



Full wwPDB EM Validation Report ⓘ

Nov 20, 2022 – 09:32 AM EST

PDB ID : 7MCS
EMDB ID : EMD-23757
Title : Cryo-electron microscopy structure of TnsC(1-503)A225V bound to DNA
Authors : Shen, Y.; Ortega, J.; Guarne, A.
Deposited on : 2021-04-02
Resolution : 3.56 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:

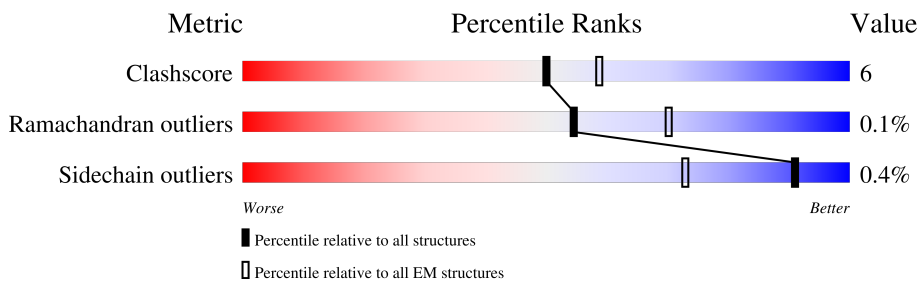
EMDB validation analysis : 0.0.1.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

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 3.56 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	523	
1	B	523	
1	C	523	
1	D	523	
1	E	523	
1	F	523	
1	G	523	
2	H	15	

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Mol	Chain	Length	Quality of chain
3	I	15	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a red segment on the left, a green segment in the middle, and a yellow segment on the right. The red segment is labeled '40%' above it. The green segment is labeled '40%' below it. The yellow segment is labeled '60%' below it.</p>

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 21600 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 Transposon Tn7 transposition protein TnsC.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	249	1967	1241	359	358	9	0	0
1	B	402	3196	2020	579	585	12	0	0
1	C	404	3208	2027	581	588	12	0	0
1	D	403	3201	2022	580	587	12	0	0
1	E	402	3196	2019	579	586	12	0	0
1	F	401	3183	2012	575	584	12	0	0
1	G	358	2842	1808	504	518	12	0	0

There are 154 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	SER	engineered mutation	UNP P05846
A	225	VAL	ALA	engineered mutation	UNP P05846
A	504	PRO	-	expression tag	UNP P05846
A	505	ASN	-	expression tag	UNP P05846
A	506	SER	-	expression tag	UNP P05846
A	507	SER	-	expression tag	UNP P05846
A	508	SER	-	expression tag	UNP P05846
A	509	VAL	-	expression tag	UNP P05846
A	510	ASP	-	expression tag	UNP P05846
A	511	LYS	-	expression tag	UNP P05846
A	512	LEU	-	expression tag	UNP P05846
A	513	ALA	-	expression tag	UNP P05846
A	514	ALA	-	expression tag	UNP P05846
A	515	ALA	-	expression tag	UNP P05846
A	516	LEU	-	expression tag	UNP P05846
A	517	GLU	-	expression tag	UNP P05846

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Chain	Residue	Modelled	Actual	Comment	Reference
A	518	HIS	-	expression tag	UNP P05846
A	519	HIS	-	expression tag	UNP P05846
A	520	HIS	-	expression tag	UNP P05846
A	521	HIS	-	expression tag	UNP P05846
A	522	HIS	-	expression tag	UNP P05846
A	523	HIS	-	expression tag	UNP P05846
B	2	GLY	SER	engineered mutation	UNP P05846
B	225	VAL	ALA	engineered mutation	UNP P05846
B	504	PRO	-	expression tag	UNP P05846
B	505	ASN	-	expression tag	UNP P05846
B	506	SER	-	expression tag	UNP P05846
B	507	SER	-	expression tag	UNP P05846
B	508	SER	-	expression tag	UNP P05846
B	509	VAL	-	expression tag	UNP P05846
B	510	ASP	-	expression tag	UNP P05846
B	511	LYS	-	expression tag	UNP P05846
B	512	LEU	-	expression tag	UNP P05846
B	513	ALA	-	expression tag	UNP P05846
B	514	ALA	-	expression tag	UNP P05846
B	515	ALA	-	expression tag	UNP P05846
B	516	LEU	-	expression tag	UNP P05846
B	517	GLU	-	expression tag	UNP P05846
B	518	HIS	-	expression tag	UNP P05846
B	519	HIS	-	expression tag	UNP P05846
B	520	HIS	-	expression tag	UNP P05846
B	521	HIS	-	expression tag	UNP P05846
B	522	HIS	-	expression tag	UNP P05846
B	523	HIS	-	expression tag	UNP P05846
C	2	GLY	SER	engineered mutation	UNP P05846
C	225	VAL	ALA	engineered mutation	UNP P05846
C	504	PRO	-	expression tag	UNP P05846
C	505	ASN	-	expression tag	UNP P05846
C	506	SER	-	expression tag	UNP P05846
C	507	SER	-	expression tag	UNP P05846
C	508	SER	-	expression tag	UNP P05846
C	509	VAL	-	expression tag	UNP P05846
C	510	ASP	-	expression tag	UNP P05846
C	511	LYS	-	expression tag	UNP P05846
C	512	LEU	-	expression tag	UNP P05846
C	513	ALA	-	expression tag	UNP P05846
C	514	ALA	-	expression tag	UNP P05846
C	515	ALA	-	expression tag	UNP P05846

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Chain	Residue	Modelled	Actual	Comment	Reference
C	516	LEU	-	expression tag	UNP P05846
C	517	GLU	-	expression tag	UNP P05846
C	518	HIS	-	expression tag	UNP P05846
C	519	HIS	-	expression tag	UNP P05846
C	520	HIS	-	expression tag	UNP P05846
C	521	HIS	-	expression tag	UNP P05846
C	522	HIS	-	expression tag	UNP P05846
C	523	HIS	-	expression tag	UNP P05846
D	2	GLY	SER	engineered mutation	UNP P05846
D	225	VAL	ALA	engineered mutation	UNP P05846
D	504	PRO	-	expression tag	UNP P05846
D	505	ASN	-	expression tag	UNP P05846
D	506	SER	-	expression tag	UNP P05846
D	507	SER	-	expression tag	UNP P05846
D	508	SER	-	expression tag	UNP P05846
D	509	VAL	-	expression tag	UNP P05846
D	510	ASP	-	expression tag	UNP P05846
D	511	LYS	-	expression tag	UNP P05846
D	512	LEU	-	expression tag	UNP P05846
D	513	ALA	-	expression tag	UNP P05846
D	514	ALA	-	expression tag	UNP P05846
D	515	ALA	-	expression tag	UNP P05846
D	516	LEU	-	expression tag	UNP P05846
D	517	GLU	-	expression tag	UNP P05846
D	518	HIS	-	expression tag	UNP P05846
D	519	HIS	-	expression tag	UNP P05846
D	520	HIS	-	expression tag	UNP P05846
D	521	HIS	-	expression tag	UNP P05846
D	522	HIS	-	expression tag	UNP P05846
D	523	HIS	-	expression tag	UNP P05846
E	2	GLY	SER	engineered mutation	UNP P05846
E	225	VAL	ALA	engineered mutation	UNP P05846
E	504	PRO	-	expression tag	UNP P05846
E	505	ASN	-	expression tag	UNP P05846
E	506	SER	-	expression tag	UNP P05846
E	507	SER	-	expression tag	UNP P05846
E	508	SER	-	expression tag	UNP P05846
E	509	VAL	-	expression tag	UNP P05846
E	510	ASP	-	expression tag	UNP P05846
E	511	LYS	-	expression tag	UNP P05846
E	512	LEU	-	expression tag	UNP P05846
E	513	ALA	-	expression tag	UNP P05846

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Chain	Residue	Modelled	Actual	Comment	Reference
E	514	ALA	-	expression tag	UNP P05846
E	515	ALA	-	expression tag	UNP P05846
E	516	LEU	-	expression tag	UNP P05846
E	517	GLU	-	expression tag	UNP P05846
E	518	HIS	-	expression tag	UNP P05846
E	519	HIS	-	expression tag	UNP P05846
E	520	HIS	-	expression tag	UNP P05846
E	521	HIS	-	expression tag	UNP P05846
E	522	HIS	-	expression tag	UNP P05846
E	523	HIS	-	expression tag	UNP P05846
F	2	GLY	SER	engineered mutation	UNP P05846
F	225	VAL	ALA	engineered mutation	UNP P05846
F	504	PRO	-	expression tag	UNP P05846
F	505	ASN	-	expression tag	UNP P05846
F	506	SER	-	expression tag	UNP P05846
F	507	SER	-	expression tag	UNP P05846
F	508	SER	-	expression tag	UNP P05846
F	509	VAL	-	expression tag	UNP P05846
F	510	ASP	-	expression tag	UNP P05846
F	511	LYS	-	expression tag	UNP P05846
F	512	LEU	-	expression tag	UNP P05846
F	513	ALA	-	expression tag	UNP P05846
F	514	ALA	-	expression tag	UNP P05846
F	515	ALA	-	expression tag	UNP P05846
F	516	LEU	-	expression tag	UNP P05846
F	517	GLU	-	expression tag	UNP P05846
F	518	HIS	-	expression tag	UNP P05846
F	519	HIS	-	expression tag	UNP P05846
F	520	HIS	-	expression tag	UNP P05846
F	521	HIS	-	expression tag	UNP P05846
F	522	HIS	-	expression tag	UNP P05846
F	523	HIS	-	expression tag	UNP P05846
G	2	GLY	SER	engineered mutation	UNP P05846
G	225	VAL	ALA	engineered mutation	UNP P05846
G	504	PRO	-	expression tag	UNP P05846
G	505	ASN	-	expression tag	UNP P05846
G	506	SER	-	expression tag	UNP P05846
G	507	SER	-	expression tag	UNP P05846
G	508	SER	-	expression tag	UNP P05846
G	509	VAL	-	expression tag	UNP P05846
G	510	ASP	-	expression tag	UNP P05846
G	511	LYS	-	expression tag	UNP P05846

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Chain	Residue	Modelled	Actual	Comment	Reference
G	512	LEU	-	expression tag	UNP P05846
G	513	ALA	-	expression tag	UNP P05846
G	514	ALA	-	expression tag	UNP P05846
G	515	ALA	-	expression tag	UNP P05846
G	516	LEU	-	expression tag	UNP P05846
G	517	GLU	-	expression tag	UNP P05846
G	518	HIS	-	expression tag	UNP P05846
G	519	HIS	-	expression tag	UNP P05846
G	520	HIS	-	expression tag	UNP P05846
G	521	HIS	-	expression tag	UNP P05846
G	522	HIS	-	expression tag	UNP P05846
G	523	HIS	-	expression tag	UNP P05846

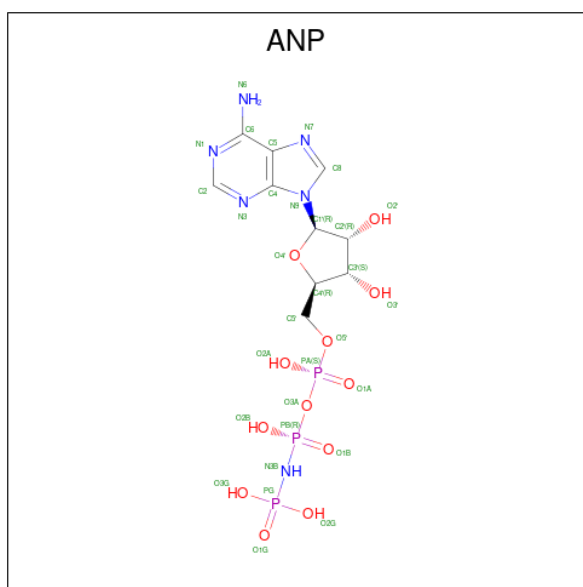
- Molecule 2 is a DNA chain called DNA (5'-D(P*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*C)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	H	15	306	142	59	90	15	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(P*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*G)-3').

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
3	I	15	309	143	61	90	15	0	0

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
4	B	1	Total	C	N	O	P	0
			31	10	6	12	3	
4	C	1	Total	C	N	O	P	0
			31	10	6	12	3	
4	D	1	Total	C	N	O	P	0
			31	10	6	12	3	
4	E	1	Total	C	N	O	P	0
			31	10	6	12	3	
4	F	1	Total	C	N	O	P	0
			31	10	6	12	3	
4	G	1	Total	C	N	O	P	0
			31	10	6	12	3	

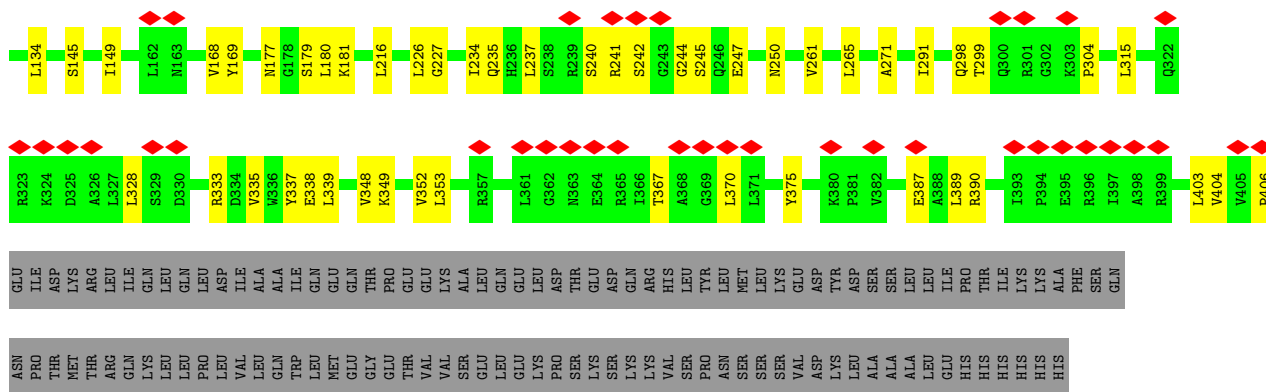
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

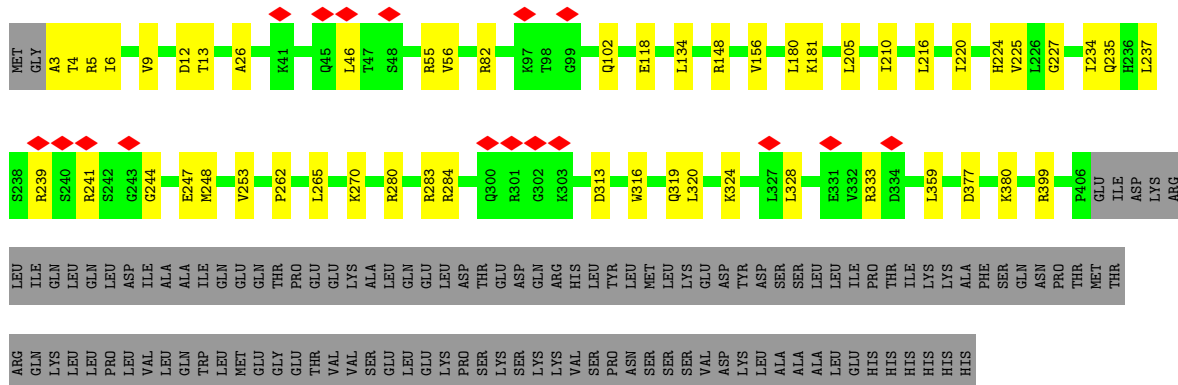
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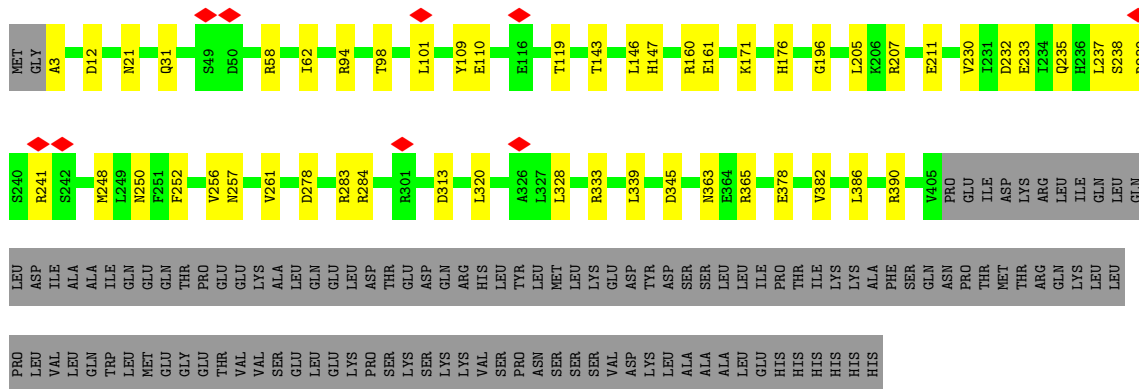
Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
5	G	1	1	1	0



• Molecule 1: Transposon Tn7 transposition protein TnsC

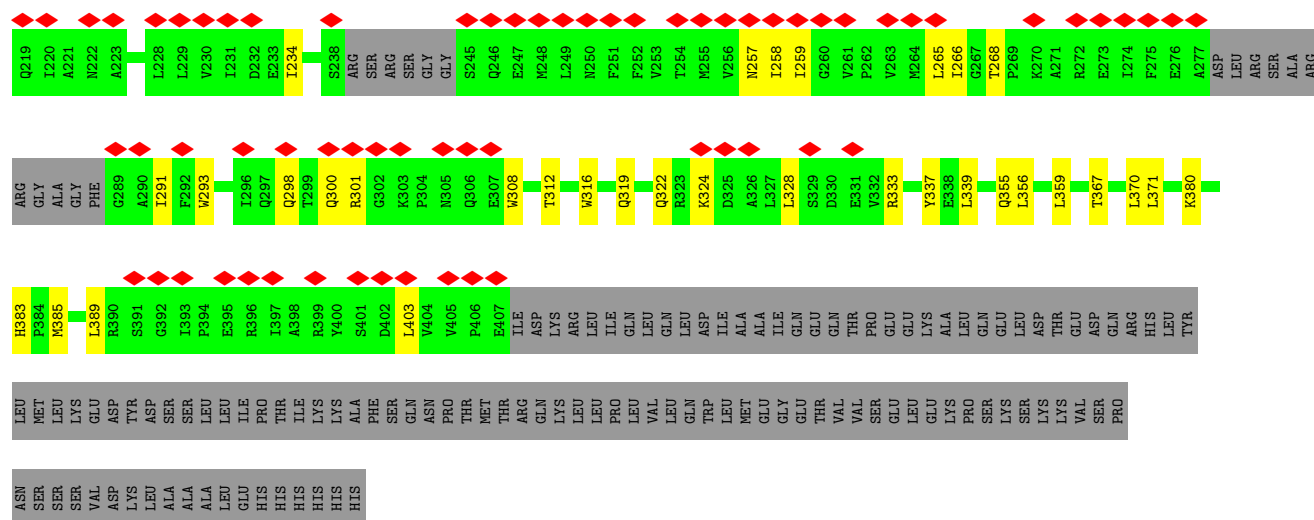


• Molecule 1: Transposon Tn7 transposition protein TnsC



• Molecule 1: Transposon Tn7 transposition protein TnsC

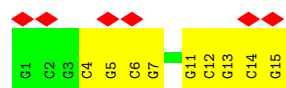




● Molecule 2: DNA (5'-D(P*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*C)-3')



● Molecule 3: DNA (5'-D(P*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*GP*CP*G)-3')



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	155988	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 ($e^-/\text{\AA}^2$)	78	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2750	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.105	Depositor
Minimum map value	-0.058	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.0145	Depositor
Map size (Å)	256.5, 256.5, 256.5	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.855, 0.855, 0.855	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, ANP

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	0.27	0/2002	0.57	0/2701
1	B	0.27	0/3254	0.57	1/4402 (0.0%)
1	C	0.31	0/3266	0.60	1/4419 (0.0%)
1	D	0.28	0/3258	0.56	0/4407
1	E	0.27	0/3253	0.56	0/4400
1	F	0.26	0/3240	0.56	0/4383
1	G	0.27	0/2892	0.60	2/3917 (0.1%)
2	H	0.58	0/342	0.87	0/525
3	I	0.54	0/346	0.80	0/532
All	All	0.29	0/21853	0.58	4/29686 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	G	22	PRO	CA-N-CD	-8.39	99.75	111.50
1	G	173	ASP	CB-CG-OD1	6.00	123.70	118.30
1	C	134	LEU	CA-CB-CG	5.62	128.23	115.30
1	B	101	LEU	CA-CB-CG	5.15	127.16	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1967	0	1977	26	0
1	B	3196	0	3235	43	0
1	C	3208	0	3247	34	0
1	D	3201	0	3240	42	0
1	E	3196	0	3235	43	0
1	F	3183	0	3217	44	0
1	G	2842	0	2892	52	0
2	H	306	0	166	7	0
3	I	309	0	166	9	0
4	B	31	0	13	1	0
4	C	31	0	13	1	0
4	D	31	0	13	0	0
4	E	31	0	13	3	0
4	F	31	0	13	0	0
4	G	31	0	13	4	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
All	All	21600	0	21453	260	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 6.

All (260) 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:A:157:ILE:HD11	1:A:166:GLN:HE21	1.45	0.81
1:B:387:GLU:HG2	1:B:390:ARG:HH22	1.50	0.77
1:C:181:LYS:HD3	1:C:210:ILE:HG13	1.69	0.75
1:G:141:GLY:HA3	4:G:601:ANP:H8	1.70	0.72
1:E:283:ARG:NH1	1:F:345:ASP:OD2	2.25	0.70
1:C:55:ARG:NH1	1:C:359:LEU:O	2.26	0.69
1:G:380:LYS:HA	1:G:383:HIS:HD2	1.58	0.68
1:F:346:ILE:HA	1:F:349:LYS:HG2	1.74	0.67
1:F:87:ILE:HG12	1:F:228:LEU:HD22	1.77	0.66
1:E:54:SER:HB3	1:E:57:ILE:HD12	1.78	0.66
1:G:312:THR:HG21	1:G:333:ARG:HG3	1.77	0.64
1:G:234:ILE:HG12	1:G:265:LEU:HD12	1.80	0.64
1:B:102:GLN:NE2	1:C:4:THR:O	2.29	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:237:LEU:HD22	1:B:245:SER:HA	1.81	0.62
1:E:28:PRO:O	1:E:88:ARG:NH2	2.33	0.62
1:D:62:ILE:HD11	1:D:320:LEU:HB2	1.82	0.61
1:F:357:ARG:HE	1:F:361:LEU:HD13	1.66	0.61
1:E:258:ILE:HD11	1:F:173:ASP:HB2	1.83	0.61
1:A:134:LEU:HD23	1:A:266:ILE:HG12	1.84	0.60
1:D:207:ARG:HH22	3:I:11:DG:H4'	1.66	0.60
1:D:233:GLU:OE1	1:D:235:GLN:NE2	2.34	0.60
1:F:156:VAL:HG11	1:F:195:LEU:HB2	1.82	0.60
1:D:313:ASP:OD1	1:D:333:ARG:NH2	2.34	0.60
4:B:601:ANP:O3G	4:B:601:ANP:O1B	2.19	0.60
1:C:118:GLU:O	1:D:31:GLN:NE2	2.35	0.59
1:D:207:ARG:HH21	3:I:12:DC:H5'	1.67	0.59
1:E:241:ARG:NH1	3:I:5:DG:OP2	2.34	0.59
1:B:126:ARG:HG2	1:B:127:SER:H	1.67	0.59
1:E:55:ARG:HH21	1:E:360:ALA:HA	1.68	0.59
1:D:109:TYR:OH	1:E:26:ALA:O	2.18	0.58
3:I:12:DC:H2''	3:I:13:DG:C8	2.38	0.58
1:C:82:ARG:NH2	1:D:378:GLU:OE2	2.37	0.58
1:G:142:LYS:HE2	1:G:266:ILE:HB	1.84	0.58
1:F:109:TYR:HE2	1:G:154:PRO:HG3	1.68	0.57
1:A:211:GLU:HG2	1:B:181:LYS:HG2	1.86	0.57
1:D:58:ARG:HD2	1:D:320:LEU:HD23	1.86	0.57
1:B:241:ARG:HG3	1:B:242:SER:H	1.68	0.57
1:D:257:ASN:OD1	1:D:284:ARG:NH2	2.38	0.57
1:C:3:ALA:O	1:C:5:ARG:NH2	2.38	0.56
1:F:226:LEU:HD23	1:F:261:VAL:HG11	1.87	0.56
1:F:146:LEU:HD22	1:F:266:ILE:HD11	1.86	0.56
1:G:4:THR:HG21	1:G:156:VAL:HG13	1.85	0.56
1:F:118:GLU:OE1	1:G:31:GLN:NE2	2.38	0.56
1:C:4:THR:HG22	1:C:156:VAL:HG13	1.88	0.55
1:A:235:GLN:OE1	1:A:268:THR:N	2.38	0.55
1:G:35:ASN:O	1:G:39:SER:N	2.38	0.55
1:A:82:ARG:O	1:A:86:MET:HG3	2.07	0.55
1:B:234:ILE:HD12	1:B:265:LEU:HD22	1.89	0.55
1:F:238:SER:OG	1:F:270:LYS:NZ	2.40	0.55
1:G:298:GLN:HA	1:G:308:TRP:HZ3	1.71	0.55
1:D:160:ARG:NH1	1:D:161:GLU:OE2	2.40	0.55
1:G:134:LEU:HB3	1:G:266:ILE:HG22	1.89	0.55
1:A:253:VAL:HG11	1:A:280:ARG:HH21	1.73	0.54
1:E:58:ARG:NH2	1:E:364:GLU:OE2	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:257:ASN:OD1	1:G:258:ILE:N	2.41	0.53
1:F:308:TRP:HZ3	1:F:333:ARG:HD3	1.73	0.53
1:E:109:TYR:OH	1:F:26:ALA:O	2.19	0.53
1:B:389:LEU:HD11	1:B:403:LEU:HD13	1.90	0.53
1:E:181:LYS:HG3	1:E:213:MET:HE1	1.91	0.53
1:B:315:LEU:HD11	1:B:348:VAL:HG12	1.91	0.53
1:G:258:ILE:HG22	1:G:259:ILE:HG23	1.92	0.52
1:F:46:LEU:HD11	1:F:318:LEU:HD23	1.91	0.52
1:B:30:LEU:HD23	1:C:56:VAL:HG12	1.91	0.52
1:B:113:GLN:HG2	1:C:9:VAL:HG23	1.92	0.52
1:D:207:ARG:NH2	3:I:12:DC:H5'	2.24	0.52
1:F:62:ILE:HD11	1:F:356:LEU:HD13	1.91	0.52
1:F:328:LEU:HD22	1:F:333:ARG:CZ	2.39	0.52
1:B:235:GLN:HA	1:B:271:ALA:HB2	1.91	0.52
1:C:4:THR:HG23	1:C:6:ILE:HG12	1.91	0.52
1:F:15:VAL:HG22	1:G:57:ILE:HG23	1.92	0.52
1:A:236:HIS:HA	1:A:239:ARG:HB3	1.92	0.51
1:C:283:ARG:NE	1:D:345:ASP:OD2	2.35	0.51
1:G:188:PHE:HB3	1:G:200:GLU:HB3	1.92	0.51
1:E:62:ILE:HD12	1:E:355:GLN:HB3	1.93	0.51
1:E:226:LEU:HD23	1:E:261:VAL:HG11	1.92	0.51
1:D:283:ARG:NE	1:E:345:ASP:OD2	2.31	0.51
1:G:38:ALA:HA	1:G:41:LYS:HE3	1.91	0.51
1:C:224:HIS:O	1:C:225:VAL:C	2.49	0.51
1:F:109:TYR:O	1:F:113:GLN:NE2	2.44	0.51
1:F:46:LEU:HD11	1:F:318:LEU:HA	1.94	0.50
1:E:330:ASP:HA	1:E:333:ARG:HG2	1.93	0.50
1:B:53:LYS:O	1:B:58:ARG:NH1	2.45	0.50
2:H:11:DC:H2''	2:H:12:DG:C8	2.47	0.50
1:D:238:SER:HB3	1:D:241:ARG:HH21	1.77	0.50
1:B:244:GLY:HA2	1:B:247:GLU:HG2	1.94	0.49
1:A:15:VAL:HG11	1:A:18:TYR:HD2	1.77	0.49
1:B:177:ASN:O	1:B:240:SER:OG	2.25	0.49
1:D:339:LEU:HA	1:D:390:ARG:HH21	1.77	0.49
1:B:349:LYS:HA	1:B:352:VAL:HG22	1.94	0.49
1:D:207:ARG:NH2	3:I:11:DG:O3'	2.45	0.49
4:E:601:ANP:O2A	4:E:601:ANP:O2B	2.30	0.49
1:G:389:LEU:HD22	1:G:403:LEU:HD21	1.95	0.49
1:D:382:VAL:HG22	1:D:386:LEU:HD23	1.93	0.49
1:D:146:LEU:HD21	1:D:230:VAL:HG11	1.94	0.49
1:B:109:TYR:OH	1:C:26:ALA:O	2.29	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:254:THR:O	1:F:258:ILE:HG12	2.13	0.49
1:F:215:ALA:O	1:F:219:GLN:HG2	2.14	0.48
1:G:51:LEU:HD21	1:G:322:GLN:HA	1.95	0.48
1:G:138:SER:HB3	1:G:268:THR:HG21	1.95	0.48
1:B:134:LEU:HD12	1:B:291:ILE:HB	1.94	0.48
1:A:229:LEU:HB3	1:A:263:VAL:HG12	1.94	0.48
1:F:397:ILE:HD13	1:F:405:VAL:HA	1.94	0.48
4:C:601:ANP:O1B	4:C:601:ANP:O3G	2.30	0.48
1:D:250:ASN:HD22	1:E:239:ARG:HH12	1.60	0.48
1:A:131:SER:O	1:A:287:GLY:N	2.35	0.48
1:C:319:GLN:HB2	1:C:324:LYS:HE2	1.95	0.48
1:D:110:GLU:HG2	1:E:9:VAL:HG21	1.95	0.48
1:C:313:ASP:OD1	1:C:333:ARG:NH2	2.41	0.47
3:I:14:DC:H2''	3:I:15:DG:C8	2.49	0.47
1:B:404:VAL:HG22	1:B:406:PRO:HD2	1.96	0.47
2:H:9:DC:H2''	2:H:10:DG:C8	2.49	0.47
1:G:55:ARG:HG2	1:G:58:ARG:HH22	1.78	0.47
1:B:367:THR:HG22	1:B:370:LEU:HG	1.96	0.47
1:D:98:THR:HB	1:D:101:LEU:HD23	1.97	0.47
1:D:143:THR:O	1:D:147:HIS:ND1	2.48	0.47
1:E:69:TYR:CD2	4:E:601:ANP:H2	2.49	0.47
1:E:312:THR:HG22	1:E:328:LEU:HD11	1.97	0.47
1:D:211:GLU:HA	1:E:182:GLU:OE2	2.14	0.47
1:F:58:ARG:CZ	1:F:322:GLN:HE22	2.28	0.47
1:G:203:TYR:HB3	1:G:216:LEU:HD21	1.96	0.47
1:C:247:GLU:HG2	1:D:176:HIS:HD2	1.78	0.47
1:E:253:VAL:HA	1:E:256:VAL:HG12	1.96	0.47
1:A:18:TYR:HB3	1:A:24:ILE:HG21	1.97	0.46
1:A:185:LEU:O	1:A:189:ARG:HG3	2.15	0.46
1:G:141:GLY:HA3	4:G:601:ANP:C8	2.43	0.46
1:G:143:THR:OG1	4:G:601:ANP:N3B	2.48	0.46
1:G:291:ILE:HD11	1:G:293:TRP:HE1	1.80	0.46
1:D:328:LEU:HD23	1:D:333:ARG:HH21	1.80	0.46
1:E:216:LEU:O	1:E:220:ILE:HG12	2.15	0.46
1:G:380:LYS:HA	1:G:383:HIS:CD2	2.45	0.46
1:D:232:ASP:OD1	1:D:233:GLU:N	2.48	0.46
1:B:216:LEU:HD21	1:C:205:LEU:HD21	1.98	0.46
1:D:58:ARG:O	1:D:62:ILE:HG12	2.16	0.46
1:G:189:ARG:HD2	1:G:200:GLU:HG2	1.97	0.46
1:A:146:LEU:HD11	1:A:230:VAL:HG21	1.97	0.46
1:D:147:HIS:HB3	1:D:171:LYS:HE3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:55:ARG:HB3	1:G:359:LEU:HD21	1.97	0.46
1:G:298:GLN:HA	1:G:308:TRP:CZ3	2.51	0.46
1:F:133:LEU:HD13	1:F:135:ILE:HG13	1.98	0.46
1:F:279:LEU:HD12	1:G:385:MET:HB3	1.98	0.46
1:C:316:TRP:CD2	1:C:328:LEU:HD13	2.51	0.45
1:E:346:ILE:HD13	1:E:349:LYS:HE3	1.96	0.45
1:B:250:ASN:HD21	1:C:239:ARG:HH21	1.65	0.45
1:E:46:LEU:HD22	1:E:320:LEU:HD11	1.98	0.45
1:E:279:LEU:HD13	1:F:403:LEU:HD21	1.99	0.45
1:G:6:ILE:HG22	1:G:157:ILE:HD12	1.98	0.45
1:G:88:ARG:O	1:G:92:VAL:HG23	2.15	0.45
1:G:319:GLN:HG2	1:G:355:GLN:NE2	2.32	0.45
1:B:145:SER:O	1:B:149:ILE:HG12	2.16	0.45
1:C:180:LEU:HD13	1:C:248:MET:HG3	1.99	0.45
1:G:367:THR:HG23	1:G:370:LEU:H	1.82	0.45
1:C:216:LEU:HD21	1:D:205:LEU:HD21	1.99	0.45
1:C:227:GLY:O	1:C:262:PRO:HG2	2.17	0.45
1:E:107:ASN:OD1	1:E:111:ARG:NH2	2.50	0.45
3:I:4:DC:H2 ^o	3:I:5:DG:C8	2.52	0.45
1:A:82:ARG:HH21	1:B:353:LEU:HD22	1.82	0.45
1:A:206:LYS:HD3	1:A:207:ARG:N	2.31	0.45
1:C:253:VAL:HG11	1:C:280:ARG:NH2	2.32	0.45
1:B:168:VAL:HG13	1:B:169:TYR:HD2	1.82	0.45
1:D:252:PHE:O	1:D:256:VAL:HG23	2.17	0.44
1:B:177:ASN:ND2	1:B:179:SER:OG	2.50	0.44
1:B:94:ARG:HH22	1:B:227:GLY:HA2	1.82	0.44
1:G:316:TRP:CE2	1:G:328:LEU:HD13	2.52	0.44
1:C:46:LEU:HD13	1:C:320:LEU:HD21	1.98	0.44
1:C:253:VAL:HG13	1:C:284:ARG:HD3	1.98	0.44
1:G:4:THR:HG23	1:G:6:ILE:H	1.82	0.44
1:A:232:ASP:OD1	1:A:233:GLU:N	2.51	0.44
1:F:211:GLU:OE2	1:G:177:ASN:ND2	2.51	0.44
2:H:13:DC:H2 ^o	2:H:14:DG:C8	2.53	0.44
1:B:335:VAL:HA	1:B:338:GLU:HB3	2.00	0.44
1:E:118:GLU:HG3	1:F:31:GLN:NE2	2.32	0.44
1:A:178:GLY:N	1:A:182:GLU:OE2	2.48	0.43
1:F:241:ARG:HA	1:F:241:ARG:HD3	1.84	0.43
1:G:339:LEU:HD12	1:G:371:LEU:HG	2.00	0.43
1:F:206:LYS:O	1:F:208:HIS:ND1	2.30	0.43
1:E:280:ARG:NH1	1:F:235:GLN:HE21	2.16	0.43
1:E:393:ILE:HG22	1:E:395:GLU:H	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:234:ILE:HD13	1:F:265:LEU:HD22	2.00	0.43
1:A:83:ILE:HD12	1:A:149:ILE:HG21	2.00	0.43
1:B:339:LEU:HD22	1:B:375:TYR:CD2	2.54	0.43
1:F:328:LEU:HD22	1:F:333:ARG:NH2	2.33	0.43
1:G:31:GLN:HG2	1:G:32:GLU:H	1.83	0.43
1:B:73:LEU:H	1:B:76:HIS:CE1	2.36	0.43
1:E:323:ARG:HH12	1:E:363:ASN:HA	1.84	0.43
1:F:175:SER:HB2	1:F:182:GLU:HB2	2.01	0.43
1:E:18:TYR:OH	1:F:64:ARG:NH1	2.52	0.43
1:G:32:GLU:OE1	1:G:34:VAL:HG12	2.18	0.43
1:C:244:GLY:HA3	1:D:239:ARG:HH22	1.84	0.43
1:E:380:LYS:HE2	1:E:380:LYS:HB2	1.78	0.43
1:D:143:THR:HG22	1:D:147:HIS:CE1	2.54	0.43
1:D:278:ASP:N	1:D:278:ASP:OD1	2.52	0.43
1:F:58:ARG:NH2	1:F:322:GLN:HE22	2.17	0.42
1:B:110:GLU:O	1:B:114:THR:OG1	2.35	0.42
1:B:299:THR:HG22	1:B:304:PRO:HA	2.00	0.42
1:D:12:ASP:N	1:D:12:ASP:OD1	2.51	0.42
1:E:113:GLN:O	1:F:11:ARG:NE	2.52	0.42
1:F:113:GLN:HG2	1:G:9:VAL:HB	2.01	0.42
1:B:83:ILE:HD12	1:B:149:ILE:HG21	2.01	0.42
1:D:94:ARG:HH12	1:D:261:VAL:HG12	1.85	0.42
1:D:119:THR:HG23	1:E:152:THR:HG23	2.01	0.42
1:A:279:LEU:HD23	1:B:403:LEU:HD21	2.01	0.42
1:C:234:ILE:HD13	1:C:234:ILE:HA	1.88	0.42
1:D:3:ALA:N	1:D:196:GLY:H	2.17	0.42
1:D:237:LEU:HD21	1:D:248:MET:SD	2.60	0.42
1:F:82:ARG:HH12	1:G:356:LEU:HD22	1.84	0.42
1:E:343:VAL:O	1:E:347:VAL:HG23	2.20	0.42
1:G:308:TRP:CH2	1:G:337:TYR:HB2	2.55	0.42
1:B:112:VAL:HA	1:B:117:LEU:O	2.20	0.42
1:G:22:PRO:HD2	1:G:23:PHE:N	2.34	0.42
2:H:1:DC:H2"	2:H:2:DG:C8	2.55	0.42
1:E:269:PRO:HB3	1:E:272:ARG:NH1	2.34	0.42
1:F:184:CYS:SG	1:F:214:LEU:HD23	2.60	0.42
1:G:316:TRP:CD1	1:G:328:LEU:HB2	2.55	0.42
1:G:12:ASP:OD1	1:G:13:THR:N	2.53	0.42
1:A:235:GLN:HA	1:A:271:ALA:HB2	2.01	0.41
1:B:298:GLN:HB2	1:B:337:TYR:CE2	2.55	0.41
1:E:93:GLY:O	1:E:126:ARG:HG3	2.20	0.41
1:C:377:ASP:HA	1:C:380:LYS:HE3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:211:GLU:HG3	1:F:182:GLU:HG3	2.02	0.41
1:E:365:ARG:NH2	1:E:367:THR:HB	2.35	0.41
1:G:10:TYR:HA	1:G:25:GLU:OE1	2.20	0.41
1:A:237:LEU:HD13	1:A:248:MET:SD	2.61	0.41
1:B:226:LEU:HD23	1:B:261:VAL:HG11	2.01	0.41
1:B:328:LEU:HD11	1:B:333:ARG:NH2	2.35	0.41
1:D:241:ARG:HA	1:D:241:ARG:HD3	1.80	0.41
1:B:18:TYR:HB3	1:B:24:ILE:HG21	2.02	0.41
1:C:234:ILE:HG13	1:C:265:LEU:HD23	2.03	0.41
1:F:306:GLN:HA	1:F:309:ILE:HG22	2.02	0.41
1:G:168:VAL:HG13	1:G:169:TYR:HD2	1.86	0.41
1:G:210:ILE:HA	1:G:213:MET:HG2	2.00	0.41
1:A:253:VAL:HA	1:A:256:VAL:HG22	2.02	0.41
1:A:120:PHE:HD2	1:A:122:PHE:CE1	2.39	0.41
1:C:12:ASP:OD1	1:C:13:THR:N	2.54	0.41
1:E:103:LYS:HB2	1:E:103:LYS:HE3	1.85	0.41
1:D:239:ARG:HA	1:D:239:ARG:HD2	1.70	0.41
1:A:259:ILE:HG22	1:A:261:VAL:HG23	2.03	0.41
1:C:237:LEU:HD21	1:C:248:MET:SD	2.60	0.41
1:E:269:PRO:O	1:E:272:ARG:NH1	2.54	0.41
1:F:18:TYR:CD1	1:F:24:ILE:HG12	2.56	0.41
1:G:145:SER:O	1:G:149:ILE:HG12	2.20	0.41
2:H:3:DC:H2''	2:H:4:DG:C8	2.56	0.41
2:H:6:DG:H2''	2:H:7:DC:C6	2.55	0.41
1:B:58:ARG:O	1:B:62:ILE:HG12	2.21	0.41
1:B:121:ARG:HH12	1:C:148:ARG:NH1	2.19	0.40
1:D:363:ASN:O	1:D:365:ARG:N	2.52	0.40
1:A:24:ILE:HG23	1:A:88:ARG:HG2	2.02	0.40
1:C:239:ARG:NH1	1:C:270:LYS:HZ3	2.19	0.40
1:E:141:GLY:HA3	4:E:601:ANP:H8	2.03	0.40
1:E:178:GLY:HA3	1:E:240:SER:HB2	2.03	0.40
1:F:312:THR:OG1	1:F:333:ARG:NH1	2.54	0.40
1:A:84:SER:HB2	1:A:88:ARG:NH2	2.36	0.40
1:E:198:ASN:O	1:E:202:ARG:HG3	2.21	0.40
1:G:300:GLN:O	1:G:301:ARG:NE	2.49	0.40
2:H:5:DC:H2''	2:H:6:DG:C8	2.56	0.40
1:C:216:LEU:O	1:C:220:ILE:HG12	2.20	0.40
1:G:142:LYS:N	4:G:601:ANP:O2B	2.54	0.40
3:I:6:DC:H2''	3:I:7:DG:C8	2.57	0.40
1:B:55:ARG:HG2	1:B:58:ARG:HH21	1.87	0.40
1:B:180:LEU:HD11	1:B:247:GLU:HG3	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:91:TYR:OH	1:G:167:VAL:O	2.38	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	245/523 (47%)	231 (94%)	14 (6%)	0	100	100
1	B	400/523 (76%)	374 (94%)	26 (6%)	0	100	100
1	C	402/523 (77%)	386 (96%)	15 (4%)	1 (0%)	47	80
1	D	401/523 (77%)	384 (96%)	17 (4%)	0	100	100
1	E	400/523 (76%)	378 (94%)	22 (6%)	0	100	100
1	F	399/523 (76%)	392 (98%)	7 (2%)	0	100	100
1	G	350/523 (67%)	339 (97%)	10 (3%)	1 (0%)	41	74
All	All	2597/3661 (71%)	2484 (96%)	111 (4%)	2 (0%)	54	84

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	41	LYS
1	C	235	GLN

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	210/458 (46%)	208 (99%)	2 (1%)	76	89
1	B	346/458 (76%)	346 (100%)	0	100	100
1	C	347/458 (76%)	344 (99%)	3 (1%)	78	90
1	D	346/458 (76%)	345 (100%)	1 (0%)	92	97
1	E	346/458 (76%)	346 (100%)	0	100	100
1	F	344/458 (75%)	344 (100%)	0	100	100
1	G	311/458 (68%)	309 (99%)	2 (1%)	86	94
All	All	2250/3206 (70%)	2242 (100%)	8 (0%)	91	97

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	ARG
1	A	201	ARG
1	C	102	GLN
1	C	241	ARG
1	C	399	ARG
1	D	21	ASN
1	G	171	LYS
1	G	324	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	GLN
1	B	177	ASN
1	D	250	ASN
1	G	355	GLN
1	G	383	HIS

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 monosaccharides 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
4	ANP	E	601	5	29,33,33	1.06	3 (10%)	31,52,52	1.15	2 (6%)
4	ANP	B	601	5	29,33,33	1.09	4 (13%)	31,52,52	1.18	2 (6%)
4	ANP	G	601	-	29,33,33	1.06	4 (13%)	31,52,52	1.10	2 (6%)
4	ANP	D	601	5	29,33,33	1.02	3 (10%)	31,52,52	1.12	2 (6%)
4	ANP	F	601	5	29,33,33	1.08	4 (13%)	31,52,52	1.14	3 (9%)
4	ANP	C	601	-	29,33,33	1.06	4 (13%)	31,52,52	1.14	2 (6%)

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
4	ANP	E	601	5	-	9/14/38/38	0/3/3/3
4	ANP	B	601	5	-	7/14/38/38	0/3/3/3
4	ANP	G	601	-	-	5/14/38/38	0/3/3/3
4	ANP	D	601	5	-	5/14/38/38	0/3/3/3
4	ANP	F	601	5	-	4/14/38/38	0/3/3/3
4	ANP	C	601	-	-	7/14/38/38	0/3/3/3

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	601	ANP	PB-O3A	-2.87	1.55	1.59
4	F	601	ANP	PB-O3A	-2.66	1.55	1.59
4	B	601	ANP	PB-O3A	-2.46	1.56	1.59
4	B	601	ANP	PG-N3B	2.43	1.69	1.63
4	B	601	ANP	PG-O1G	2.37	1.49	1.46
4	G	601	ANP	PG-O1G	2.36	1.49	1.46
4	F	601	ANP	PG-O1G	2.35	1.49	1.46
4	C	601	ANP	PG-N3B	2.34	1.69	1.63
4	G	601	ANP	PG-N3B	2.34	1.69	1.63
4	D	601	ANP	PB-O3A	-2.33	1.56	1.59
4	C	601	ANP	PG-O1G	2.30	1.49	1.46
4	C	601	ANP	PB-O3A	-2.29	1.56	1.59
4	G	601	ANP	PB-O3A	-2.27	1.56	1.59
4	E	601	ANP	PG-O1G	2.25	1.49	1.46
4	F	601	ANP	PG-N3B	2.23	1.69	1.63
4	G	601	ANP	PB-O1B	2.21	1.49	1.46
4	D	601	ANP	PG-N3B	2.20	1.69	1.63
4	E	601	ANP	PG-N3B	2.19	1.69	1.63
4	D	601	ANP	PG-O1G	2.18	1.49	1.46
4	F	601	ANP	PB-O1B	2.16	1.49	1.46
4	B	601	ANP	PB-O1B	2.15	1.49	1.46
4	C	601	ANP	PB-O1B	2.09	1.49	1.46

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	601	ANP	PB-O3A-PA	-4.42	117.03	132.62
4	E	601	ANP	PB-O3A-PA	-4.39	117.14	132.62
4	F	601	ANP	PB-O3A-PA	-3.94	118.75	132.62
4	C	601	ANP	PB-O3A-PA	-3.78	119.29	132.62
4	G	601	ANP	PB-O3A-PA	-3.68	119.66	132.62
4	D	601	ANP	PB-O3A-PA	-3.54	120.14	132.62
4	D	601	ANP	C5-C6-N6	2.27	123.81	120.35
4	E	601	ANP	C5-C6-N6	2.27	123.80	120.35
4	G	601	ANP	C5-C6-N6	2.26	123.78	120.35
4	B	601	ANP	C5-C6-N6	2.25	123.78	120.35
4	C	601	ANP	C5-C6-N6	2.23	123.75	120.35
4	F	601	ANP	C5-C6-N6	2.23	123.73	120.35
4	F	601	ANP	O1G-PG-N3B	-2.05	108.75	111.77

There are no chirality outliers.

All (37) torsion outliers are listed below:

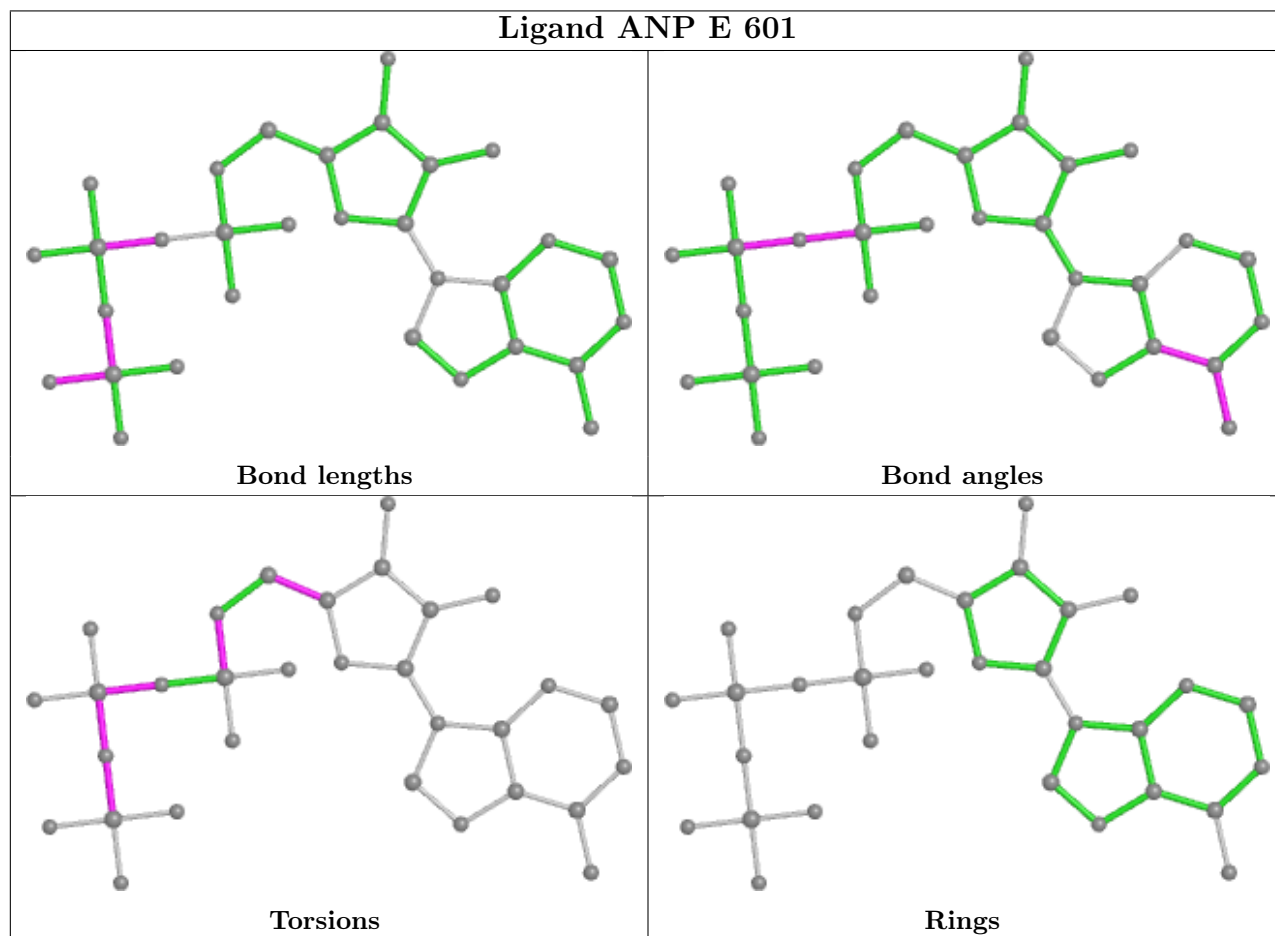
Mol	Chain	Res	Type	Atoms
4	B	601	ANP	PB-N3B-PG-O1G
4	B	601	ANP	C5'-O5'-PA-O1A
4	B	601	ANP	C5'-O5'-PA-O2A
4	B	601	ANP	O4'-C4'-C5'-O5'
4	B	601	ANP	C3'-C4'-C5'-O5'
4	C	601	ANP	PB-N3B-PG-O1G
4	C	601	ANP	PG-N3B-PB-O3A
4	C	601	ANP	C5'-O5'-PA-O1A
4	C	601	ANP	C5'-O5'-PA-O2A
4	C	601	ANP	O4'-C4'-C5'-O5'
4	C	601	ANP	C3'-C4'-C5'-O5'
4	D	601	ANP	PB-N3B-PG-O1G
4	D	601	ANP	PG-N3B-PB-O1B
4	D	601	ANP	PG-N3B-PB-O3A
4	E	601	ANP	PB-N3B-PG-O1G
4	E	601	ANP	PG-N3B-PB-O1B
4	E	601	ANP	PG-N3B-PB-O3A
4	E	601	ANP	PA-O3A-PB-O1B
4	E	601	ANP	PA-O3A-PB-O2B
4	E	601	ANP	C5'-O5'-PA-O1A
4	E	601	ANP	O4'-C4'-C5'-O5'
4	F	601	ANP	PB-N3B-PG-O1G
4	F	601	ANP	PG-N3B-PB-O1B
4	F	601	ANP	PG-N3B-PB-O3A
4	F	601	ANP	C5'-O5'-PA-O1A
4	G	601	ANP	C5'-O5'-PA-O1A
4	G	601	ANP	C5'-O5'-PA-O2A
4	G	601	ANP	O4'-C4'-C5'-O5'
4	G	601	ANP	C3'-C4'-C5'-O5'
4	D	601	ANP	O4'-C4'-C5'-O5'
4	E	601	ANP	C5'-O5'-PA-O3A
4	B	601	ANP	C4'-C5'-O5'-PA
4	D	601	ANP	C3'-C4'-C5'-O5'
4	E	601	ANP	C3'-C4'-C5'-O5'
4	B	601	ANP	C5'-O5'-PA-O3A
4	C	601	ANP	C5'-O5'-PA-O3A
4	G	601	ANP	C5'-O5'-PA-O3A

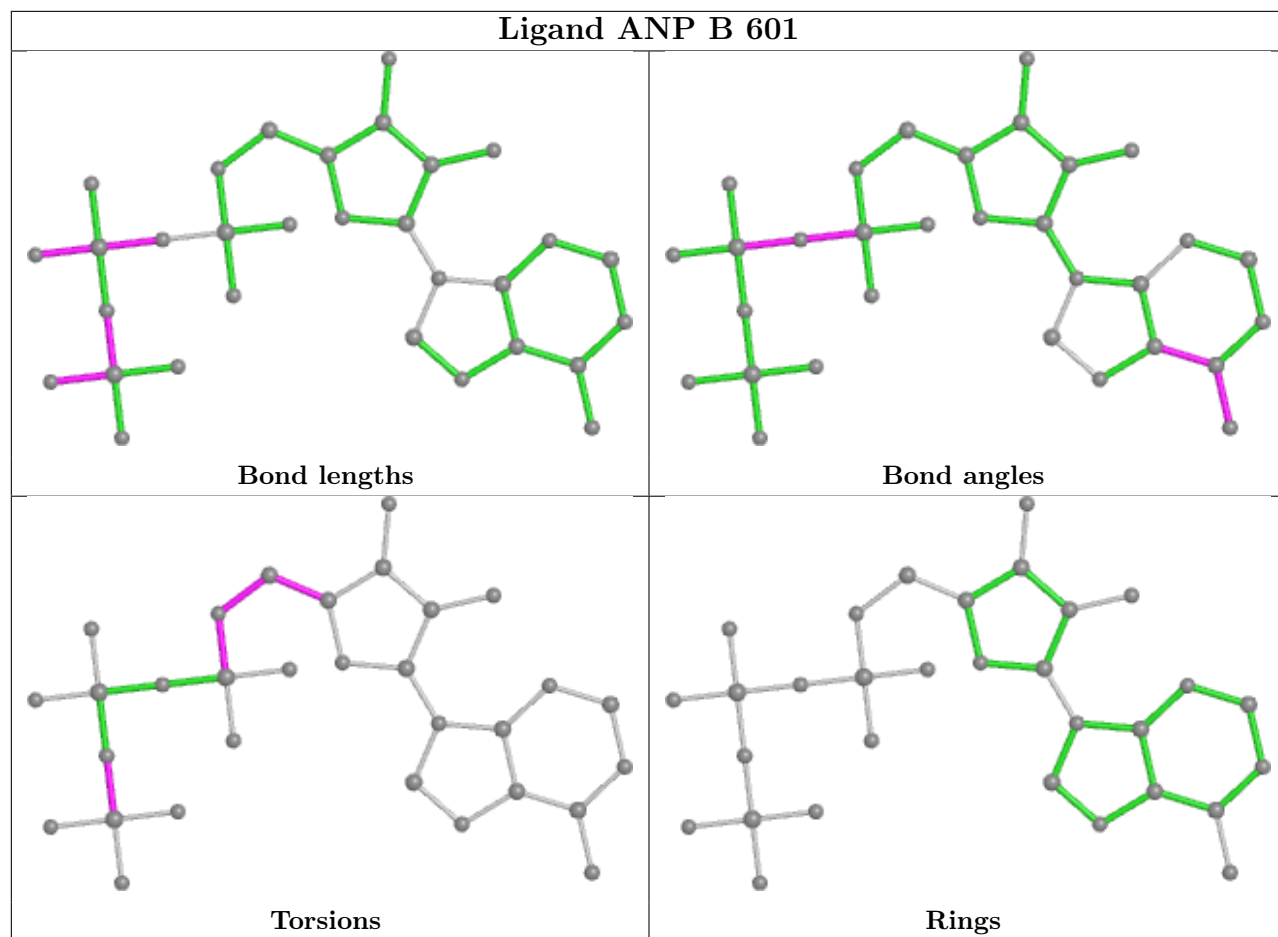
There are no ring outliers.

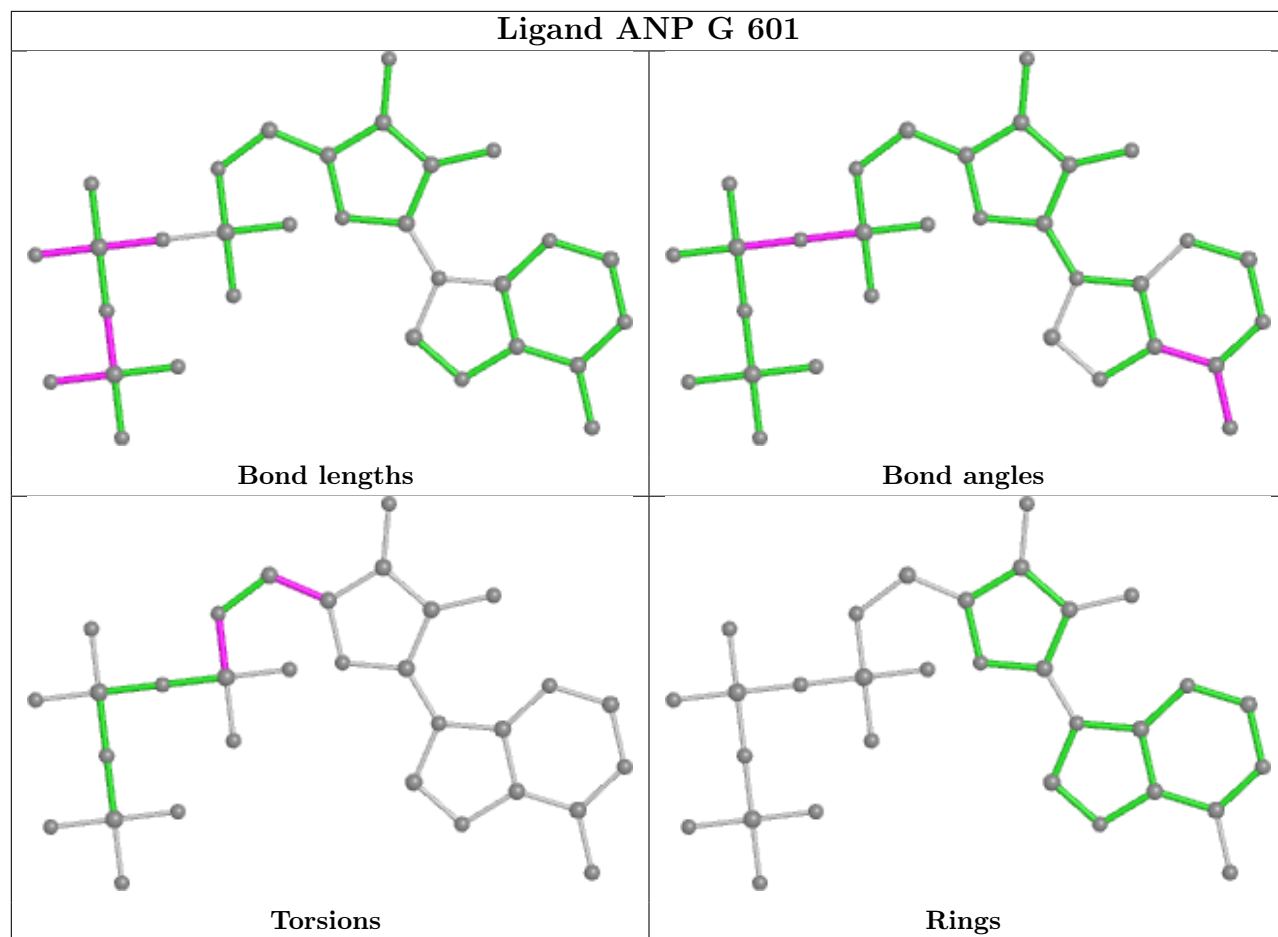
4 monomers are involved in 9 short contacts:

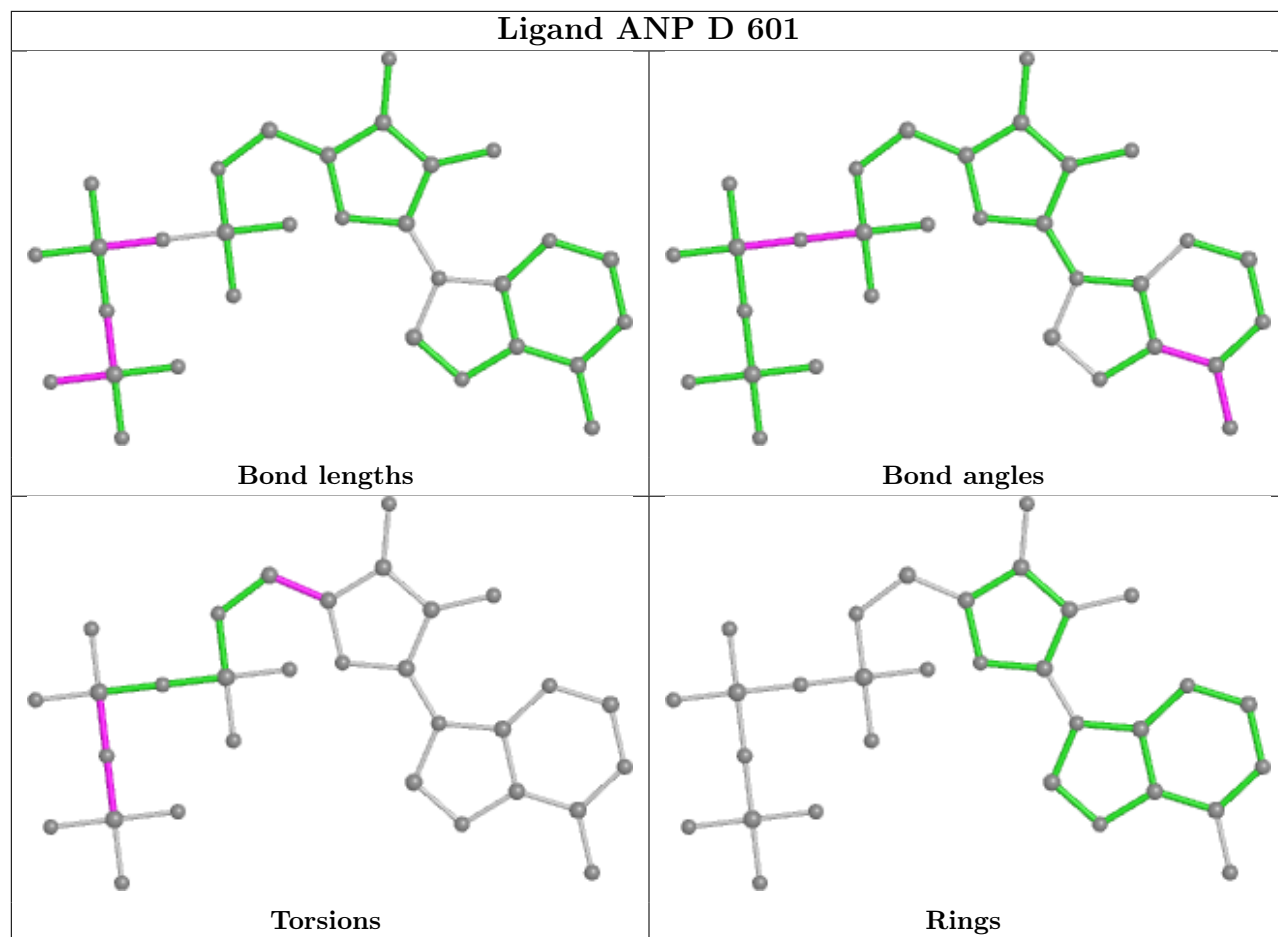
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	601	ANP	3	0
4	B	601	ANP	1	0
4	G	601	ANP	4	0
4	C	601	ANP	1	0

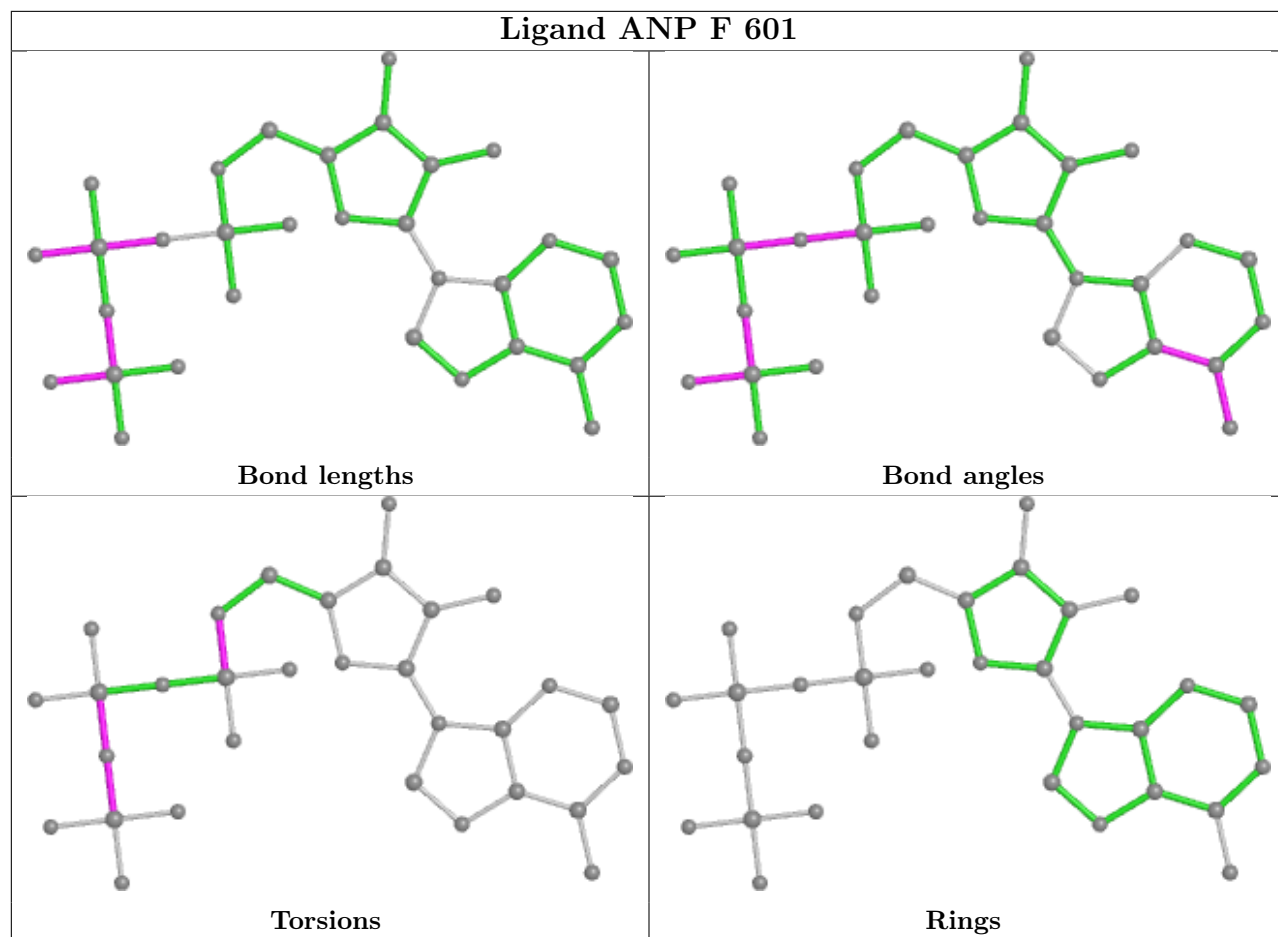
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

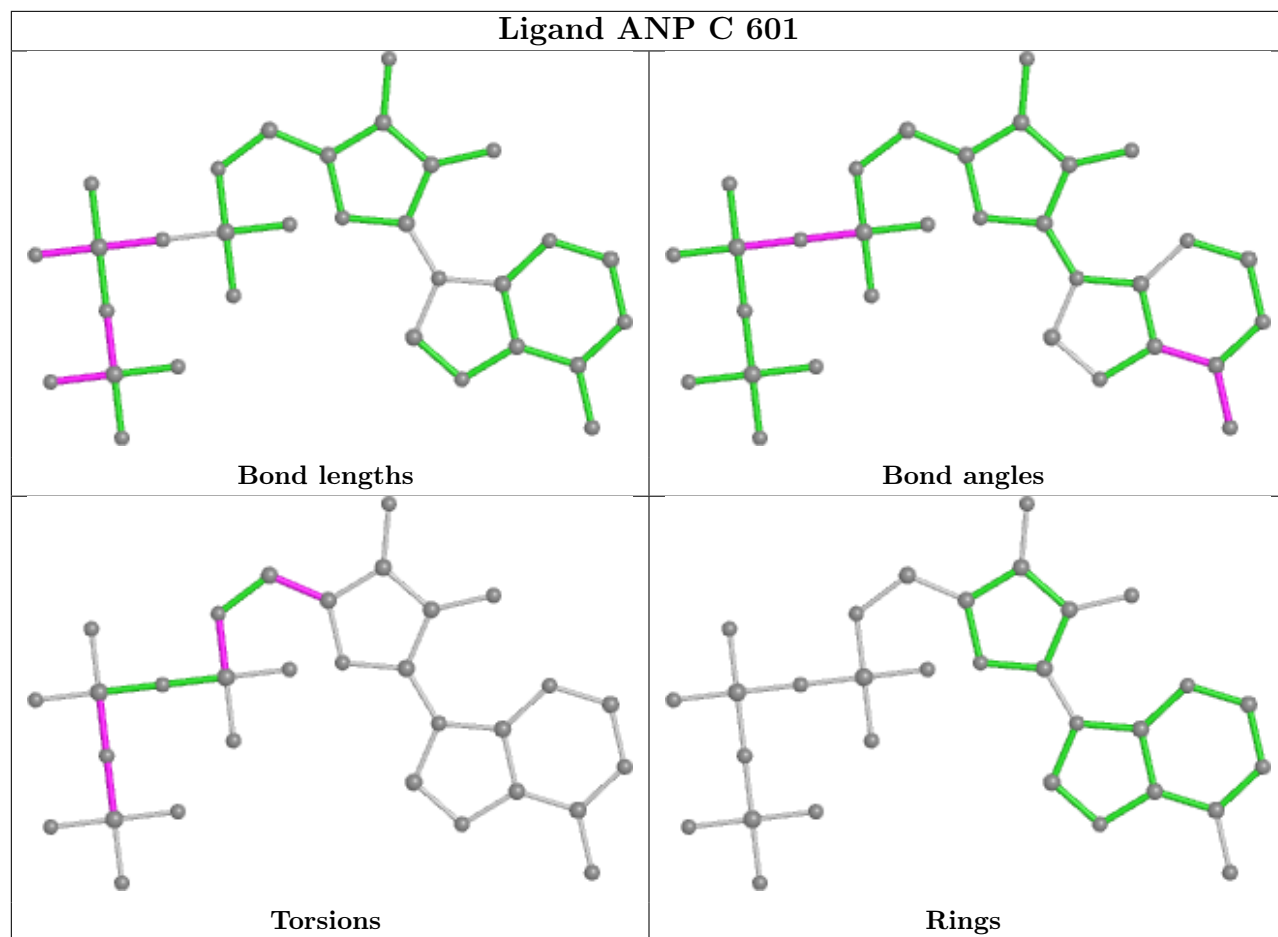












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

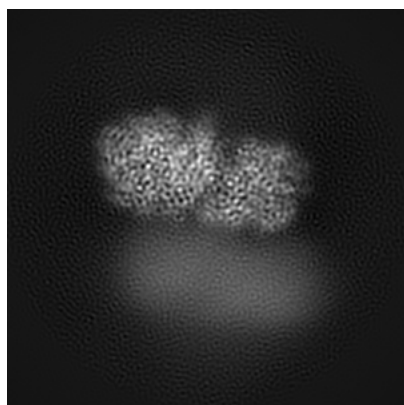
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-23757. These allow visual inspection of the internal detail of the map and identification of artifacts.

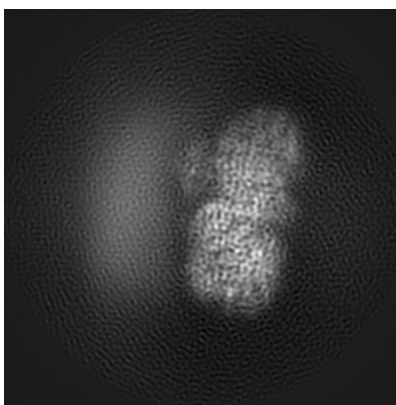
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

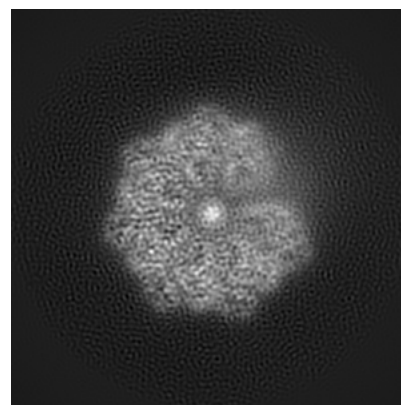
6.1.1 Primary map



X



Y

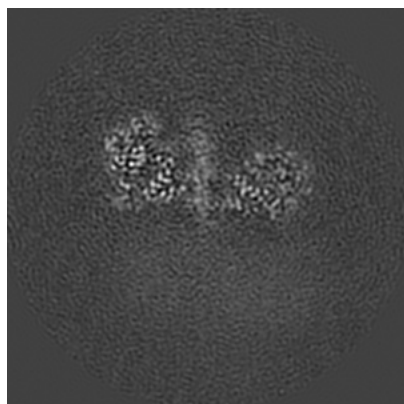


Z

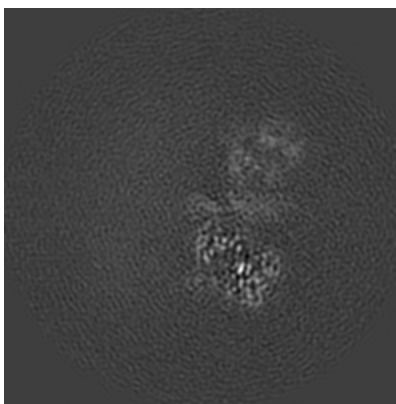
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

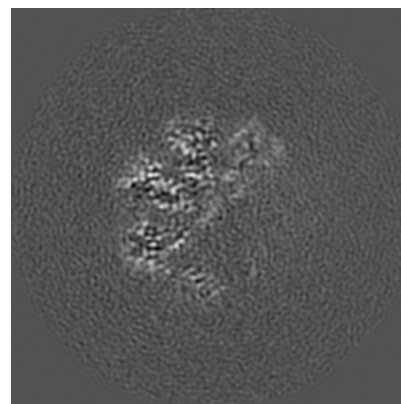
6.2.1 Primary map



X Index: 150



Y Index: 150

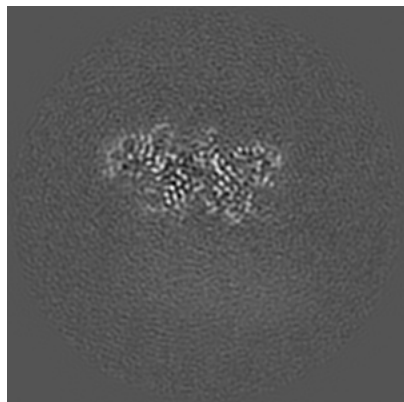


Z Index: 150

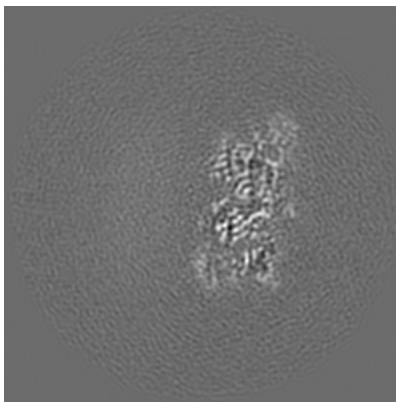
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

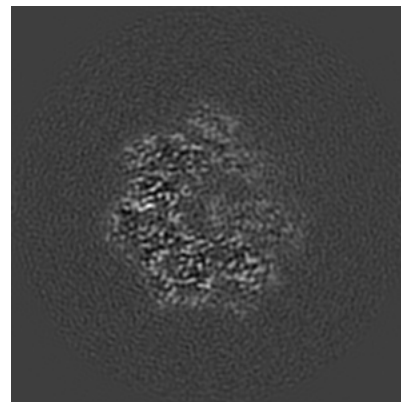
6.3.1 Primary map



X Index: 111



Y Index: 105



Z Index: 177

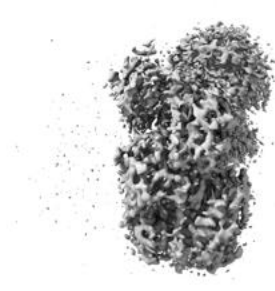
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0145. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

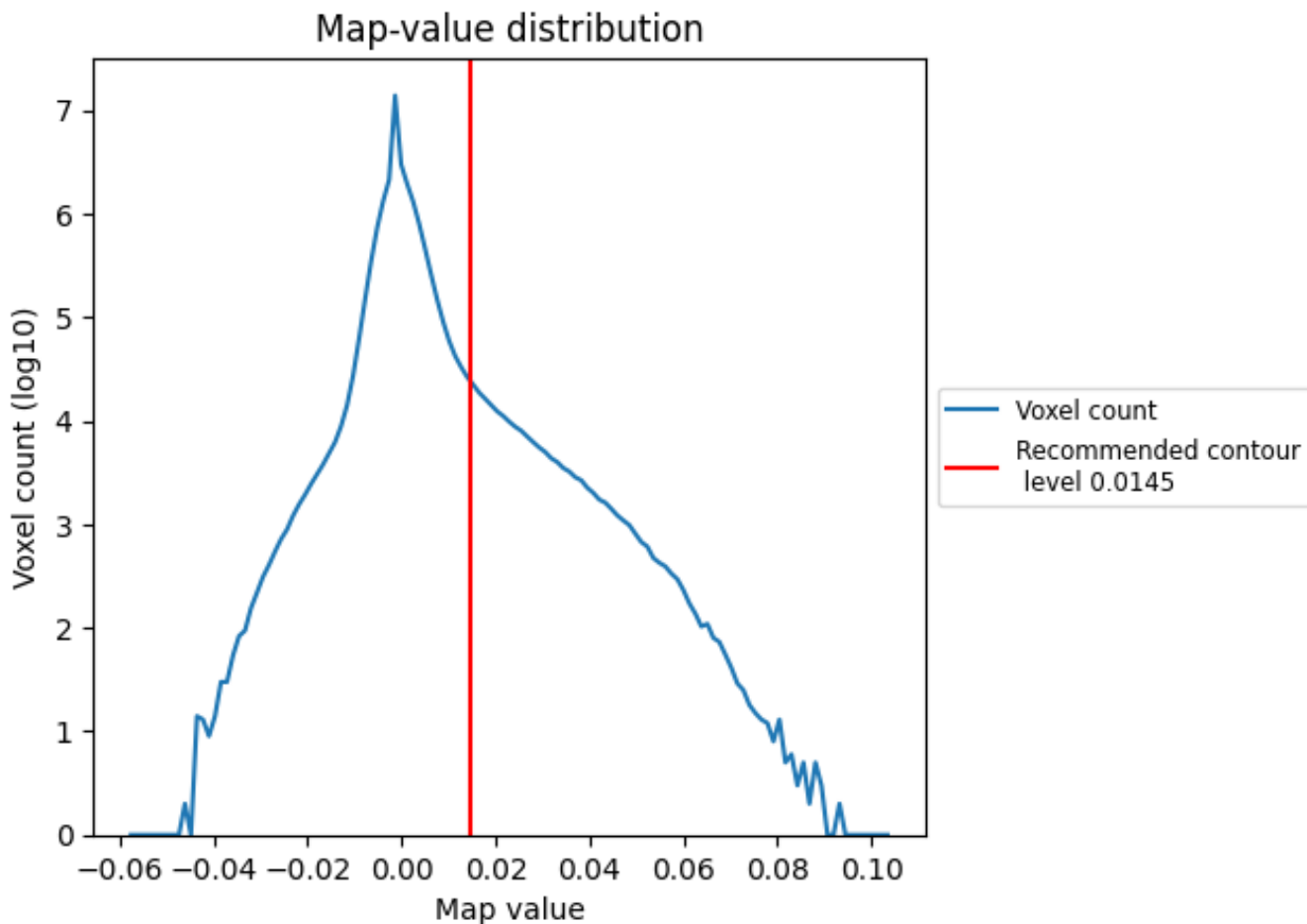
6.5 Mask visualisation

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

7 Map analysis [i](#)

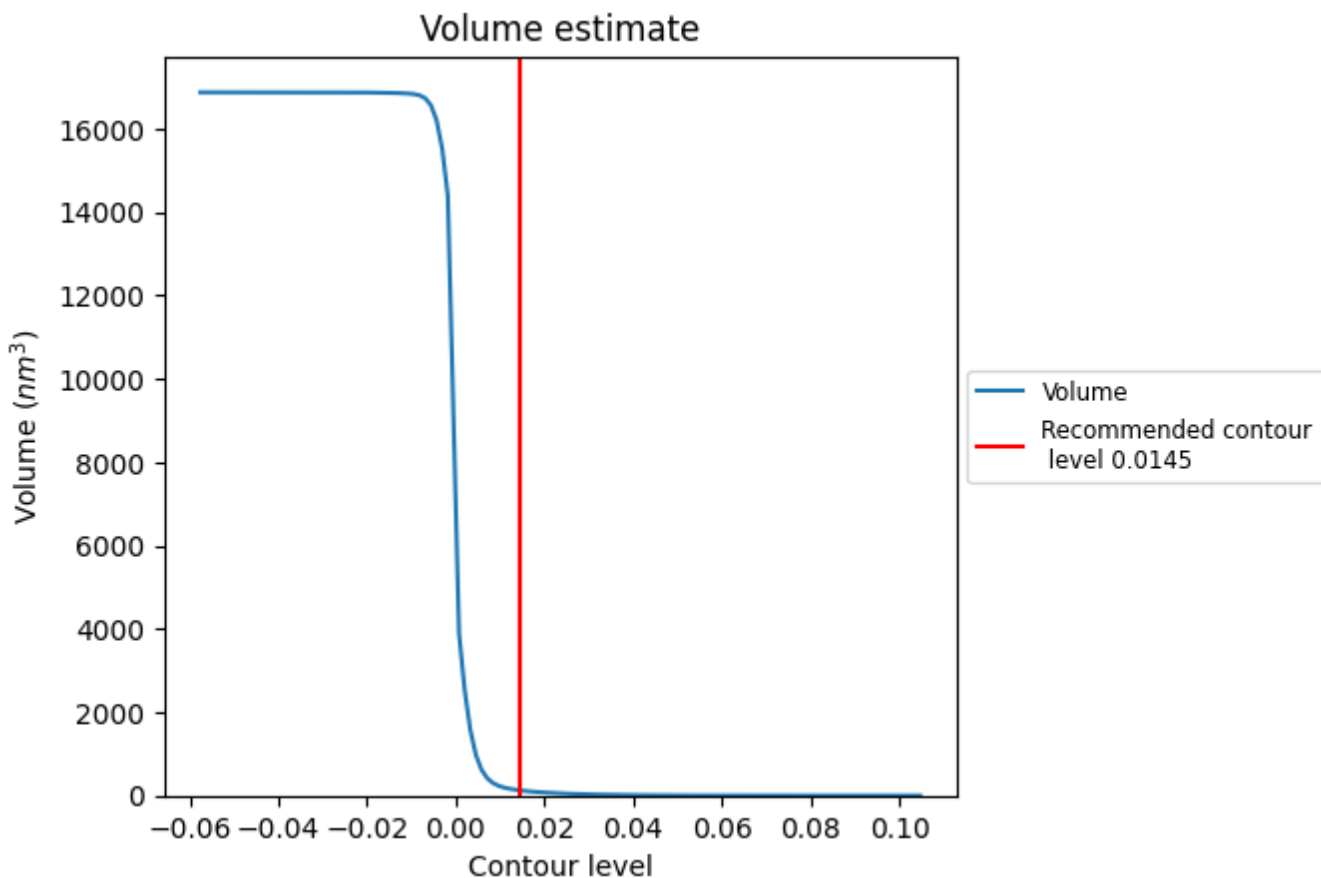
This section contains the results of statistical analysis of the map.

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.

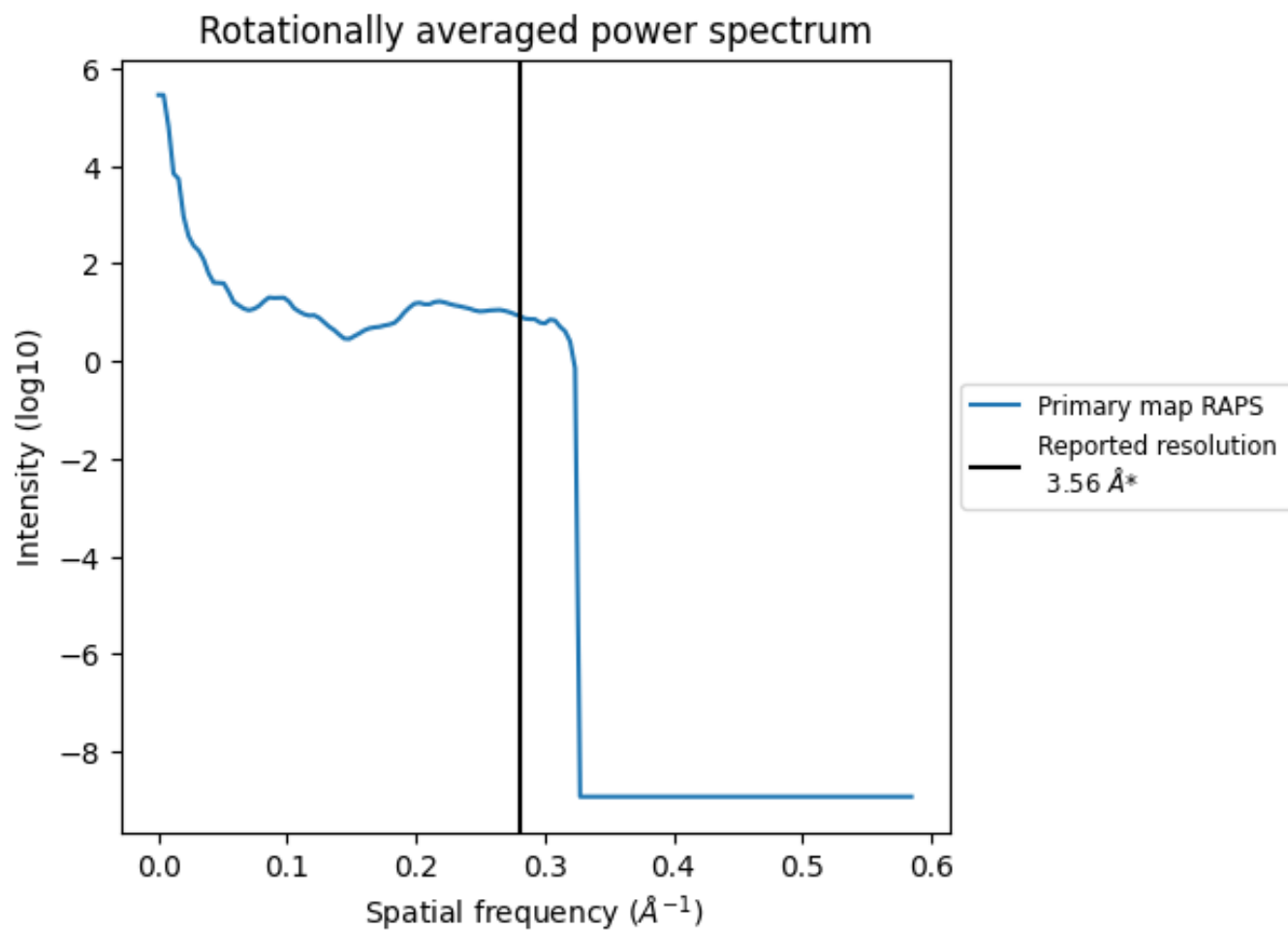
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 124 nm^3 ; this corresponds to an approximate mass of 112 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](#)

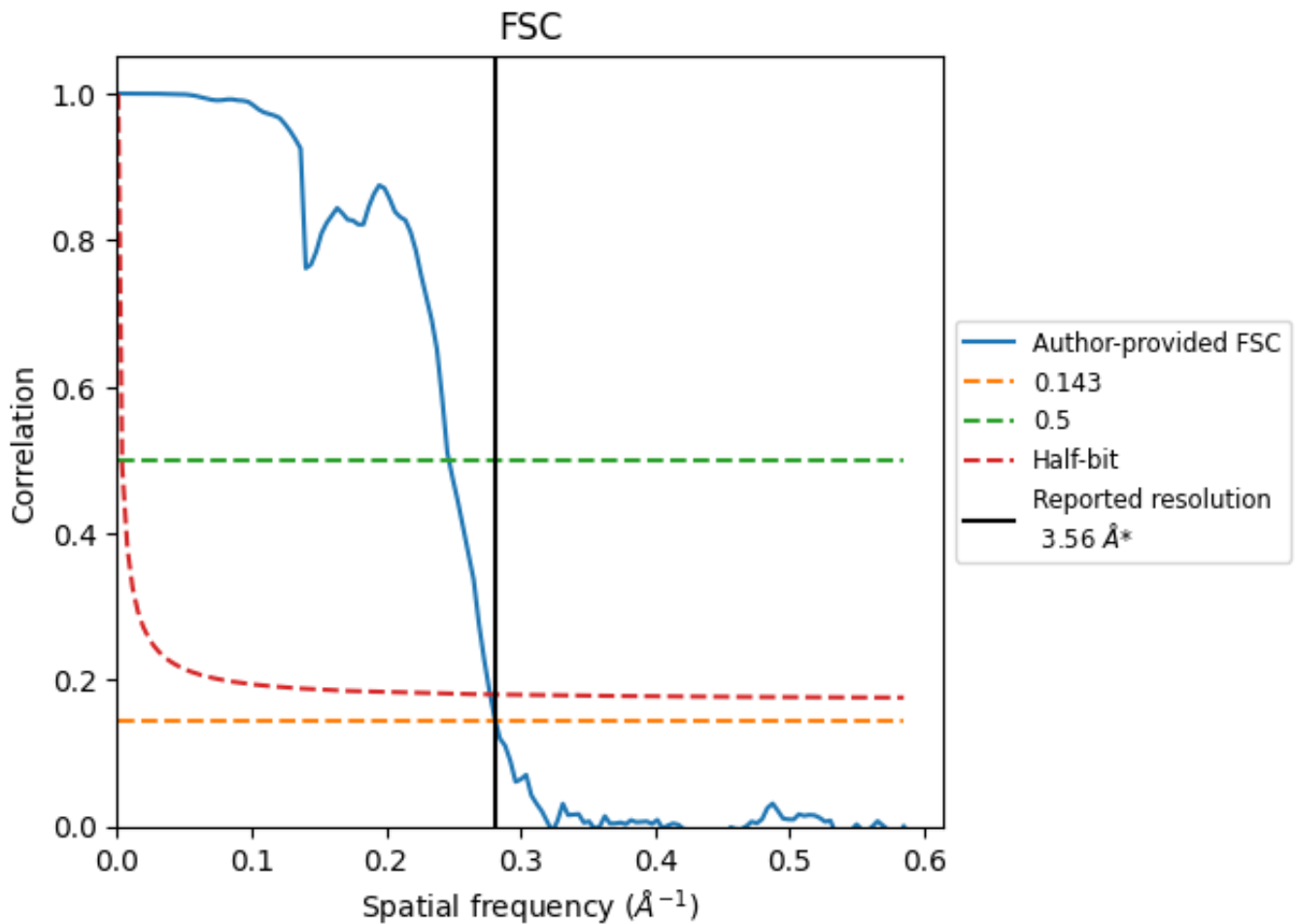


*Reported resolution corresponds to spatial frequency of 0.281 Å⁻¹

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 0.281 Å⁻¹

8.2 Resolution estimates [i](#)

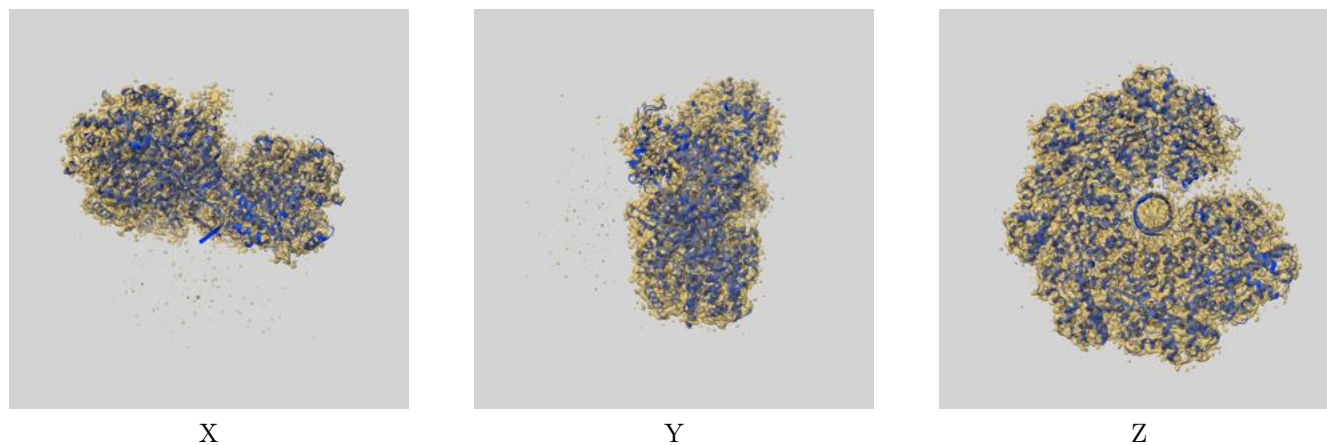
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.56	-	-
Author-provided FSC curve	3.55	4.05	3.60
Unmasked-calculated*	-	-	-

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

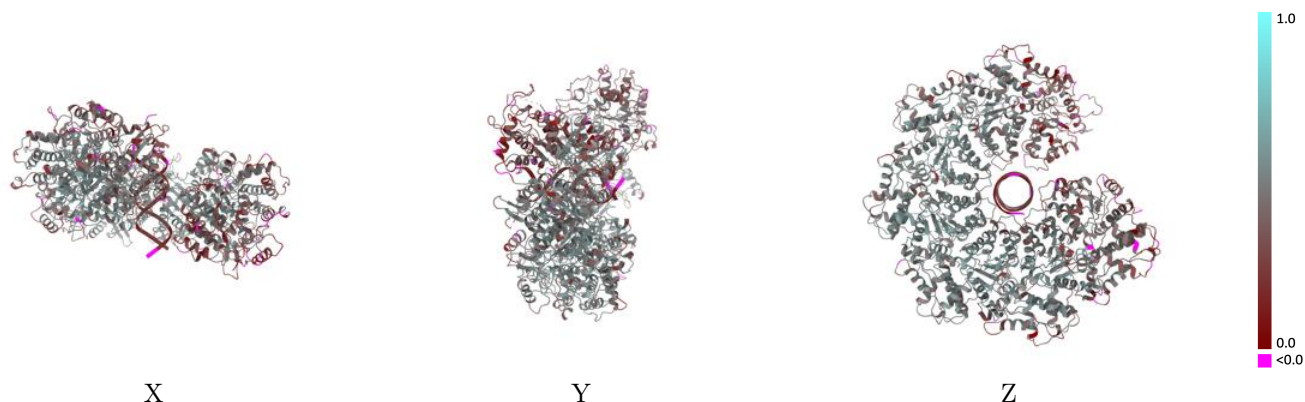
This section contains information regarding the fit between EMDB map EMD-23757 and PDB model 7MCS. Per-residue inclusion information can be found in section 3 on page 11.

9.1 Map-model overlay [i](#)



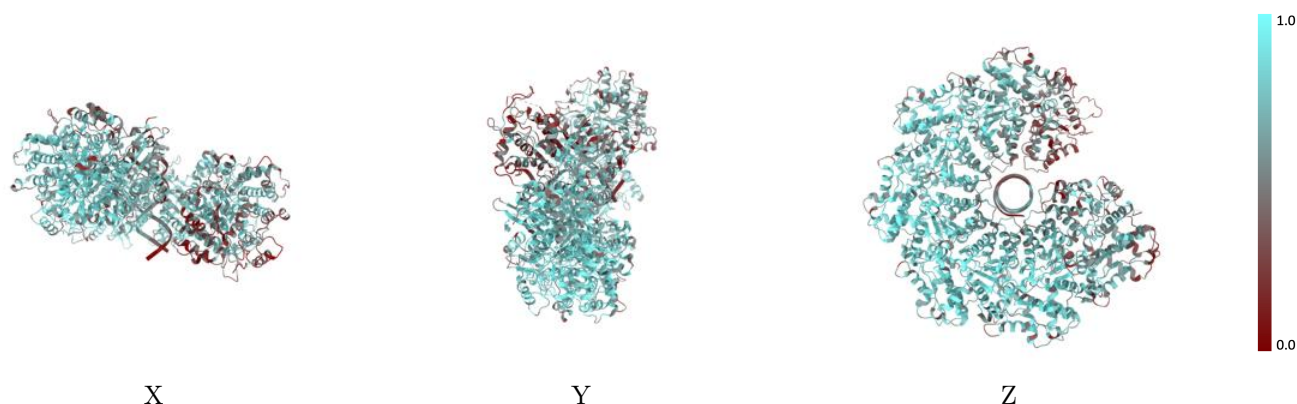
The images above show the 3D surface view of the map at the recommended contour level 0.0145 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



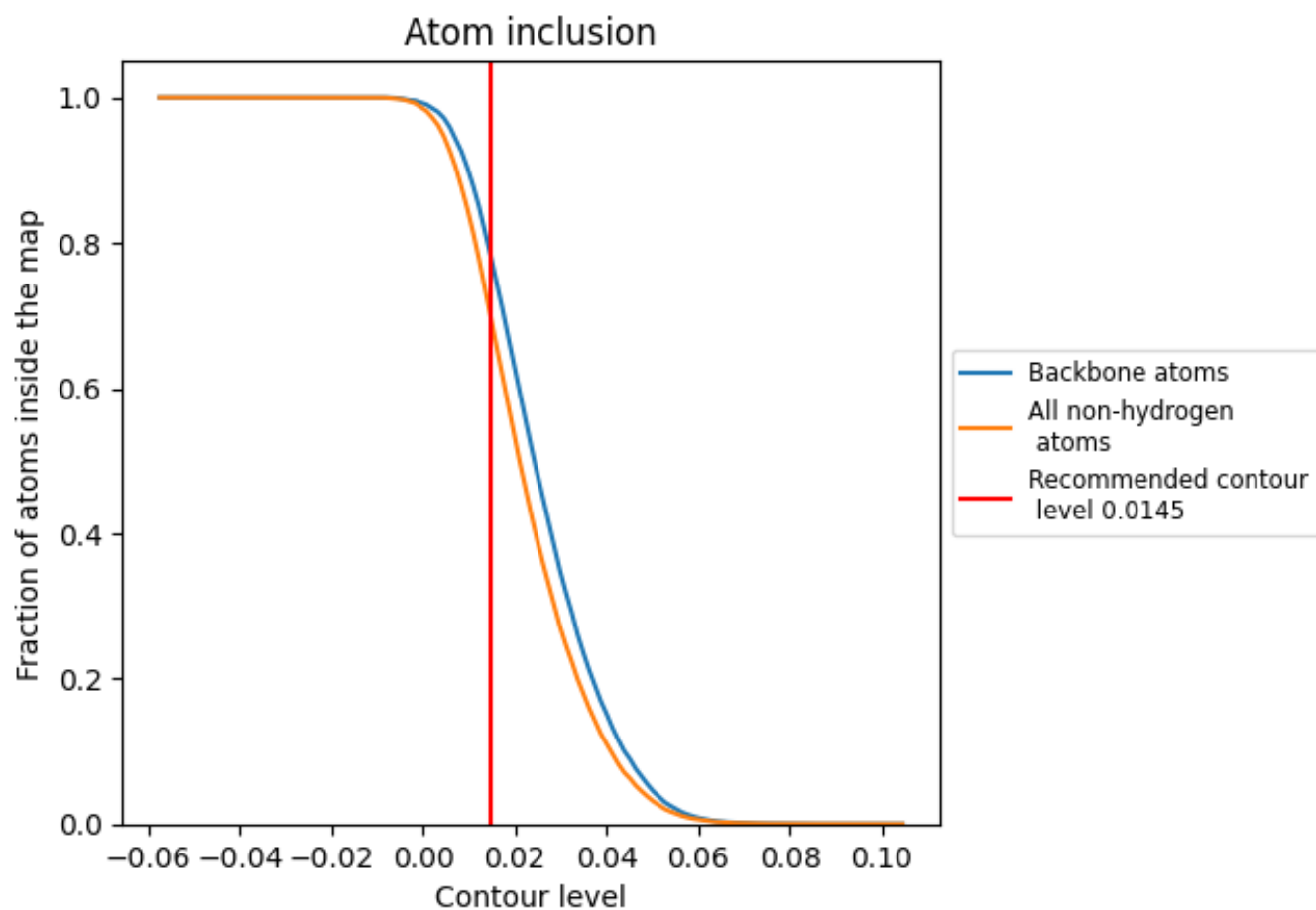
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0145).





















9.4 Atom inclusion [i](#)



At the recommended contour level, 79% of all backbone atoms, 70% 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 (0.0145) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7039	 0.4390
A	 0.5886	 0.3790
B	 0.7025	 0.4350
C	 0.8057	 0.4920
D	 0.8221	 0.5080
E	 0.8046	 0.4960
F	 0.7107	 0.4460
G	 0.4697	 0.3300
H	 0.4510	 0.1900
I	 0.4693	 0.1920

