



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 02:54 PM UTC

PDB ID : 7RY3 / pdb_00007ry3
EMDB ID : EMD-24732
Title : Multidrug Efflux pump AdeJ with TP-6076 bound
Authors : Zhang, Z.
Deposited on : 2021-08-24
Resolution : 2.91 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

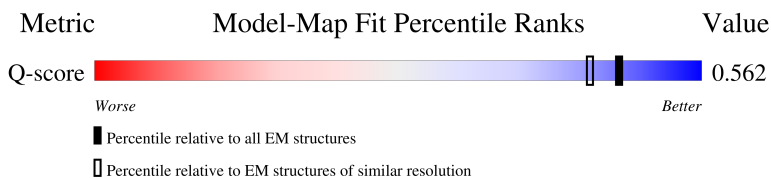
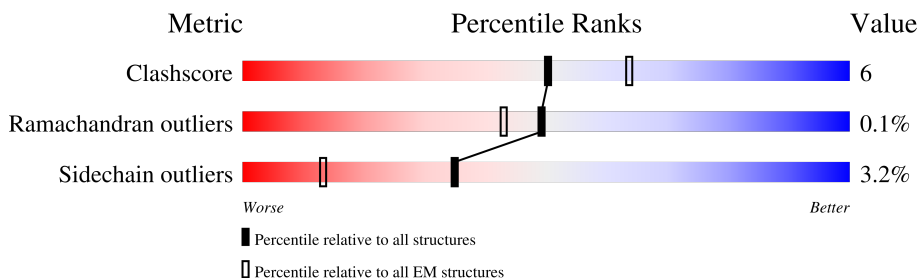
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.91 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	12972 (2.41 - 3.41)

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	1058	 78% 20% ..
1	B	1058	 79% 19% ..
1	C	1058	 81% 17% ..

2 Entry composition [i](#)

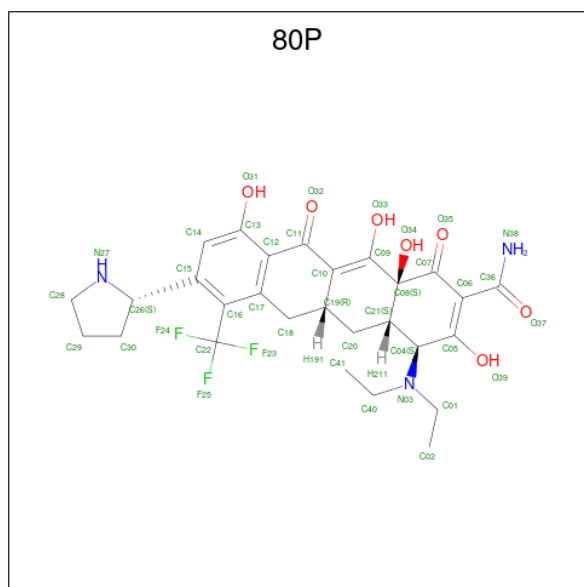
There are 2 unique types of molecules in this entry. The entry contains 23996 atoms, of which 32 are hydrogens and 0 are deuteriums.

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 Efflux pump membrane transporter.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1047	Total	C	N	O	S	0	0
			7979	5135	1324	1486	34		
1	B	1045	Total	C	N	O	S	0	0
			7958	5120	1321	1483	34		
1	C	1048	Total	C	N	O	S	0	0
			7986	5140	1325	1487	34		

- Molecule 2 is (4S,4aS,5aR,12aS)-4-(diethylamino)-3,10,12,12a-tetrahydroxy-1,11-dioxo-8-[(2S)-pyrrolidin-2-yl]-7-(trifluoromethyl)-1,4,4a,5,5a,6,11,12a-octahydrotetracene-2-carboxamide (CCD ID: 80P) (formula: C₂₈H₃₂F₃N₃O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						AltConf
2	C	1	Total	C	F	H	N	O	0
			73	28	3	32	3	7	

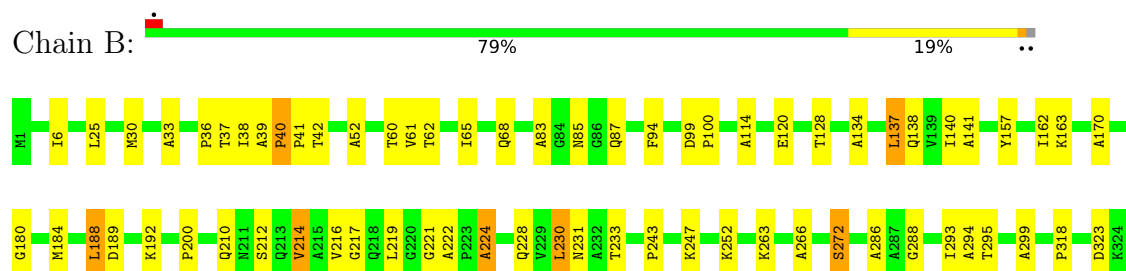
3 Residue-property plots

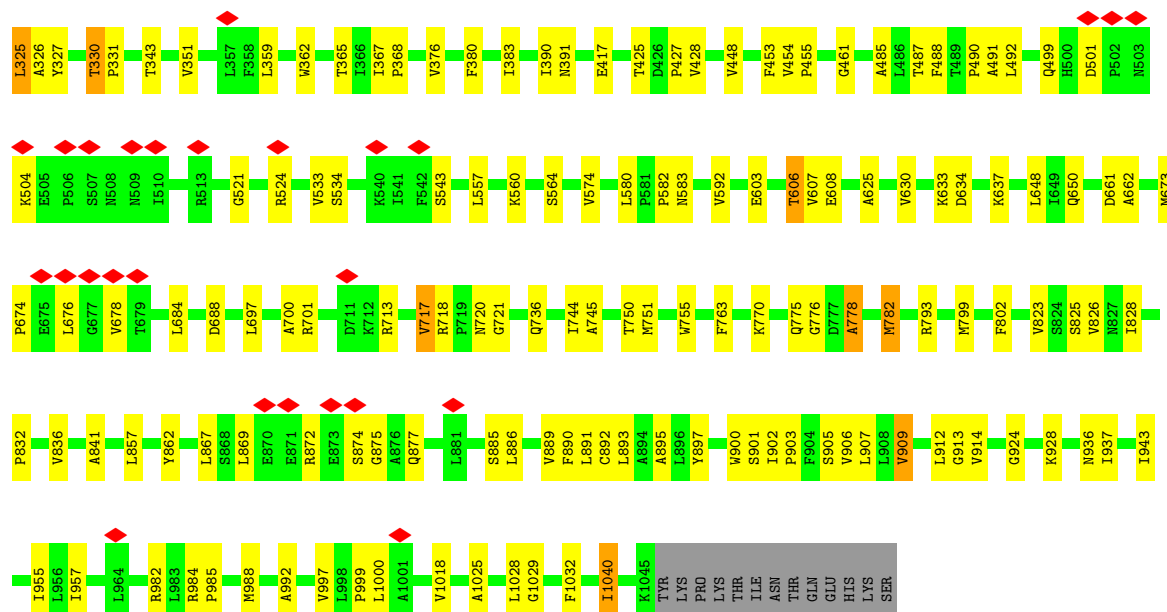
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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



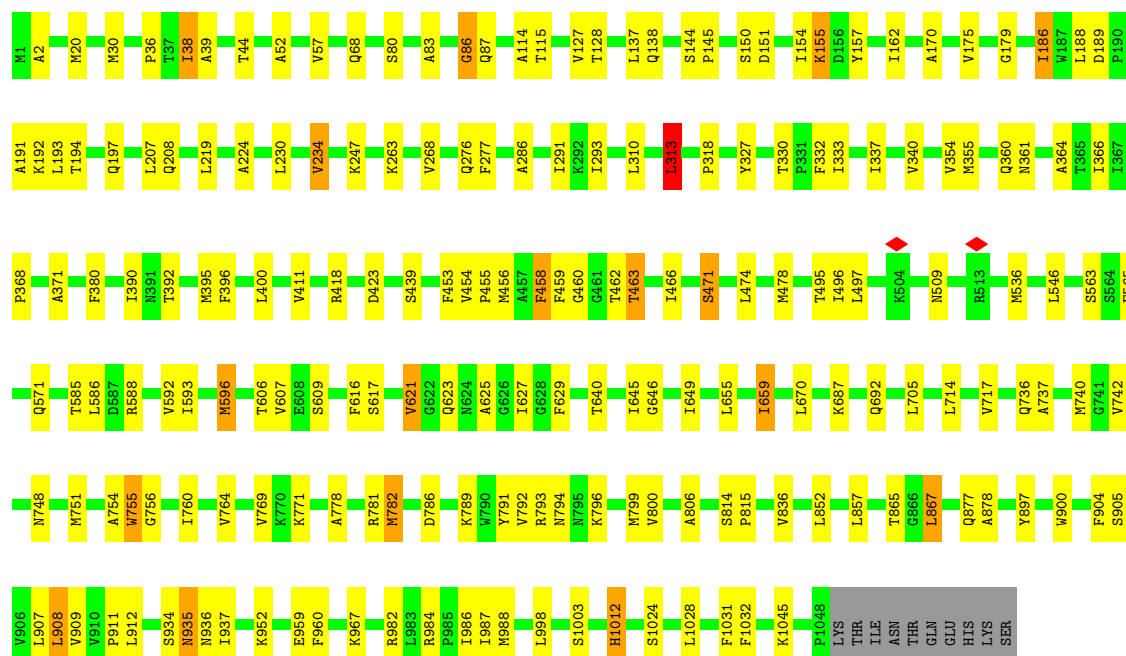
• Molecule 1: Efflux pump membrane transporter





• Molecule 1: Efflux pump membrane transporter

Chain C: 81% 17%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	68346	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 ($e^-/\text{\AA}^2$)	36	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	3.100	Depositor
Minimum map value	-1.725	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.217	Depositor
Recommended contour level	0.11	Depositor
Map size (Å)	132.84001, 132.84001, 160.92001	wwPDB
Map dimensions	149, 123, 123	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.08, 1.08, 1.08	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 80P

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	0.89	0/8137	1.08	62/11064 (0.6%)
1	B	0.70	0/8115	1.06	64/11035 (0.6%)
1	C	0.87	0/8145	1.10	68/11076 (0.6%)
All	All	0.82	0/24397	1.08	194/33175 (0.6%)

There are no bond length outliers.

All (194) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	376	VAL	N-CA-C	12.20	122.13	110.42
1	C	796	LYS	N-CA-C	-11.78	97.52	113.30
1	A	172	VAL	N-CA-C	9.77	120.59	110.72
1	A	276	GLN	N-CA-C	9.50	121.72	111.36
1	A	236	ALA	N-CA-C	9.38	121.88	108.00
1	C	495	THR	N-CA-C	9.16	121.03	111.14
1	C	371	ALA	N-CA-C	8.86	120.71	111.14
1	B	210	GLN	N-CA-C	8.73	120.88	111.36
1	C	760	ILE	N-CA-C	8.70	118.60	110.42
1	B	606	THR	N-CA-C	8.66	120.41	110.97
1	C	360	GLN	N-CA-C	8.65	120.70	111.28
1	B	984	ARG	CA-C-N	-8.48	111.08	119.64
1	B	984	ARG	C-N-CA	-8.48	111.08	119.64
1	B	170	ALA	N-CA-C	8.36	121.14	111.11
1	C	162	ILE	N-CA-C	8.36	121.54	111.09
1	A	694	HIS	N-CA-C	-8.34	102.70	113.12
1	B	85	ASN	N-CA-C	-8.27	103.29	112.72
1	B	37	THR	N-CA-C	8.23	122.84	109.76
1	B	720	ASN	N-CA-C	8.00	120.08	111.36
1	A	266	ALA	N-CA-C	7.98	120.25	108.60
1	B	61	VAL	N-CA-C	7.90	118.91	111.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	83	ALA	N-CA-C	7.87	121.37	109.41
1	C	361	ASN	N-CA-C	7.87	121.64	110.23
1	B	224	ALA	N-CA-C	7.86	120.67	110.53
1	B	216	VAL	N-CA-C	7.83	119.64	112.43
1	B	52	ALA	N-CA-C	7.75	121.95	110.48
1	B	1032	PHE	N-CA-C	7.74	119.80	111.36
1	A	235	ASN	N-CA-C	7.71	122.94	112.68
1	B	272	SER	N-CA-C	7.71	121.37	110.50
1	B	222	ALA	N-CA-C	7.66	124.04	113.16
1	A	321	LEU	N-CA-C	7.65	121.50	110.10
1	A	604	LYS	N-CA-C	7.60	119.56	111.28
1	B	134	ALA	N-CA-C	7.55	120.98	111.69
1	C	877	GLN	N-CA-C	7.47	119.42	111.28
1	C	621	VAL	CB-CA-C	-7.46	103.89	111.80
1	B	162	ILE	N-CA-C	7.42	117.91	111.90
1	C	984	ARG	CA-C-N	-7.39	112.06	120.04
1	C	984	ARG	C-N-CA	-7.39	112.06	120.04
1	B	461	GLY	N-CA-C	-7.38	103.15	111.63
1	C	234	VAL	N-CA-C	7.30	119.35	108.54
1	B	875	GLY	N-CA-C	7.29	121.47	112.73
1	A	439	SER	N-CA-C	7.25	120.23	111.82
1	C	935	ASN	N-CA-C	7.22	121.36	111.54
1	A	229	VAL	N-CA-C	7.18	117.98	110.72
1	B	661	ASP	N-CA-C	7.14	119.07	111.28
1	C	621	VAL	N-CA-C	7.13	117.50	107.37
1	B	188	LEU	N-CA-C	7.12	120.54	110.50
1	B	874	SER	N-CA-C	7.11	119.03	111.28
1	B	607	VAL	N-CA-C	7.09	118.79	108.58
1	B	330	THR	N-CA-C	7.08	123.22	113.16
1	C	782	MET	N-CA-C	7.08	120.04	111.40
1	A	900	TRP	N-CA-C	-7.08	105.27	114.04
1	B	36	PRO	N-CA-C	-7.07	100.15	111.03
1	B	62	THR	N-CA-C	7.05	118.62	111.07
1	C	115	THR	N-CA-C	6.97	118.88	111.28
1	C	276	GLN	N-CA-C	6.96	119.89	111.82
1	C	607	VAL	N-CA-C	6.96	118.82	108.46
1	C	756	GLY	N-CA-C	6.95	121.67	112.77
1	C	155	LYS	N-CA-C	-6.94	103.80	111.36
1	A	188	LEU	N-CA-C	6.82	120.98	110.20
1	A	230	LEU	N-CA-C	6.81	120.61	109.24
1	C	188	LEU	N-CA-C	6.73	120.22	109.24
1	C	471	SER	N-CA-C	6.71	118.25	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	148	ASN	N-CA-C	-6.71	104.23	112.88
1	C	224	ALA	N-CA-C	6.71	119.60	110.55
1	A	319	THR	N-CA-C	6.70	118.24	111.07
1	B	454	VAL	N-CA-C	6.70	122.33	112.61
1	B	120	GLU	N-CA-C	6.67	119.12	111.11
1	C	692	GLN	N-CA-C	-6.63	102.25	110.41
1	A	609	SER	N-CA-C	6.63	119.01	109.14
1	B	60	THR	N-CA-C	6.62	118.58	111.36
1	B	778	ALA	N-CA-C	6.59	118.46	111.28
1	A	825	SER	N-CA-C	6.55	119.37	109.41
1	C	998	LEU	CA-C-N	-6.51	111.83	119.05
1	C	998	LEU	C-N-CA	-6.51	111.83	119.05
1	C	1045	LYS	N-CA-C	6.49	119.65	110.50
1	A	756	GLY	N-CA-C	-6.48	101.20	111.18
1	A	782	MET	N-CA-C	6.47	119.65	111.69
1	C	987	ILE	N-CA-C	-6.47	104.45	110.53
1	C	52	ALA	N-CA-C	6.45	120.07	110.52
1	A	625	ALA	N-CA-C	6.39	119.51	110.50
1	B	877	GLN	N-CA-C	6.38	118.23	111.28
1	A	134	ALA	N-CA-C	6.37	121.22	112.04
1	A	412	VAL	N-CA-C	6.37	117.15	110.72
1	A	58	GLU	N-CA-C	6.36	118.00	111.14
1	A	1035	VAL	N-CA-C	6.35	116.52	110.42
1	C	38	ILE	N-CA-C	6.32	117.10	110.72
1	B	83	ALA	N-CA-C	6.29	118.97	109.41
1	C	208	GLN	N-CA-C	-6.25	103.05	112.04
1	C	771	LYS	N-CA-C	6.25	119.65	110.59
1	A	293	ILE	N-CA-C	6.24	118.61	109.63
1	C	806	ALA	N-CA-C	6.21	119.61	109.24
1	C	934	SER	N-CA-C	6.21	117.75	107.20
1	B	491	ALA	N-CA-C	-6.19	103.12	112.04
1	A	404	LEU	N-CA-C	6.17	118.09	111.36
1	A	757	GLY	N-CA-C	6.15	122.80	113.02
1	C	44	THR	N-CA-C	6.15	118.97	109.07
1	B	901	SER	N-CA-C	6.15	119.25	111.69
1	A	217	GLY	N-CA-C	6.12	122.86	112.22
1	B	582	PRO	N-CA-C	6.09	120.77	111.15
1	C	836	VAL	N-CA-C	6.09	118.39	109.63
1	C	609	SER	N-CA-C	6.08	118.01	108.96
1	C	453	PHE	N-CA-C	6.07	117.98	111.36
1	C	286	ALA	N-CA-C	6.07	118.29	109.07
1	C	865	THR	N-CA-C	6.06	118.39	109.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	458	PHE	N-CA-C	6.06	118.85	111.82
1	B	6	ILE	N-CA-C	-6.05	104.61	110.72
1	A	174	GLU	N-CA-C	6.04	119.93	110.32
1	A	658	ILE	N-CA-C	6.04	116.16	110.30
1	B	65	ILE	N-CA-C	6.04	116.81	110.72
1	C	2	ALA	N-CA-C	-6.01	105.95	113.28
1	A	277	PHE	N-CA-C	5.98	119.14	109.40
1	A	1027	PHE	N-CA-C	5.97	120.56	113.28
1	B	721	GLY	N-CA-C	5.95	120.10	112.13
1	A	736	GLN	N-CA-C	-5.92	104.76	112.23
1	C	509	ASN	N-CA-C	-5.92	100.58	108.34
1	A	762	ASP	N-CA-C	5.91	119.03	109.40
1	B	217	GLY	N-CA-C	5.87	123.59	112.62
1	A	454	VAL	CA-C-N	-5.86	113.58	119.56
1	A	454	VAL	C-N-CA	-5.86	113.58	119.56
1	C	1032	PHE	N-CA-C	5.85	117.65	111.28
1	C	439	SER	N-CA-C	5.83	117.31	111.07
1	B	180	GLY	N-CA-C	5.82	121.66	111.25
1	A	1015	GLY	N-CA-C	5.80	119.23	112.50
1	B	453	PHE	N-CA-C	5.78	117.66	111.36
1	A	273	ASP	N-CA-C	-5.77	104.91	111.14
1	C	454	VAL	N-CA-C	5.74	120.92	112.61
1	C	852	LEU	CA-C-N	-5.71	112.70	119.84
1	C	852	LEU	C-N-CA	-5.71	112.70	119.84
1	A	170	ALA	N-CA-C	5.71	118.47	109.96
1	A	48	THR	N-CA-C	5.65	118.11	108.90
1	B	128	THR	N-CA-C	5.57	119.32	109.96
1	C	800	VAL	CA-C-N	-5.56	114.60	120.66
1	C	800	VAL	C-N-CA	-5.56	114.60	120.66
1	B	825	SER	N-CA-C	5.55	117.62	109.24
1	B	40	PRO	CA-C-N	-5.54	114.22	119.76
1	B	40	PRO	C-N-CA	-5.54	114.22	119.76
1	C	794	ASN	N-CA-C	5.54	119.74	112.92
1	C	170	ALA	N-CA-C	5.51	118.32	110.10
1	C	460	GLY	N-CA-C	5.51	121.73	112.85
1	C	36	PRO	CA-C-O	-5.49	115.51	121.77
1	C	87	GLN	N-CA-C	5.46	118.14	109.24
1	C	39	ALA	N-CA-C	5.45	117.42	109.84
1	B	214	VAL	N-CA-C	5.44	116.41	108.58
1	B	288	GLY	N-CA-C	5.43	119.21	110.90
1	A	836	VAL	N-CA-C	5.41	117.36	108.97
1	B	718	ARG	N-CA-C	5.40	118.30	110.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	313	LEU	N-CA-C	5.39	119.12	112.54
1	A	83	ALA	N-CA-C	5.38	118.35	109.85
1	A	968	GLY	N-CA-C	5.38	122.74	115.32
1	A	589	THR	N-CA-C	-5.37	105.46	112.23
1	A	645	ILE	N-CA-C	5.37	116.10	110.62
1	A	130	THR	N-CA-C	5.36	117.97	109.50
1	B	745	ALA	N-CA-C	-5.36	105.88	112.90
1	A	834	PRO	N-CA-C	5.35	121.72	113.75
1	C	128	THR	N-CA-C	5.35	118.45	109.95
1	A	568	GLU	N-CA-C	5.33	118.61	110.20
1	C	86	GLY	N-CA-C	-5.33	107.86	115.64
1	C	411	VAL	N-CA-C	5.32	115.53	110.42
1	A	357	LEU	N-CA-C	-5.32	105.15	111.69
1	B	326	ALA	N-CA-C	5.29	117.12	111.36
1	B	775	GLN	N-CA-C	-5.28	101.04	109.07
1	A	237	GLN	N-CA-C	5.28	117.88	110.23
1	A	407	ASP	N-CA-C	5.27	117.03	111.28
1	A	752	ARG	N-CA-C	-5.27	105.13	112.45
1	C	755	TRP	N-CA-C	5.27	116.83	111.14
1	A	223	PRO	N-CA-C	-5.26	101.62	112.47
1	C	967	LYS	N-CA-C	-5.25	103.11	110.35
1	B	87	GLN	N-CA-C	5.24	118.03	109.59
1	B	228	GLN	N-CA-C	-5.23	102.54	110.28
1	A	607	VAL	N-CA-C	5.22	115.64	108.12
1	C	645	ILE	N-CA-C	5.21	116.30	111.45
1	B	230	LEU	N-CA-C	5.20	117.44	109.07
1	B	212	SER	N-CA-C	5.19	117.39	110.53
1	C	852	LEU	N-CA-C	5.17	120.51	113.16
1	A	654	ALA	N-CA-C	-5.17	106.93	113.18
1	B	266	ALA	N-CA-C	5.16	116.66	108.96
1	B	603	GLU	N-CA-C	-5.16	102.01	109.59
1	B	39	ALA	CA-C-N	-5.16	115.07	120.38
1	B	39	ALA	C-N-CA	-5.16	115.07	120.38
1	B	782	MET	N-CA-C	5.15	116.98	111.36
1	B	286	ALA	N-CA-C	5.15	115.84	108.74
1	A	320	GLY	N-CA-C	5.14	120.25	113.37
1	A	162	ILE	N-CA-C	5.13	117.83	112.90
1	A	563	SER	N-CA-C	5.13	117.47	109.52
1	A	120	GLU	N-CA-C	5.12	116.86	111.28
1	C	230	LEU	N-CA-C	5.11	117.24	108.90
1	A	488	PHE	N-CA-C	5.11	116.93	111.36
1	C	900	TRP	N-CA-C	-5.11	107.05	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	210	GLN	N-CA-C	5.07	119.45	113.16
1	C	80	SER	N-CA-C	5.05	117.48	109.50
1	A	744	ILE	N-CA-C	-5.04	104.68	111.44
1	B	1029	GLY	N-CA-C	5.01	118.74	112.73
1	B	625	ALA	N-CA-C	5.01	117.97	109.76

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	7979	0	8112	106	0
1	B	7958	0	8090	98	0
1	C	7986	0	8119	77	0
2	C	41	32	0	1	0
All	All	23964	32	24321	269	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 (269) 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:606:THR:HG21	1:B:648:LEU:HD22	1.52	0.90
1:C:737:ALA:HB1	1:C:742:VAL:HG23	1.61	0.83
1:A:582:PRO:HB3	1:A:725:THR:HG22	1.62	0.81
1:A:367:ILE:HD12	1:A:492:LEU:HB3	1.63	0.78
1:A:956:LEU:HD12	1:A:986:ILE:HG12	1.65	0.78
1:A:671:PRO:HG2	1:A:674:PRO:HA	1.67	0.77
1:B:713:ARG:HG2	1:B:836:VAL:HG21	1.67	0.76
1:B:606:THR:CG2	1:B:648:LEU:HD22	2.16	0.74
1:B:343:THR:HG21	1:B:1000:LEU:HG	1.70	0.72
1:A:193:LEU:HD12	1:A:265:VAL:HB	1.73	0.71
1:C:617:SER:HB2	1:C:625:ALA:HB1	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:340:VAL:HG11	1:C:395:MET:HB3	1.74	0.69
1:B:137:LEU:CD2	1:B:293:ILE:HD11	2.23	0.68
1:A:617:SER:HB2	1:A:625:ALA:HB1	1.73	0.68
1:A:762:ASP:HB3	1:A:769:VAL:HG23	1.75	0.68
1:B:68:GLN:HG3	1:B:114:ALA:HB2	1.75	0.67
1:B:367:ILE:HB	1:B:368:PRO:HD3	1.77	0.67
1:B:634:ASP:HB2	1:B:637:LYS:HG2	1.77	0.65
1:A:912:LEU:HD22	1:A:1028:LEU:HD23	1.78	0.64
1:C:737:ALA:HB1	1:C:742:VAL:CG2	2.26	0.64
1:B:343:THR:HG23	1:B:999:PRO:HB2	1.80	0.64
1:C:68:GLN:HG3	1:C:114:ALA:HB2	1.79	0.63
1:B:359:LEU:HD22	1:B:417:GLU:OE2	1.98	0.62
1:C:646:GLY:HA2	1:C:649:ILE:HG12	1.81	0.62
1:C:277:PHE:HE1	1:C:616:PHE:HB2	1.63	0.62
1:B:184:MET:HE1	1:B:243:PRO:HG3	1.81	0.62
1:A:367:ILE:HB	1:A:368:PRO:HD3	1.81	0.61
1:A:41:PRO:HG2	1:A:94:PHE:HB2	1.82	0.60
1:B:905:SER:HB3	1:B:1040:ILE:HG12	1.83	0.60
1:C:38:ILE:HG22	1:C:462:THR:HG23	1.82	0.59
1:A:340:VAL:HG13	1:A:395:MET:HG2	1.83	0.59
1:C:277:PHE:CE1	1:C:616:PHE:HB2	2.37	0.59
1:A:182:TYR:HB3	1:A:270:LEU:HD11	1.84	0.59
1:C:247:LYS:HB3	1:C:263:LYS:HB3	1.83	0.59
1:B:701:ARG:HD3	1:B:826:VAL:HG11	1.84	0.58
1:C:157:TYR:CZ	1:C:318:PRO:HD3	2.38	0.58
1:A:75:LEU:HD21	1:A:78:ILE:HD11	1.86	0.58
1:B:678:VAL:HG21	1:B:867:LEU:HB3	1.85	0.58
1:A:580:LEU:HB3	1:A:581:PRO:HD2	1.84	0.58
1:B:30:MET:HE1	1:B:383:ILE:HG13	1.85	0.58
1:C:310:LEU:HD23	1:C:313:LEU:HD11	1.86	0.57
1:A:912:LEU:CD1	1:A:1029:GLY:HA2	2.35	0.57
1:B:293:ILE:HG22	1:B:294:ALA:O	2.05	0.57
1:B:869:LEU:HD12	1:B:872:ARG:HH12	1.70	0.57
1:B:362:TRP:HA	1:B:365:THR:HG22	1.87	0.57
1:B:914:VAL:HG13	1:B:943:ILE:HD12	1.87	0.56
1:A:912:LEU:HD13	1:A:1029:GLY:HA2	1.88	0.56
1:B:219:LEU:HD23	1:C:755:TRP:HZ3	1.71	0.56
1:A:537:LEU:HD21	1:A:1035:VAL:HA	1.88	0.55
1:A:11:PHE:CD1	1:B:895:ALA:HB1	2.41	0.55
1:C:459:PHE:HE1	1:C:878:ALA:HA	1.71	0.55
1:B:247:LYS:HB3	1:B:263:LYS:HB3	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:597:THR:HG23	1:A:610:ILE:HG21	1.89	0.54
1:A:2:ALA:CB	1:A:435:MET:HG2	2.38	0.54
1:C:354:VAL:HG12	1:C:988:MET:HG3	1.89	0.54
1:A:340:VAL:CG1	1:A:395:MET:HG2	2.38	0.54
1:B:138:GLN:HB2	1:B:325:LEU:HD21	1.89	0.53
1:A:686:LEU:HB3	1:A:826:VAL:HG22	1.90	0.53
1:B:137:LEU:HG	1:B:293:ILE:HD11	1.90	0.53
1:C:137:LEU:HD11	1:C:330:THR:HG22	1.89	0.53
1:B:42:THR:H	1:B:295:THR:HG21	1.73	0.53
1:C:778:ALA:O	1:C:782:MET:HG2	2.07	0.53
1:B:912:LEU:HD23	1:B:1028:LEU:HB3	1.91	0.53
1:C:596:MET:HA	1:C:655:LEU:HD21	1.91	0.53
1:A:10:ILE:O	1:A:14:VAL:HG23	2.09	0.53
1:B:913:GLY:HA2	1:B:1025:ALA:HB2	1.89	0.53
1:A:174:GLU:HB2	1:A:292:LYS:HB2	1.89	0.53
1:A:592:VAL:HG13	1:A:659:ILE:HG21	1.91	0.53
1:C:333:ILE:O	1:C:337:ILE:HG12	2.09	0.53
1:C:791:TYR:HB3	1:C:799:MET:HG3	1.90	0.53
1:B:137:LEU:HG	1:B:293:ILE:CG1	2.39	0.53
1:B:380:PHE:HA	1:B:383:ILE:HG12	1.91	0.52
1:C:30:MET:HG2	1:C:390:ILE:HG13	1.90	0.52
1:C:186:ILE:HG23	1:C:268:VAL:HG22	1.91	0.52
1:B:592:VAL:HG11	1:B:662:ALA:HB2	1.90	0.52
1:A:780:SER:O	1:A:786:ASP:HB3	2.08	0.52
1:A:162:ILE:HG23	1:A:313:LEU:HD13	1.91	0.52
1:B:574:VAL:HG13	1:B:630:VAL:HB	1.91	0.52
1:A:912:LEU:O	1:A:915:ILE:HG22	2.10	0.52
1:C:355:MET:HE3	1:C:368:PRO:HG2	1.91	0.52
1:C:563:SER:HA	1:C:935:ASN:HB3	1.91	0.52
1:A:698:ILE:HD12	1:A:701:ARG:HH21	1.73	0.52
1:B:137:LEU:HG	1:B:293:ILE:HG13	1.93	0.51
1:B:985:PRO:HA	1:B:988:MET:HG2	1.92	0.51
1:A:230:LEU:HG	1:B:583:ASN:HB3	1.92	0.51
1:B:455:PRO:HG2	1:B:885:SER:HB2	1.92	0.51
1:B:189:ASP:HB3	1:B:192:LYS:HB2	1.92	0.51
1:C:764:VAL:HG22	1:C:769:VAL:HG12	1.93	0.51
1:C:982:ARG:O	1:C:986:ILE:HG12	2.11	0.51
1:B:485:ALA:O	1:B:490:PRO:HD3	2.11	0.51
1:A:557:LEU:HB3	1:A:918:ILE:HG12	1.93	0.51
1:B:501:ASP:HB2	1:B:504:LYS:HG2	1.91	0.51
1:B:997:VAL:HG21	1:B:1018:VAL:HB	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:MET:HB3	1:B:390:ILE:HD11	1.93	0.51
1:B:219:LEU:HD23	1:C:755:TRP:CZ3	2.45	0.51
1:B:889:VAL:O	1:B:893:LEU:HG	2.10	0.51
1:B:700:ALA:HA	1:B:857:LEU:HD13	1.92	0.51
1:A:240:LEU:HD22	1:A:245:GLN:HB3	1.93	0.50
1:B:137:LEU:CD2	1:B:293:ILE:CD1	2.88	0.50
1:C:592:VAL:HG13	1:C:659:ILE:HD12	1.93	0.50
1:A:34:GLN:HB2	1:A:333:ILE:HG23	1.94	0.50
1:A:390:ILE:HG23	1:A:395:MET:CE	2.42	0.50
1:A:184:MET:HE1	1:A:243:PRO:HG3	1.94	0.50
1:B:676:LEU:HD13	1:B:867:LEU:HD22	1.93	0.50
1:A:485:ALA:O	1:A:490:PRO:HD3	2.12	0.49
1:B:763:PHE:CE1	1:B:770:LYS:HB2	2.46	0.49
1:A:337:ILE:O	1:A:340:VAL:HG12	2.13	0.49
1:A:14:VAL:HG13	1:B:891:LEU:HB3	1.95	0.49
1:A:755:TRP:HZ3	1:C:219:LEU:HD23	1.78	0.49
1:B:157:TYR:CZ	1:B:318:PRO:HD3	2.48	0.49
1:A:778:ALA:O	1:A:782:MET:HG2	2.13	0.49
1:B:521:GLY:O	1:B:524:ARG:HG2	2.13	0.49
1:A:69:MET:HE1	1:A:107:VAL:HA	1.95	0.48
1:A:396:PHE:CE1	1:A:1011:GLN:HG2	2.48	0.48
1:C:364:ALA:HB2	1:C:497:LEU:HD21	1.96	0.48
1:B:38:ILE:HG13	1:B:40:PRO:HD3	1.95	0.48
1:B:924:GLY:O	1:B:928:LYS:HB2	2.14	0.48
1:A:121:ASP:O	1:A:125:GLN:HG2	2.13	0.47
1:A:762:ASP:HB3	1:A:769:VAL:CG2	2.43	0.47
1:A:802:PHE:HA	1:A:805:PHE:CZ	2.49	0.47
1:B:359:LEU:CD2	1:B:417:GLU:OE2	2.62	0.47
1:C:456:MET:HB2	1:C:471:SER:HB2	1.96	0.47
1:A:224:ALA:HB1	1:B:782:MET:HE1	1.96	0.47
1:A:69:MET:HE3	1:A:72:LEU:HD11	1.96	0.47
1:C:593:ILE:HA	1:C:596:MET:HE3	1.95	0.47
1:B:33:ALA:O	1:B:391:ASN:HA	2.15	0.47
1:B:137:LEU:CG	1:B:293:ILE:HD11	2.45	0.47
1:C:327:TYR:HB2	2:C:1101:80P:C29	2.45	0.47
1:C:585:THR:HG22	1:C:588:ARG:HG3	1.97	0.47
1:C:175:VAL:HG22	1:C:291:ILE:CD1	2.45	0.47
1:A:912:LEU:HD11	1:A:1032:PHE:CG	2.51	0.46
1:B:219:LEU:O	1:B:231:ASN:HA	2.14	0.46
1:C:908:LEU:O	1:C:911:PRO:HD2	2.14	0.46
1:A:219:LEU:HD23	1:B:755:TRP:HZ3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1003:SER:HB3	1:C:1012:HIS:HD2	1.81	0.46
1:A:17:LEU:HA	1:A:20:MET:HE3	1.98	0.46
1:A:343:THR:HG23	1:A:999:PRO:HB2	1.97	0.46
1:A:14:VAL:O	1:A:18:VAL:HG23	2.16	0.46
1:B:717:VAL:HG12	1:B:828:ILE:HG22	1.97	0.46
1:C:197:GLN:HA	1:C:799:MET:HE3	1.96	0.46
1:C:474:LEU:O	1:C:478:MET:HG2	2.16	0.46
1:A:403:GLY:HA3	1:A:993:PHE:HA	1.97	0.46
1:A:952:LYS:HE3	1:A:952:LYS:HB3	1.70	0.46
1:A:26:THR:HA	1:A:29:LYS:HZ2	1.81	0.46
1:B:137:LEU:HG	1:B:293:ILE:CD1	2.46	0.46
1:C:616:PHE:HD1	1:C:621:VAL:HG22	1.80	0.46
1:C:904:PHE:HA	1:C:907:LEU:HB2	1.98	0.46
1:B:137:LEU:HD21	1:B:293:ILE:HD11	1.97	0.46
1:B:832:PRO:HB3	1:B:841:ALA:HB2	1.98	0.46
1:C:57:VAL:HG21	1:C:86:GLY:HA2	1.97	0.45
1:A:912:LEU:CD1	1:A:1032:PHE:HB2	2.47	0.45
1:C:466:ILE:HG21	1:C:937:ILE:HD13	1.98	0.45
1:A:163:LYS:HE3	1:A:163:LYS:HB3	1.72	0.45
1:A:876:ALA:O	1:A:879:PRO:HD2	2.17	0.45
1:A:438:ILE:HG23	1:A:442:LEU:HG	1.97	0.45
1:B:137:LEU:HD23	1:B:293:ILE:CD1	2.46	0.45
1:B:221:GLY:O	1:B:224:ALA:HB2	2.17	0.45
1:B:330:THR:O	1:B:331:PRO:C	2.58	0.45
1:C:30:MET:HE1	1:C:380:PHE:HB3	1.98	0.45
1:C:952:LYS:HB3	1:C:952:LYS:HE2	1.42	0.45
1:A:580:LEU:CB	1:A:581:PRO:HD2	2.46	0.45
1:A:1007:GLY:O	1:A:1011:GLN:HG3	2.16	0.45
1:A:912:LEU:HD11	1:A:1032:PHE:HB2	1.99	0.45
1:C:418:ARG:HH12	1:C:960:PHE:HE2	1.65	0.44
1:B:247:LYS:HE2	1:B:263:LYS:HE2	1.98	0.44
1:B:137:LEU:HD21	1:B:299:ALA:HB1	2.00	0.44
1:B:778:ALA:O	1:B:782:MET:HG2	2.18	0.44
1:A:17:LEU:HD23	1:A:20:MET:CE	2.48	0.44
1:A:562:PRO:O	1:A:934:SER:HB2	2.18	0.44
1:B:427:PRO:HD3	1:B:499:GLN:HB2	2.00	0.44
1:A:678:VAL:HG22	1:A:679:THR:HG23	1.99	0.44
1:B:560:LYS:HA	1:B:560:LYS:HD3	1.81	0.43
1:C:337:ILE:O	1:C:340:VAL:HG22	2.19	0.43
1:C:616:PHE:CD1	1:C:621:VAL:HG22	2.52	0.43
1:A:142:PHE:HB3	1:A:321:LEU:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:ILE:HG23	1:A:367:ILE:HD13	2.00	0.43
1:C:340:VAL:HG21	1:C:395:MET:HE2	2.01	0.43
1:A:955:ILE:HA	1:A:955:ILE:HD13	1.74	0.43
1:B:890:PHE:HD1	1:B:907:LEU:HD11	1.83	0.43
1:C:191:ALA:O	1:C:194:THR:HG22	2.19	0.43
1:B:684:LEU:HD11	1:B:862:TYR:HB2	2.00	0.43
1:C:175:VAL:HG22	1:C:291:ILE:HD13	2.00	0.43
1:B:141:ALA:O	1:B:323:ASP:HA	2.17	0.43
1:B:906:VAL:O	1:B:909:VAL:HG22	2.19	0.43
1:B:564:SER:O	1:B:936:ASN:HA	2.19	0.43
1:C:144:SER:HA	1:C:145:PRO:HD2	1.92	0.43
1:C:909:VAL:HG23	1:C:912:LEU:HD12	2.01	0.43
1:A:225:VAL:HG12	1:B:782:MET:HE2	2.01	0.43
1:B:890:PHE:CE1	1:B:903:PRO:HB2	2.54	0.43
1:C:150:SER:O	1:C:154:ILE:HG13	2.18	0.43
1:C:455:PRO:HA	1:C:458:PHE:CD2	2.54	0.43
1:A:136:PHE:CD1	1:A:292:LYS:HG3	2.54	0.43
1:A:701:ARG:HG3	1:A:826:VAL:HG21	2.01	0.43
1:A:1030:ILE:HG13	1:A:1031:PHE:N	2.34	0.43
1:B:188:LEU:HB2	1:B:776:GLY:HA2	2.01	0.43
1:A:930:ASP:HA	1:A:931:PRO:HD2	1.90	0.42
1:B:906:VAL:HG11	1:B:955:ILE:HD11	2.01	0.42
1:C:536:MET:HE2	1:C:1031:PHE:HB3	2.01	0.42
1:A:872:ARG:HA	1:A:872:ARG:HD3	1.85	0.42
1:B:487:THR:C	1:B:490:PRO:HD2	2.44	0.42
1:A:450:THR:HG23	1:A:475:VAL:HG13	2.01	0.42
1:A:10:ILE:HD12	1:B:900:TRP:NE1	2.35	0.42
1:C:332:PHE:HZ	1:C:571:GLN:O	2.02	0.42
1:C:400:LEU:HD23	1:C:400:LEU:HA	1.80	0.42
1:A:842:MET:HE2	1:A:872:ARG:HH21	1.84	0.42
1:C:627:ILE:HD12	1:C:629:PHE:HE2	1.83	0.42
1:A:349:ILE:O	1:A:353:ILE:HG12	2.19	0.42
1:C:786:ASP:HA	1:C:789:LYS:HD3	2.01	0.42
1:A:557:LEU:HD23	1:A:557:LEU:HA	1.76	0.42
1:C:1024:SER:O	1:C:1028:LEU:HB2	2.20	0.42
1:A:586:LEU:HD13	1:A:623:GLN:HA	2.02	0.42
1:A:428:VAL:HG22	1:A:494:ALA:HB1	2.02	0.41
1:C:189:ASP:HB3	1:C:192:LYS:HB2	2.02	0.41
1:B:608:GLU:HB2	1:B:633:LYS:HG2	2.01	0.41
1:C:463:THR:HG23	1:C:565:PHE:HE1	1.86	0.41
1:C:705:LEU:HD23	1:C:717:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:580:LEU:HB3	1:A:581:PRO:CD	2.50	0.41
1:C:155:LYS:HD2	1:C:179:GLY:HA3	2.02	0.41
1:C:586:LEU:HD13	1:C:623:GLN:HA	2.02	0.41
1:C:736:GLN:O	1:C:740:MET:HG2	2.21	0.41
1:A:452:VAL:HG22	1:A:944:ILE:HG23	2.01	0.41
1:A:791:TYR:HB3	1:A:799:MET:HE3	2.01	0.41
1:A:850:GLY:O	1:A:853:PRO:HD2	2.20	0.41
1:B:41:PRO:HG2	1:B:94:PHE:HB2	2.02	0.41
1:B:886:LEU:O	1:B:889:VAL:HG12	2.21	0.41
1:A:222:ALA:HB3	1:A:223:PRO:HD3	2.01	0.41
1:B:957:ILE:CD1	1:B:982:ARG:HB2	2.50	0.41
1:A:360:GLN:HG2	1:A:515:PHE:CG	2.55	0.41
1:A:696:LYS:HD2	1:A:696:LYS:HA	1.85	0.41
1:A:912:LEU:HD11	1:A:1032:PHE:CD2	2.56	0.41
1:B:351:VAL:HG22	1:B:992:ALA:HB1	2.02	0.41
1:B:448:VAL:HG12	1:B:892:CYS:HB3	2.02	0.41
1:B:557:LEU:HD23	1:B:557:LEU:HA	1.91	0.41
1:A:908:LEU:O	1:A:911:PRO:HD2	2.21	0.41
1:B:793:ARG:HB2	1:B:799:MET:HE2	2.02	0.41
1:B:897:TYR:HB3	1:B:902:ILE:HD12	2.03	0.41
1:B:902:ILE:HB	1:B:903:PRO:HD3	2.03	0.41
1:C:155:LYS:HD2	1:C:179:GLY:CA	2.50	0.41
1:B:200:PRO:HG2	1:B:750:THR:HG23	2.03	0.41
1:B:673:MET:HA	1:B:674:PRO:HD3	1.84	0.41
1:C:366:ILE:HD11	1:C:496:ILE:HG21	2.02	0.41
1:A:118:LEU:HD23	1:A:118:LEU:HA	1.82	0.41
1:A:437:GLN:H	1:A:437:GLN:HG3	1.60	0.41
1:A:455:PRO:HG3	1:A:888:ILE:HG13	2.03	0.41
1:C:748:ASN:HA	1:C:751:MET:HE2	2.02	0.41
1:A:27:LEU:HD23	1:A:27:LEU:HA	1.91	0.41
1:A:36:PRO:HD2	1:A:393:LEU:HD23	2.03	0.41
1:A:246:PHE:O	1:A:262:LEU:HD23	2.21	0.41
1:A:360:GLN:HG2	1:A:515:PHE:CD1	2.56	0.41
1:A:915:ILE:HD12	1:A:915:ILE:HA	1.91	0.41
1:B:163:LYS:HE2	1:B:163:LYS:HB3	1.68	0.41
1:A:744:ILE:HA	1:A:747:ILE:HD12	2.04	0.40
1:C:897:TYR:HE1	1:C:959:GLU:HG2	1.85	0.40
1:A:15:ILE:O	1:A:19:ILE:HG13	2.21	0.40
1:A:36:PRO:CD	1:A:393:LEU:HD23	2.51	0.40
1:A:836:VAL:CG1	1:A:840:ASP:HB2	2.51	0.40
1:A:912:LEU:HD11	1:A:1032:PHE:CB	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:MET:HE2	1:C:20:MET:HB2	1.68	0.40
1:A:119:PRO:HG3	1:C:769:VAL:HG21	2.02	0.40
1:B:99:ASP:HA	1:B:100:PRO:HD2	1.96	0.40
1:B:138:GLN:HA	1:B:327:TYR:O	2.20	0.40
1:B:688:ASP:HB2	1:B:697:LEU:HD13	2.02	0.40
1:C:792:VAL:HG12	1:C:793:ARG:N	2.36	0.40
1:A:216:VAL:HG12	1:B:751:MET:HB3	2.01	0.40
1:A:724:ASP:HA	1:A:815:PRO:HD3	2.03	0.40
1:B:488:PHE:CE2	1:B:492:LEU:HD11	2.56	0.40
1:C:814:SER:HA	1:C:815:PRO:HD3	1.91	0.40
1:C:137:LEU:C	1:C:138:GLN:HG3	2.47	0.40
1:C:867:LEU:HD12	1:C:867:LEU:HA	1.98	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	1045/1058 (99%)	1028 (98%)	16 (2%)	1 (0%)	48	76
1	B	1043/1058 (99%)	1027 (98%)	16 (2%)	0	100	100
1	C	1046/1058 (99%)	1028 (98%)	17 (2%)	1 (0%)	48	76
All	All	3134/3174 (99%)	3083 (98%)	49 (2%)	2 (0%)	49	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	754	ALA
1	A	223	PRO

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	863/874 (99%)	831 (96%)	32 (4%)	30	63
1	B	861/874 (98%)	837 (97%)	24 (3%)	38	70
1	C	864/874 (99%)	837 (97%)	27 (3%)	35	67
All	All	2588/2622 (99%)	2505 (97%)	83 (3%)	35	66

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ILE
1	A	68	GLN
1	A	162	ILE
1	A	177	VAL
1	A	214	VAL
1	A	230	LEU
1	A	312	GLU
1	A	438	ILE
1	A	496	ILE
1	A	497	LEU
1	A	510	ILE
1	A	511	PHE
1	A	570	ASP
1	A	606	THR
1	A	612	THR
1	A	630	VAL
1	A	640	THR
1	A	678	VAL
1	A	683	ASN
1	A	684	LEU
1	A	687	LYS
1	A	696	LYS
1	A	760	ILE
1	A	777	ASP
1	A	818	GLU
1	A	823	VAL

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Mol	Chain	Res	Type
1	A	831	THR
1	A	918	ILE
1	A	955	ILE
1	A	959	GLU
1	A	975	THR
1	A	981	MET
1	B	25	LEU
1	B	137	LEU
1	B	140	ILE
1	B	214	VAL
1	B	230	LEU
1	B	233	THR
1	B	252	LYS
1	B	272	SER
1	B	325	LEU
1	B	425	THR
1	B	428	VAL
1	B	533	VAL
1	B	534	SER
1	B	543	SER
1	B	580	LEU
1	B	650	GLN
1	B	717	VAL
1	B	736	GLN
1	B	744	ILE
1	B	802	PHE
1	B	823	VAL
1	B	909	VAL
1	B	937	ILE
1	B	1040	ILE
1	C	127	VAL
1	C	151	ASP
1	C	186	ILE
1	C	193	LEU
1	C	207	LEU
1	C	234	VAL
1	C	293	ILE
1	C	313	LEU
1	C	392	THR
1	C	396	PHE
1	C	423	ASP
1	C	463	THR

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Mol	Chain	Res	Type
1	C	546	LEU
1	C	596	MET
1	C	606	THR
1	C	640	THR
1	C	659	ILE
1	C	670	LEU
1	C	687	LYS
1	C	714	LEU
1	C	781	ARG
1	C	857	LEU
1	C	867	LEU
1	C	905	SER
1	C	908	LEU
1	C	936	ASN
1	C	1012	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	213	GLN
1	A	231	ASN
1	A	274	ASN
1	A	308	GLN
1	A	424	HIS
1	A	469	GLN
1	A	508	ASN
1	A	827	ASN
1	A	977	HIS
1	B	274	ASN
1	B	342	HIS
1	B	650	GLN
1	B	683	ASN
1	B	827	ASN
1	B	935	ASN
1	C	70	ASN
1	C	93	ASN
1	C	360	GLN
1	C	519	ASN
1	C	530	GLN
1	C	683	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

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	80P	C	1101	-	45,45,45	4.84	22 (48%)	51,72,72	2.70	20 (39%)

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	80P	C	1101	-	-	14/22/87/87	0/5/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1101	80P	C08-C09	14.87	1.65	1.52
2	C	1101	80P	C21-C04	14.50	1.67	1.54
2	C	1101	80P	C04-N03	11.58	1.58	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1101	80P	C08-C21	10.68	1.62	1.53
2	C	1101	80P	C06-C36	7.35	1.62	1.47
2	C	1101	80P	C08-C07	6.62	1.63	1.55
2	C	1101	80P	C04-C05	6.17	1.63	1.51
2	C	1101	80P	C18-C17	5.57	1.59	1.51
2	C	1101	80P	C15-C26	5.15	1.59	1.52
2	C	1101	80P	C06-C07	5.10	1.60	1.46
2	C	1101	80P	C18-C19	4.59	1.65	1.52
2	C	1101	80P	C16-C17	4.42	1.47	1.40
2	C	1101	80P	C22-C16	3.80	1.60	1.51
2	C	1101	80P	C12-C13	3.78	1.47	1.41
2	C	1101	80P	C12-C11	3.64	1.55	1.46
2	C	1101	80P	C10-C11	3.27	1.54	1.46
2	C	1101	80P	C16-C15	3.01	1.45	1.40
2	C	1101	80P	C40-N03	2.91	1.55	1.47
2	C	1101	80P	C36-N38	2.60	1.40	1.33
2	C	1101	80P	C01-N03	2.25	1.53	1.47
2	C	1101	80P	C14-C15	2.24	1.43	1.39
2	C	1101	80P	C20-C19	2.20	1.58	1.53

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1101	80P	C11-C10-C09	8.62	125.62	118.80
2	C	1101	80P	C17-C18-C19	7.26	122.73	113.12
2	C	1101	80P	C36-C06-C07	-6.12	113.73	120.97
2	C	1101	80P	C12-C11-C10	4.94	127.54	118.28
2	C	1101	80P	O33-C09-C08	-4.17	107.33	113.37
2	C	1101	80P	O32-C11-C10	-3.83	114.49	121.14
2	C	1101	80P	C18-C17-C16	3.64	125.55	119.75
2	C	1101	80P	C08-C21-C04	3.60	116.56	111.64
2	C	1101	80P	F23-C22-C16	3.58	120.66	112.32
2	C	1101	80P	C13-C12-C17	3.43	122.72	119.03
2	C	1101	80P	C21-C08-C07	-3.43	107.12	111.05
2	C	1101	80P	O33-C09-C10	3.21	129.75	123.52
2	C	1101	80P	F25-C22-C16	3.13	119.62	112.32
2	C	1101	80P	O37-C36-N38	-2.88	116.43	122.89
2	C	1101	80P	O32-C11-C12	-2.27	117.94	122.06
2	C	1101	80P	C19-C20-C21	2.25	114.21	110.38
2	C	1101	80P	C07-C08-C09	2.20	112.46	109.88
2	C	1101	80P	O39-C05-C06	-2.12	119.38	122.93
2	C	1101	80P	C14-C13-C12	-2.10	118.56	120.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1101	80P	C15-C26-N27	2.00	115.81	112.59

There are no chirality outliers.

All (14) torsion outliers are listed below:

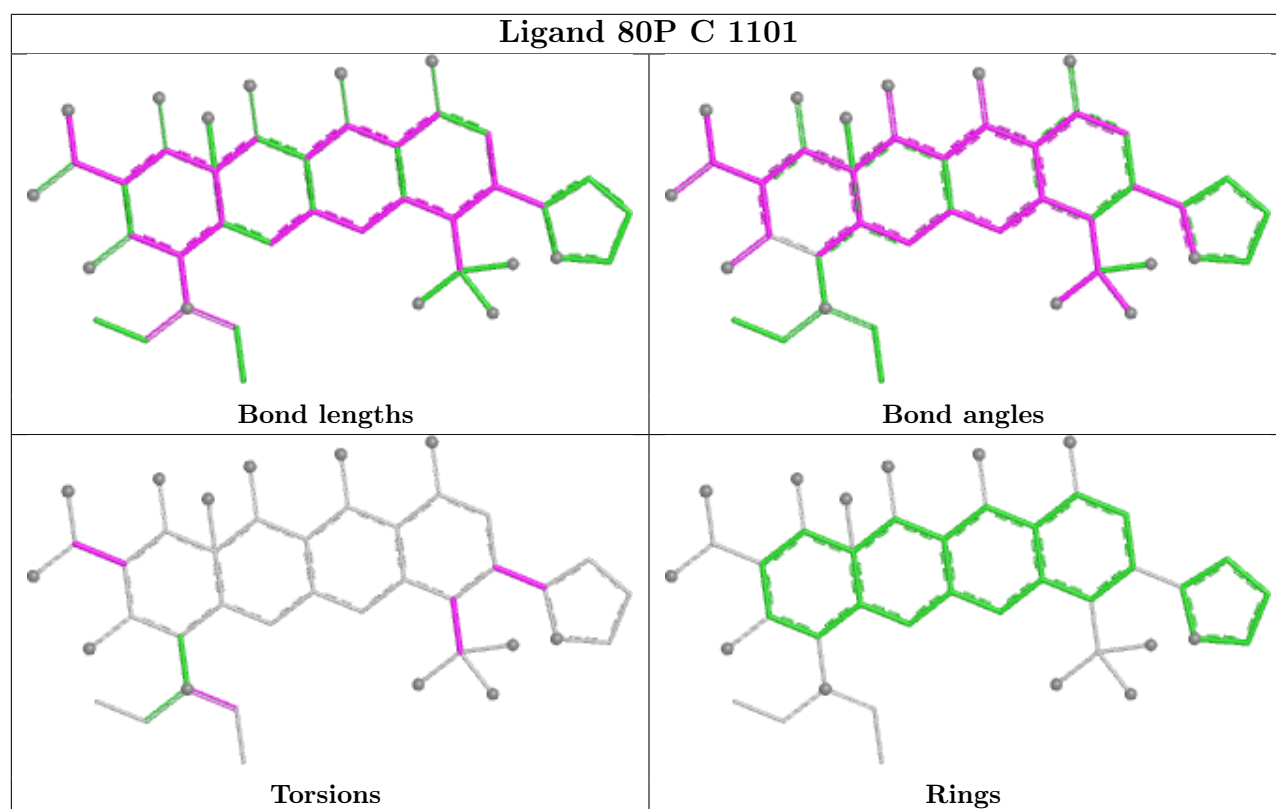
Mol	Chain	Res	Type	Atoms
2	C	1101	80P	C05-C06-C36-N38
2	C	1101	80P	C05-C06-C36-O37
2	C	1101	80P	C07-C06-C36-N38
2	C	1101	80P	C07-C06-C36-O37
2	C	1101	80P	C14-C15-C26-N27
2	C	1101	80P	C16-C15-C26-N27
2	C	1101	80P	C14-C15-C26-C30
2	C	1101	80P	C15-C16-C22-F23
2	C	1101	80P	C41-C40-N03-C04
2	C	1101	80P	C17-C16-C22-F23
2	C	1101	80P	C41-C40-N03-C01
2	C	1101	80P	C15-C16-C22-F24
2	C	1101	80P	C15-C16-C22-F25
2	C	1101	80P	C17-C16-C22-F24

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1101	80P	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.

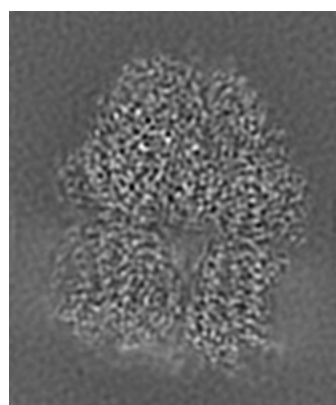
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-24732. These allow visual inspection of the internal detail of the map and identification of artifacts.

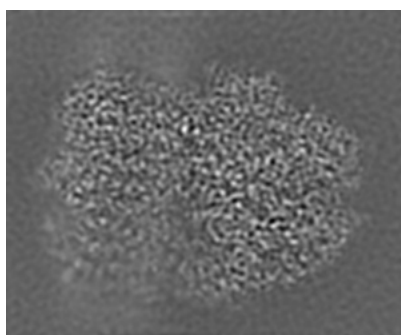
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

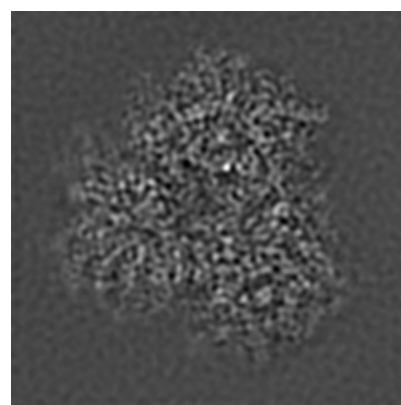
6.1.1 Primary map



X



Y

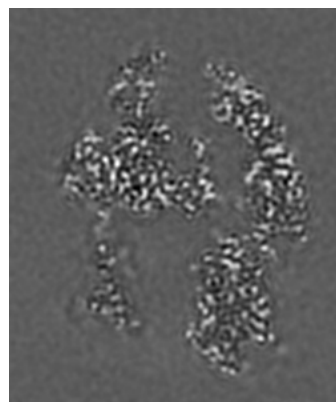


Z

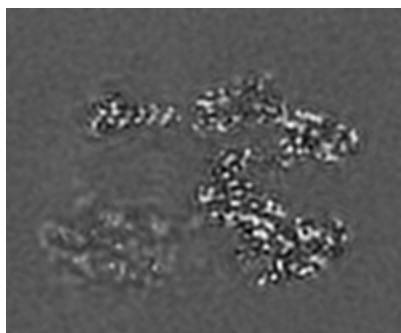
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

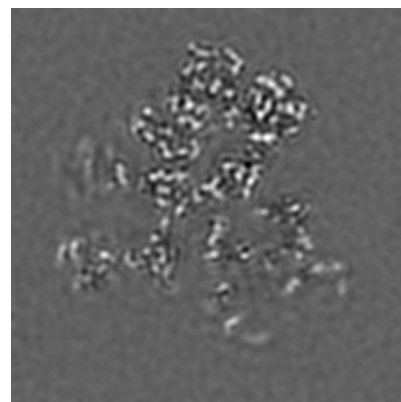
6.2.1 Primary map



X Index: 61



Y Index: 61

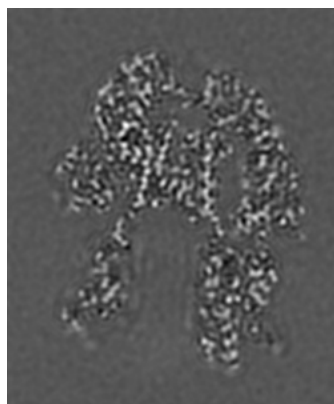


Z Index: 74

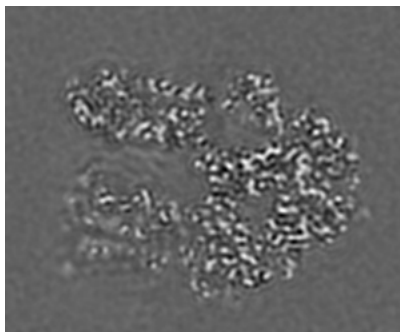
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

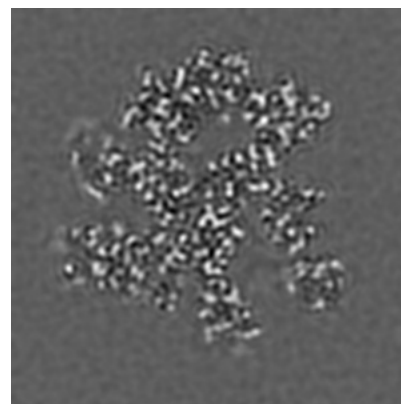
6.3.1 Primary map



X Index: 66



Y Index: 48

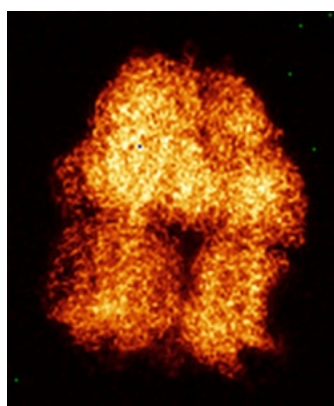


Z Index: 82

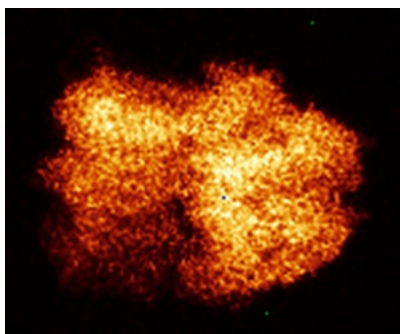
The images above show the largest variance slices of the map in three orthogonal directions.

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

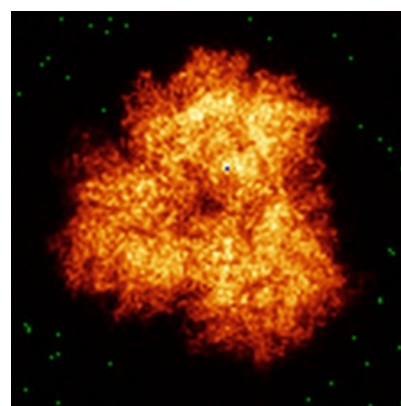
6.4.1 Primary map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

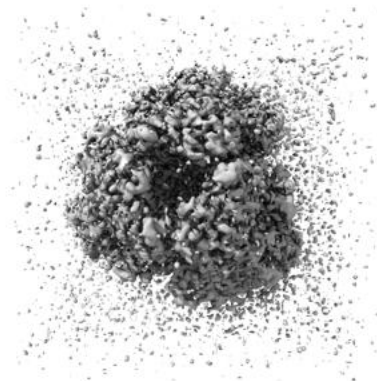
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.11. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

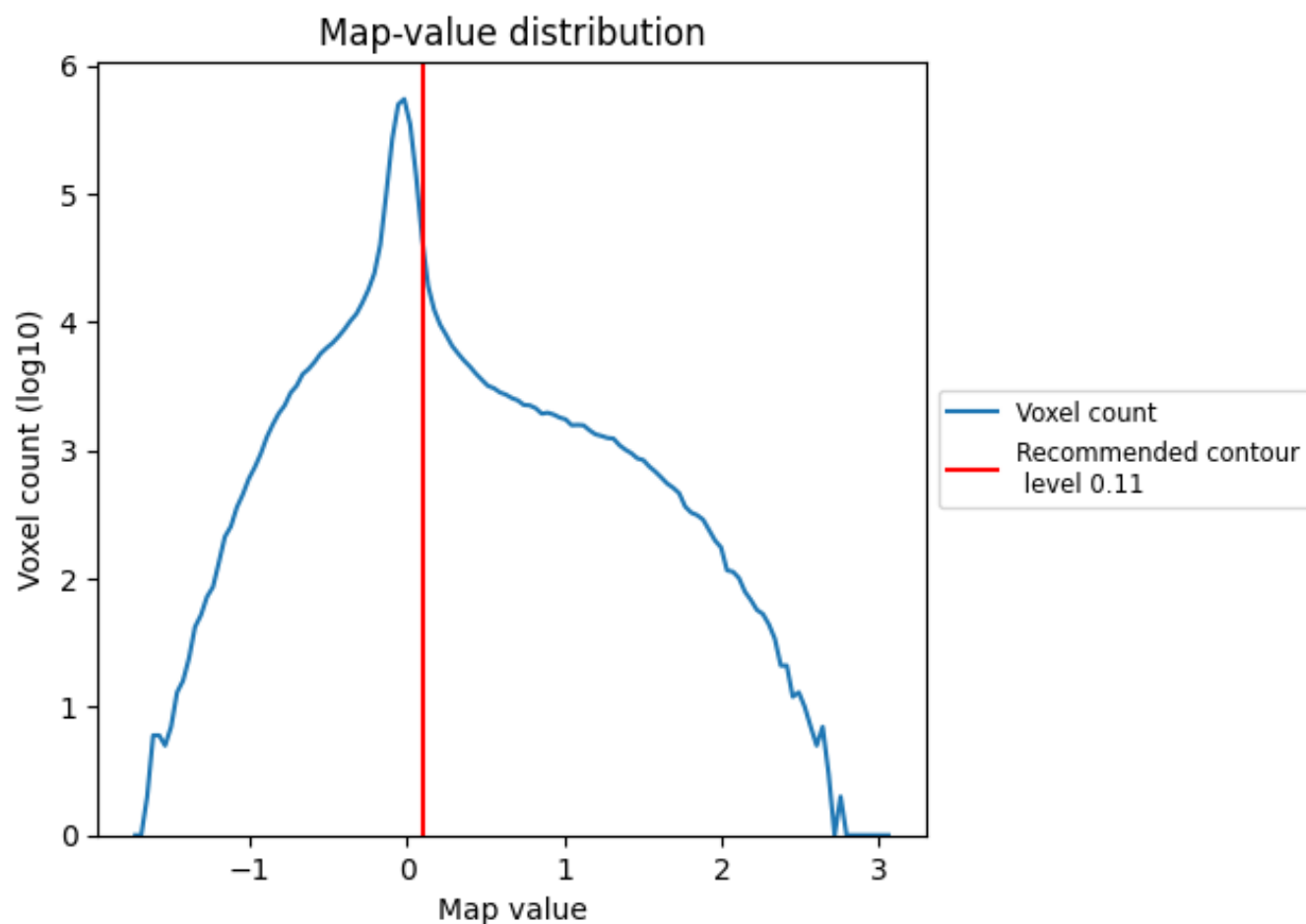
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

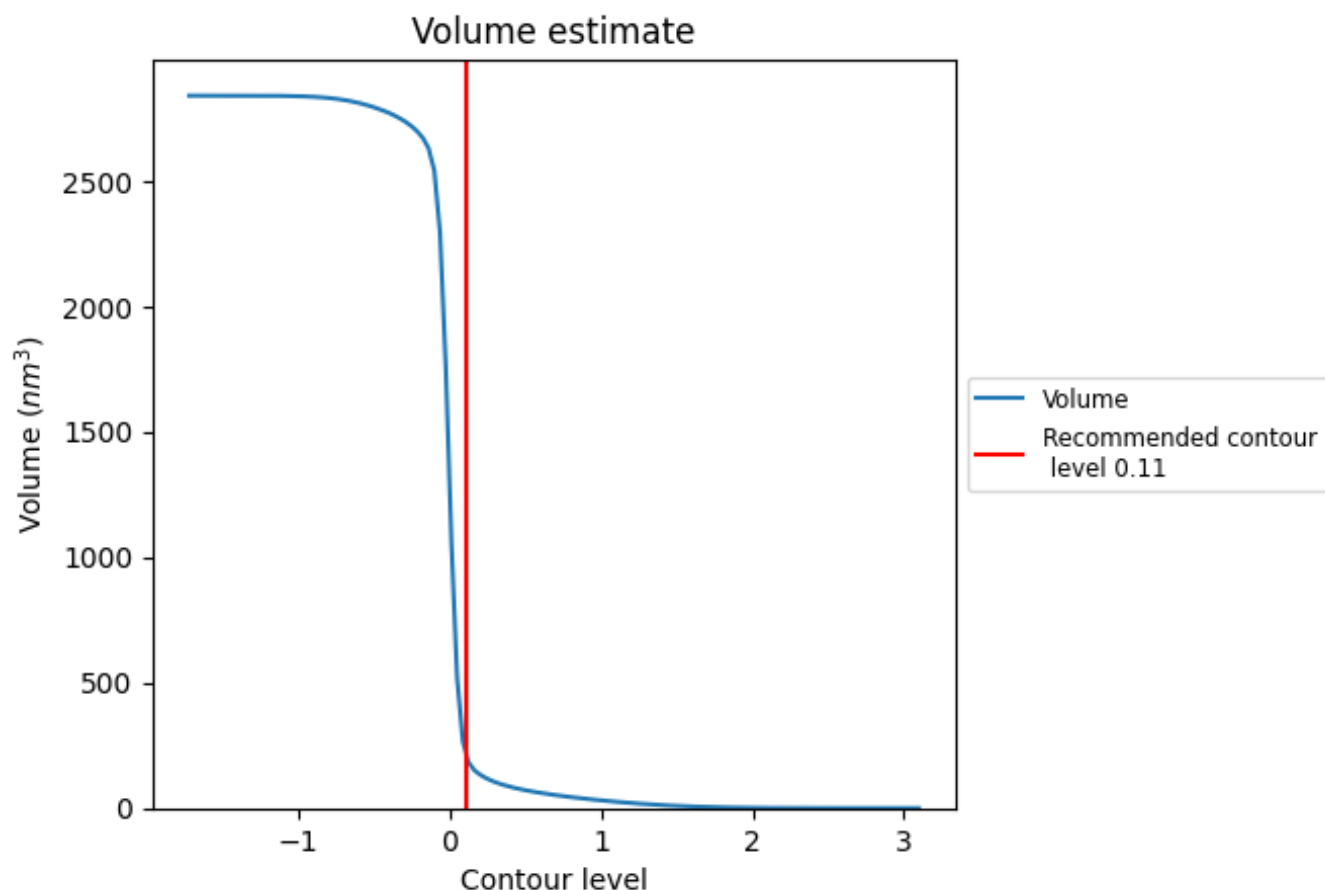
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

7.2 Volume estimate [i](#)



The volume at the recommended contour level is 210 nm³; this corresponds to an approximate mass of 190 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](#)

This section was not generated. The rotationally averaged power spectrum is only generated for cubic maps.

8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

This section contains information regarding the fit between EMD map EMD-24732 and PDB model 7RY3. Per-residue inclusion information can be found in section 3 on page 4.

9.1 Map-model overlay [i](#)

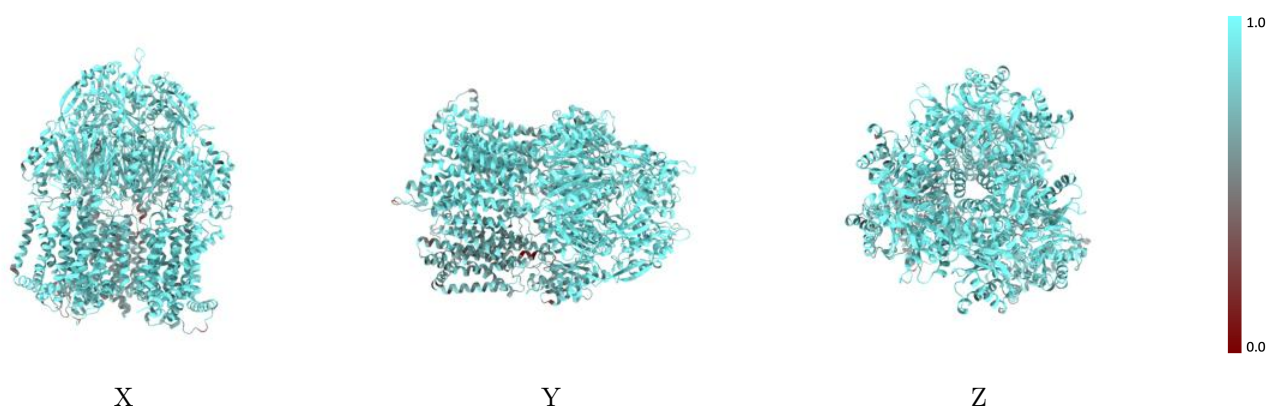
This section was not generated.

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



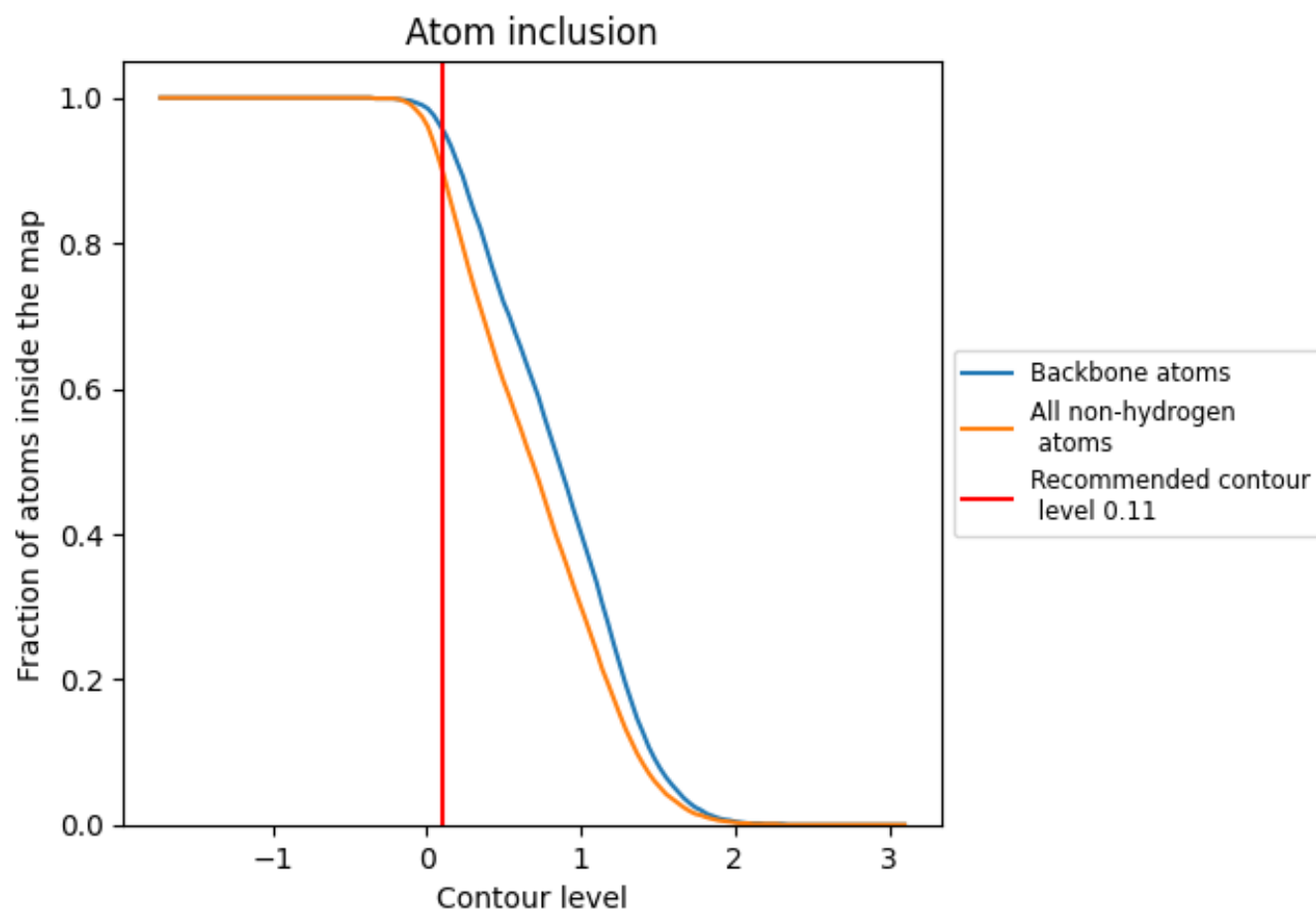
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.11).

9.4 Atom inclusion [i](#)



At the recommended contour level, 95% 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 (0.11) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8950	0.5620
A	0.9340	0.6000
B	0.8430	0.5090
C	0.9070	0.5760

