



Full wwPDB EM Validation Report ⓘ

Mar 17, 2026 – 10:59 PM UTC

PDB ID : 8JYL / pdb_00008jyl
EMDB ID : EMD-36725
Title : Acyl-ACP Synthetase structure bound to C10-AMS
Authors : Huang, H.; Chang, S.; Huang, M.; Zhang, H.; Zhou, C.; Zhang, X.; Feng, Y.
Deposited on : 2023-07-03
Resolution : 2.33 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

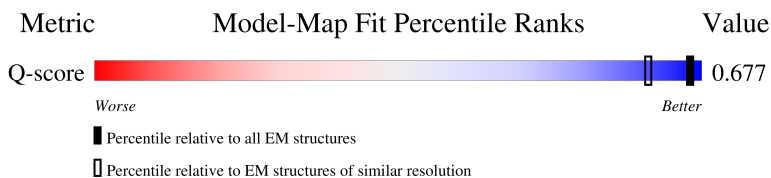
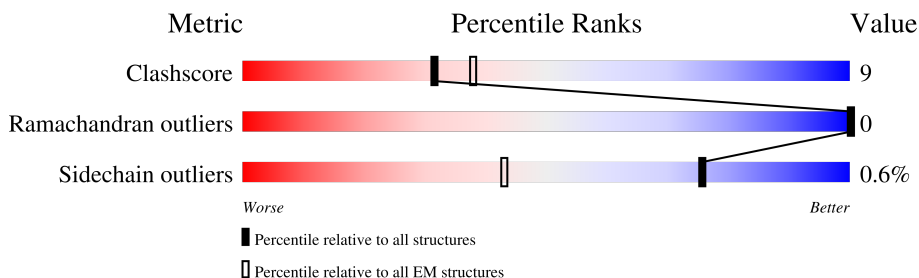
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.33 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	4434 (1.83 - 2.83)

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	532	5% 78% 21% .
1	B	532	6% 78% 21% .
1	C	532	5% 78% 21% .
1	D	532	5% 78% 21% .

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	532	5% 78% 21%
1	F	532	6% 78% 21%

2 Entry composition [i](#)

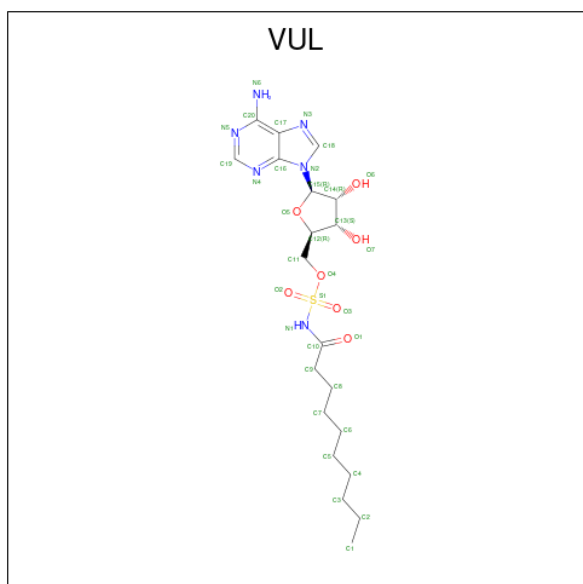
There are 3 unique types of molecules in this entry. The entry contains 25434 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Acyl-acyl carrier protein synthetase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	527	Total	C	N	O	S	0	0
			4204	2696	705	785	18		
1	B	527	Total	C	N	O	S	0	0
			4204	2696	705	785	18		
1	C	527	Total	C	N	O	S	0	0
			4204	2696	705	785	18		
1	D	527	Total	C	N	O	S	0	0
			4204	2696	705	785	18		
1	E	527	Total	C	N	O	S	0	0
			4204	2696	705	785	18		
1	F	527	Total	C	N	O	S	0	0
			4204	2696	705	785	18		

- Molecule 2 is [(2R,3S,4R,5R)-5-(6-aminopurin-9-yl)-3,4-bis(oxidanyl)oxolan-2-yl]methyl N-decanoylsulfamate (CCD ID: VUL) (formula: C₂₀H₃₂N₆O₇S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	S	0
			34	20	6	7	1	
2	B	1	Total	C	N	O	S	0
			34	20	6	7	1	
2	C	1	Total	C	N	O	S	0
			34	20	6	7	1	
2	D	1	Total	C	N	O	S	0
			34	20	6	7	1	
2	E	1	Total	C	N	O	S	0
			34	20	6	7	1	
2	F	1	Total	C	N	O	S	0
			34	20	6	7	1	

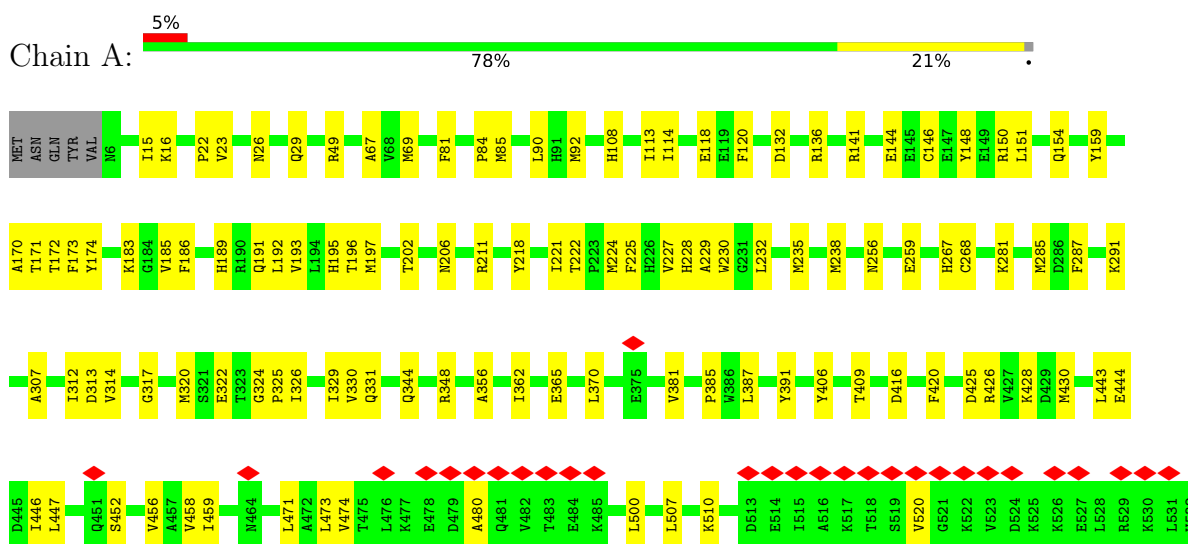
- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	Mg	0
			1	1	
3	B	1	Total	Mg	0
			1	1	
3	C	1	Total	Mg	0
			1	1	
3	D	1	Total	Mg	0
			1	1	
3	E	1	Total	Mg	0
			1	1	
3	F	1	Total	Mg	0
			1	1	

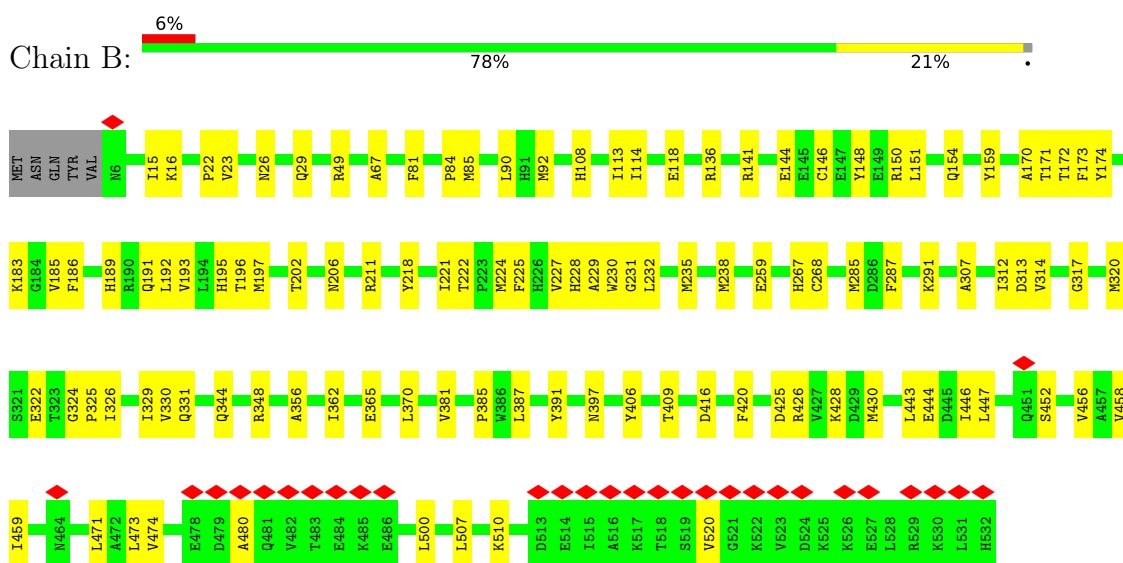
3 Residue-property plots

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.

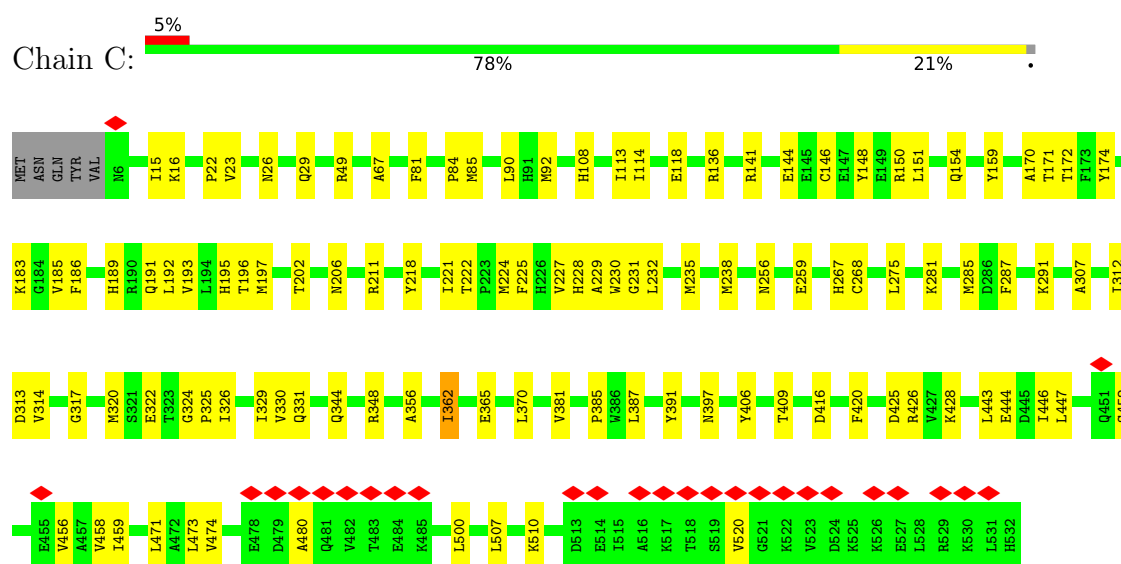
• Molecule 1: Acyl-acyl carrier protein synthetase



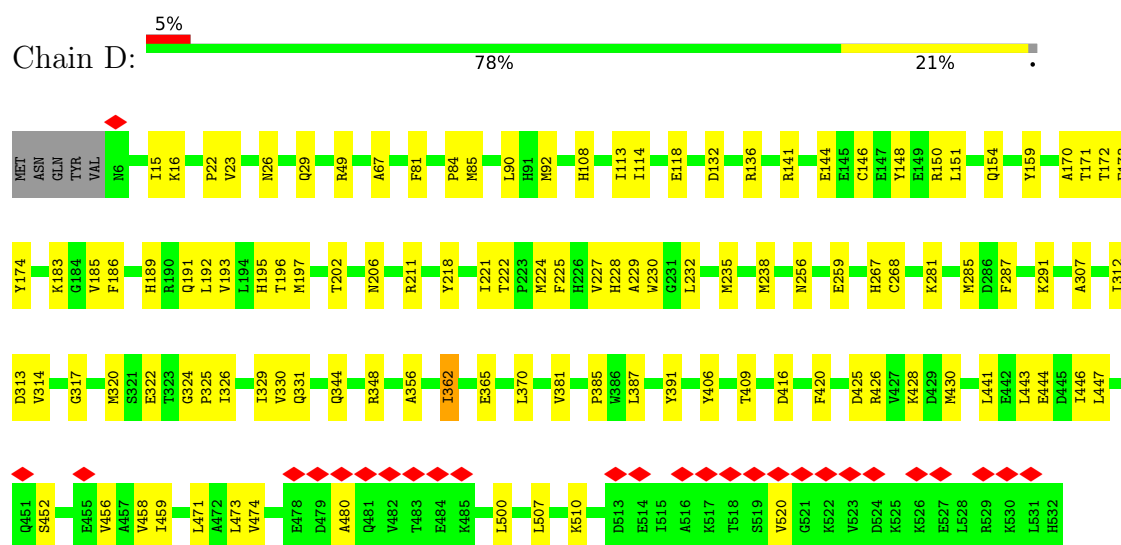
• Molecule 1: Acyl-acyl carrier protein synthetase



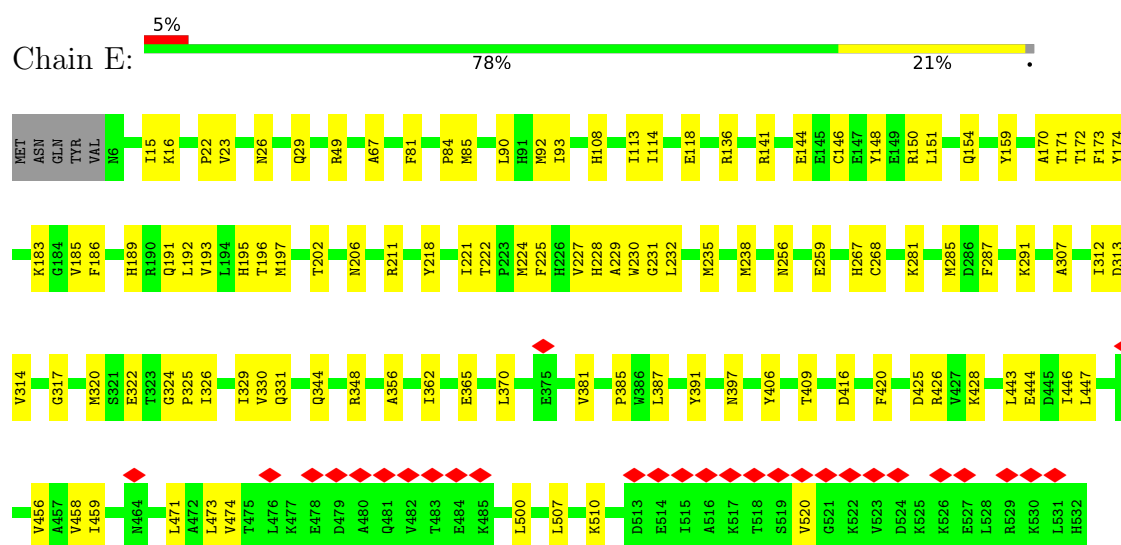
• Molecule 1: Acyl-acyl carrier protein synthetase



• Molecule 1: Acyl-acyl carrier protein synthetase



• Molecule 1: Acyl-acyl carrier protein synthetase



Chain F:

6%

78%

21%

•

Value	Category
MET	Green
ASN	Green
GLN	Green
TYR	Green
VAL	Green
N6	Green
I15	Green
K16	Green
P22	Green
V23	Green
N26	Green
Q29	Green
R49	Green
A67	Green
F81	Green
P84	Green
M85	Green
L90	Green
H91	Green
M92	Green
H108	Green
I113	Green
I114	Green
E118	Green
D132	Green
R136	Green
R141	Green
E144	Green
E145	Green
C146	Green
E147	Green
Y148	Green
E149	Green
R150	Green
L151	Green
Q154	Green
Y159	Green
A170	Green
T171	Green
T172	Green
F173	Green
Y174	Green
K183	Green
G184	Green
V185	Green
F186	Green
H189	Green
R190	Green
Q191	Green
L192	Green
V193	Green
L194	Green
H195	Green
T196	Green
M197	Green
T202	Green
N206	Green
R211	Green
Y218	Green
I221	Green
T222	Green
M224	Green
F225	Green
H226	Green
V227	Green
H228	Green
A229	Green
W230	Green
G231	Green
L232	Green
M235	Green
M238	Green
N256	Green
E259	Green
H267	Green
C268	Green
K281	Green
M285	Green
D286	Green
F287	Green
K291	Green
A307	Green
I312	Green
D313	Green
V314	Green
G317	Green
K320	Green
S321	Green
E322	Green
T323	Green
G324	Green
P325	Green
T326	Green
T329	Green
V330	Green
Q331	Green
Q344	Green
R348	Green
A356	Green
I362	Green
E365	Green
L370	Green
V381	Green
P385	Green
V386	Green
L387	Green
Y391	Green
Y406	Green
T409	Green
D416	Green
F420	Green
D425	Green
R426	Green
V427	Green
K428	Green
L443	Green
E444	Green
D445	Green
L446	Green
L447	Green
Q453	Green
S452	Green
V456	Green

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	212195	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	6.301	Depositor
Minimum map value	-4.023	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.155	Depositor
Recommended contour level	0.43	Depositor
Map size (Å)	279.0, 279.0, 279.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.93, 0.93, 0.93	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, VUL

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	0.23	0/4299	0.35	0/5829
1	B	0.24	0/4299	0.35	0/5829
1	C	0.24	0/4299	0.35	0/5829
1	D	0.24	0/4299	0.35	0/5829
1	E	0.23	0/4299	0.35	0/5829
1	F	0.24	0/4299	0.35	0/5829
All	All	0.24	0/25794	0.35	0/34974

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
1	E	0	1
1	F	0	1
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	324	GLY	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	B	324	GLY	Peptide
1	C	324	GLY	Peptide
1	D	324	GLY	Peptide
1	E	324	GLY	Peptide
1	F	324	GLY	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4204	0	4208	77	0
1	B	4204	0	4208	77	0
1	C	4204	0	4208	78	0
1	D	4204	0	4208	77	0
1	E	4204	0	4208	78	0
1	F	4204	0	4208	77	0
2	A	34	0	0	1	0
2	B	34	0	0	1	0
2	C	34	0	0	1	0
2	D	34	0	0	1	0
2	E	34	0	0	1	0
2	F	34	0	0	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
All	All	25434	0	25248	443	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 9.

All (443) 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:B:443:LEU:HD23	1:B:458:VAL:HG11	1.33	1.10
1:A:443:LEU:HD23	1:A:458:VAL:HG11	1.33	1.09
1:D:443:LEU:HD23	1:D:458:VAL:HG11	1.33	1.09
1:F:443:LEU:HD23	1:F:458:VAL:HG11	1.33	1.09
1:C:443:LEU:HD23	1:C:458:VAL:HG11	1.33	1.06
1:E:443:LEU:HD23	1:E:458:VAL:HG11	1.33	1.05
1:D:227:VAL:HG22	1:D:325:PRO:HD3	1.44	1.00
1:C:227:VAL:HG22	1:C:325:PRO:HD3	1.44	1.00
1:A:227:VAL:HG22	1:A:325:PRO:HD3	1.44	0.99
1:E:227:VAL:HG22	1:E:325:PRO:HD3	1.44	0.99
1:B:227:VAL:HG22	1:B:325:PRO:HD3	1.44	0.99
1:F:227:VAL:HG22	1:F:325:PRO:HD3	1.44	0.98
1:C:443:LEU:HD23	1:C:458:VAL:CG1	2.07	0.84
1:E:443:LEU:HD23	1:E:458:VAL:CG1	2.07	0.84
1:B:443:LEU:HD23	1:B:458:VAL:CG1	2.07	0.83
1:A:443:LEU:HD23	1:A:458:VAL:CG1	2.07	0.83
1:D:443:LEU:HD23	1:D:458:VAL:CG1	2.08	0.83
1:F:443:LEU:HD23	1:F:458:VAL:CG1	2.07	0.83
1:E:446:ILE:HD12	1:E:500:LEU:HD13	1.73	0.71
1:A:446:ILE:HD12	1:A:500:LEU:HD13	1.73	0.71
1:C:446:ILE:HD12	1:C:500:LEU:HD13	1.73	0.70
1:D:446:ILE:HD12	1:D:500:LEU:HD13	1.73	0.70
1:C:230:TRP:HA	1:C:325:PRO:HG3	1.74	0.70
1:A:230:TRP:HA	1:A:325:PRO:HG3	1.73	0.70
1:F:230:TRP:HA	1:F:325:PRO:HG3	1.73	0.70
1:F:446:ILE:HD12	1:F:500:LEU:HD13	1.73	0.70
1:B:230:TRP:HA	1:B:325:PRO:HG3	1.73	0.70
1:B:446:ILE:HD12	1:B:500:LEU:HD13	1.73	0.70
1:D:230:TRP:HA	1:D:325:PRO:HG3	1.73	0.69
1:E:230:TRP:HA	1:E:325:PRO:HG3	1.73	0.68
1:A:92:MET:HE1	1:A:171:THR:HB	1.78	0.66
1:C:92:MET:HE1	1:C:171:THR:HB	1.78	0.66
1:B:92:MET:HE1	1:B:171:THR:HB	1.77	0.65
1:D:92:MET:HE1	1:D:171:THR:HB	1.78	0.65
1:E:92:MET:HE1	1:E:171:THR:HB	1.78	0.65
1:A:227:VAL:HG11	1:A:325:PRO:HB3	1.79	0.65
1:B:227:VAL:HG11	1:B:325:PRO:HB3	1.79	0.64
1:E:259:GLU:HB2	1:E:285:MET:HE3	1.80	0.64
1:F:259:GLU:HB2	1:F:285:MET:HE3	1.80	0.64
1:F:92:MET:HE1	1:F:171:THR:HB	1.77	0.64
1:F:227:VAL:HG11	1:F:325:PRO:HB3	1.79	0.64
1:D:259:GLU:HB2	1:D:285:MET:HE3	1.80	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:227:VAL:HG11	1:D:325:PRO:HB3	1.79	0.63
1:C:227:VAL:HG11	1:C:325:PRO:HB3	1.79	0.63
1:E:227:VAL:HG11	1:E:325:PRO:HB3	1.79	0.63
1:C:259:GLU:HB2	1:C:285:MET:HE3	1.80	0.63
1:A:259:GLU:HB2	1:A:285:MET:HE3	1.80	0.62
1:B:259:GLU:HB2	1:B:285:MET:HE3	1.80	0.62
1:B:456:VAL:HG22	1:B:474:VAL:HG22	1.84	0.60
1:E:172:THR:HG22	1:E:186:PHE:HB3	1.84	0.60
1:E:456:VAL:HG22	1:E:474:VAL:HG22	1.84	0.60
1:F:456:VAL:HG22	1:F:474:VAL:HG22	1.84	0.60
1:F:172:THR:HG22	1:F:186:PHE:HB3	1.84	0.60
1:D:456:VAL:HG22	1:D:474:VAL:HG22	1.84	0.59
1:A:456:VAL:HG22	1:A:474:VAL:HG22	1.84	0.59
1:D:172:THR:HG22	1:D:186:PHE:HB3	1.84	0.59
1:B:344:GLN:HG2	1:B:348:ARG:HE	1.67	0.59
1:C:172:THR:HG22	1:C:186:PHE:HB3	1.84	0.59
1:B:172:THR:HG22	1:B:186:PHE:HB3	1.84	0.59
1:C:344:GLN:HG2	1:C:348:ARG:HE	1.67	0.59
1:D:344:GLN:HG2	1:D:348:ARG:HE	1.67	0.59
1:E:344:GLN:HG2	1:E:348:ARG:HE	1.67	0.59
1:F:344:GLN:HG2	1:F:348:ARG:HE	1.67	0.59
1:B:193:VAL:O	1:B:197:MET:HG2	2.03	0.59
1:D:193:VAL:O	1:D:197:MET:HG2	2.03	0.59
1:A:172:THR:HG22	1:A:186:PHE:HB3	1.84	0.58
1:A:344:GLN:HG2	1:A:348:ARG:HE	1.67	0.58
1:C:456:VAL:HG22	1:C:474:VAL:HG22	1.84	0.58
1:A:193:VAL:O	1:A:197:MET:HG2	2.03	0.58
1:C:90:LEU:HG	1:C:92:MET:HE2	1.86	0.58
1:E:193:VAL:O	1:E:197:MET:HG2	2.03	0.57
1:F:193:VAL:O	1:F:197:MET:HG2	2.03	0.57
1:C:193:VAL:O	1:C:197:MET:HG2	2.03	0.57
1:F:90:LEU:HG	1:F:92:MET:HE2	1.86	0.57
1:A:90:LEU:HG	1:A:92:MET:HE2	1.86	0.57
1:C:426:ARG:HG2	1:C:428:LYS:HG3	1.87	0.57
1:F:426:ARG:HG2	1:F:428:LYS:HG3	1.87	0.56
1:B:90:LEU:HG	1:B:92:MET:HE2	1.86	0.56
1:E:90:LEU:HG	1:E:92:MET:HE2	1.86	0.56
1:F:447:LEU:HD11	1:F:507:LEU:HD21	1.88	0.56
1:B:426:ARG:HG2	1:B:428:LYS:HG3	1.87	0.56
1:C:447:LEU:HD11	1:C:507:LEU:HD21	1.88	0.56
1:D:426:ARG:HG2	1:D:428:LYS:HG3	1.87	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:LEU:HD11	1:A:507:LEU:HD21	1.88	0.56
1:D:90:LEU:HG	1:D:92:MET:HE2	1.86	0.56
1:E:447:LEU:HD11	1:E:507:LEU:HD21	1.88	0.56
1:E:426:ARG:HG2	1:E:428:LYS:HG3	1.87	0.56
1:A:426:ARG:HG2	1:A:428:LYS:HG3	1.87	0.55
1:B:81:PHE:O	1:B:85:MET:HB2	2.07	0.55
1:D:81:PHE:O	1:D:85:MET:HB2	2.07	0.55
1:A:67:ALA:HB3	1:A:114:ILE:HG12	1.89	0.55
1:A:81:PHE:O	1:A:85:MET:HB2	2.07	0.55
1:A:365:GLU:OE1	1:A:365:GLU:N	2.36	0.55
1:A:206:ASN:O	1:B:211:ARG:NH2	2.34	0.55
1:C:81:PHE:O	1:C:85:MET:HB2	2.07	0.55
1:B:67:ALA:HB3	1:B:114:ILE:HG12	1.89	0.55
1:D:67:ALA:HB3	1:D:114:ILE:HG12	1.89	0.55
1:B:447:LEU:HD11	1:B:507:LEU:HD21	1.88	0.54
1:A:185:VAL:HG11	1:A:322:GLU:HG3	1.89	0.54
1:D:447:LEU:HD11	1:D:507:LEU:HD21	1.88	0.54
1:E:81:PHE:O	1:E:85:MET:HB2	2.07	0.54
1:C:67:ALA:HB3	1:C:114:ILE:HG12	1.89	0.54
1:E:227:VAL:CG1	1:E:325:PRO:HB3	2.37	0.54
1:F:227:VAL:CG1	1:F:325:PRO:HB3	2.37	0.54
1:A:317:GLY:O	2:A:601:VUL:N6	2.41	0.54
1:E:317:GLY:O	2:E:601:VUL:N6	2.41	0.54
1:F:317:GLY:O	2:F:601:VUL:N6	2.41	0.54
1:B:227:VAL:CG1	1:B:325:PRO:HB3	2.37	0.54
1:A:227:VAL:CG1	1:A:325:PRO:HB3	2.37	0.54
1:B:317:GLY:O	2:B:601:VUL:N6	2.41	0.54
1:D:317:GLY:O	2:D:601:VUL:N6	2.41	0.54
1:C:317:GLY:O	2:C:601:VUL:N6	2.41	0.54
1:D:227:VAL:CG1	1:D:325:PRO:HB3	2.37	0.54
1:E:67:ALA:HB3	1:E:114:ILE:HG12	1.89	0.54
1:C:185:VAL:HG11	1:C:322:GLU:HG3	1.89	0.54
1:C:443:LEU:HD21	1:C:507:LEU:HD13	1.90	0.54
1:F:185:VAL:HG11	1:F:322:GLU:HG3	1.89	0.54
1:D:185:VAL:HG11	1:D:322:GLU:HG3	1.89	0.53
1:E:211:ARG:NH2	1:F:206:ASN:O	2.34	0.53
1:F:81:PHE:O	1:F:85:MET:HB2	2.07	0.53
1:B:185:VAL:HG11	1:B:322:GLU:HG3	1.89	0.53
1:C:227:VAL:CG1	1:C:325:PRO:HB3	2.37	0.53
1:E:206:ASN:O	1:F:211:ARG:NH2	2.34	0.53
1:F:67:ALA:HB3	1:F:114:ILE:HG12	1.89	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:443:LEU:HD21	1:A:507:LEU:HD13	1.90	0.53
1:D:443:LEU:HD21	1:D:507:LEU:HD13	1.90	0.53
1:F:443:LEU:HD21	1:F:507:LEU:HD13	1.90	0.53
1:E:185:VAL:HG11	1:E:322:GLU:HG3	1.89	0.53
1:E:443:LEU:HD21	1:E:507:LEU:HD13	1.90	0.53
1:C:365:GLU:OE1	1:C:365:GLU:N	2.36	0.52
1:B:443:LEU:HD21	1:B:507:LEU:HD13	1.90	0.52
1:B:211:ARG:O	1:B:291:LYS:NZ	2.41	0.52
1:A:425:ASP:CG	1:A:426:ARG:H	2.18	0.52
1:C:425:ASP:CG	1:C:426:ARG:H	2.18	0.51
1:E:211:ARG:O	1:E:291:LYS:NZ	2.41	0.51
1:B:224:MET:HA	1:B:229:ALA:HB2	1.92	0.51
1:E:365:GLU:OE1	1:E:365:GLU:N	2.36	0.51
1:A:211:ARG:NH2	1:B:206:ASN:O	2.34	0.51
1:A:224:MET:HA	1:A:229:ALA:HB2	1.93	0.51
1:B:221:ILE:HB	1:B:268:CYS:HB3	1.93	0.51
1:F:425:ASP:CG	1:F:426:ARG:H	2.18	0.51
1:B:425:ASP:CG	1:B:426:ARG:H	2.18	0.51
1:F:224:MET:HA	1:F:229:ALA:HB2	1.93	0.51
1:A:221:ILE:HB	1:A:268:CYS:HB3	1.93	0.51
1:E:224:MET:HA	1:E:229:ALA:HB2	1.93	0.51
1:D:224:MET:HA	1:D:229:ALA:HB2	1.92	0.51
1:E:320:MET:HE1	1:E:409:THR:HA	1.93	0.51
1:D:221:ILE:HB	1:D:268:CYS:HB3	1.93	0.50
1:A:320:MET:HE1	1:A:409:THR:HA	1.93	0.50
1:C:206:ASN:O	1:D:211:ARG:NH2	2.35	0.50
1:D:425:ASP:CG	1:D:426:ARG:H	2.18	0.50
1:C:221:ILE:HB	1:C:268:CYS:HB3	1.93	0.50
1:C:224:MET:HA	1:C:229:ALA:HB2	1.92	0.50
1:B:170:ALA:HB2	1:B:189:HIS:CD2	2.47	0.50
1:B:365:GLU:OE1	1:B:365:GLU:N	2.36	0.50
1:C:227:VAL:HG22	1:C:325:PRO:CD	2.31	0.50
1:C:320:MET:HE1	1:C:409:THR:HA	1.93	0.50
1:E:425:ASP:CG	1:E:426:ARG:H	2.18	0.50
1:C:170:ALA:HB2	1:C:189:HIS:CD2	2.47	0.50
1:D:320:MET:HE1	1:D:409:THR:HA	1.93	0.50
1:D:365:GLU:OE1	1:D:365:GLU:N	2.36	0.50
1:E:221:ILE:HB	1:E:268:CYS:HB3	1.93	0.50
1:F:227:VAL:HG22	1:F:325:PRO:CD	2.31	0.50
1:A:287:PHE:HB3	1:A:312:ILE:HG12	1.94	0.49
1:C:211:ARG:NH2	1:D:206:ASN:O	2.35	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:221:ILE:HB	1:F:268:CYS:HB3	1.93	0.49
1:F:287:PHE:HB3	1:F:312:ILE:HG12	1.94	0.49
1:F:320:MET:HE1	1:F:409:THR:HA	1.93	0.49
1:A:170:ALA:HB2	1:A:189:HIS:CD2	2.47	0.49
1:F:365:GLU:OE1	1:F:365:GLU:N	2.36	0.49
1:E:287:PHE:HB3	1:E:312:ILE:HG12	1.94	0.49
1:B:227:VAL:HG22	1:B:325:PRO:CD	2.31	0.49
1:C:287:PHE:HB3	1:C:312:ILE:HG12	1.94	0.49
1:D:287:PHE:HB3	1:D:312:ILE:HG12	1.95	0.49
1:B:287:PHE:HB3	1:B:312:ILE:HG12	1.94	0.49
1:D:170:ALA:HB2	1:D:189:HIS:CD2	2.47	0.49
1:F:170:ALA:HB2	1:F:189:HIS:CD2	2.47	0.49
1:F:195:HIS:HE1	1:F:326:ILE:H	1.61	0.49
1:B:320:MET:HE1	1:B:409:THR:HA	1.93	0.49
1:E:170:ALA:HB2	1:E:189:HIS:CD2	2.47	0.49
1:E:195:HIS:HE1	1:E:326:ILE:H	1.61	0.49
1:D:291:LYS:HG2	1:D:313:ASP:HB3	1.96	0.48
1:B:381:VAL:HB	1:B:406:TYR:HB3	1.96	0.48
1:C:195:HIS:HE1	1:C:326:ILE:H	1.61	0.48
1:F:291:LYS:HG2	1:F:313:ASP:HB3	1.96	0.48
1:E:291:LYS:HG2	1:E:313:ASP:HB3	1.96	0.48
1:B:291:LYS:HG2	1:B:313:ASP:HB3	1.96	0.48
1:C:381:VAL:HB	1:C:406:TYR:HB3	1.96	0.48
1:A:381:VAL:HB	1:A:406:TYR:HB3	1.96	0.48
1:A:444:GLU:HG2	1:A:456:VAL:HG12	1.96	0.48
1:B:22:PRO:HA	1:B:238:MET:SD	2.54	0.48
1:A:22:PRO:HA	1:A:238:MET:SD	2.54	0.48
1:A:291:LYS:HG2	1:A:313:ASP:HB3	1.96	0.48
1:B:444:GLU:HG2	1:B:456:VAL:HG12	1.96	0.48
1:C:291:LYS:HG2	1:C:313:ASP:HB3	1.96	0.48
1:C:444:GLU:HG2	1:C:456:VAL:HG12	1.96	0.48
1:D:195:HIS:HE1	1:D:326:ILE:H	1.61	0.48
1:F:381:VAL:HB	1:F:406:TYR:HB3	1.96	0.47
1:A:49:ARG:HG2	1:A:159:TYR:HB3	1.97	0.47
1:F:49:ARG:HG2	1:F:159:TYR:HB3	1.97	0.47
1:B:222:THR:HB	1:B:267:HIS:CD2	2.50	0.47
1:D:381:VAL:HB	1:D:406:TYR:HB3	1.96	0.47
1:A:196:THR:HG23	1:A:235:MET:HG2	1.96	0.47
1:B:430:MET:HE3	1:B:430:MET:HB3	1.77	0.47
1:D:430:MET:HB3	1:D:430:MET:HE3	1.77	0.47
1:A:195:HIS:HE1	1:A:326:ILE:H	1.61	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:HIS:HE1	1:B:326:ILE:H	1.61	0.47
1:C:49:ARG:HG2	1:C:159:TYR:HB3	1.97	0.47
1:C:326:ILE:HG21	1:C:329:ILE:HD11	1.96	0.47
1:E:381:VAL:HB	1:E:406:TYR:HB3	1.96	0.47
1:F:227:VAL:CG2	1:F:325:PRO:HD3	2.30	0.47
1:F:326:ILE:HG21	1:F:329:ILE:HD11	1.96	0.47
1:B:144:GLU:H	1:B:144:GLU:CD	2.23	0.47
1:C:144:GLU:CD	1:C:144:GLU:H	2.23	0.47
1:C:196:THR:HG23	1:C:235:MET:HG2	1.96	0.47
1:D:144:GLU:CD	1:D:144:GLU:H	2.23	0.47
1:D:196:THR:HG23	1:D:235:MET:HG2	1.96	0.47
1:E:144:GLU:CD	1:E:144:GLU:H	2.23	0.47
1:E:222:THR:HB	1:E:267:HIS:CD2	2.50	0.47
1:F:22:PRO:HA	1:F:238:MET:SD	2.54	0.47
1:F:222:THR:HB	1:F:267:HIS:CD2	2.50	0.47
1:F:444:GLU:HG2	1:F:456:VAL:HG12	1.96	0.47
1:C:22:PRO:HA	1:C:238:MET:SD	2.54	0.47
1:A:144:GLU:CD	1:A:144:GLU:H	2.23	0.47
1:B:49:ARG:HG2	1:B:159:TYR:HB3	1.97	0.47
1:D:22:PRO:HA	1:D:238:MET:SD	2.54	0.47
1:E:22:PRO:HA	1:E:238:MET:SD	2.54	0.47
1:E:49:ARG:HG2	1:E:159:TYR:HB3	1.97	0.47
1:E:196:THR:HG23	1:E:235:MET:HG2	1.96	0.47
1:F:144:GLU:CD	1:F:144:GLU:H	2.23	0.47
1:A:222:THR:HB	1:A:267:HIS:CD2	2.50	0.47
1:A:326:ILE:HG21	1:A:329:ILE:HD11	1.97	0.47
1:D:222:THR:HB	1:D:267:HIS:CD2	2.50	0.47
1:D:444:GLU:HG2	1:D:456:VAL:HG12	1.96	0.47
1:E:227:VAL:CG2	1:E:325:PRO:HD3	2.30	0.47
1:F:443:LEU:CD2	1:F:458:VAL:HG11	2.25	0.47
1:B:196:THR:HG23	1:B:235:MET:HG2	1.96	0.46
1:E:326:ILE:HG21	1:E:329:ILE:HD11	1.97	0.46
1:E:444:GLU:HG2	1:E:456:VAL:HG12	1.96	0.46
1:B:191:GLN:HG2	1:B:385:PRO:O	2.16	0.46
1:B:326:ILE:HG21	1:B:329:ILE:HD11	1.96	0.46
1:D:326:ILE:HG21	1:D:329:ILE:HD11	1.97	0.46
1:C:191:GLN:HG2	1:C:385:PRO:O	2.16	0.46
1:B:16:LYS:HB3	1:B:85:MET:HG2	1.98	0.46
1:C:222:THR:HB	1:C:267:HIS:CD2	2.50	0.46
1:D:49:ARG:HG2	1:D:159:TYR:HB3	1.97	0.46
1:A:16:LYS:HB3	1:A:85:MET:HG2	1.98	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:227:VAL:HG22	1:D:325:PRO:CD	2.31	0.46
1:E:26:ASN:HB3	1:E:29:GLN:HB2	1.98	0.46
1:F:416:ASP:OD1	1:F:420:PHE:N	2.48	0.46
1:A:191:GLN:HG2	1:A:385:PRO:O	2.16	0.46
1:B:206:ASN:OD1	1:B:331:GLN:NE2	2.50	0.45
1:D:191:GLN:HG2	1:D:385:PRO:O	2.16	0.45
1:E:191:GLN:HG2	1:E:385:PRO:O	2.16	0.45
1:F:196:THR:HG23	1:F:235:MET:HG2	1.96	0.45
1:B:92:MET:HB3	1:B:225:PHE:HA	1.99	0.45
1:C:227:VAL:CG2	1:C:325:PRO:HD3	2.30	0.45
1:F:206:ASN:OD1	1:F:331:GLN:NE2	2.50	0.45
1:D:26:ASN:HB3	1:D:29:GLN:HB2	1.98	0.45
1:E:206:ASN:OD1	1:E:331:GLN:NE2	2.50	0.45
1:F:26:ASN:HB3	1:F:29:GLN:HB2	1.98	0.45
1:F:191:GLN:HG2	1:F:385:PRO:O	2.16	0.45
1:C:416:ASP:OD1	1:C:420:PHE:N	2.48	0.45
1:F:16:LYS:HB3	1:F:85:MET:HG2	1.98	0.45
1:F:92:MET:HB3	1:F:225:PHE:HA	1.98	0.45
1:D:92:MET:HB3	1:D:225:PHE:HA	1.99	0.45
1:E:416:ASP:OD1	1:E:420:PHE:N	2.48	0.45
1:C:206:ASN:OD1	1:C:331:GLN:NE2	2.50	0.45
1:E:16:LYS:HB3	1:E:85:MET:HG2	1.98	0.45
1:C:15:ILE:HD11	1:C:192:LEU:HD12	1.99	0.45
1:D:15:ILE:HD11	1:D:192:LEU:HD12	1.99	0.45
1:D:206:ASN:OD1	1:D:331:GLN:NE2	2.50	0.45
1:A:191:GLN:NE2	1:A:387:LEU:O	2.50	0.45
1:B:15:ILE:HD11	1:B:192:LEU:HD12	1.99	0.45
1:C:26:ASN:HB3	1:C:29:GLN:HB2	1.98	0.45
1:F:15:ILE:HD11	1:F:192:LEU:HD12	1.99	0.45
1:A:92:MET:HB3	1:A:225:PHE:HA	1.99	0.44
1:C:307:ALA:HB3	1:C:314:VAL:HG11	2.00	0.44
1:E:15:ILE:HD11	1:E:192:LEU:HD12	1.99	0.44
1:F:330:VAL:HG11	1:F:348:ARG:HA	2.00	0.44
1:A:206:ASN:OD1	1:A:331:GLN:NE2	2.50	0.44
1:C:16:LYS:HB3	1:C:85:MET:HG2	1.98	0.44
1:B:330:VAL:HG11	1:B:348:ARG:HA	2.00	0.44
1:D:16:LYS:HB3	1:D:85:MET:HG2	1.98	0.44
1:E:459:ILE:HD11	1:E:471:LEU:HD23	1.99	0.44
1:A:15:ILE:HD11	1:A:192:LEU:HD12	1.99	0.44
1:A:26:ASN:HB3	1:A:29:GLN:HB2	1.98	0.44
1:A:430:MET:HE3	1:A:430:MET:HB3	1.77	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:92:MET:HB3	1:E:225:PHE:HA	1.99	0.44
1:E:330:VAL:HG11	1:E:348:ARG:HA	2.00	0.44
1:B:26:ASN:HB3	1:B:29:GLN:HB2	1.98	0.44
1:C:92:MET:HB3	1:C:225:PHE:HA	1.99	0.44
1:D:307:ALA:HB3	1:D:314:VAL:HG11	2.00	0.44
1:B:397:ASN:ND2	1:D:132:ASP:OD2	2.48	0.44
1:C:330:VAL:HG11	1:C:348:ARG:HA	2.00	0.44
1:D:330:VAL:HG11	1:D:348:ARG:HA	2.00	0.44
1:A:330:VAL:HG11	1:A:348:ARG:HA	2.00	0.44
1:D:459:ILE:HD11	1:D:471:LEU:HD23	1.99	0.44
1:B:191:GLN:NE2	1:B:387:LEU:O	2.50	0.44
1:F:183:LYS:HB3	1:F:391:TYR:CD1	2.53	0.44
1:F:459:ILE:HD11	1:F:471:LEU:HD23	1.99	0.44
1:A:307:ALA:HB3	1:A:314:VAL:HG11	2.00	0.44
1:C:183:LYS:HB3	1:C:391:TYR:CD1	2.53	0.43
1:D:227:VAL:CG2	1:D:325:PRO:HD3	2.30	0.43
1:F:118:GLU:OE1	1:F:141:ARG:HG2	2.18	0.43
1:A:23:VAL:HG13	1:B:202:THR:HA	2.00	0.43
1:A:416:ASP:OD1	1:A:420:PHE:N	2.48	0.43
1:C:459:ILE:HD11	1:C:471:LEU:HD23	1.99	0.43
1:F:307:ALA:HB3	1:F:314:VAL:HG11	2.00	0.43
1:B:307:ALA:HB3	1:B:314:VAL:HG11	2.00	0.43
1:D:183:LYS:HB3	1:D:391:TYR:CD1	2.53	0.43
1:D:191:GLN:NE2	1:D:387:LEU:O	2.50	0.43
1:C:23:VAL:HG13	1:D:202:THR:HA	2.00	0.43
1:D:232:LEU:HD13	1:D:235:MET:HE3	2.01	0.43
1:E:118:GLU:OE1	1:E:141:ARG:HG2	2.18	0.43
1:A:118:GLU:OE1	1:A:141:ARG:HG2	2.18	0.43
1:A:183:LYS:HB3	1:A:391:TYR:CD1	2.53	0.43
1:D:416:ASP:OD1	1:D:420:PHE:N	2.48	0.43
1:E:232:LEU:HD13	1:E:235:MET:HE3	2.01	0.43
1:F:146:CYS:SG	1:F:148:TYR:HB3	2.59	0.43
1:A:146:CYS:SG	1:A:148:TYR:HB3	2.59	0.43
1:A:150:ARG:O	1:A:154:GLN:HG2	2.19	0.43
1:E:307:ALA:HB3	1:E:314:VAL:HG11	2.00	0.43
1:A:202:THR:HA	1:B:23:VAL:HG13	2.00	0.43
1:B:183:LYS:HB3	1:B:391:TYR:CD1	2.53	0.43
1:B:232:LEU:HD13	1:B:235:MET:HE3	2.01	0.43
1:D:473:LEU:HD23	1:D:510:LYS:HB3	2.01	0.43
1:E:473:LEU:HD23	1:E:510:LYS:HB3	2.01	0.43
1:B:146:CYS:SG	1:B:148:TYR:HB3	2.59	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:ARG:O	1:B:154:GLN:HG2	2.19	0.43
1:B:416:ASP:OD1	1:B:420:PHE:N	2.48	0.43
1:C:146:CYS:SG	1:C:148:TYR:HB3	2.59	0.43
1:D:443:LEU:HD11	1:D:507:LEU:HD22	2.01	0.43
1:C:362:ILE:HG22	1:C:370:LEU:HB2	2.01	0.43
1:D:146:CYS:SG	1:D:148:TYR:HB3	2.59	0.43
1:E:146:CYS:SG	1:E:148:TYR:HB3	2.59	0.43
1:A:459:ILE:HD11	1:A:471:LEU:HD23	1.99	0.43
1:C:238:MET:HE3	1:D:356:ALA:HB1	2.01	0.43
1:C:356:ALA:HB1	1:D:238:MET:HE3	2.01	0.43
1:E:183:LYS:HB3	1:E:391:TYR:CD1	2.53	0.43
1:E:238:MET:HE3	1:F:356:ALA:HB1	2.01	0.43
1:E:443:LEU:HD11	1:E:507:LEU:HD22	2.01	0.43
1:F:232:LEU:HD13	1:F:235:MET:HE3	2.01	0.43
1:B:118:GLU:OE1	1:B:141:ARG:HG2	2.18	0.42
1:B:362:ILE:HG22	1:B:370:LEU:HB2	2.01	0.42
1:B:443:LEU:HD11	1:B:507:LEU:HD22	2.01	0.42
1:B:459:ILE:HD11	1:B:471:LEU:HD23	1.99	0.42
1:C:118:GLU:OE1	1:C:141:ARG:HG2	2.18	0.42
1:C:218:TYR:OH	1:C:267:HIS:HE1	2.03	0.42
1:E:150:ARG:O	1:E:154:GLN:HG2	2.19	0.42
1:E:218:TYR:OH	1:E:267:HIS:HE1	2.03	0.42
1:E:356:ALA:HB1	1:F:238:MET:HE3	2.01	0.42
1:E:362:ILE:HG22	1:E:370:LEU:HB2	2.01	0.42
1:F:150:ARG:O	1:F:154:GLN:HG2	2.19	0.42
1:F:443:LEU:HD11	1:F:507:LEU:HD22	2.01	0.42
1:A:362:ILE:HG22	1:A:370:LEU:HB2	2.01	0.42
1:A:473:LEU:HD23	1:A:510:LYS:HB3	2.01	0.42
1:C:150:ARG:O	1:C:154:GLN:HG2	2.19	0.42
1:D:211:ARG:O	1:D:291:LYS:NZ	2.40	0.42
1:D:218:TYR:OH	1:D:267:HIS:HE1	2.02	0.42
1:D:362:ILE:HG22	1:D:370:LEU:HB2	2.01	0.42
1:F:218:TYR:OH	1:F:267:HIS:HE1	2.03	0.42
1:F:362:ILE:HG22	1:F:370:LEU:HB2	2.01	0.42
1:D:150:ARG:O	1:D:154:GLN:HG2	2.19	0.42
1:F:191:GLN:NE2	1:F:387:LEU:O	2.50	0.42
1:A:227:VAL:CG2	1:A:325:PRO:HD3	2.30	0.42
1:C:202:THR:HA	1:D:23:VAL:HG13	2.00	0.42
1:F:473:LEU:HD23	1:F:510:LYS:HB3	2.00	0.42
1:B:136:ARG:HH22	1:B:151:LEU:HD21	1.85	0.42
1:E:202:THR:HA	1:F:23:VAL:HG13	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:TYR:OH	1:A:267:HIS:HE1	2.03	0.42
1:A:232:LEU:HD13	1:A:235:MET:HE3	2.01	0.42
1:B:218:TYR:OH	1:B:267:HIS:HE1	2.03	0.42
1:D:118:GLU:OE1	1:D:141:ARG:HG2	2.18	0.42
1:E:23:VAL:HG13	1:F:202:THR:HA	2.00	0.42
1:A:238:MET:HE3	1:B:356:ALA:HB1	2.01	0.42
1:C:232:LEU:HD13	1:C:235:MET:HE3	2.00	0.42
1:C:473:LEU:HD23	1:C:510:LYS:HB3	2.00	0.42
1:D:113:ILE:HG12	1:D:136:ARG:HB3	2.02	0.42
1:F:197:MET:SD	1:F:238:MET:HE2	2.60	0.42
1:A:356:ALA:HB1	1:B:238:MET:HE3	2.01	0.42
1:B:84:PRO:HB3	1:B:170:ALA:HB1	2.02	0.42
1:C:197:MET:SD	1:C:238:MET:HE2	2.60	0.42
1:A:173:PHE:HZ	1:A:227:VAL:HG23	1.85	0.41
1:B:173:PHE:HZ	1:B:227:VAL:HG23	1.85	0.41
1:C:191:GLN:NE2	1:C:387:LEU:O	2.50	0.41
1:D:136:ARG:HH22	1:D:151:LEU:HD21	1.85	0.41
1:E:191:GLN:NE2	1:E:387:LEU:O	2.50	0.41
1:E:197:MET:SD	1:E:238:MET:HE2	2.60	0.41
1:B:452:SER:HB2	1:B:480:ALA:HB1	2.02	0.41
1:B:473:LEU:HD23	1:B:510:LYS:HB3	2.00	0.41
1:C:113:ILE:HG12	1:C:136:ARG:HB3	2.02	0.41
1:D:84:PRO:HB3	1:D:170:ALA:HB1	2.02	0.41
1:D:173:PHE:HZ	1:D:227:VAL:HG23	1.85	0.41
1:A:136:ARG:HH22	1:A:151:LEU:HD21	1.85	0.41
1:A:197:MET:SD	1:A:238:MET:HE2	2.60	0.41
1:A:443:LEU:HD11	1:A:507:LEU:HD22	2.01	0.41
1:A:452:SER:HB2	1:A:480:ALA:HB1	2.02	0.41
1:B:108:HIS:HD2	1:B:174:TYR:OH	2.04	0.41
1:C:443:LEU:HD11	1:C:507:LEU:HD22	2.01	0.41
1:E:108:HIS:HD2	1:E:174:TYR:OH	2.04	0.41
1:E:173:PHE:HZ	1:E:227:VAL:HG23	1.85	0.41
1:C:452:SER:HB2	1:C:480:ALA:HB1	2.02	0.41
1:B:113:ILE:HG12	1:B:136:ARG:HB3	2.02	0.41
1:B:197:MET:SD	1:B:238:MET:HE2	2.60	0.41
1:B:227:VAL:CG2	1:B:325:PRO:HD3	2.30	0.41
1:F:231:GLY:O	1:F:235:MET:HG3	2.21	0.41
1:F:256:ASN:HB2	1:F:281:LYS:NZ	2.36	0.41
1:A:84:PRO:HB3	1:A:170:ALA:HB1	2.02	0.41
1:B:231:GLY:O	1:B:235:MET:HG3	2.21	0.41
1:C:320:MET:O	1:C:325:PRO:HA	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:231:GLY:O	1:E:235:MET:HG3	2.21	0.41
1:E:256:ASN:HB2	1:E:281:LYS:NZ	2.36	0.41
1:E:320:MET:O	1:E:325:PRO:HA	2.21	0.41
1:F:320:MET:O	1:F:325:PRO:HA	2.21	0.41
1:F:113:ILE:HG12	1:F:136:ARG:HB3	2.02	0.41
1:A:256:ASN:HB2	1:A:281:LYS:NZ	2.36	0.41
1:C:108:HIS:HD2	1:C:174:TYR:OH	2.04	0.41
1:D:197:MET:SD	1:D:238:MET:HE2	2.60	0.41
1:E:84:PRO:HB3	1:E:170:ALA:HB1	2.02	0.41
1:E:136:ARG:HH22	1:E:151:LEU:HD21	1.85	0.41
1:F:136:ARG:HH22	1:F:151:LEU:HD21	1.85	0.41
1:A:108:HIS:HD2	1:A:174:TYR:OH	2.04	0.41
1:D:108:HIS:HD2	1:D:174:TYR:OH	2.04	0.41
1:D:452:SER:HB2	1:D:480:ALA:HB1	2.02	0.41
1:A:132:ASP:OD2	1:E:397:ASN:ND2	2.49	0.40
1:A:113:ILE:HG12	1:A:136:ARG:HB3	2.02	0.40
1:C:231:GLY:O	1:C:235:MET:HG3	2.21	0.40
1:C:256:ASN:HB2	1:C:281:LYS:NZ	2.36	0.40
1:C:443:LEU:CD2	1:C:458:VAL:HG11	2.25	0.40
1:A:227:VAL:HG22	1:A:325:PRO:CD	2.31	0.40
1:C:84:PRO:HB3	1:C:170:ALA:HB1	2.02	0.40
1:C:211:ARG:O	1:C:291:LYS:NZ	2.40	0.40
1:E:67:ALA:HB1	1:E:93:ILE:HD11	2.04	0.40
1:E:113:ILE:HG12	1:E:136:ARG:HB3	2.02	0.40
1:E:227:VAL:HG22	1:E:325:PRO:CD	2.31	0.40
1:C:136:ARG:HH22	1:C:151:LEU:HD21	1.85	0.40
1:D:256:ASN:HB2	1:D:281:LYS:NZ	2.36	0.40
1:F:108:HIS:HD2	1:F:174:TYR:OH	2.04	0.40
1:F:452:SER:HB2	1:F:480:ALA:HB1	2.02	0.40
1:A:69:MET:HE2	1:A:120:PHE:HB3	2.04	0.40
1:C:397:ASN:ND2	1:F:132:ASP:OD2	2.50	0.40
1:F:84:PRO:HB3	1:F:170:ALA:HB1	2.02	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	525/532 (99%)	515 (98%)	10 (2%)	0	100	100
1	B	525/532 (99%)	515 (98%)	10 (2%)	0	100	100
1	C	525/532 (99%)	516 (98%)	9 (2%)	0	100	100
1	D	525/532 (99%)	515 (98%)	10 (2%)	0	100	100
1	E	525/532 (99%)	515 (98%)	10 (2%)	0	100	100
1	F	525/532 (99%)	515 (98%)	10 (2%)	0	100	100
All	All	3150/3192 (99%)	3091 (98%)	59 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	462/467 (99%)	460 (100%)	2 (0%)	84	89
1	B	462/467 (99%)	460 (100%)	2 (0%)	84	89
1	C	462/467 (99%)	458 (99%)	4 (1%)	70	80
1	D	462/467 (99%)	458 (99%)	4 (1%)	70	80
1	E	462/467 (99%)	460 (100%)	2 (0%)	84	89
1	F	462/467 (99%)	459 (99%)	3 (1%)	78	86
All	All	2772/2802 (99%)	2755 (99%)	17 (1%)	76	86

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	228	HIS
1	A	520	VAL
1	B	228	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	520	VAL
1	C	228	HIS
1	C	275	LEU
1	C	362	ILE
1	C	520	VAL
1	D	228	HIS
1	D	362	ILE
1	D	441	LEU
1	D	520	VAL
1	E	228	HIS
1	E	520	VAL
1	F	228	HIS
1	F	362	ILE
1	F	520	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (59) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	108	HIS
1	A	160	ASN
1	A	195	HIS
1	A	228	HIS
1	A	260	GLN
1	A	267	HIS
1	A	331	GLN
1	A	344	GLN
1	B	29	GLN
1	B	50	GLN
1	B	108	HIS
1	B	160	ASN
1	B	195	HIS
1	B	228	HIS
1	B	260	GLN
1	B	267	HIS
1	B	331	GLN
1	B	344	GLN
1	C	29	GLN
1	C	50	GLN
1	C	108	HIS
1	C	160	ASN
1	C	195	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	228	HIS
1	C	260	GLN
1	C	267	HIS
1	C	331	GLN
1	C	344	GLN
1	D	50	GLN
1	D	108	HIS
1	D	160	ASN
1	D	195	HIS
1	D	228	HIS
1	D	260	GLN
1	D	267	HIS
1	D	331	GLN
1	D	344	GLN
1	D	397	ASN
1	E	50	GLN
1	E	108	HIS
1	E	160	ASN
1	E	195	HIS
1	E	228	HIS
1	E	260	GLN
1	E	267	HIS
1	E	331	GLN
1	E	344	GLN
1	E	448	HIS
1	F	29	GLN
1	F	50	GLN
1	F	108	HIS
1	F	160	ASN
1	F	195	HIS
1	F	228	HIS
1	F	260	GLN
1	F	267	HIS
1	F	331	GLN
1	F	344	GLN

5.3.3 RNA ⓘ

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 12 ligands modelled in this entry, 6 are monoatomic - leaving 6 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
2	VUL	B	601	3	36,36,36	6.72	13 (36%)	46,50,50	2.16	10 (21%)
2	VUL	E	601	3	36,36,36	6.72	13 (36%)	46,50,50	2.16	10 (21%)
2	VUL	D	601	3	36,36,36	6.73	13 (36%)	46,50,50	2.16	10 (21%)
2	VUL	F	601	3	36,36,36	6.73	13 (36%)	46,50,50	2.15	10 (21%)
2	VUL	C	601	3	36,36,36	6.73	13 (36%)	46,50,50	2.16	10 (21%)
2	VUL	A	601	3	36,36,36	6.73	13 (36%)	46,50,50	2.16	10 (21%)

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
2	VUL	B	601	3	-	5/23/40/40	0/3/3/3
2	VUL	E	601	3	-	5/23/40/40	0/3/3/3
2	VUL	D	601	3	-	5/23/40/40	0/3/3/3
2	VUL	F	601	3	-	5/23/40/40	0/3/3/3
2	VUL	C	601	3	-	5/23/40/40	0/3/3/3
2	VUL	A	601	3	-	5/23/40/40	0/3/3/3

All (78) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	601	VUL	O3-S1	25.91	1.65	1.42
2	A	601	VUL	O3-S1	25.91	1.65	1.42
2	E	601	VUL	O3-S1	25.91	1.65	1.42
2	C	601	VUL	O2-S1	25.88	1.65	1.42
2	D	601	VUL	O2-S1	25.88	1.65	1.42
2	B	601	VUL	O2-S1	25.85	1.65	1.42
2	F	601	VUL	O2-S1	25.85	1.65	1.42
2	C	601	VUL	O3-S1	25.85	1.65	1.42
2	D	601	VUL	O3-S1	25.85	1.65	1.42
2	A	601	VUL	O2-S1	25.83	1.65	1.42
2	B	601	VUL	O3-S1	25.81	1.65	1.42
2	E	601	VUL	O2-S1	25.74	1.65	1.42
2	A	601	VUL	O5-C15	8.77	1.62	1.42
2	E	601	VUL	O5-C15	8.77	1.62	1.42
2	C	601	VUL	O5-C15	8.76	1.62	1.42
2	D	601	VUL	O5-C15	8.76	1.62	1.42
2	B	601	VUL	O5-C15	8.75	1.62	1.42
2	F	601	VUL	O5-C15	8.73	1.62	1.42
2	A	601	VUL	C14-C15	-7.22	1.30	1.53
2	E	601	VUL	C14-C15	-7.22	1.30	1.53
2	C	601	VUL	C14-C15	-7.22	1.30	1.53
2	D	601	VUL	C14-C15	-7.22	1.30	1.53
2	F	601	VUL	C14-C15	-7.22	1.30	1.53
2	B	601	VUL	C14-C15	-7.19	1.30	1.53
2	E	601	VUL	S1-N1	7.13	1.71	1.59
2	F	601	VUL	S1-N1	7.13	1.71	1.59
2	B	601	VUL	S1-N1	7.12	1.71	1.59
2	C	601	VUL	S1-N1	7.10	1.71	1.59
2	D	601	VUL	S1-N1	7.10	1.71	1.59
2	A	601	VUL	S1-N1	7.10	1.71	1.59
2	B	601	VUL	O5-C12	-6.37	1.30	1.45
2	A	601	VUL	O5-C12	-6.36	1.30	1.45
2	E	601	VUL	O5-C12	-6.36	1.30	1.45
2	D	601	VUL	O5-C12	-6.36	1.30	1.45
2	F	601	VUL	O5-C12	-6.33	1.30	1.45
2	C	601	VUL	O5-C12	-6.32	1.30	1.45
2	F	601	VUL	C20-N6	4.23	1.45	1.34
2	A	601	VUL	C20-N6	4.19	1.44	1.34
2	E	601	VUL	C20-N6	4.19	1.44	1.34
2	B	601	VUL	C20-N6	4.19	1.44	1.34
2	D	601	VUL	C20-N6	4.19	1.44	1.34
2	C	601	VUL	C20-N6	4.17	1.44	1.34

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	VUL	C18-N2	-2.92	1.32	1.37
2	D	601	VUL	C18-N2	-2.92	1.32	1.37
2	A	601	VUL	C18-N2	-2.90	1.32	1.37
2	E	601	VUL	C18-N2	-2.90	1.32	1.37
2	B	601	VUL	C18-N2	-2.90	1.32	1.37
2	B	601	VUL	O7-C13	-2.88	1.35	1.43
2	F	601	VUL	C18-N2	-2.87	1.32	1.37
2	A	601	VUL	O7-C13	-2.86	1.35	1.43
2	E	601	VUL	O7-C13	-2.86	1.35	1.43
2	C	601	VUL	O7-C13	-2.86	1.35	1.43
2	D	601	VUL	O7-C13	-2.86	1.35	1.43
2	F	601	VUL	O7-C13	-2.83	1.35	1.43
2	B	601	VUL	O4-S1	2.75	1.65	1.60
2	C	601	VUL	O4-S1	2.74	1.65	1.60
2	D	601	VUL	O4-S1	2.74	1.65	1.60
2	C	601	VUL	O6-C14	2.74	1.49	1.43
2	D	601	VUL	O6-C14	2.74	1.49	1.43
2	A	601	VUL	O4-S1	2.73	1.65	1.60
2	E	601	VUL	O4-S1	2.73	1.65	1.60
2	F	601	VUL	O4-S1	2.73	1.65	1.60
2	B	601	VUL	O6-C14	2.72	1.49	1.43
2	A	601	VUL	O6-C14	2.72	1.49	1.43
2	E	601	VUL	O6-C14	2.72	1.49	1.43
2	F	601	VUL	O6-C14	2.72	1.49	1.43
2	E	601	VUL	C10-N1	2.45	1.45	1.39
2	C	601	VUL	C10-N1	2.45	1.45	1.39
2	D	601	VUL	C10-N1	2.45	1.45	1.39
2	B	601	VUL	C10-N1	2.45	1.45	1.39
2	F	601	VUL	C10-N1	2.45	1.45	1.39
2	A	601	VUL	C10-N1	2.44	1.45	1.39
2	C	601	VUL	O1-C10	-2.42	1.18	1.23
2	D	601	VUL	O1-C10	-2.42	1.18	1.23
2	B	601	VUL	O1-C10	-2.40	1.18	1.23
2	A	601	VUL	O1-C10	-2.39	1.18	1.23
2	E	601	VUL	O1-C10	-2.39	1.18	1.23
2	F	601	VUL	O1-C10	-2.38	1.18	1.23

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	601	VUL	O3-S1-O2	-7.35	109.84	120.85
2	B	601	VUL	O3-S1-O2	-7.35	109.84	120.85

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	VUL	O3-S1-O2	-7.35	109.85	120.85
2	D	601	VUL	O3-S1-O2	-7.35	109.85	120.85
2	A	601	VUL	O3-S1-O2	-7.35	109.85	120.85
2	F	601	VUL	O3-S1-O2	-7.34	109.86	120.85
2	C	601	VUL	C17-C16-N4	-5.55	119.08	126.72
2	E	601	VUL	C17-C16-N4	-5.51	119.12	126.72
2	A	601	VUL	C17-C16-N4	-5.51	119.13	126.72
2	B	601	VUL	C17-C16-N4	-5.51	119.13	126.72
2	D	601	VUL	C17-C16-N4	-5.50	119.14	126.72
2	F	601	VUL	C17-C16-N4	-5.50	119.14	126.72
2	D	601	VUL	N5-C19-N4	-5.01	121.00	128.58
2	C	601	VUL	N5-C19-N4	-4.99	121.03	128.58
2	B	601	VUL	N5-C19-N4	-4.99	121.03	128.58
2	A	601	VUL	N5-C19-N4	-4.99	121.03	128.58
2	E	601	VUL	N5-C19-N4	-4.99	121.03	128.58
2	F	601	VUL	N5-C19-N4	-4.97	121.06	128.58
2	C	601	VUL	C19-N4-C16	4.19	122.06	111.83
2	D	601	VUL	C19-N4-C16	4.17	122.02	111.83
2	E	601	VUL	C19-N4-C16	4.17	122.01	111.83
2	A	601	VUL	C19-N4-C16	4.17	122.01	111.83
2	B	601	VUL	C19-N4-C16	4.17	122.00	111.83
2	F	601	VUL	C19-N4-C16	4.16	121.99	111.83
2	C	601	VUL	N4-C16-N2	4.07	134.09	127.17
2	E	601	VUL	N4-C16-N2	4.05	134.05	127.17
2	A	601	VUL	N4-C16-N2	4.04	134.04	127.17
2	D	601	VUL	N4-C16-N2	4.04	134.04	127.17
2	B	601	VUL	N4-C16-N2	4.04	134.04	127.17
2	F	601	VUL	N4-C16-N2	4.01	133.99	127.17
2	C	601	VUL	C16-C17-N3	-3.34	106.76	110.58
2	D	601	VUL	C16-C17-N3	-3.32	106.79	110.58
2	B	601	VUL	C16-C17-N3	-3.31	106.80	110.58
2	F	601	VUL	C16-C17-N3	-3.31	106.80	110.58
2	A	601	VUL	C16-C17-N3	-3.30	106.81	110.58
2	E	601	VUL	C16-C17-N3	-3.30	106.81	110.58
2	C	601	VUL	C17-N3-C18	3.28	108.61	103.45
2	D	601	VUL	C17-N3-C18	3.28	108.61	103.45
2	B	601	VUL	C17-N3-C18	3.27	108.58	103.45
2	F	601	VUL	C17-N3-C18	3.27	108.58	103.45
2	A	601	VUL	C17-N3-C18	3.26	108.58	103.45
2	E	601	VUL	C17-N3-C18	3.26	108.58	103.45
2	C	601	VUL	N2-C18-N3	-3.22	109.37	113.94
2	D	601	VUL	N2-C18-N3	-3.22	109.37	113.94

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	VUL	N2-C18-N3	-3.21	109.39	113.94
2	A	601	VUL	N2-C18-N3	-3.20	109.40	113.94
2	E	601	VUL	N2-C18-N3	-3.20	109.40	113.94
2	F	601	VUL	N2-C18-N3	-3.18	109.42	113.94
2	C	601	VUL	C16-N2-C18	2.56	108.42	105.74
2	D	601	VUL	C16-N2-C18	2.54	108.41	105.74
2	B	601	VUL	C16-N2-C18	2.53	108.40	105.74
2	A	601	VUL	C16-N2-C18	2.52	108.39	105.74
2	E	601	VUL	C16-N2-C18	2.52	108.39	105.74
2	F	601	VUL	C16-N2-C18	2.47	108.33	105.74
2	C	601	VUL	C20-C17-N3	2.28	136.49	132.09
2	D	601	VUL	C20-C17-N3	2.28	136.49	132.09
2	F	601	VUL	C20-C17-N3	2.27	136.47	132.09
2	B	601	VUL	C20-C17-N3	2.27	136.46	132.09
2	A	601	VUL	C20-C17-N3	2.26	136.46	132.09
2	E	601	VUL	C20-C17-N3	2.26	136.46	132.09

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	VUL	C5-C6-C7-C8
2	B	601	VUL	C5-C6-C7-C8
2	C	601	VUL	C5-C6-C7-C8
2	D	601	VUL	C5-C6-C7-C8
2	E	601	VUL	C5-C6-C7-C8
2	F	601	VUL	C5-C6-C7-C8
2	A	601	VUL	C2-C3-C4-C5
2	B	601	VUL	C2-C3-C4-C5
2	C	601	VUL	C2-C3-C4-C5
2	D	601	VUL	C2-C3-C4-C5
2	E	601	VUL	C2-C3-C4-C5
2	F	601	VUL	C2-C3-C4-C5
2	A	601	VUL	C4-C5-C6-C7
2	C	601	VUL	C4-C5-C6-C7
2	E	601	VUL	C4-C5-C6-C7
2	B	601	VUL	C4-C5-C6-C7
2	F	601	VUL	C4-C5-C6-C7
2	D	601	VUL	C4-C5-C6-C7
2	A	601	VUL	C1-C2-C3-C4
2	B	601	VUL	C1-C2-C3-C4
2	C	601	VUL	C1-C2-C3-C4

Continued on next page...

Continued from previous page...

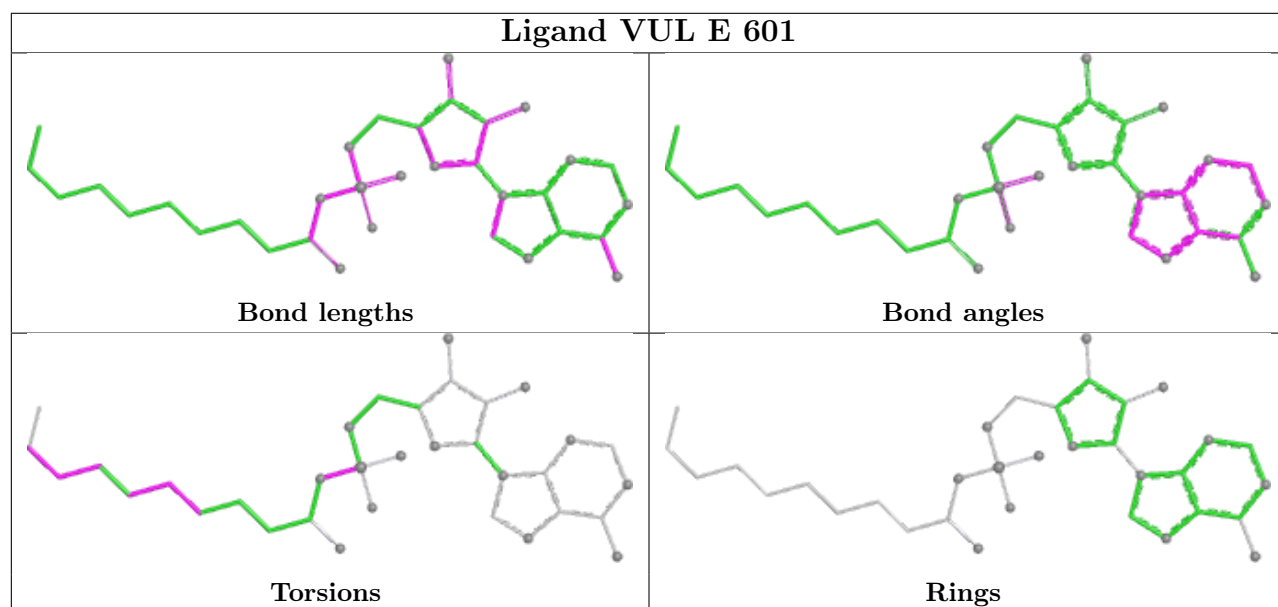
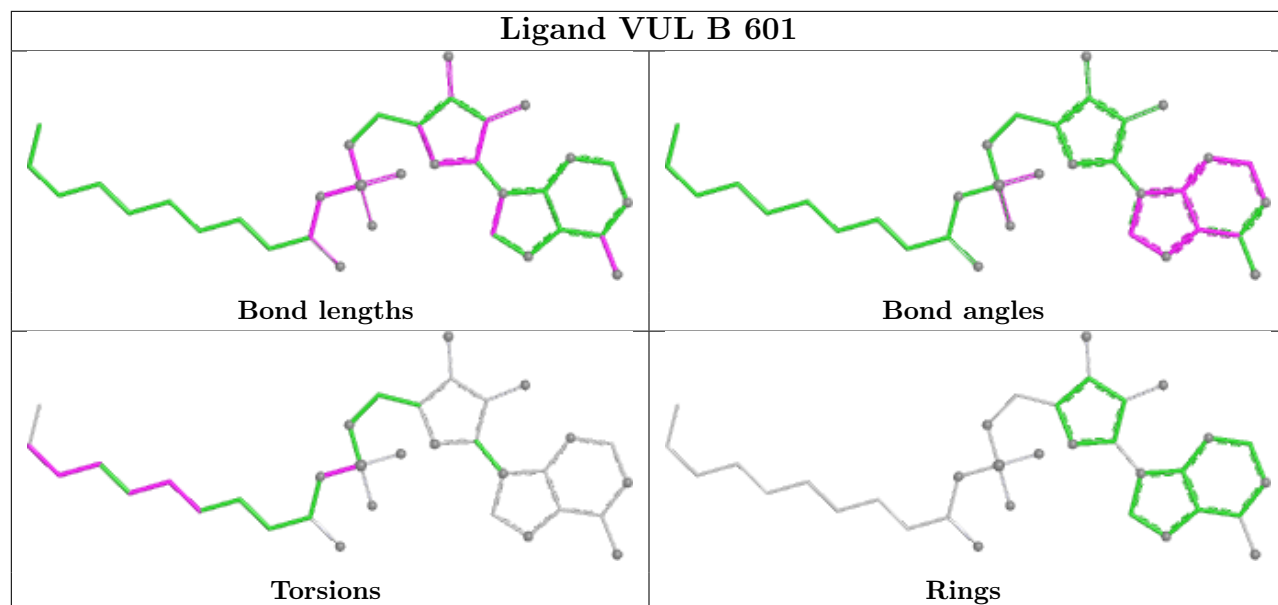
Mol	Chain	Res	Type	Atoms
2	D	601	VUL	C1-C2-C3-C4
2	E	601	VUL	C1-C2-C3-C4
2	F	601	VUL	C1-C2-C3-C4
2	A	601	VUL	C10-N1-S1-O2
2	B	601	VUL	C10-N1-S1-O2
2	C	601	VUL	C10-N1-S1-O2
2	D	601	VUL	C10-N1-S1-O2
2	E	601	VUL	C10-N1-S1-O2
2	F	601	VUL	C10-N1-S1-O2

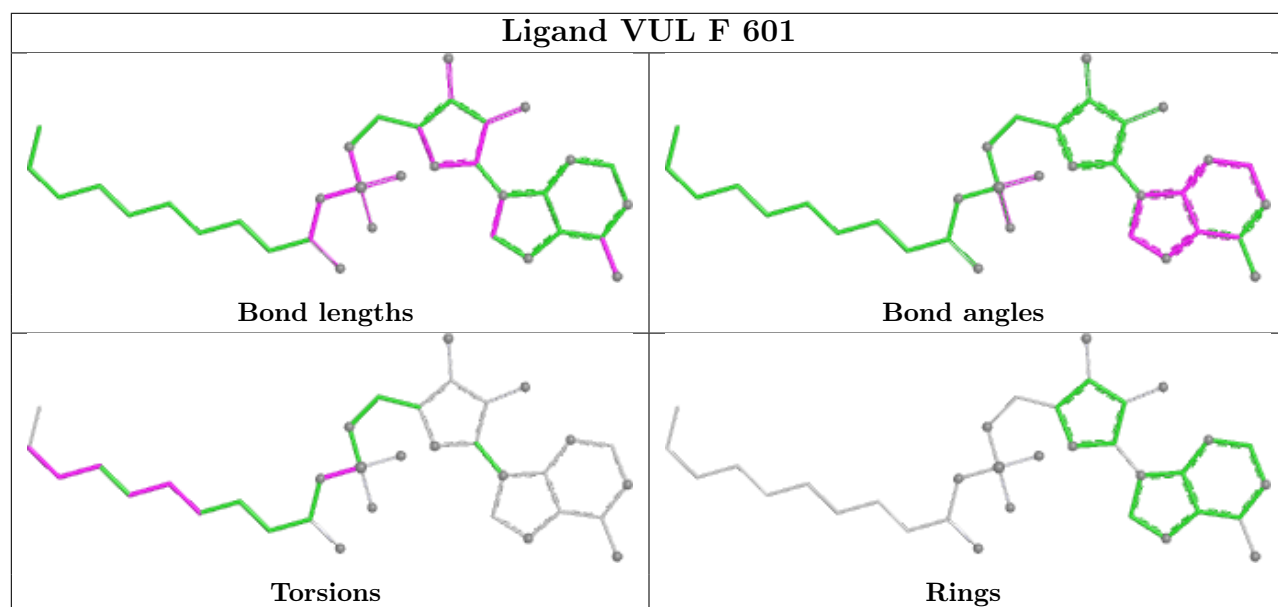
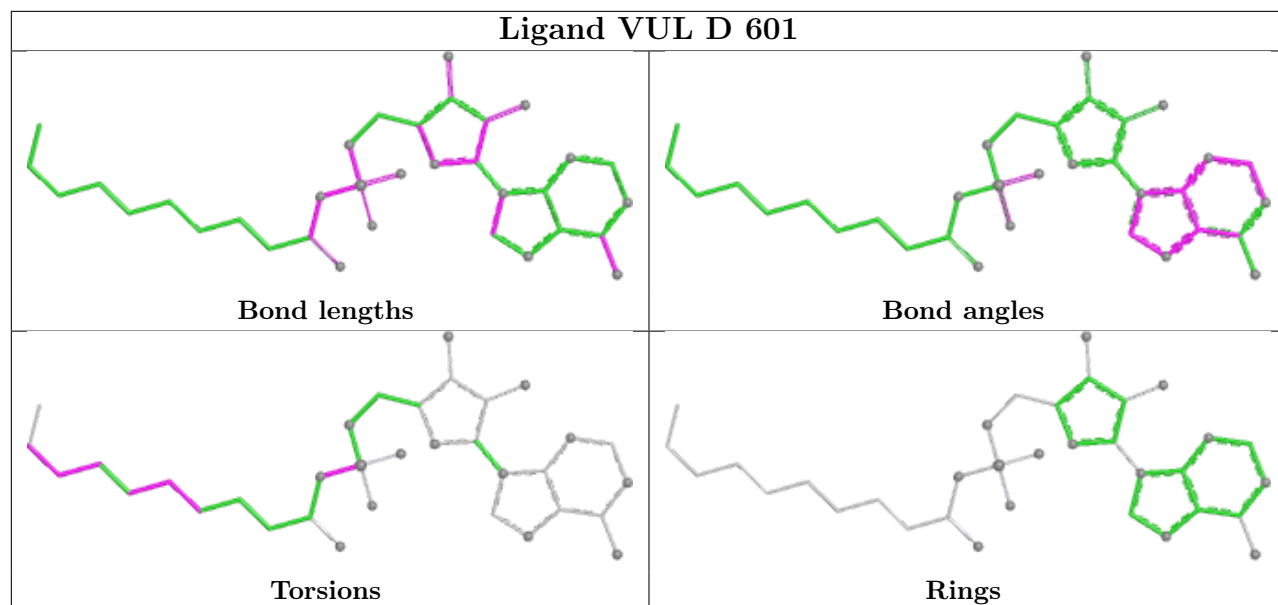
There are no ring outliers.

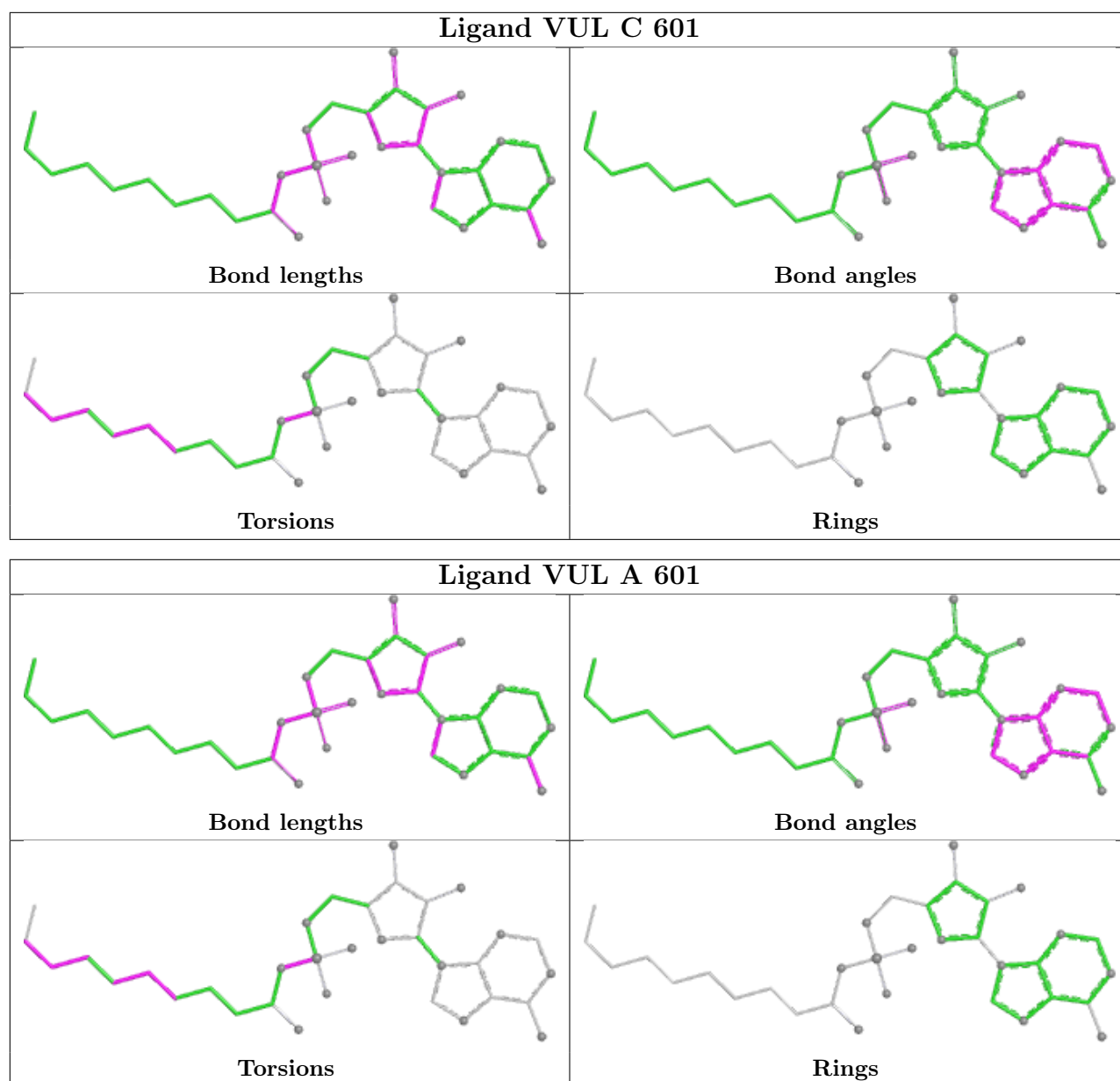
6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	VUL	1	0
2	E	601	VUL	1	0
2	D	601	VUL	1	0
2	F	601	VUL	1	0
2	C	601	VUL	1	0
2	A	601	VUL	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.

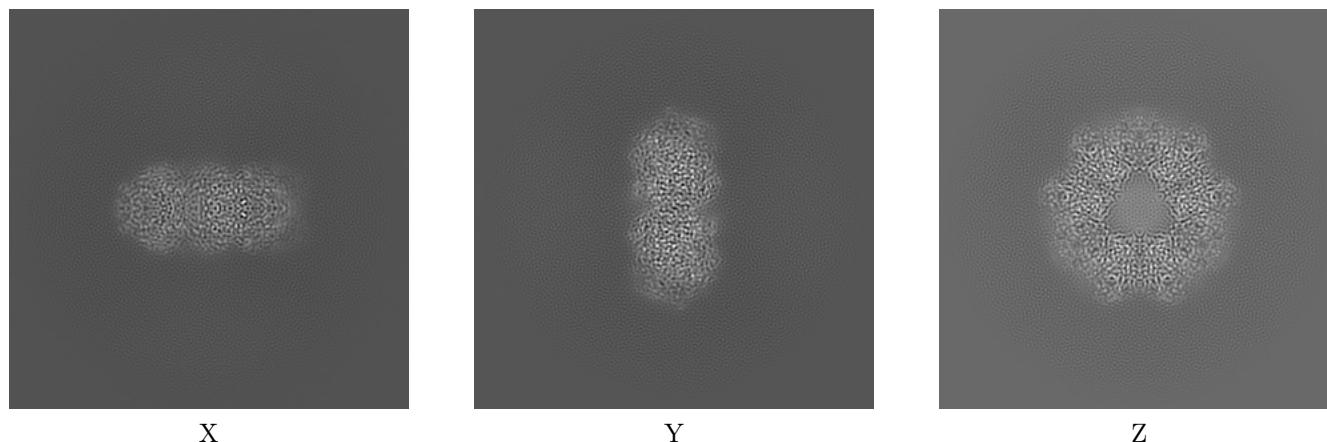
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-36725. These allow visual inspection of the internal detail of the map and identification of artifacts.

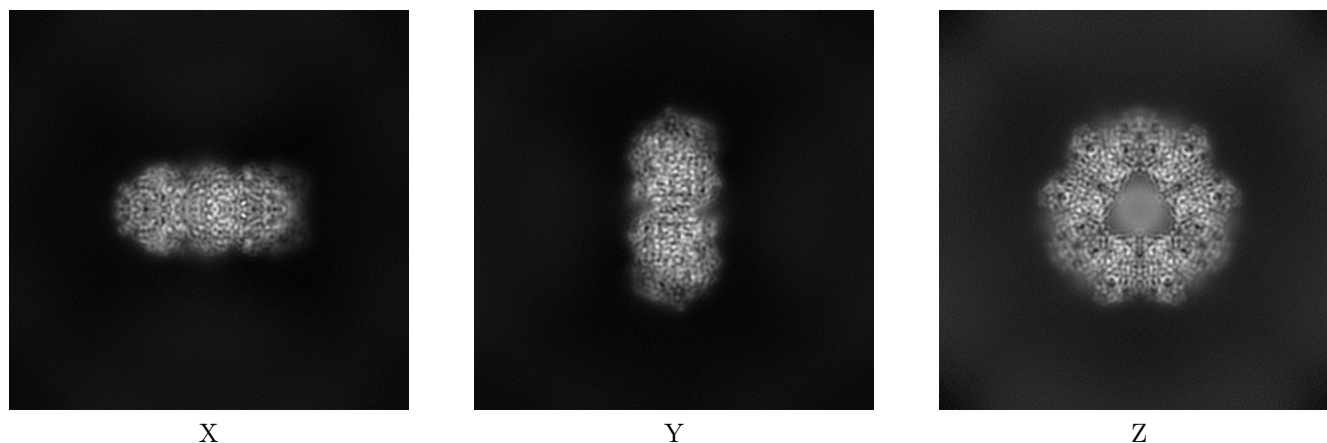
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

6.1.1 Primary map



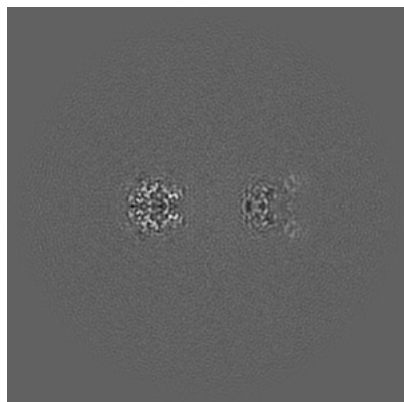
6.1.2 Raw map



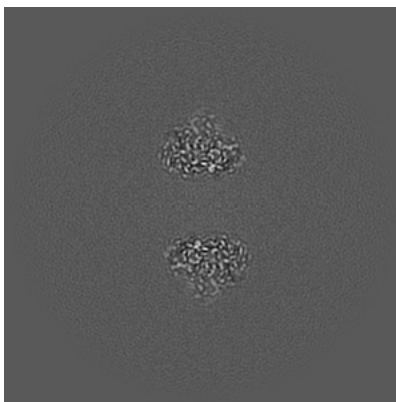
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

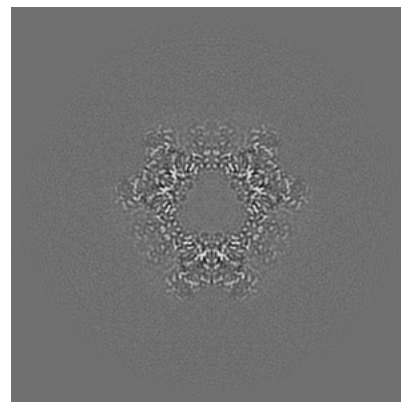
6.2.1 Primary map



X Index: 150

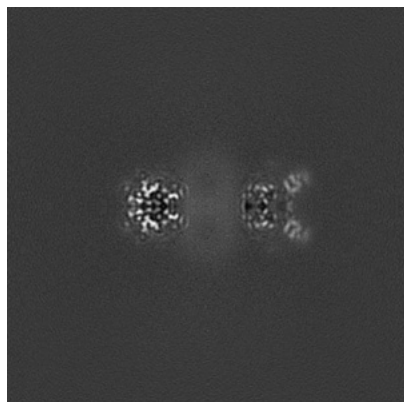


Y Index: 150

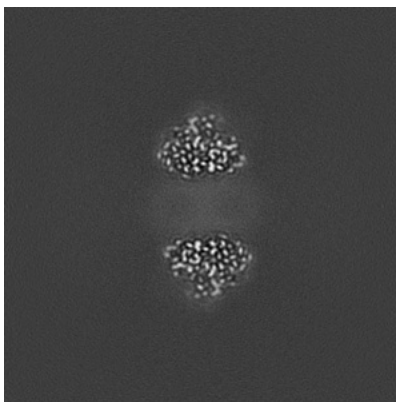


Z Index: 150

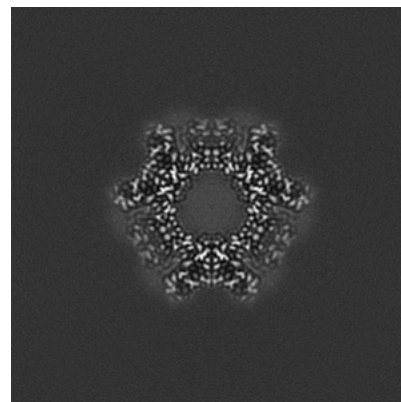
6.2.2 Raw map



X Index: 150



Y Index: 150

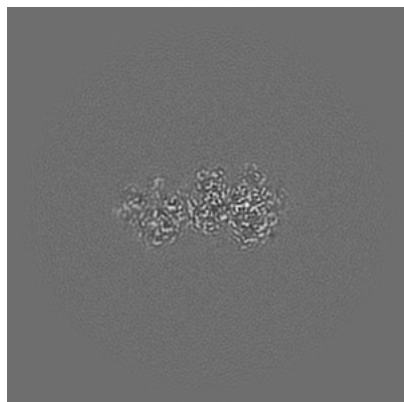


Z Index: 150

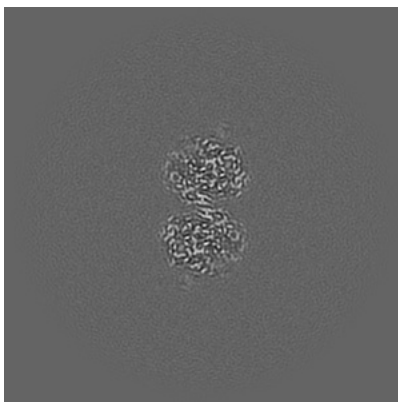
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

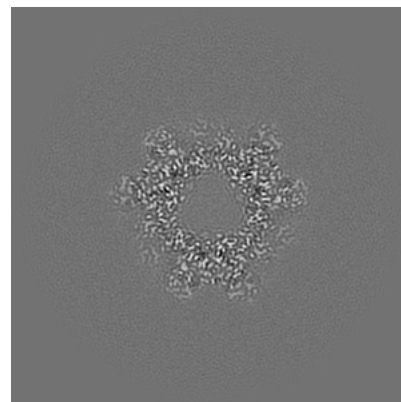
6.3.1 Primary map



X Index: 122

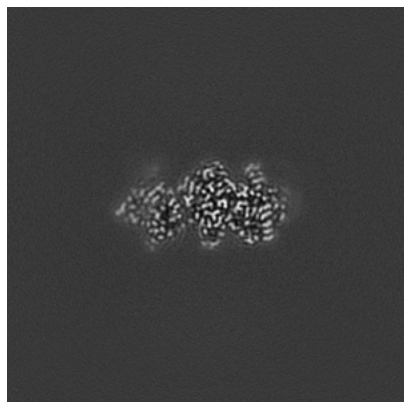


Y Index: 180

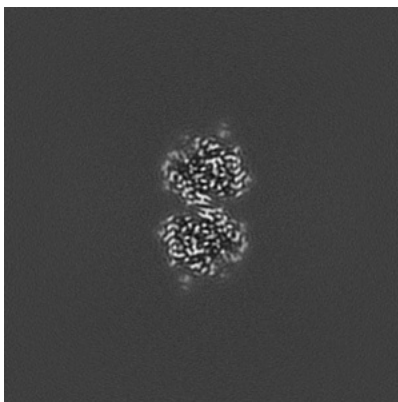


Z Index: 152

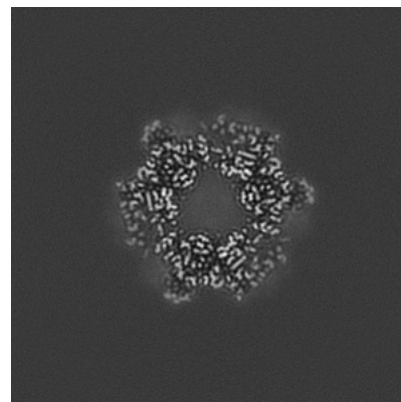
6.3.2 Raw map



X Index: 117



Y Index: 180

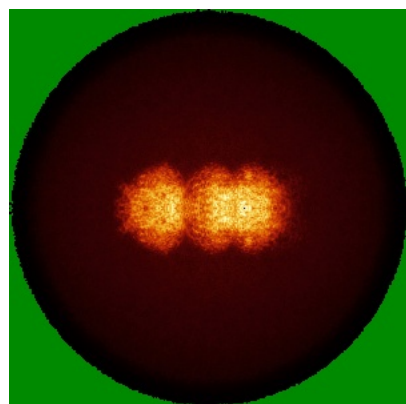


Z Index: 145

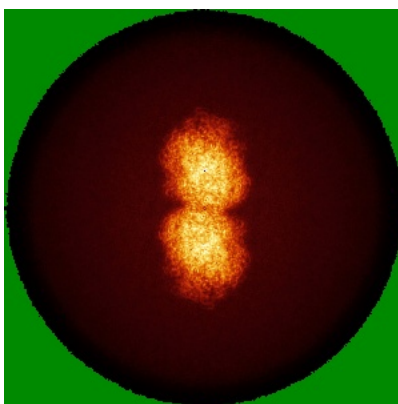
The images above show the largest variance slices of the map in three orthogonal directions.

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

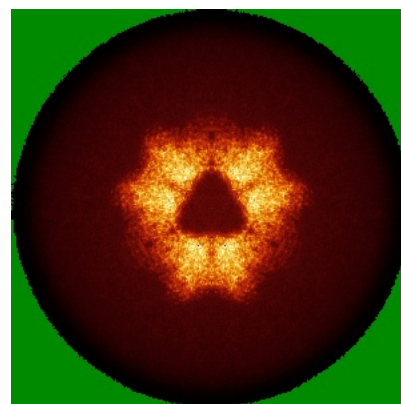
6.4.1 Primary map



X

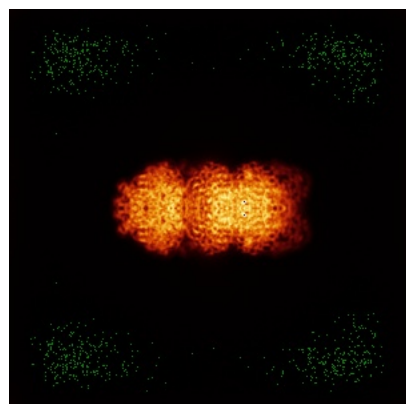


Y

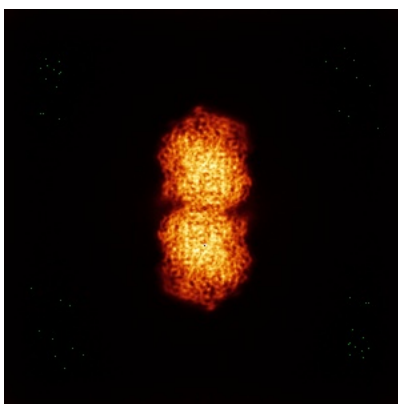


Z

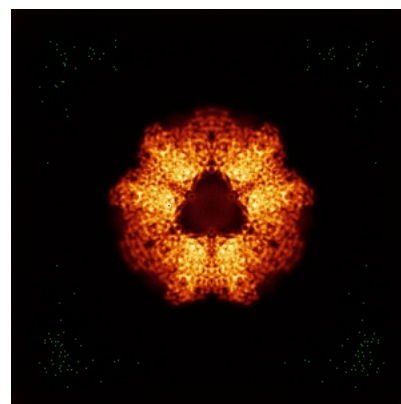
6.4.2 Raw map



X



Y

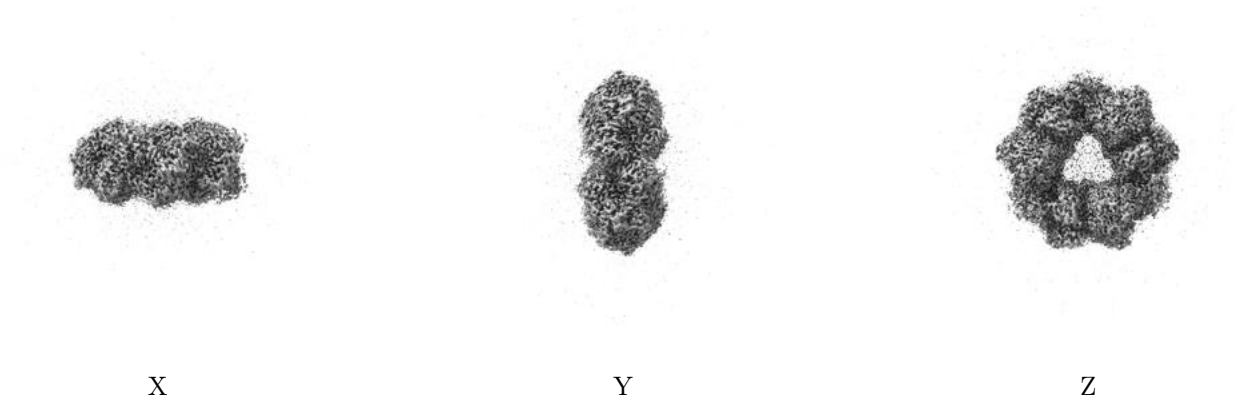


Z

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.

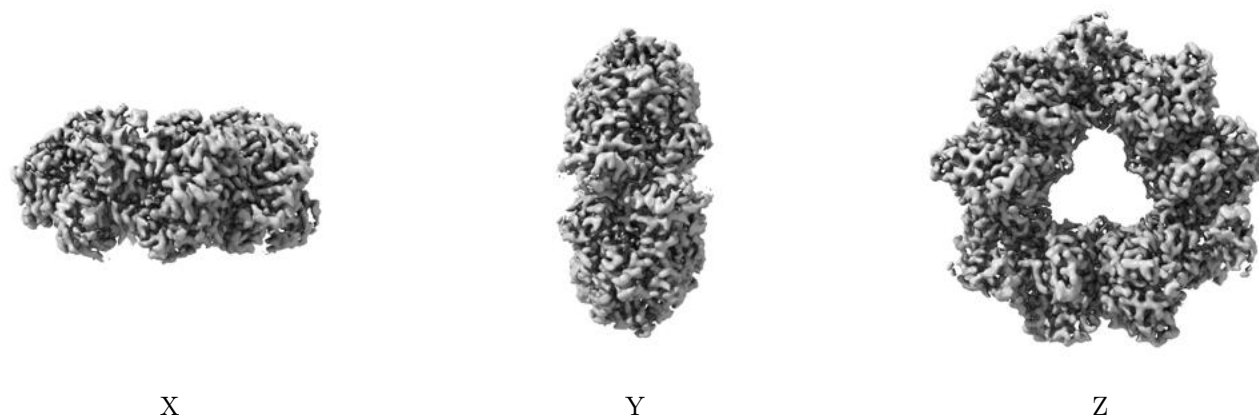
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.43. 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.

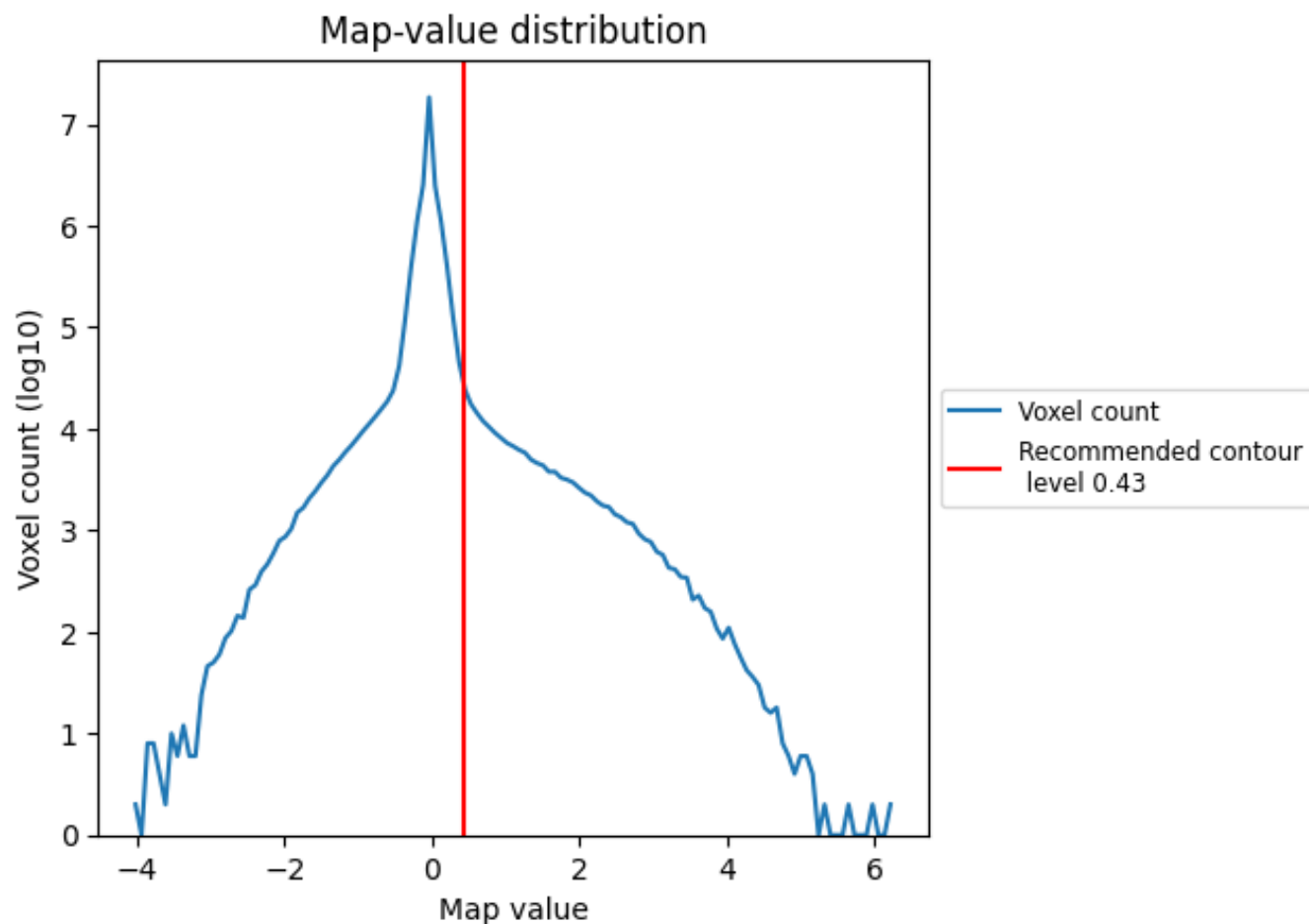
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

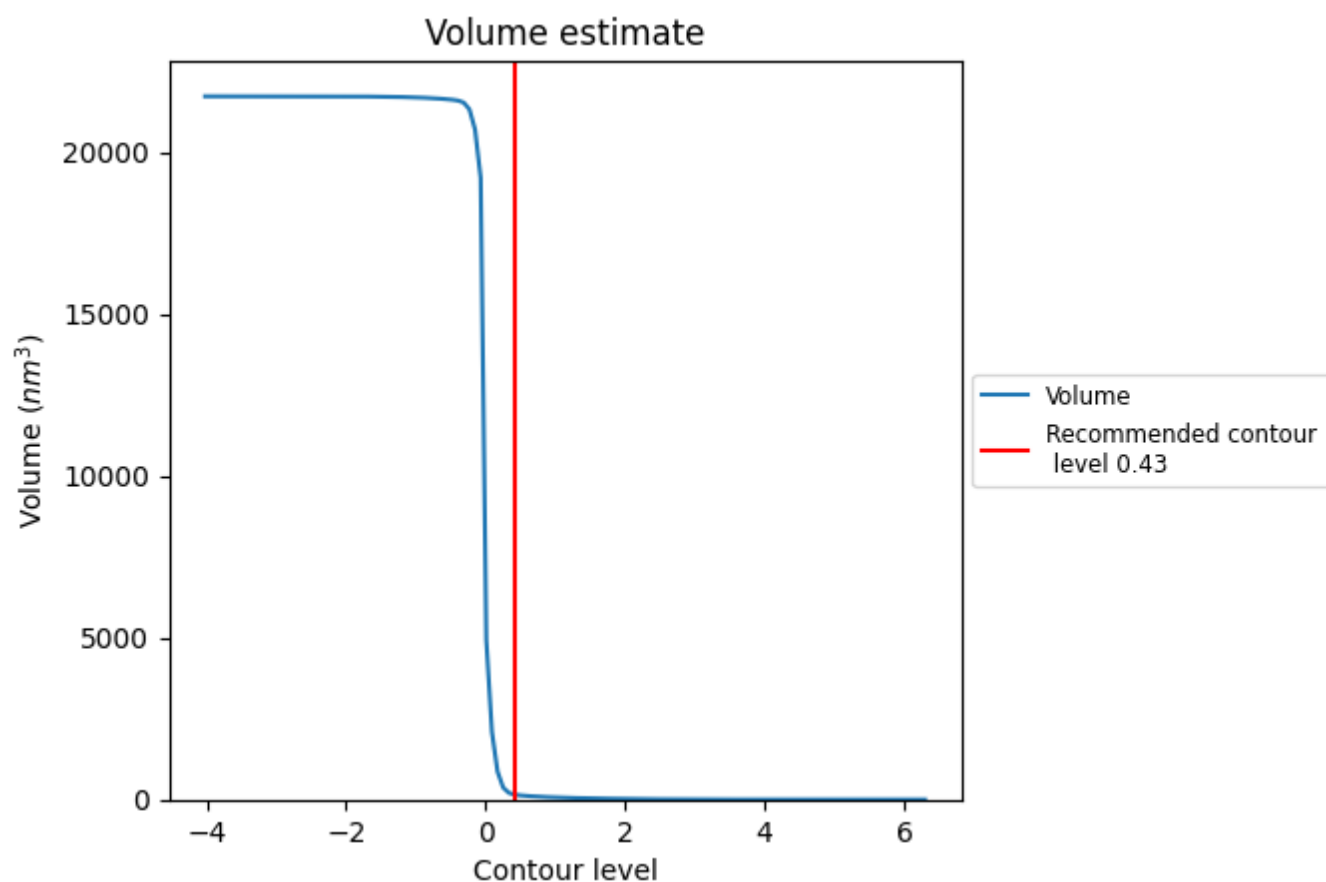
This section contains the results of statistical analysis of the map.

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.

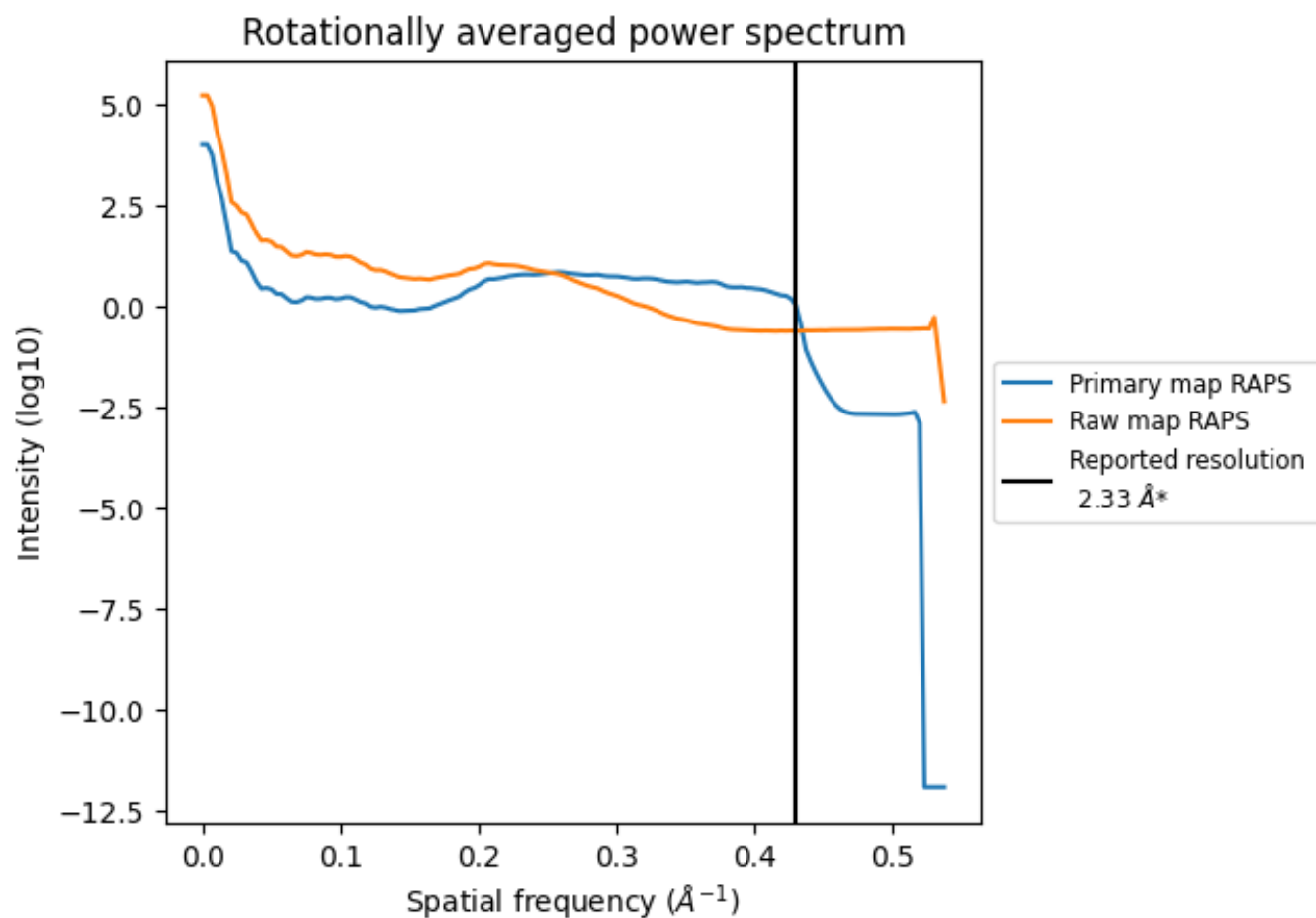
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 149 nm³; this corresponds to an approximate mass of 135 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 ⓘ

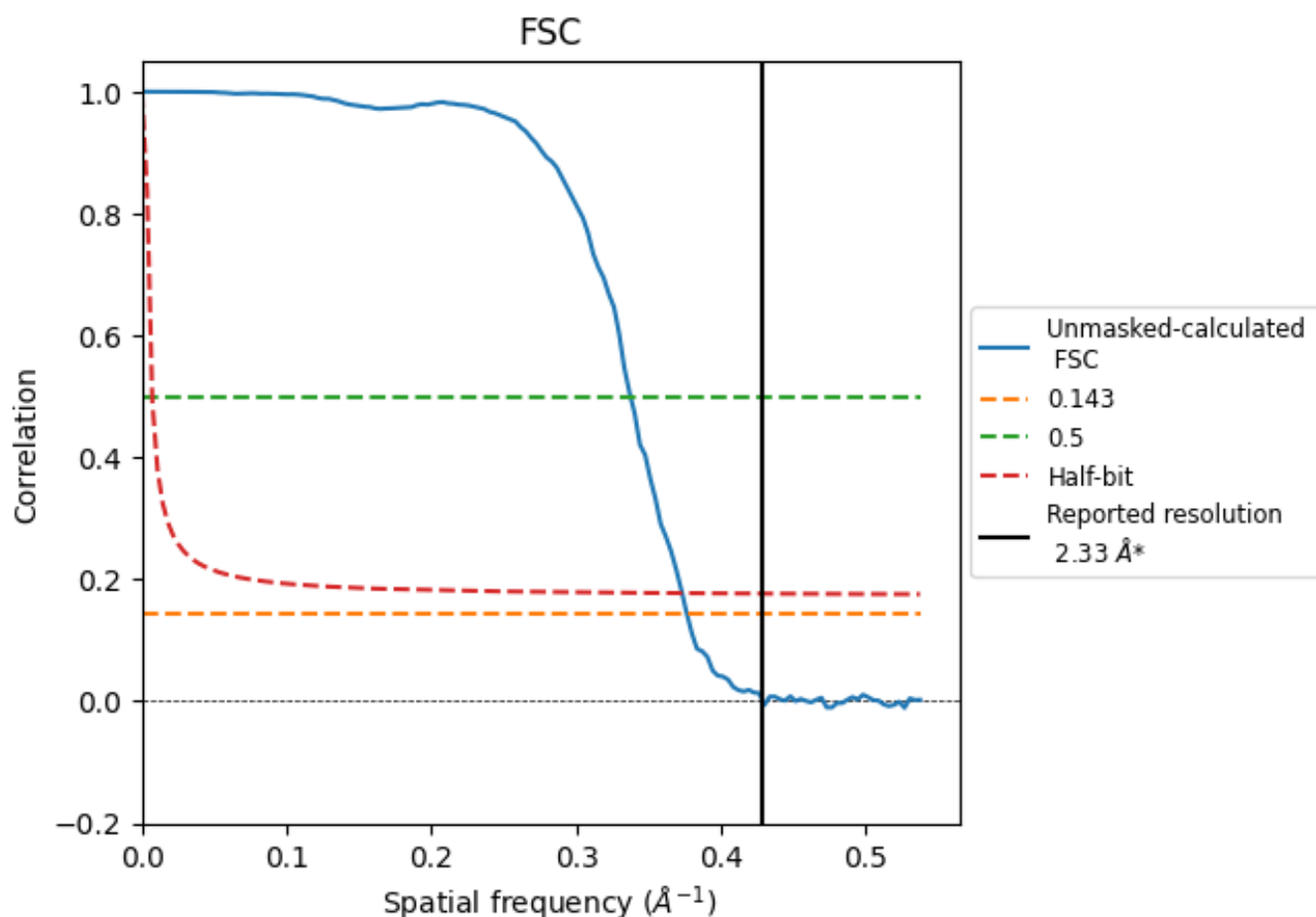


*Reported resolution corresponds to spatial frequency of 0.429 Å⁻¹

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 0.429 Å⁻¹

8.2 Resolution estimates [i](#)

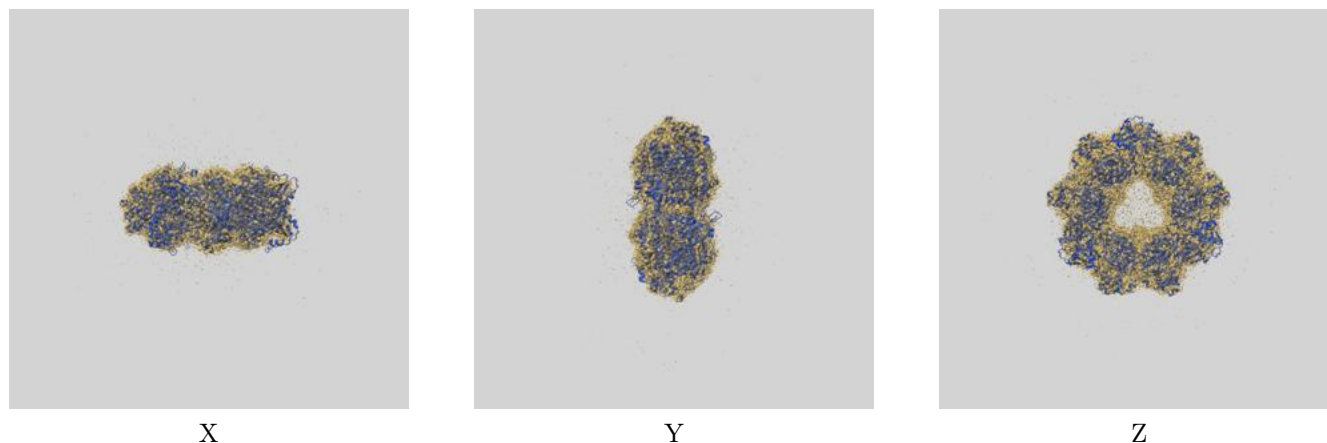
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.33	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	2.66	2.96	2.68

*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 2.66 differs from the reported value 2.33 by more than 10 %

9 Map-model fit [i](#)

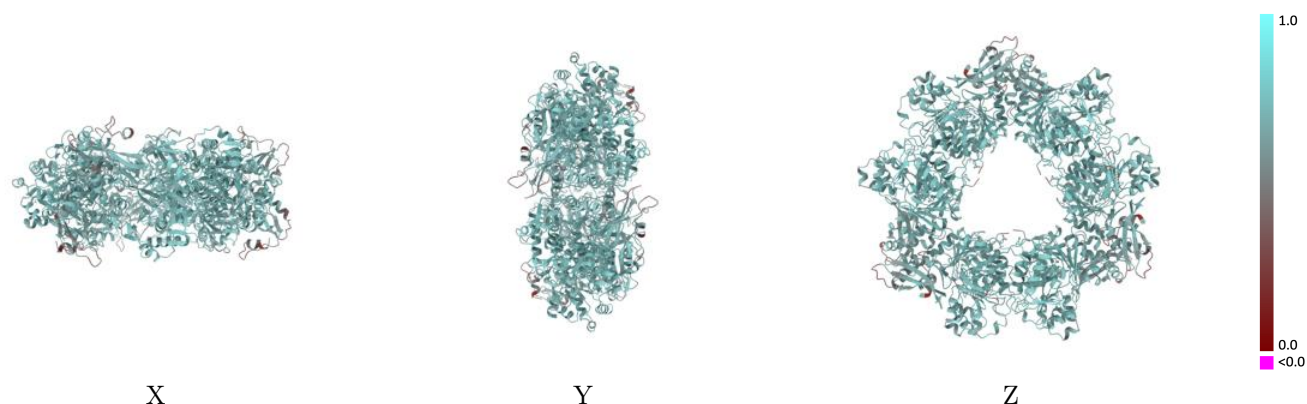
This section contains information regarding the fit between EMDB map EMD-36725 and PDB model 8JYL. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



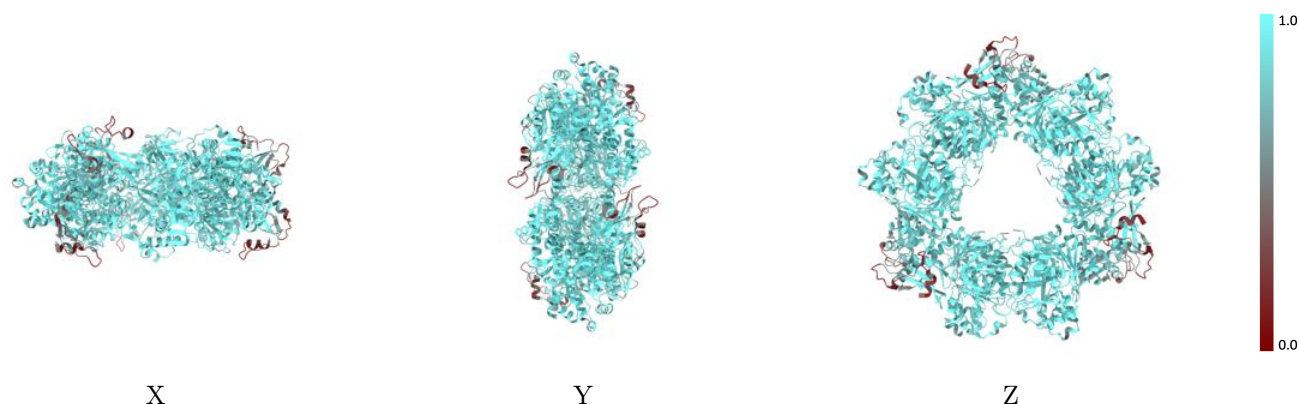
The images above show the 3D surface view of the map at the recommended contour level 0.43 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



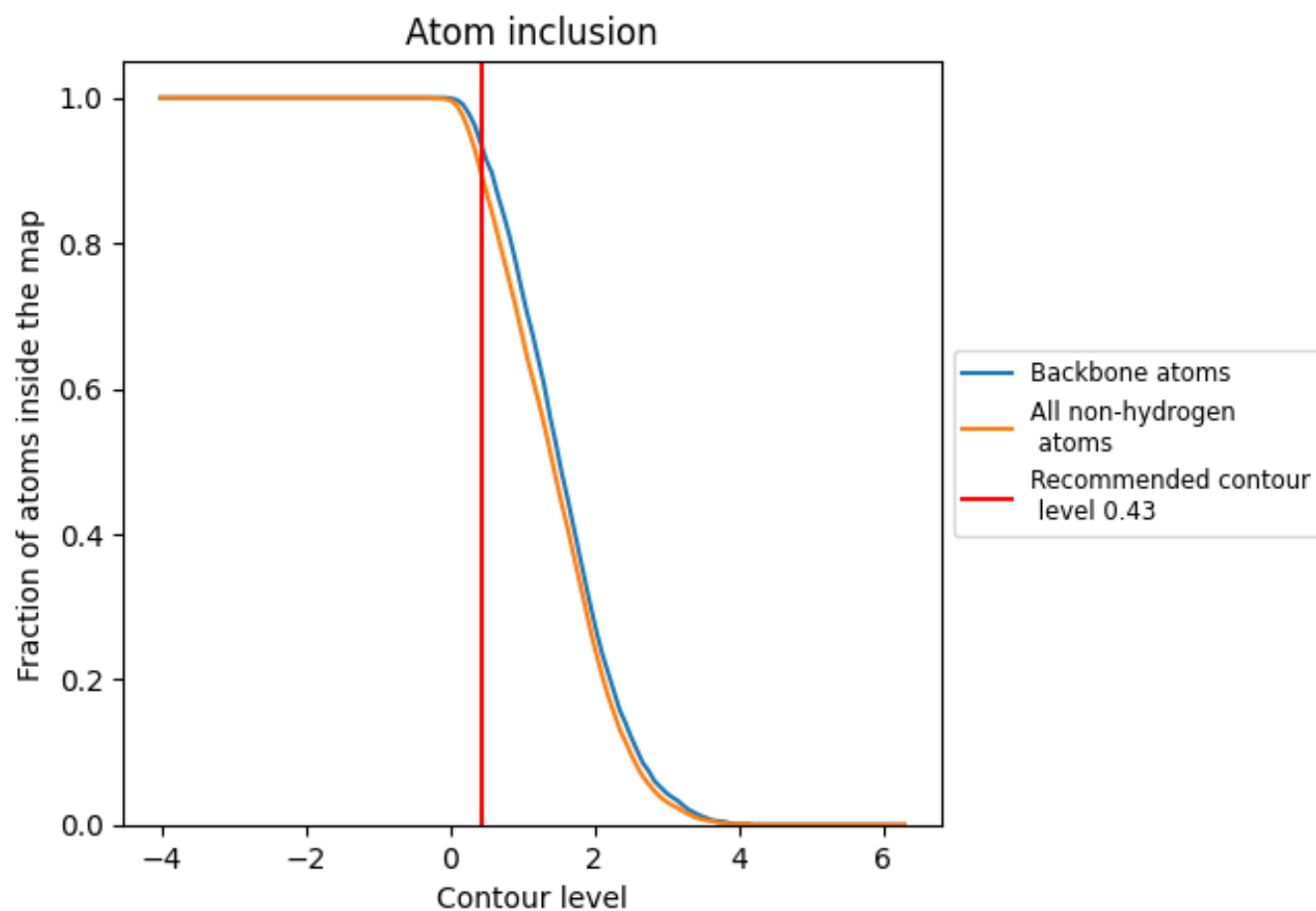
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.43).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 90% 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 (0.43) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8950	0.6770
A	0.8960	0.6780
B	0.8940	0.6760
C	0.8950	0.6760
D	0.8950	0.6770
E	0.8960	0.6780
F	0.8940	0.6760

