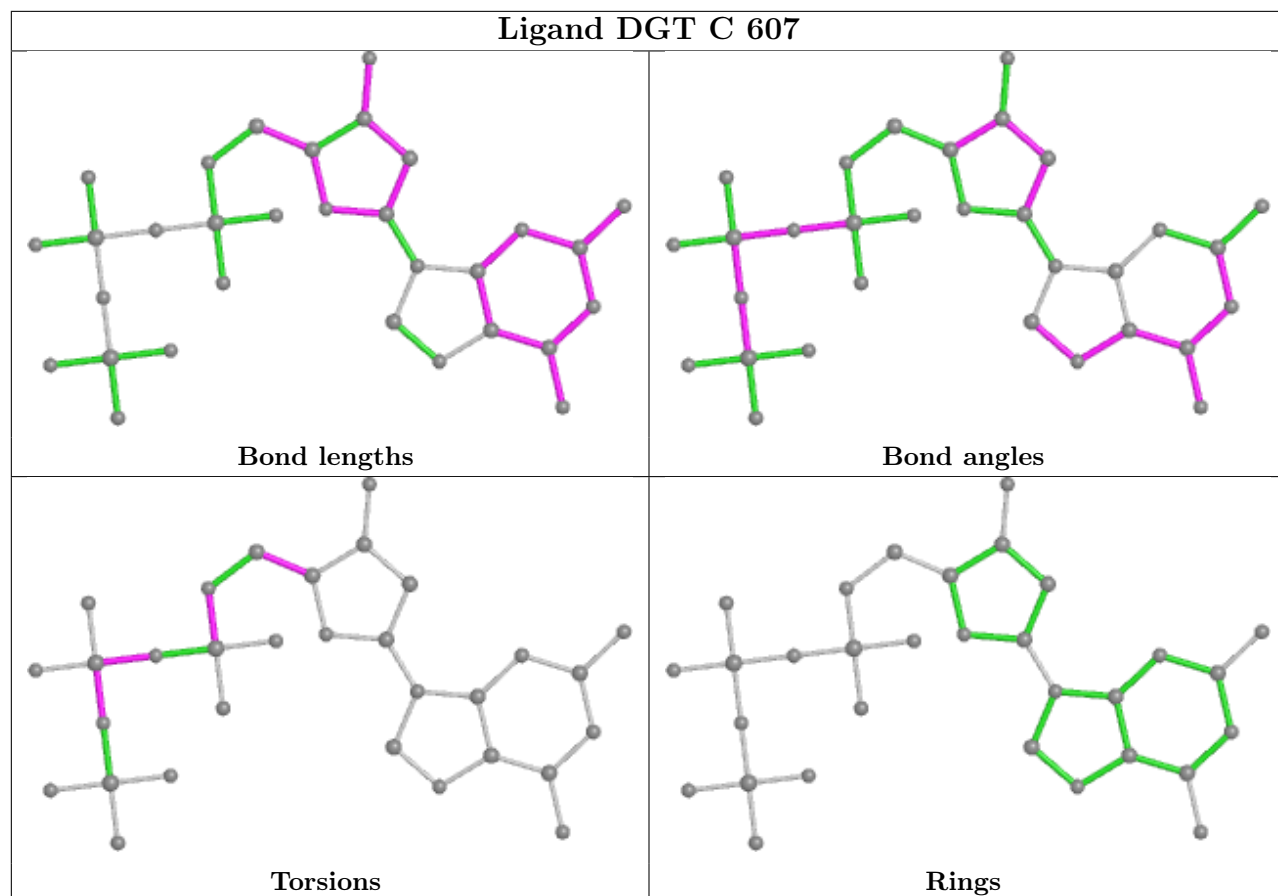


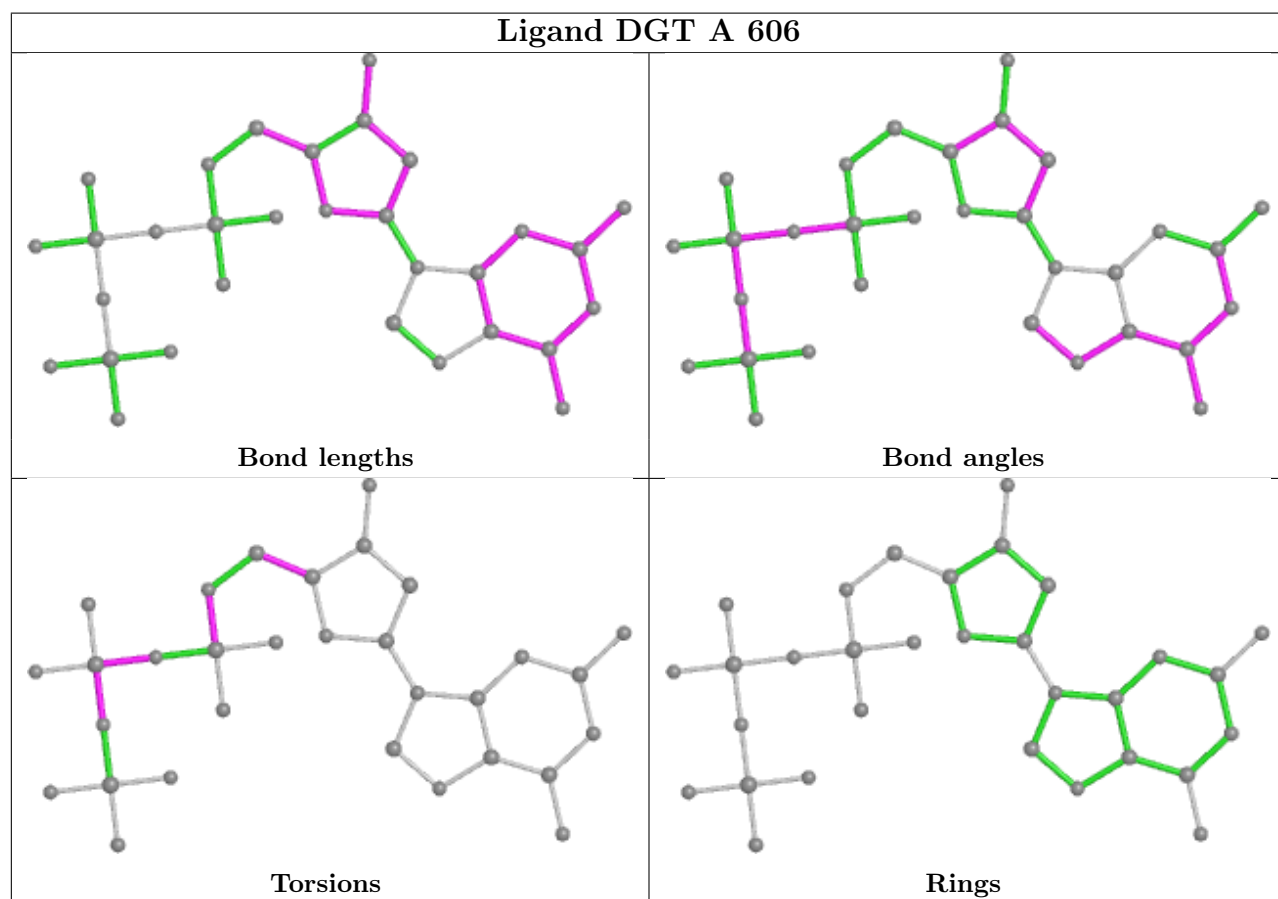
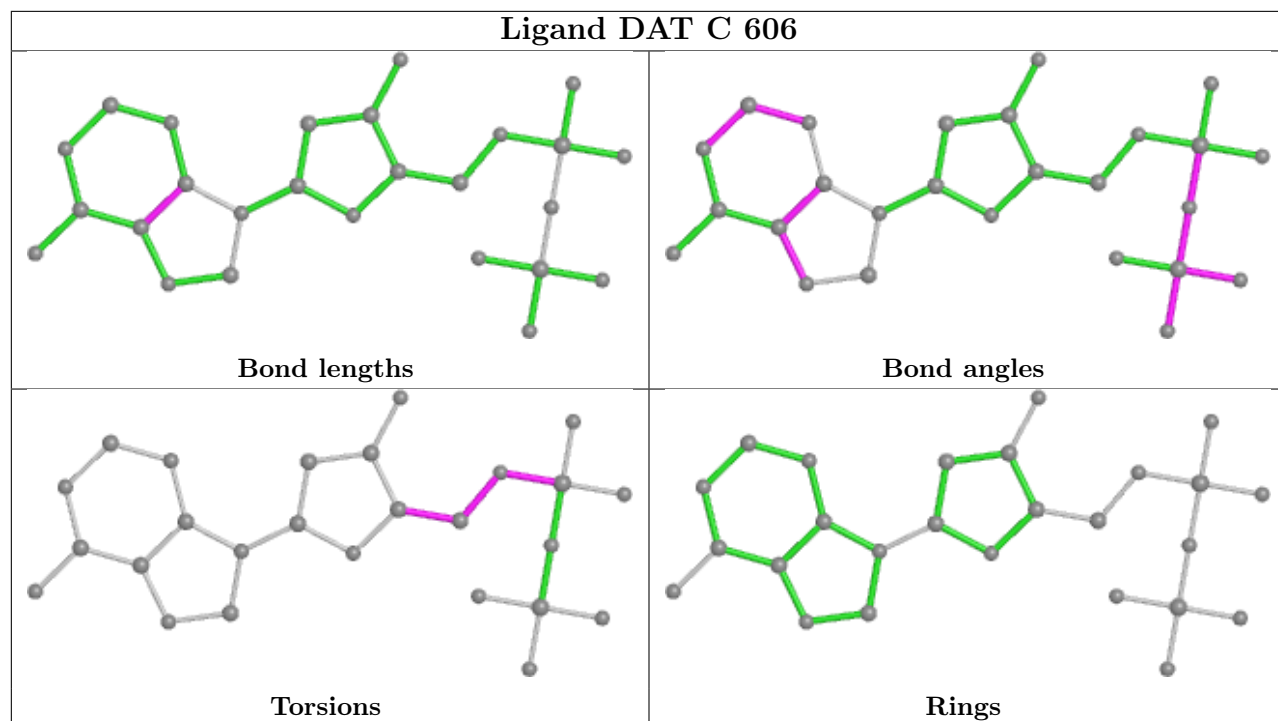












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

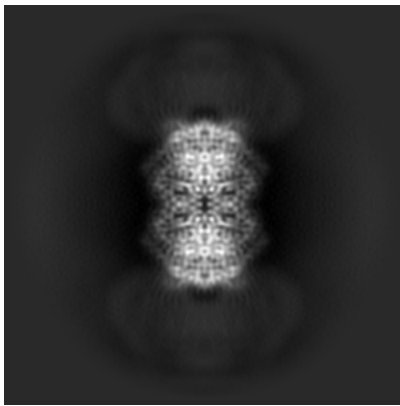
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61008. These allow visual inspection of the internal detail of the map and identification of artifacts.

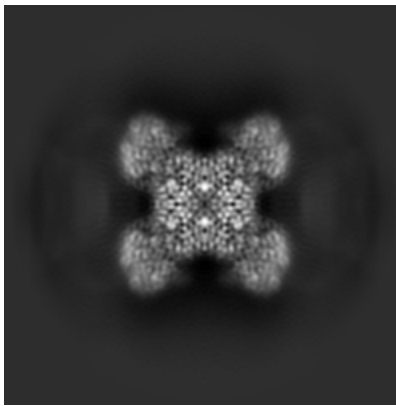
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

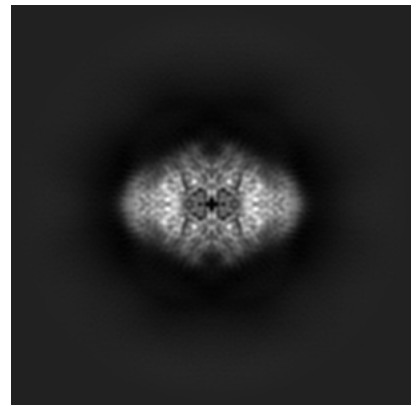
6.1.1 Primary map



X

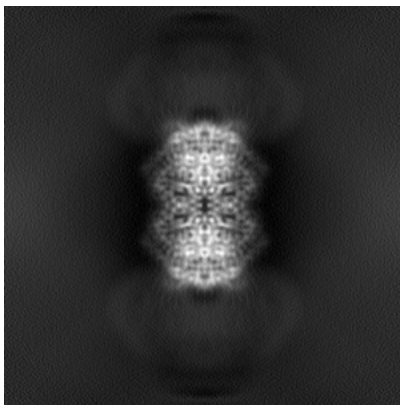


Y

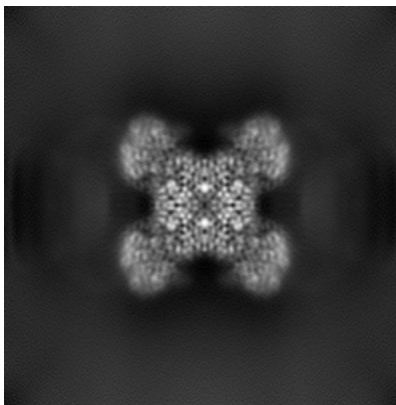


Z

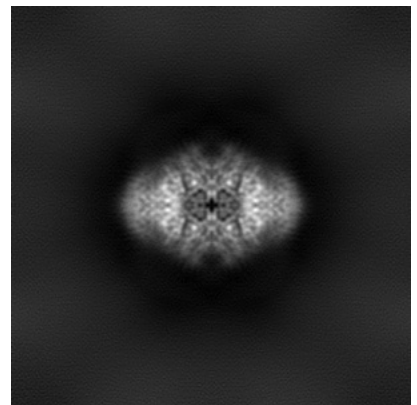
6.1.2 Raw map



X



Y

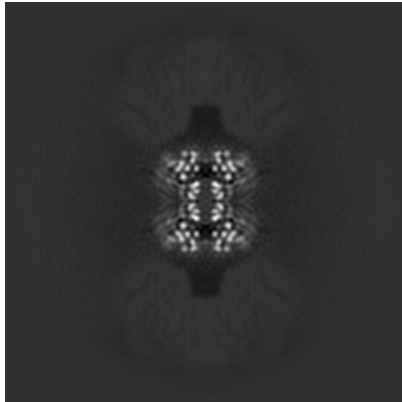


Z

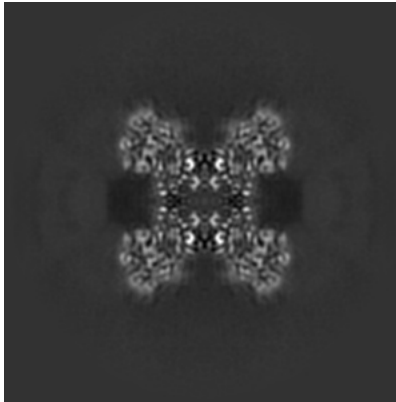
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

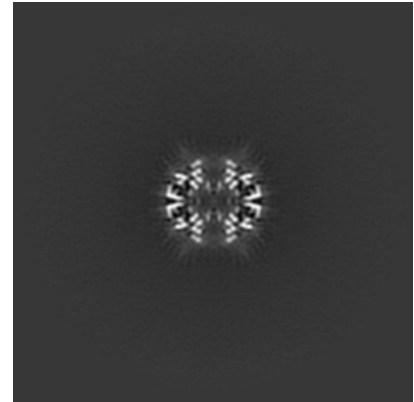
6.2.1 Primary map



X Index: 128

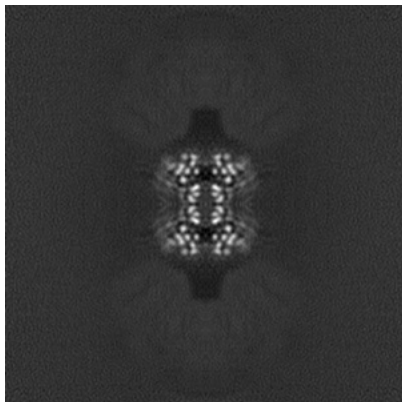


Y Index: 128

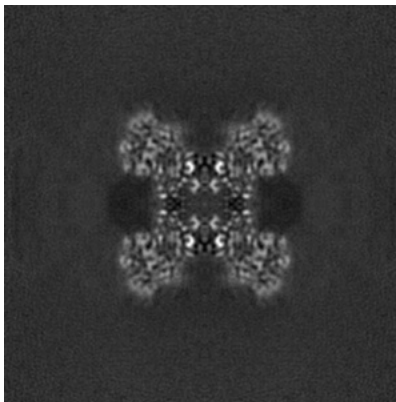


Z Index: 128

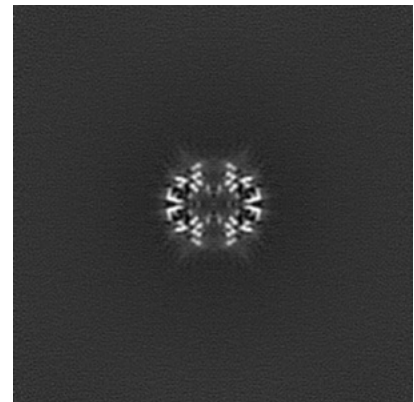
6.2.2 Raw map



X Index: 128



Y Index: 128

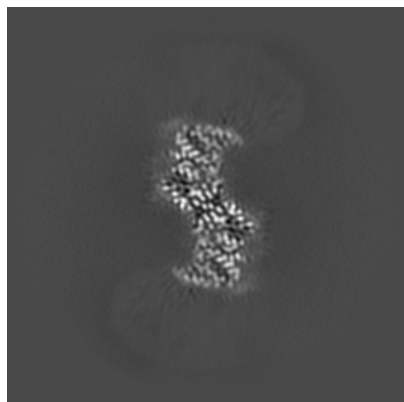


Z Index: 128

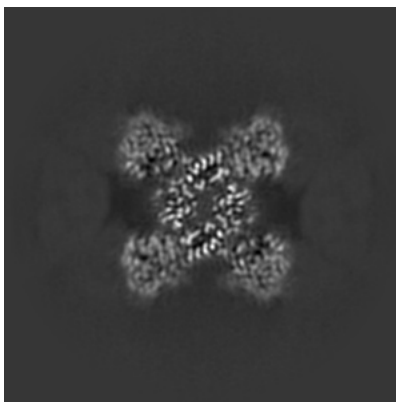
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

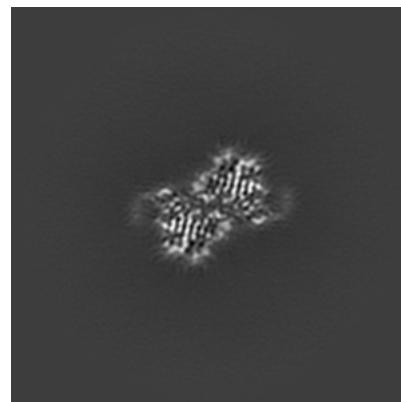
6.3.1 Primary map



X Index: 151

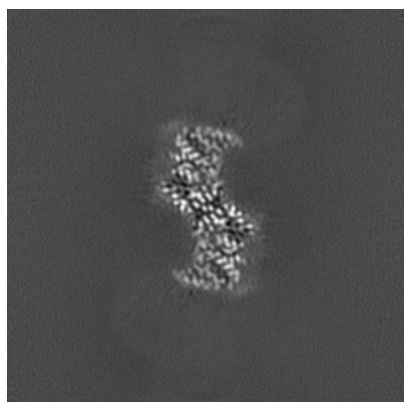


Y Index: 132

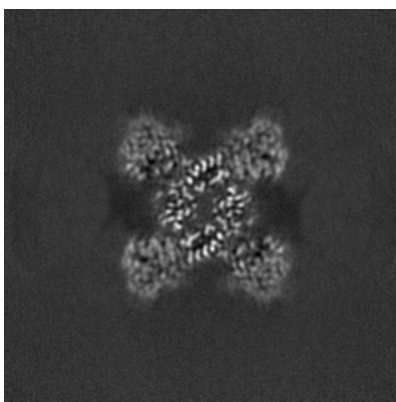


Z Index: 116

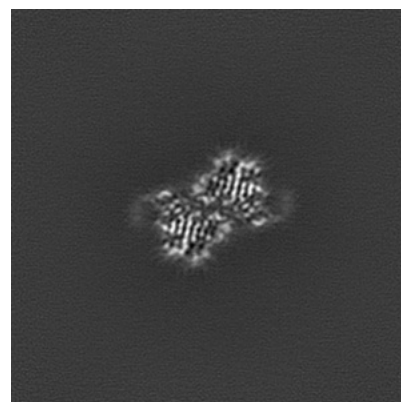
6.3.2 Raw map



X Index: 151



Y Index: 132

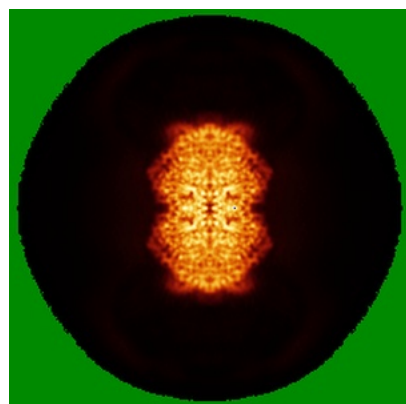


Z Index: 116

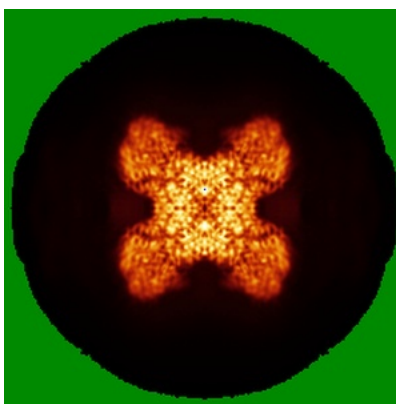
The images above show the largest variance slices of the map in three orthogonal directions.

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

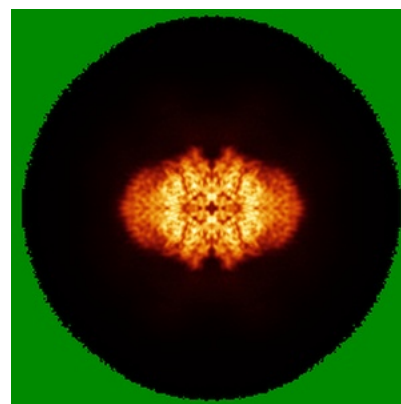
6.4.1 Primary map



X

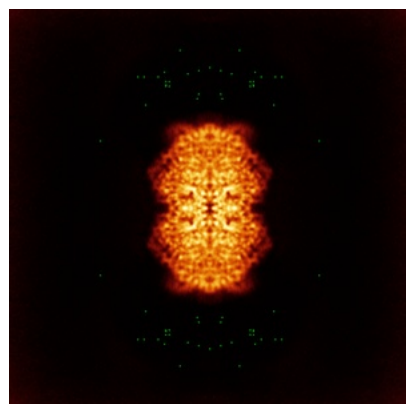


Y

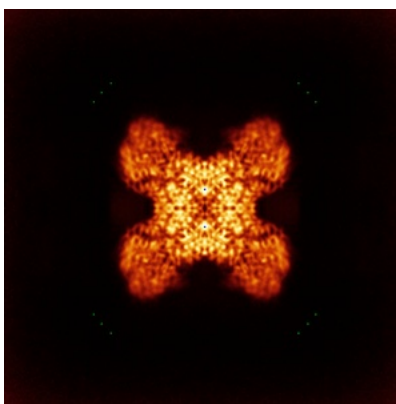


Z

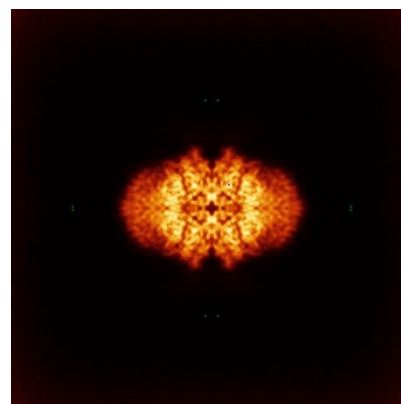
6.4.2 Raw map



X



Y

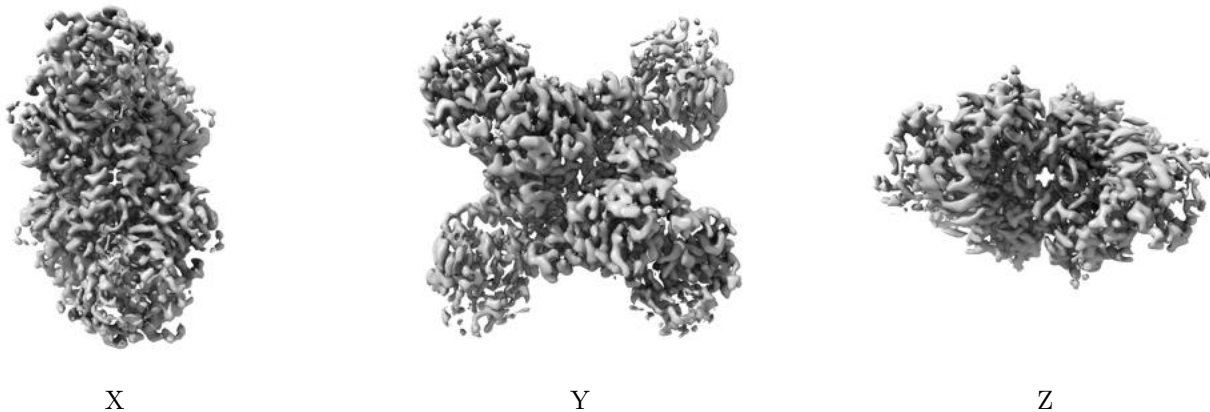


Z

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.

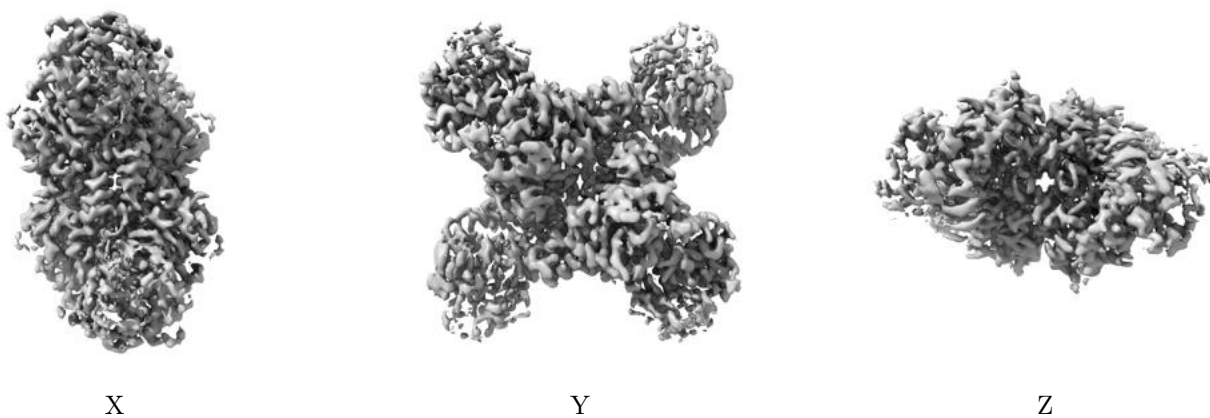
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.07. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

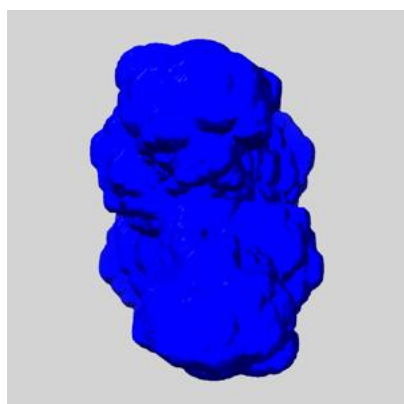
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

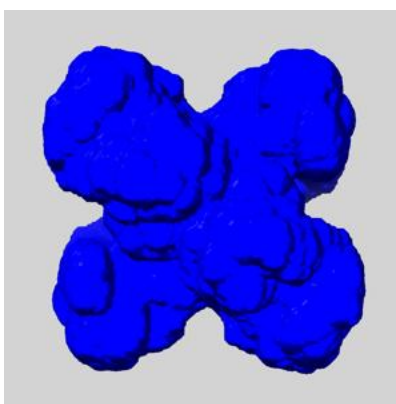
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

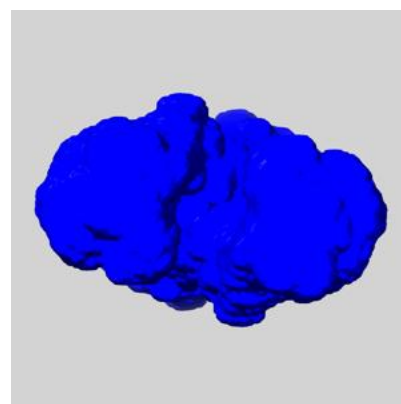
6.6.1 emd_61008_msk_1.map [i](#)



X



Y

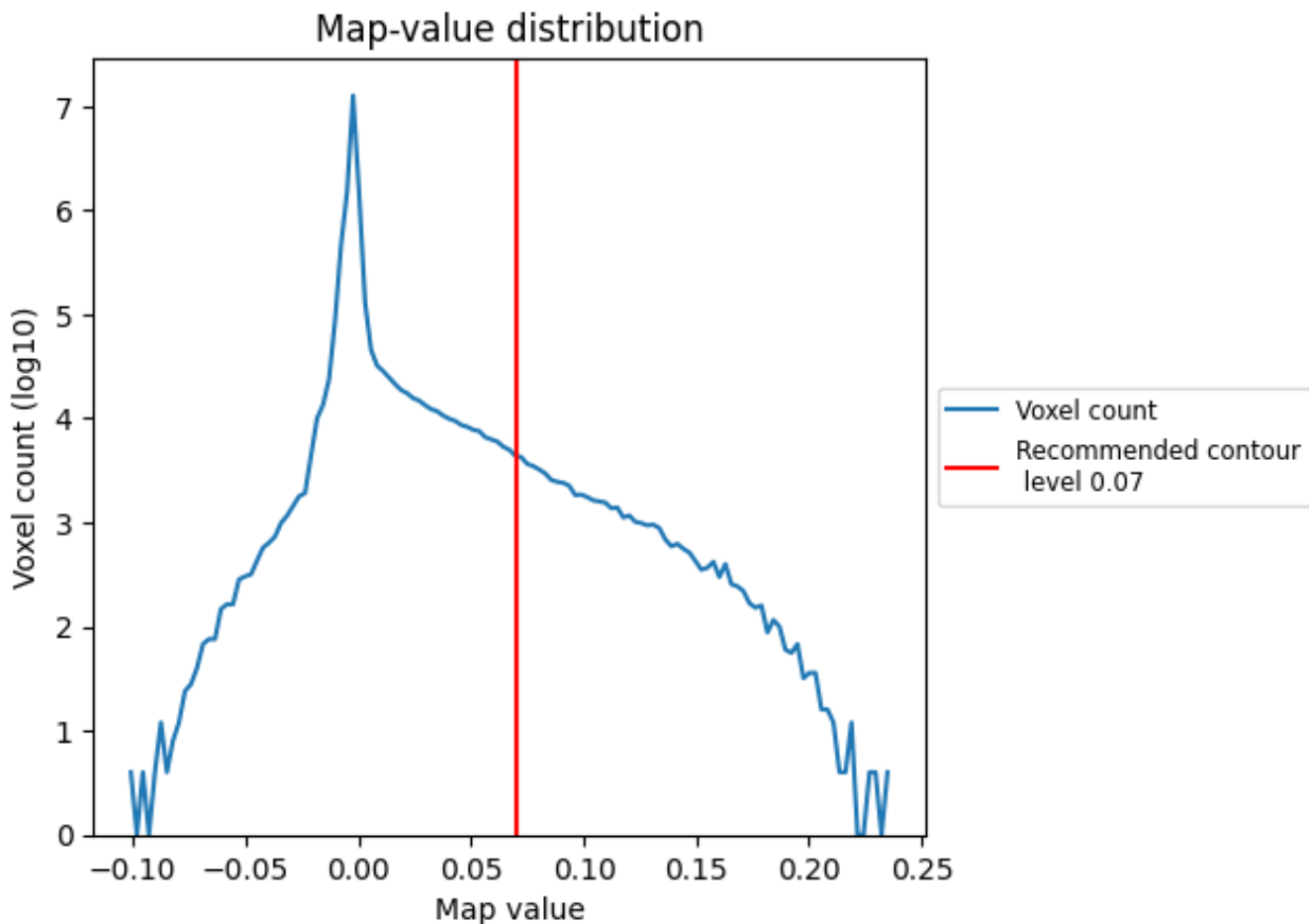


Z

7 Map analysis [i](#)

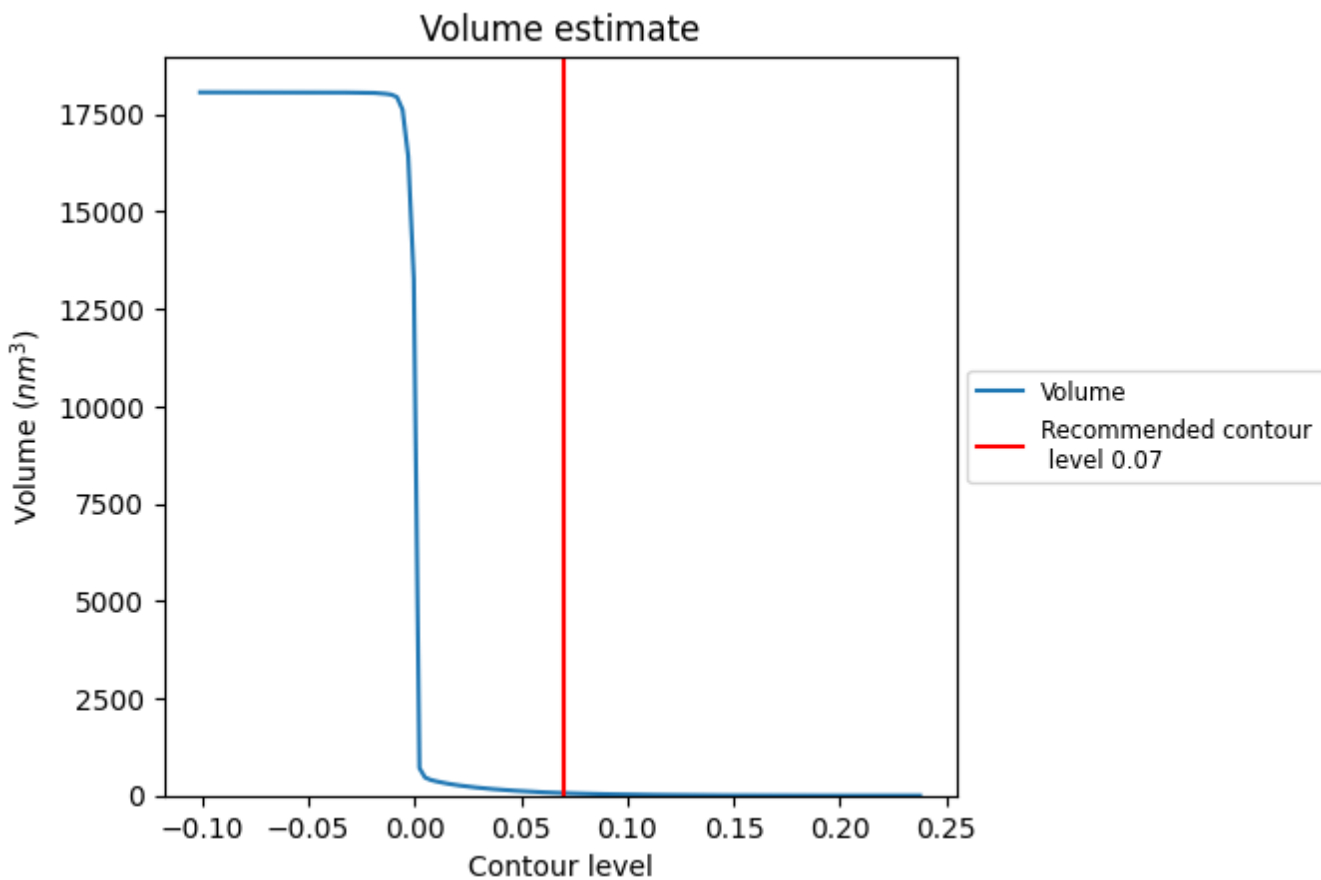
This section contains the results of statistical analysis of the map.

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.

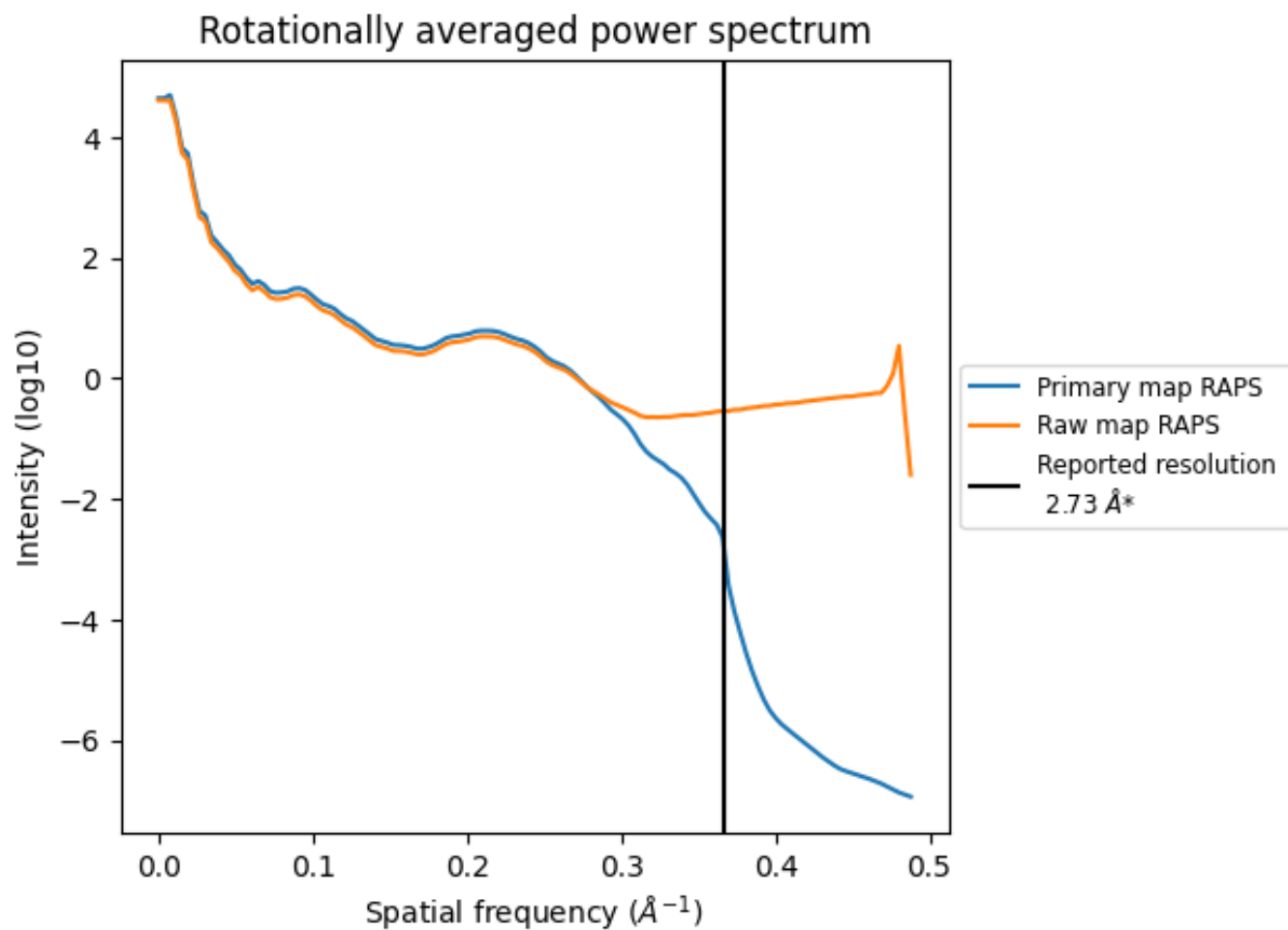
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 63 nm³; this corresponds to an approximate mass of 57 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 [i](#)

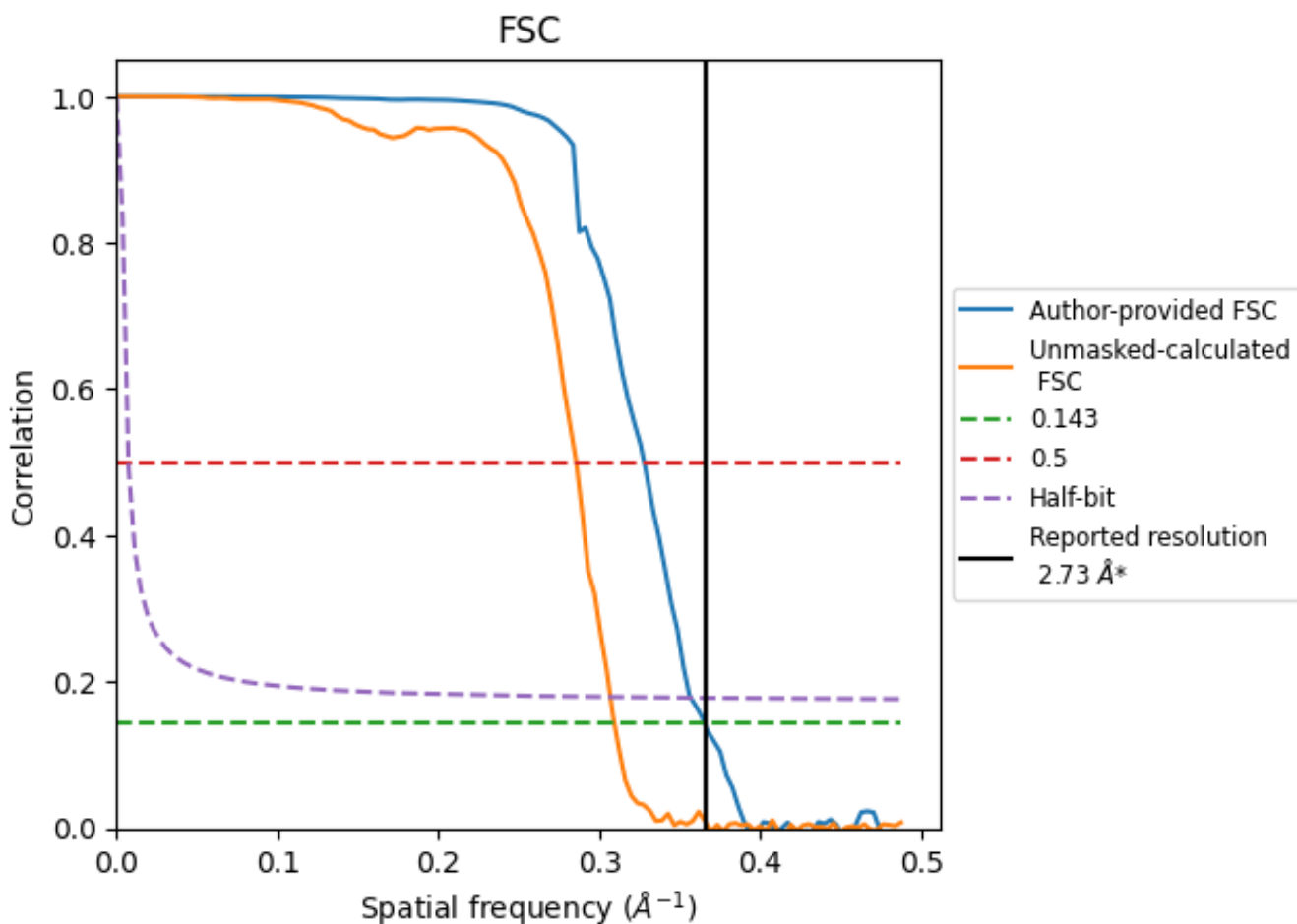


*Reported resolution corresponds to spatial frequency of 0.366 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.366 Å⁻¹

8.2 Resolution estimates [i](#)

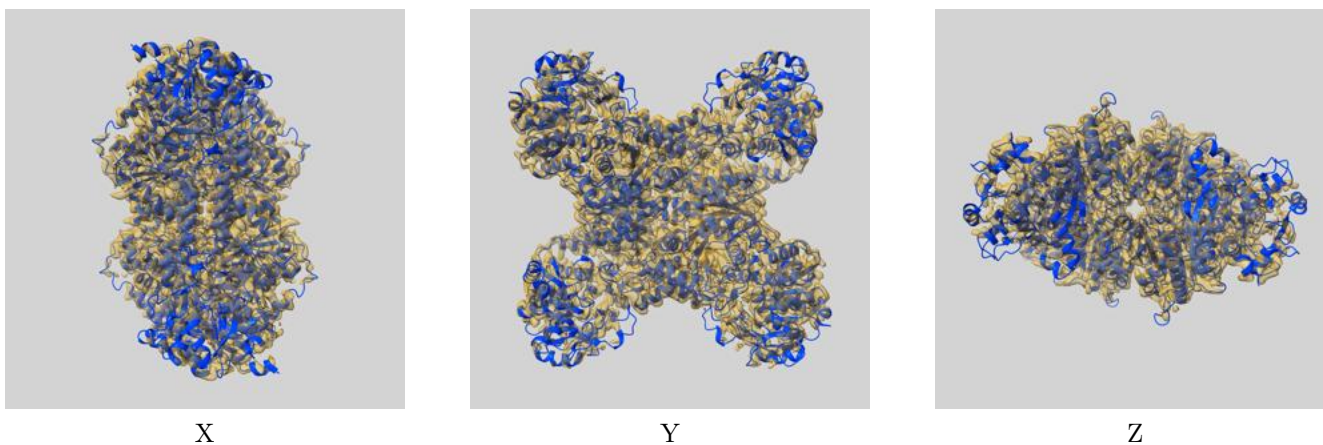
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.73	-	-
Author-provided FSC curve	2.74	3.05	2.80
Unmasked-calculated*	3.23	3.50	3.26

*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 3.23 differs from the reported value 2.73 by more than 10 %

9 Map-model fit [i](#)

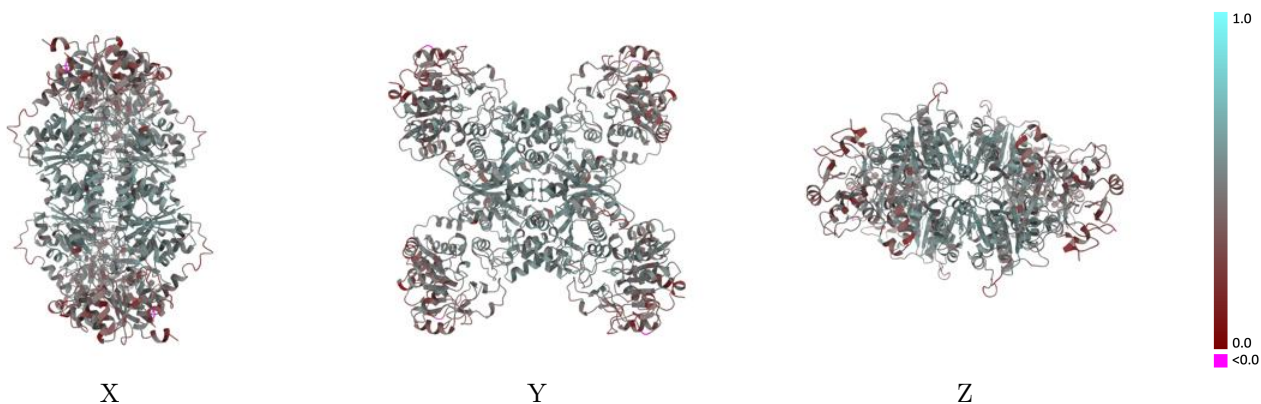
This section contains information regarding the fit between EMDB map EMD-61008 and PDB model 9IZ1. Per-residue inclusion information can be found in section 3 on page 7.

9.1 Map-model overlay [i](#)



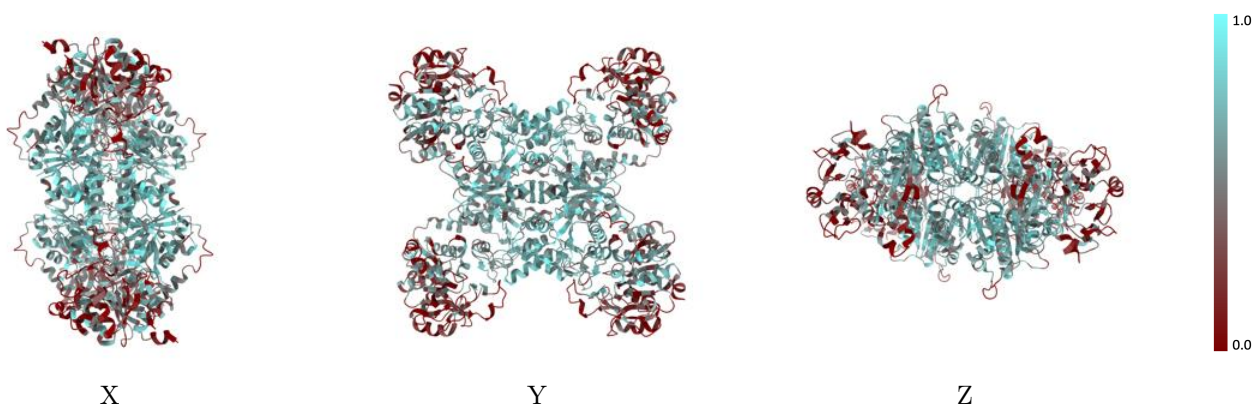
The images above show the 3D surface view of the map at the recommended contour level 0.07 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



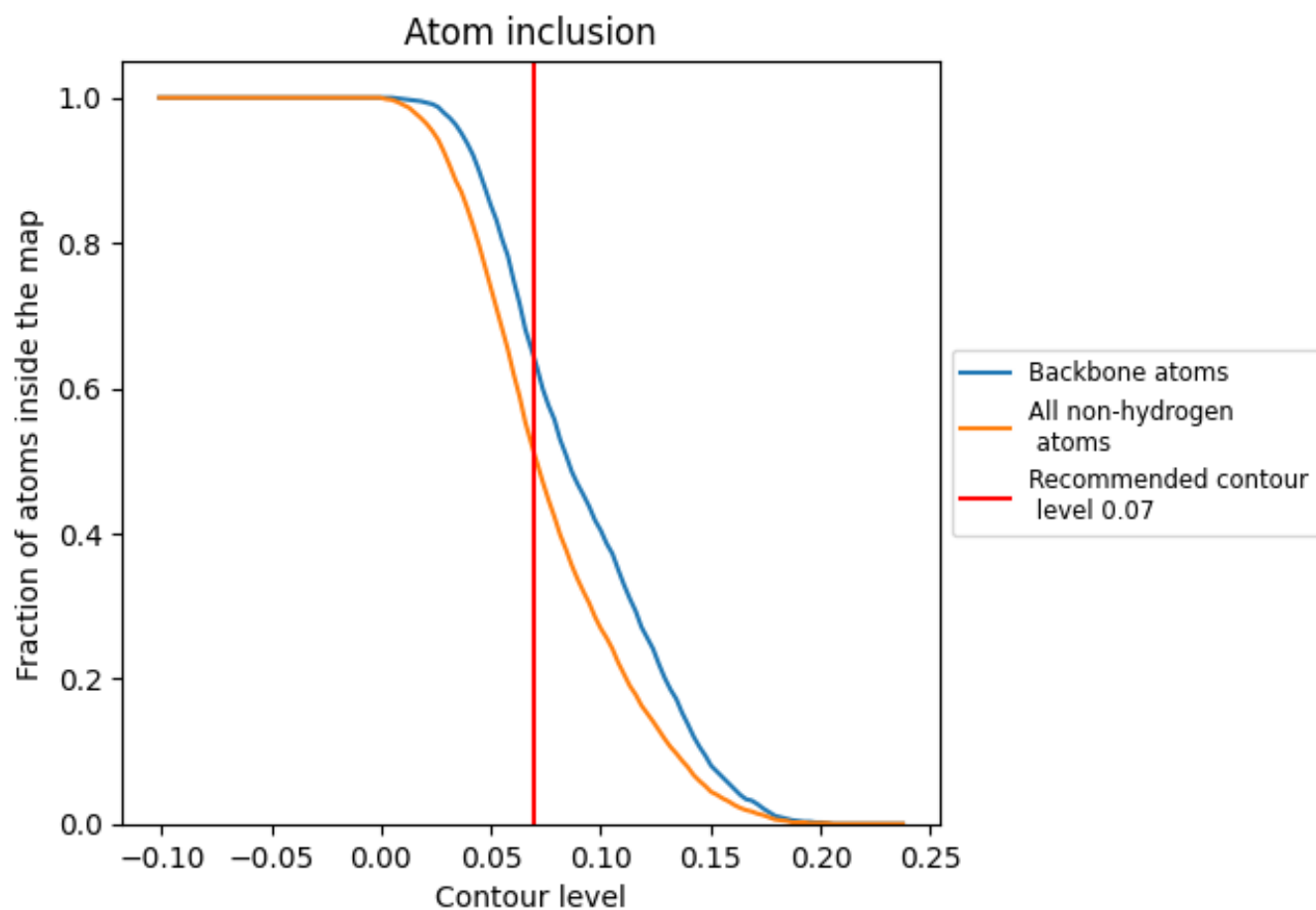
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.07).

9.4 Atom inclusion [i](#)



At the recommended contour level, 64% of all backbone atoms, 51% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.07) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	█ 0.5070	█ 0.4730
A	█ 0.5070	█ 0.4720
B	█ 0.5080	█ 0.4730
C	█ 0.5070	█ 0.4720
D	█ 0.5070	█ 0.4730

