



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:49 AM UTC

PDB ID : 1XDJ / pdb_00001xdj
Title : Crystal Structure of T. maritima Cobalamin-Independent Methionine Synthase complexed with Zn²⁺ and Homocysteine
Authors : Pejchal, R.; Ludwig, M.L.
Deposited on : 2004-09-06
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

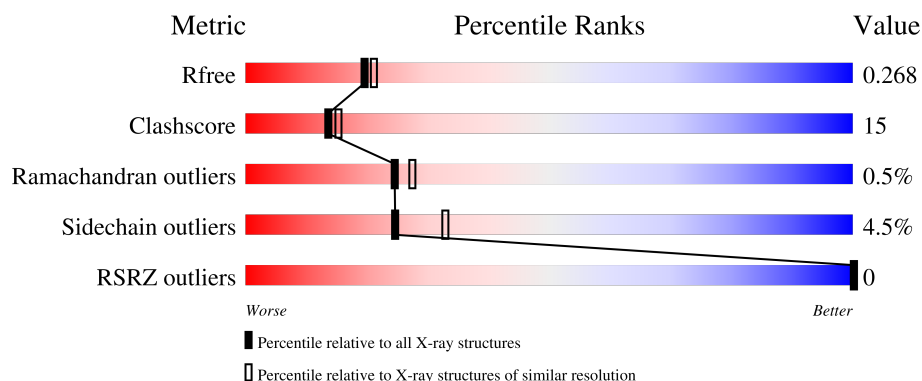
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	766	 65% 26% • 6%
1	B	766	 64% 26% • 7%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11545 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	722	Total	C	N	O	S	Se	0	0	0
			5657	3673	930	1030	7	17			
1	B	712	Total	C	N	O	S	Se	0	0	0
			5546	3604	901	1017	7	17			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-31	MET	-	expression tag	UNP Q9X112
A	-30	HIS	-	expression tag	UNP Q9X112
A	-29	HIS	-	expression tag	UNP Q9X112
A	-28	HIS	-	expression tag	UNP Q9X112
A	-27	HIS	-	expression tag	UNP Q9X112
A	-26	HIS	-	expression tag	UNP Q9X112
A	-25	HIS	-	expression tag	UNP Q9X112
A	-24	GLY	-	expression tag	UNP Q9X112
A	-23	LYS	-	expression tag	UNP Q9X112
A	-22	PRO	-	expression tag	UNP Q9X112
A	-21	ILE	-	expression tag	UNP Q9X112
A	-20	PRO	-	expression tag	UNP Q9X112
A	-19	ASN	-	expression tag	UNP Q9X112
A	-18	PRO	-	expression tag	UNP Q9X112
A	-17	LEU	-	expression tag	UNP Q9X112
A	-16	LEU	-	expression tag	UNP Q9X112
A	-15	GLY	-	expression tag	UNP Q9X112
A	-14	LEU	-	expression tag	UNP Q9X112
A	-13	ASP	-	expression tag	UNP Q9X112
A	-12	SER	-	expression tag	UNP Q9X112
A	-11	THR	-	expression tag	UNP Q9X112
A	-10	GLU	-	expression tag	UNP Q9X112
A	-9	ASN	-	expression tag	UNP Q9X112
A	-8	LEU	-	expression tag	UNP Q9X112

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	TYR	-	expression tag	UNP Q9X112
A	-6	PHE	-	expression tag	UNP Q9X112
A	-5	GLN	-	expression tag	UNP Q9X112
A	-4	GLY	-	expression tag	UNP Q9X112
A	-3	ILE	-	expression tag	UNP Q9X112
A	-2	ASP	-	expression tag	UNP Q9X112
A	-1	PRO	-	expression tag	UNP Q9X112
A	0	PHE	-	expression tag	UNP Q9X112
A	1	THR	-	expression tag	UNP Q9X112
A	38	MSE	MET	modified residue	UNP Q9X112
A	43	MSE	MET	modified residue	UNP Q9X112
A	45	MSE	MET	modified residue	UNP Q9X112
A	73	MSE	MET	modified residue	UNP Q9X112
A	93	MSE	MET	modified residue	UNP Q9X112
A	102	MSE	MET	modified residue	UNP Q9X112
A	170	MSE	MET	modified residue	UNP Q9X112
A	344	MSE	MET	modified residue	UNP Q9X112
A	421	MSE	MET	modified residue	UNP Q9X112
A	468	MSE	MET	modified residue	UNP Q9X112
A	510	MSE	MET	modified residue	UNP Q9X112
A	529	MSE	MET	modified residue	UNP Q9X112
A	537	MSE	MET	modified residue	UNP Q9X112
A	619	MSE	MET	modified residue	UNP Q9X112
A	682	MSE	MET	modified residue	UNP Q9X112
A	721	MSE	MET	modified residue	UNP Q9X112
A	728	MSE	MET	modified residue	UNP Q9X112
B	-31	MET	-	expression tag	UNP Q9X112
B	-30	HIS	-	expression tag	UNP Q9X112
B	-29	HIS	-	expression tag	UNP Q9X112
B	-28	HIS	-	expression tag	UNP Q9X112
B	-27	HIS	-	expression tag	UNP Q9X112
B	-26	HIS	-	expression tag	UNP Q9X112
B	-25	HIS	-	expression tag	UNP Q9X112
B	-24	GLY	-	expression tag	UNP Q9X112
B	-23	LYS	-	expression tag	UNP Q9X112
B	-22	PRO	-	expression tag	UNP Q9X112
B	-21	ILE	-	expression tag	UNP Q9X112
B	-20	PRO	-	expression tag	UNP Q9X112
B	-19	ASN	-	expression tag	UNP Q9X112
B	-18	PRO	-	expression tag	UNP Q9X112
B	-17	LEU	-	expression tag	UNP Q9X112
B	-16	LEU	-	expression tag	UNP Q9X112

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	GLY	-	expression tag	UNP Q9X112
B	-14	LEU	-	expression tag	UNP Q9X112
B	-13	ASP	-	expression tag	UNP Q9X112
B	-12	SER	-	expression tag	UNP Q9X112
B	-11	THR	-	expression tag	UNP Q9X112
B	-10	GLU	-	expression tag	UNP Q9X112
B	-9	ASN	-	expression tag	UNP Q9X112
B	-8	LEU	-	expression tag	UNP Q9X112
B	-7	TYR	-	expression tag	UNP Q9X112
B	-6	PHE	-	expression tag	UNP Q9X112
B	-5	GLN	-	expression tag	UNP Q9X112
B	-4	GLY	-	expression tag	UNP Q9X112
B	-3	ILE	-	expression tag	UNP Q9X112
B	-2	ASP	-	expression tag	UNP Q9X112
B	-1	PRO	-	expression tag	UNP Q9X112
B	0	PHE	-	expression tag	UNP Q9X112
B	1	THR	-	expression tag	UNP Q9X112
B	38	MSE	MET	modified residue	UNP Q9X112
B	43	MSE	MET	modified residue	UNP Q9X112
B	45	MSE	MET	modified residue	UNP Q9X112
B	73	MSE	MET	modified residue	UNP Q9X112
B	93	MSE	MET	modified residue	UNP Q9X112
B	102	MSE	MET	modified residue	UNP Q9X112
B	170	MSE	MET	modified residue	UNP Q9X112
B	344	MSE	MET	modified residue	UNP Q9X112
B	421	MSE	MET	modified residue	UNP Q9X112
B	468	MSE	MET	modified residue	UNP Q9X112
B	510	MSE	MET	modified residue	UNP Q9X112
B	529	MSE	MET	modified residue	UNP Q9X112
B	537	MSE	MET	modified residue	UNP Q9X112
B	619	MSE	MET	modified residue	UNP Q9X112
B	682	MSE	MET	modified residue	UNP Q9X112
B	721	MSE	MET	modified residue	UNP Q9X112
B	728	MSE	MET	modified residue	UNP Q9X112

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

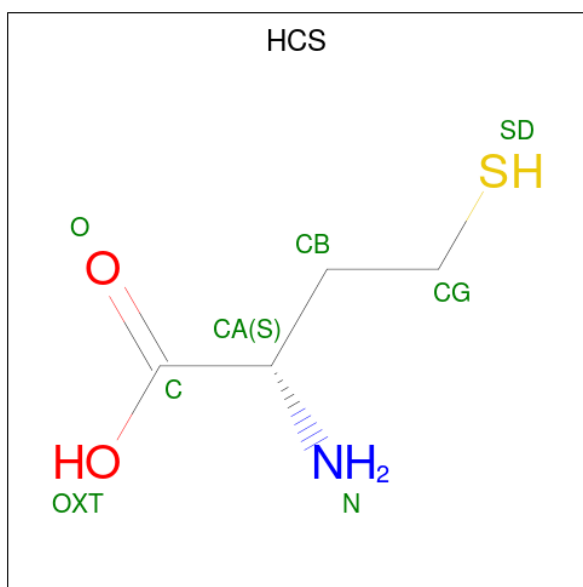


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

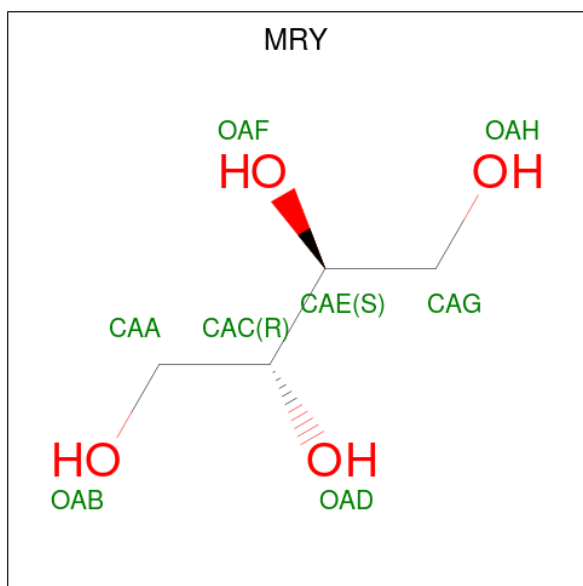
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 2-AMINO-4-MERCAPTO-BUTYRIC ACID (CCD ID: HCS) (formula: C₄H₉NO₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			8	4	1	2	1		
4	A	1	Total	C	N	O	S	0	0
			8	4	1	2	1		

- Molecule 5 is MESO-ERYTHRITOL (CCD ID: MRY) (formula: C₄H₁₀O₄).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	C O	0	0
			8	4 4		
5	B	1	Total	C O	0	0
			8	4 4		

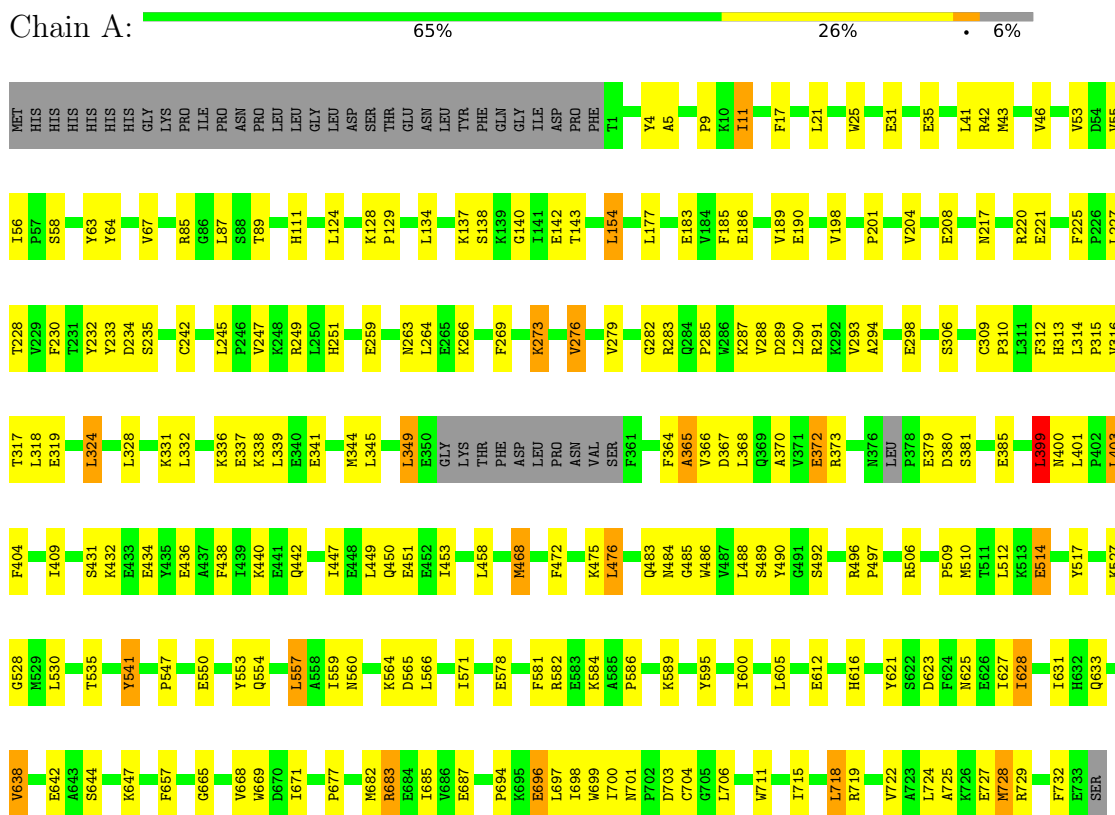
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	147	Total 147	O 147	0	0
6	B	151	Total 151	O 151	0	0

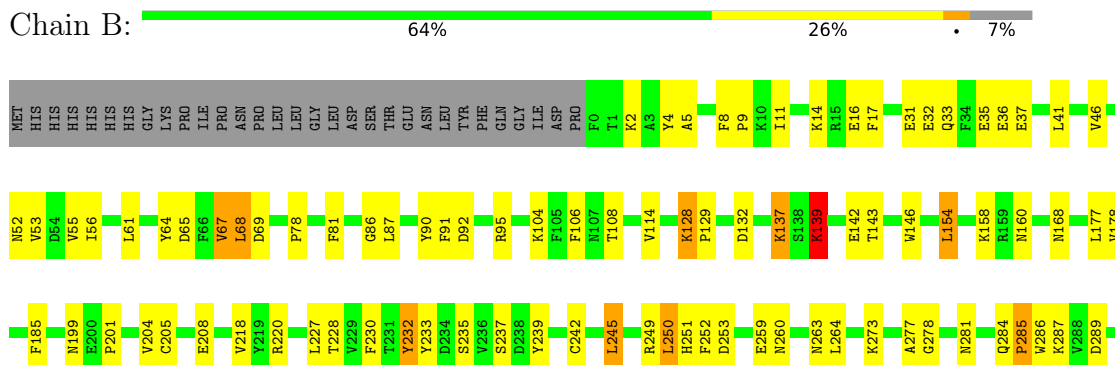
3 Residue-property plots

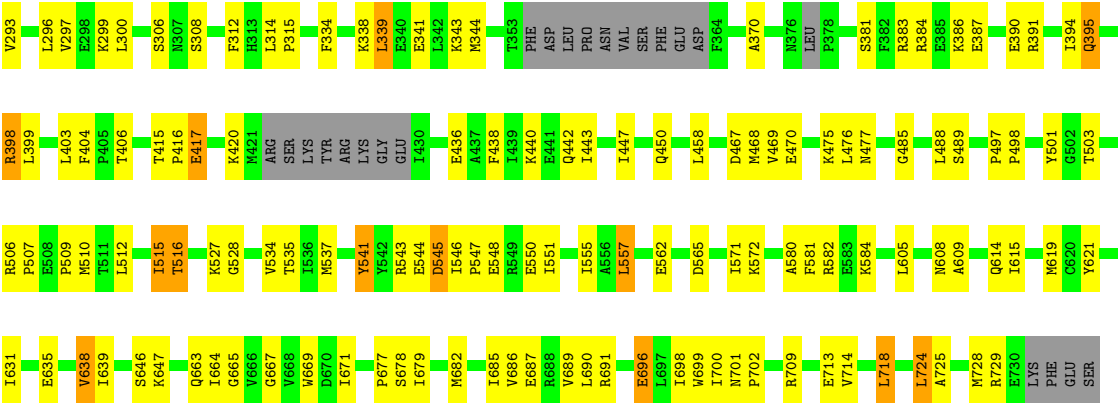
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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: 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase



- Molecule 1: 5-methyltetrahydropteroyltriglutamate--homocysteine methyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	167.05Å 159.37Å 65.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 2.20 19.90 – 2.20	Depositor EDS
% Data completeness (in resolution range)	92.9 (19.90-2.20) 93.0 (19.90-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.03 (at 2.19Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.268 0.227 , 0.268	Depositor DCC
R_{free} test set	1666 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	24.2	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 22.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.077 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11545	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HCS, ZN, SO4, MRY

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	0.44	0/5781	0.93	12/7832 (0.2%)
1	B	0.44	0/5667	0.94	16/7680 (0.2%)
All	All	0.44	0/11448	0.93	28/15512 (0.2%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	698	ILE	N-CA-C	9.36	122.54	108.23
1	A	698	ILE	N-CA-C	8.91	120.66	108.89
1	A	528	GLY	N-CA-C	-8.88	99.93	112.37
1	B	713	GLU	N-CA-C	-7.11	103.22	112.68
1	A	399	LEU	N-CA-C	6.79	118.77	111.36
1	A	541	TYR	N-CA-C	-6.75	100.54	110.52
1	A	67	VAL	N-CA-C	-6.61	103.80	111.00
1	B	541	TYR	N-CA-C	-6.55	101.71	110.55
1	B	527	LYS	N-CA-C	6.48	120.60	110.17
1	A	11	ILE	N-CA-C	6.44	119.14	111.09
1	B	528	GLY	N-CA-C	-6.36	103.93	112.14
1	A	266	LYS	N-CA-C	-6.10	106.47	114.04
1	B	139	LYS	N-CA-C	-5.88	105.42	113.30
1	B	68	LEU	N-CA-C	-5.81	104.86	111.07
1	B	469	VAL	N-CA-C	-5.70	106.66	111.56
1	A	287	LYS	N-CA-C	-5.58	102.02	110.28
1	B	220	ARG	N-CA-C	-5.50	105.37	111.36
1	B	92	ASP	N-CA-C	-5.42	105.40	112.23
1	B	67	VAL	N-CA-C	-5.33	105.93	111.58
1	B	489	SER	N-CA-C	-5.21	105.75	112.68
1	A	527	LYS	N-CA-C	5.16	117.71	110.23
1	A	220	ARG	N-CA-C	-5.14	105.36	111.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	468	MSE	N-CA-C	5.14	119.02	112.34
1	B	287	LYS	N-CA-C	-5.14	103.25	110.50
1	B	232	TYR	N-CA-C	5.09	116.03	108.60
1	B	664	ILE	N-CA-C	5.08	115.28	107.77
1	A	279	VAL	N-CA-C	5.05	118.00	113.10
1	B	277	ALA	N-CA-C	5.03	116.46	108.67

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	5657	0	5416	165	0
1	B	5546	0	5298	160	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	16	0	16	0	0
5	A	8	0	10	0	0
5	B	8	0	10	1	0
6	A	147	0	0	7	0
6	B	151	0	0	2	0
All	All	11545	0	10750	321	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 15.

All (321) 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:317:THR:HG22	1:A:319:GLU:H	1.26	0.97
1:B:228:THR:HG23	1:B:249:ARG:HG3	1.56	0.87
1:A:557:LEU:HD22	1:A:605:LEU:HD21	1.55	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:512:LEU:HD11	1:A:565:ASP:HB3	1.62	0.82
1:B:512:LEU:HD11	1:B:565:ASP:HB3	1.62	0.81
1:A:251:HIS:HA	1:A:276:VAL:HG13	1.62	0.81
1:B:4:TYR:HB2	1:B:306:SER:HB3	1.65	0.78
1:A:58:SER:HB2	1:A:143:THR:HB	1.65	0.78
1:B:677:PRO:HB2	1:B:682:MSE:HE3	1.63	0.78
1:A:557:LEU:CD2	1:A:605:LEU:HD21	2.14	0.77
1:B:296:LEU:O	1:B:300:LEU:HD13	1.85	0.76
1:B:391:ARG:O	1:B:395:GLN:HG2	1.86	0.75
1:B:619:MSE:HG3	1:B:639:ILE:HD11	1.67	0.75
1:B:450:GLN:HB3	1:B:458:LEU:HD22	1.69	0.75
1:B:399:LEU:HD11	6:B:986:HOH:O	1.87	0.74
1:A:582:ARG:HB3	1:A:621:TYR:CE2	2.24	0.72
1:B:477:ASN:HD21	1:B:507:PRO:HB3	1.55	0.72
1:A:582:ARG:HB3	1:A:621:TYR:CZ	2.25	0.72
1:B:477:ASN:ND2	1:B:507:PRO:HD3	2.05	0.72
1:A:42:ARG:O	1:A:46:VAL:HG22	1.90	0.70
1:A:290:LEU:HD22	1:A:345:LEU:HG	1.73	0.70
1:A:728:MSE:HE2	1:A:732:PHE:HE1	1.57	0.70
1:B:386:LYS:O	1:B:391:ARG:HD3	1.93	0.69
1:A:317:THR:HG23	1:A:337:GLU:OE2	1.93	0.68
1:B:245:LEU:O	1:B:273:LYS:HE3	1.94	0.68
1:B:638:VAL:HB	1:B:663:GLN:HB2	1.74	0.68
1:B:137:LYS:HD3	1:B:142:GLU:OE1	1.95	0.66
1:A:728:MSE:HE3	1:A:728:MSE:HA	1.77	0.66
1:B:686:VAL:O	1:B:690:LEU:HG	1.96	0.66
1:A:201:PRO:O	1:A:204:VAL:HG22	1.96	0.65
1:B:467:ASP:HB3	1:B:470:GLU:HB3	1.78	0.65
1:B:543:ARG:HB3	1:B:545:ASP:OD1	1.97	0.65
1:B:242:CYS:O	1:B:245:LEU:HB2	1.96	0.65
1:B:679:ILE:HG23	1:B:724:LEU:HB2	1.78	0.65
1:B:52:ASN:HA	1:B:343:LYS:HD2	1.78	0.64
1:B:114:VAL:HG11	1:B:158:LYS:HD2	1.79	0.64
1:B:239:TYR:HB2	1:B:263:ASN:HD22	1.63	0.64
1:A:475:LYS:HE2	1:A:510:MSE:O	1.98	0.64
1:B:128:LYS:HB2	1:B:128:LYS:NZ	2.13	0.63
1:B:506:ARG:NH2	1:B:562:GLU:HA	2.12	0.63
1:B:341:GLU:HA	1:B:344:MSE:HE3	1.79	0.63
1:A:4:TYR:HB2	1:A:306:SER:HB3	1.81	0.62
1:B:264:LEU:O	1:B:264:LEU:HD23	1.99	0.62
1:A:468:MSE:HE2	1:A:535:THR:HG21	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:711:TRP:O	1:A:715:ILE:HG12	1.99	0.62
1:B:250:LEU:HD13	1:B:252:PHE:HE1	1.64	0.62
1:A:668:VAL:HG23	1:A:682:MSE:HE2	1.82	0.62
1:B:546:ILE:HB	1:B:547:PRO:HD2	1.81	0.61
1:B:546:ILE:HD11	1:B:551:ILE:HG12	1.82	0.61
1:A:134:LEU:HA	1:A:137:LYS:HB2	1.82	0.61
1:A:331:LYS:HE2	1:A:364:PHE:CD1	2.36	0.60
1:B:534:VAL:HG13	1:B:584:LYS:HD3	1.83	0.60
1:A:264:LEU:HD23	1:A:264:LEU:C	2.26	0.60
1:B:14:LYS:HB2	1:B:16:GLU:OE1	2.02	0.59
1:B:501:TYR:O	1:B:543:ARG:NH1	2.30	0.59
1:A:725:ALA:O	1:A:729:ARG:HG3	2.01	0.59
1:A:512:LEU:HD23	1:A:566:LEU:HD23	1.85	0.59
1:A:550:GLU:O	1:A:554:GLN:HG3	2.01	0.59
1:B:5:ALA:HB2	1:B:53:VAL:HG11	1.86	0.58
1:B:139:LYS:NZ	1:B:139:LYS:HB2	2.19	0.58
1:A:58:SER:HB2	1:A:143:THR:CB	2.33	0.58
1:B:11:ILE:HA	1:B:17:PHE:HB3	1.85	0.58
1:A:288:VAL:HG22	1:A:289:ASP:N	2.19	0.58
1:B:232:TYR:O	1:B:233:TYR:HB2	2.03	0.58
1:A:409:ILE:HG13	1:A:704:CYS:HA	1.86	0.58
1:A:671:ILE:HD12	1:A:671:ILE:C	2.28	0.58
1:B:547:PRO:HG2	1:B:550:GLU:HB2	1.84	0.58
1:A:665:GLY:CA	1:A:701:ASN:HD21	2.17	0.57
1:A:85:ARG:H	1:A:89:THR:HG21	1.69	0.57
1:A:506:ARG:O	1:A:506:ARG:HG3	2.03	0.57
1:B:467:ASP:HB3	1:B:470:GLU:CB	2.34	0.57
1:A:438:PHE:O	1:A:442:GLN:HG2	2.03	0.57
1:B:383:ARG:HA	1:B:635:GLU:OE1	2.04	0.57
1:A:432:LYS:O	1:A:436:GLU:HB2	2.04	0.57
1:B:671:ILE:C	1:B:671:ILE:HD12	2.30	0.57
1:A:63:TYR:HD2	1:A:87:LEU:HD21	1.70	0.57
1:A:186:GLU:O	1:A:190:GLU:HG3	2.04	0.57
1:A:696:GLU:H	1:A:696:GLU:CD	2.12	0.57
1:A:373:ARG:HG3	1:A:373:ARG:HH11	1.70	0.57
1:B:4:TYR:CE1	1:B:55:VAL:HG21	2.39	0.57
1:A:242:CYS:O	1:A:245:LEU:HD13	2.04	0.56
1:A:269:PHE:CE1	1:A:273:LYS:HG3	2.39	0.56
1:B:114:VAL:HG11	1:B:158:LYS:CD	2.34	0.56
1:B:398:ARG:HD3	1:B:572:LYS:HD3	1.87	0.56
1:B:228:THR:OG1	1:B:249:ARG:HD2	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:PHE:O	1:A:189:VAL:HG23	2.06	0.56
1:A:328:LEU:O	1:A:332:LEU:HD23	2.05	0.56
1:A:154:LEU:HD13	1:A:177:LEU:HD22	1.87	0.56
1:A:294:ALA:O	1:A:298:GLU:HG3	2.06	0.56
1:B:178:VAL:CG2	1:B:218:VAL:HG13	2.36	0.56
1:B:264:LEU:HD23	1:B:264:LEU:C	2.30	0.56
1:B:609:ALA:HB3	1:B:615:ILE:HD11	1.87	0.55
1:A:581:PHE:HE2	1:A:631:ILE:HD12	1.70	0.55
1:B:128:LYS:HB2	1:B:128:LYS:HZ2	1.70	0.55
1:A:647:LYS:HZ2	1:B:691:ARG:HD3	1.70	0.55
1:B:509:PRO:HB3	1:B:512:LEU:HD12	1.87	0.55
1:A:728:MSE:HE2	1:A:732:PHE:CE1	2.40	0.55
1:B:32:GLU:O	1:B:36:GLU:HG3	2.07	0.55
1:A:647:LYS:NZ	1:B:691:ARG:HD3	2.22	0.55
1:A:449:LEU:O	1:A:453:ILE:HG12	2.07	0.55
1:B:31:GLU:O	1:B:35:GLU:HG2	2.06	0.55
1:A:727:GLU:HG3	6:A:884:HOH:O	2.07	0.54
1:A:21:LEU:HD23	1:A:492:SER:HB3	1.90	0.54
1:A:230:PHE:HA	1:A:251:HIS:HB3	1.89	0.54
1:B:178:VAL:HG22	1:B:218:VAL:HG13	1.89	0.54
1:A:547:PRO:HG2	1:A:550:GLU:HB2	1.89	0.54
1:B:33:GLN:HA	1:B:36:GLU:OE1	2.07	0.54
1:A:694:PRO:HD2	1:A:697:LEU:CD1	2.37	0.54
1:B:403:LEU:HD22	1:B:404:PHE:CD2	2.43	0.54
1:A:627:ILE:HG13	1:A:631:ILE:HD13	1.90	0.53
1:A:506:ARG:NH1	1:A:565:ASP:OD2	2.40	0.53
1:B:515:ILE:HD12	1:B:571:ILE:HD12	1.89	0.53
1:B:696:GLU:CD	1:B:696:GLU:H	2.16	0.53
1:A:403:LEU:HD13	1:A:404:PHE:CD2	2.43	0.53
1:A:628:ILE:HD12	1:A:657:PHE:CD2	2.43	0.53
1:B:250:LEU:HD13	1:B:252:PHE:CE1	2.42	0.53
1:B:678:SER:O	1:B:682:MSE:HG3	2.09	0.53
1:A:694:PRO:HD2	1:A:697:LEU:HD12	1.90	0.52
1:B:395:GLN:OE1	1:B:614:GLN:HG3	2.08	0.52
1:B:724:LEU:O	1:B:728:MSE:HG2	2.08	0.52
1:A:450:GLN:HB3	1:A:458:LEU:HD22	1.92	0.52
1:B:293:VAL:O	1:B:297:VAL:HG23	2.09	0.52
1:A:228:THR:HG23	1:A:249:ARG:HB3	1.92	0.52
1:A:4:TYR:CE1	1:A:55:VAL:HG21	2.46	0.51
1:A:317:THR:HG22	1:A:319:GLU:N	2.09	0.51
1:A:315:PRO:HB3	1:A:485:GLY:HA2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:669:TRP:CD1	1:A:677:PRO:HG2	2.46	0.51
1:A:683:ARG:HD2	1:A:687:GLU:OE1	2.10	0.51
1:B:154:LEU:HD13	1:B:177:LEU:HD22	1.91	0.51
1:A:616:HIS:CG	1:A:638:VAL:HG22	2.46	0.51
1:B:415:THR:HB	1:B:416:PRO:HD2	1.91	0.51
1:A:138:SER:C	1:A:140:GLY:H	2.18	0.51
1:A:251:HIS:CA	1:A:276:VAL:HG13	2.36	0.51
1:B:314:LEU:HA	1:B:541:TYR:OH	2.11	0.51
1:A:232:TYR:O	1:A:233:TYR:HB2	2.10	0.51
6:A:958:HOH:O	1:B:160:ASN:HB3	2.09	0.51
1:A:368:LEU:O	1:A:372:GLU:HB2	2.11	0.51
1:B:128:LYS:N	1:B:129:PRO:CD	2.74	0.51
1:A:5:ALA:HB2	1:A:53:VAL:HG11	1.93	0.51
1:B:37:GLU:O	1:B:41:LEU:HD22	2.11	0.50
1:B:381:SER:C	1:B:608:ASN:HD22	2.19	0.50
1:B:403:LEU:HD22	1:B:404:PHE:CE2	2.46	0.50
1:A:399:LEU:HB3	1:A:401:LEU:HG	1.94	0.50
1:A:530:LEU:HD13	1:A:559:ILE:HG23	1.93	0.50
1:A:732:PHE:CD1	1:A:732:PHE:N	2.80	0.50
1:A:198:VAL:HG21	1:A:227:LEU:HG	1.94	0.50
1:A:584:LYS:HB3	1:A:595:TYR:OH	2.12	0.50
1:B:384:ARG:CZ	1:B:615:ILE:HD12	2.42	0.50
1:B:235:SER:CB	1:B:259:GLU:HB2	2.42	0.50
1:A:21:LEU:HG	1:A:25:TRP:CZ2	2.47	0.50
1:A:328:LEU:CD1	1:A:332:LEU:HD21	2.42	0.50
1:A:85:ARG:H	1:A:89:THR:CG2	2.24	0.49
1:B:403:LEU:HD23	1:B:404:PHE:H	1.77	0.49
1:B:106:PHE:CD2	5:B:1054:MRY:HAG2	2.48	0.49
1:B:2:LYS:HE2	1:B:55:VAL:CG2	2.42	0.49
1:A:137:LYS:HD3	1:A:142:GLU:OE2	2.13	0.49
1:A:208:GLU:HA	1:B:208:GLU:HA	1.95	0.49
1:A:453:ILE:HD13	1:A:719:ARG:NH1	2.28	0.48
1:A:586:PRO:HD3	1:A:595:TYR:CG	2.48	0.48
1:B:725:ALA:O	1:B:729:ARG:HG2	2.14	0.48
1:A:447:ILE:O	1:A:451:GLU:HG3	2.12	0.48
1:A:506:ARG:NH2	1:A:509:PRO:HA	2.28	0.48
1:A:9:PRO:HB2	1:A:316:VAL:HG22	1.96	0.48
1:B:403:LEU:HD23	1:B:404:PHE:N	2.29	0.48
1:B:537:MSE:HG3	1:B:555:ILE:HD12	1.96	0.48
1:B:201:PRO:O	1:B:204:VAL:HG22	2.14	0.48
1:B:284:GLN:HG3	1:B:286:TRP:CD1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:373:ARG:HG3	1:A:373:ARG:NH1	2.29	0.47
1:B:8:PHE:CD2	1:B:9:PRO:HD2	2.49	0.47
1:B:56:ILE:HG22	1:B:143:THR:HG22	1.94	0.47
1:A:724:LEU:C	1:A:724:LEU:HD13	2.40	0.47
1:B:235:SER:HB3	1:B:259:GLU:HB2	1.96	0.47
1:B:582:ARG:N	1:B:621:TYR:OH	2.29	0.47
1:A:379:GLU:C	1:A:381:SER:H	2.22	0.47
1:A:700:ILE:HD13	1:A:724:LEU:HD12	1.97	0.47
1:B:537:MSE:HE1	1:B:548:GLU:CG	2.44	0.47
1:B:232:TYR:O	1:B:233:TYR:CB	2.62	0.47
1:B:535:THR:HA	1:B:580:ALA:HB3	1.97	0.47
1:A:235:SER:CB	1:A:259:GLU:HB2	2.45	0.47
1:B:436:GLU:O	1:B:440:LYS:HG3	2.14	0.47
1:A:324:LEU:HB2	6:A:828:HOH:O	2.14	0.46
1:B:137:LYS:C	1:B:139:LYS:H	2.23	0.46
1:B:467:ASP:OD1	1:B:467:ASP:C	2.58	0.46
1:A:293:VAL:HG11	1:A:345:LEU:HD11	1.95	0.46
1:A:364:PHE:O	1:A:365:ALA:C	2.57	0.46
1:A:582:ARG:HD3	1:A:627:ILE:HG21	1.97	0.46
1:A:627:ILE:HD12	1:A:631:ILE:HD11	1.96	0.46
1:B:438:PHE:O	1:B:442:GLN:HG2	2.15	0.46
1:B:714:VAL:O	1:B:718:LEU:HB2	2.15	0.46
1:A:506:ARG:NH2	1:A:565:ASP:OD2	2.46	0.46
1:A:264:LEU:HD23	1:A:264:LEU:O	2.16	0.46
1:B:398:ARG:NH1	1:B:572:LYS:NZ	2.64	0.46
1:A:242:CYS:HA	1:A:245:LEU:HD13	1.97	0.46
1:B:104:LYS:HG2	1:B:108:THR:O	2.15	0.46
1:A:669:TRP:HZ3	1:A:718:LEU:HD13	1.80	0.46
1:B:285:PRO:HD3	1:B:541:TYR:CD1	2.49	0.46
1:B:547:PRO:HG2	1:B:550:GLU:CB	2.45	0.46
1:A:289:ASP:C	1:A:291:ARG:H	2.23	0.46
1:A:367:ASP:O	1:A:370:ALA:N	2.48	0.46
1:A:644:SER:HA	1:A:685:ILE:HD13	1.98	0.46
1:A:328:LEU:HG	1:A:332:LEU:HD21	1.98	0.45
1:A:399:LEU:HD21	6:A:831:HOH:O	2.15	0.45
1:A:472:PHE:O	1:A:476:LEU:HD23	2.16	0.45
1:A:642:GLU:HG2	1:A:703:ASP:O	2.15	0.45
1:A:732:PHE:H	1:A:732:PHE:HD1	1.65	0.45
1:B:381:SER:O	1:B:608:ASN:ND2	2.44	0.45
1:A:364:PHE:O	1:A:366:VAL:HG13	2.17	0.45
1:A:683:ARG:HD2	1:A:687:GLU:CD	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:514:GLU:HG3	6:A:842:HOH:O	2.16	0.45
1:A:600:ILE:HD13	1:A:633:GLN:HB2	1.99	0.45
1:B:235:SER:HA	1:B:260:ASN:OD1	2.17	0.45
1:B:146:TRP:CD1	1:B:199:ASN:HD21	2.35	0.45
1:B:315:PRO:HB2	1:B:334:PHE:CD2	2.52	0.45
1:A:514:GLU:CD	1:A:514:GLU:H	2.24	0.45
1:B:467:ASP:OD1	1:B:468:MSE:N	2.50	0.45
1:A:571:ILE:HD12	1:A:571:ILE:N	2.32	0.44
1:A:578:GLU:OE2	1:A:578:GLU:HA	2.17	0.44
1:A:282:GLY:O	1:A:313:HIS:HB2	2.17	0.44
1:A:409:ILE:HG13	1:A:704:CYS:CA	2.47	0.44
1:A:453:ILE:HD11	1:A:715:ILE:HG23	1.99	0.44
1:A:623:ASP:OD1	1:A:625:ASN:ND2	2.51	0.44
1:B:535:THR:HA	1:B:580:ALA:CB	2.47	0.44
1:A:124:LEU:HD22	1:A:183:GLU:HG2	1.99	0.44
1:B:61:LEU:O	1:B:61:LEU:HG	2.17	0.44
1:B:557:LEU:CD2	1:B:605:LEU:HD11	2.47	0.44
1:B:14:LYS:HE3	6:B:898:HOH:O	2.17	0.44
1:B:582:ARG:HG2	1:B:582:ARG:HH21	1.82	0.44
1:A:647:LYS:HA	1:A:685:ILE:HG12	2.00	0.44
1:A:665:GLY:HA3	1:A:701:ASN:HD21	1.82	0.44
1:B:65:ASP:OD1	1:B:67:VAL:HB	2.18	0.44
1:A:288:VAL:CG2	1:A:289:ASP:N	2.81	0.44
1:B:253:ASP:OD1	1:B:253:ASP:C	2.59	0.44
1:A:484:ASN:HB3	6:A:792:HOH:O	2.18	0.44
1:B:239:TYR:HB2	1:B:263:ASN:ND2	2.32	0.44
1:B:281:ASN:HA	1:B:308:SER:OG	2.17	0.44
1:B:557:LEU:HD23	1:B:605:LEU:HD11	1.98	0.44
1:A:337:GLU:O	1:A:341:GLU:HG3	2.17	0.43
1:B:185:PHE:CD1	1:B:227:LEU:HD21	2.53	0.43
1:A:43:MSE:HA	1:A:46:VAL:CG2	2.48	0.43
1:A:251:HIS:HA	1:A:276:VAL:CG1	2.41	0.43
1:B:139:LYS:HB2	1:B:139:LYS:HZ3	1.82	0.43
1:B:390:GLU:O	1:B:394:ILE:HG12	2.19	0.43
1:A:63:TYR:CD2	1:A:87:LEU:HD21	2.51	0.43
1:A:111:HIS:HB2	1:A:490:TYR:HB3	2.00	0.43
1:B:230:PHE:HA	1:B:251:HIS:HB3	2.00	0.43
1:B:417:GLU:OE2	1:B:438:PHE:CE1	2.71	0.43
1:B:443:ILE:O	1:B:447:ILE:HG12	2.18	0.43
1:B:646:SER:O	1:B:647:LYS:HB2	2.19	0.43
1:A:483:GLN:O	1:A:496:ARG:NH1	2.46	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:PHE:O	1:B:95:ARG:HG2	2.19	0.43
1:B:296:LEU:HD12	1:B:299:LYS:HE2	2.01	0.43
1:A:668:VAL:CG2	1:A:682:MSE:HE2	2.49	0.43
1:B:259:GLU:CD	1:B:259:GLU:H	2.26	0.43
1:B:537:MSE:HG3	1:B:555:ILE:CD1	2.49	0.43
1:A:11:ILE:HA	1:A:17:PHE:HB3	2.00	0.42
1:B:315:PRO:HD3	1:B:541:TYR:OH	2.19	0.42
1:A:56:ILE:HG22	1:A:143:THR:HG22	2.00	0.42
1:A:338:LYS:HD3	1:A:338:LYS:HA	1.83	0.42
1:A:560:ASN:HD21	1:A:564:LYS:HE2	1.83	0.42
1:B:205:CYS:O	1:B:237:SER:HB3	2.19	0.42
1:B:370:ALA:HB3	1:B:503:THR:HG21	2.01	0.42
1:B:543:ARG:HE	1:B:545:ASP:CG	2.28	0.42
1:A:368:LEU:HD12	1:A:368:LEU:N	2.35	0.42
1:A:497:PRO:HG2	1:A:541:TYR:CE2	2.54	0.42
1:A:719:ARG:HG2	1:A:719:ARG:HH11	1.84	0.42
1:B:11:ILE:HA	1:B:17:PHE:CB	2.50	0.42
1:B:68:LEU:HG	1:B:90:TYR:CE2	2.55	0.42
1:B:334:PHE:O	1:B:338:LYS:HG2	2.19	0.42
1:B:669:TRP:HB2	1:B:682:MSE:HE2	2.02	0.42
1:A:317:THR:CG2	1:A:336:LYS:HE2	2.48	0.42
1:B:78:PRO:HB2	1:B:81:PHE:CE1	2.54	0.42
1:B:185:PHE:CE1	1:B:227:LEU:HD21	2.53	0.42
1:B:315:PRO:HB3	1:B:485:GLY:HA2	2.01	0.42
1:A:486:TRP:CD2	1:A:496:ARG:HG3	2.54	0.42
1:A:317:THR:HG22	1:A:318:LEU:N	2.34	0.42
1:A:341:GLU:HA	1:A:344:MSE:HE3	2.02	0.42
1:A:665:GLY:HA2	1:A:699:TRP:HB2	2.01	0.42
1:B:685:ILE:O	1:B:689:VAL:HG13	2.20	0.42
1:A:310:PRO:HB3	1:A:312:PHE:CE1	2.55	0.42
1:B:581:PHE:HB3	1:B:621:TYR:OH	2.19	0.42
1:A:31:GLU:O	1:A:35:GLU:HG2	2.20	0.41
1:A:247:VAL:HA	6:A:943:HOH:O	2.19	0.41
1:A:128:LYS:N	1:A:129:PRO:CD	2.83	0.41
1:A:283:ARG:HG2	1:A:309:CYS:SG	2.60	0.41
1:B:475:LYS:HD2	1:B:510:MSE:HB3	2.01	0.41
1:B:515:ILE:CG2	1:B:516:THR:N	2.83	0.41
1:B:286:TRP:HD1	1:B:544:GLU:OE2	2.04	0.41
1:A:225:PHE:HB2	1:A:227:LEU:HD13	2.03	0.41
1:A:233:TYR:O	1:A:234:ASP:HB2	2.20	0.41
1:A:289:ASP:C	1:A:291:ARG:N	2.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:406:THR:HA	1:B:700:ILE:O	2.20	0.41
1:B:581:PHE:CE2	1:B:631:ILE:HD13	2.55	0.41
1:B:582:ARG:HG2	1:B:582:ARG:NH2	2.36	0.41
1:A:293:VAL:CG1	1:A:345:LEU:HD11	2.51	0.41
1:A:431:SER:OG	1:A:434:GLU:HG3	2.21	0.41
1:A:453:ILE:HG22	1:A:722:VAL:HG21	2.02	0.41
1:B:69:ASP:HB3	1:B:128:LYS:NZ	2.36	0.41
1:A:217:ASN:O	1:A:221:GLU:HG3	2.21	0.41
1:A:399:LEU:O	1:A:400:ASN:HB2	2.20	0.41
1:A:450:GLN:CB	1:A:458:LEU:HD22	2.49	0.41
1:B:537:MSE:HE1	1:B:548:GLU:HG2	2.03	0.41
1:B:381:SER:C	1:B:608:ASN:ND2	2.78	0.41
1:B:701:ASN:HB2	1:B:702:PRO:CD	2.51	0.41
1:A:56:ILE:HD12	1:A:56:ILE:N	2.35	0.41
1:A:345:LEU:O	1:A:349:LEU:HB2	2.21	0.41
1:B:245:LEU:HD12	1:B:245:LEU:HA	1.94	0.41
1:B:515:ILE:HD12	1:B:571:ILE:CD1	2.51	0.41
1:A:553:TYR:O	1:A:557:LEU:HD23	2.20	0.41
1:B:55:VAL:C	1:B:56:ILE:HD12	2.46	0.41
1:B:665:GLY:HA2	1:B:699:TRP:HB2	2.03	0.41
1:A:489:SER:O	1:A:490:TYR:HB3	2.20	0.40
1:B:278:GLY:HA2	1:B:306:SER:O	2.21	0.40
1:B:284:GLN:C	1:B:314:LEU:HD21	2.46	0.40
1:B:339:LEU:HD12	1:B:339:LEU:HA	1.93	0.40
1:B:497:PRO:HA	1:B:498:PRO:HD3	1.94	0.40
1:B:667:GLY:HA2	1:B:701:ASN:O	2.21	0.40
1:A:440:LYS:HG2	1:A:517:TYR:CD2	2.57	0.40
1:A:314:LEU:HA	1:A:541:TYR:OH	2.20	0.40
1:A:589:LYS:O	1:B:168:ASN:HB3	2.21	0.40
1:B:86:GLY:O	1:B:87:LEU:C	2.65	0.40
1:B:534:VAL:CG1	1:B:584:LYS:HD3	2.49	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 X-ray entries followed by that with respect to entries of similar resolution.

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	716/766 (94%)	673 (94%)	39 (5%)	4 (1%)	21	23
1	B	704/766 (92%)	667 (95%)	34 (5%)	3 (0%)	30	34
All	All	1420/1532 (93%)	1340 (94%)	73 (5%)	7 (0%)	24	27

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	365	ALA
1	B	289	ASP
1	B	420	LYS
1	A	64	TYR
1	A	380	ASP
1	B	64	TYR
1	A	385	GLU

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 X-ray entries followed by that with respect to entries of similar resolution.

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	576/670 (86%)	552 (96%)	24 (4%)	26	36
1	B	565/670 (84%)	538 (95%)	27 (5%)	23	30
All	All	1141/1340 (85%)	1090 (96%)	51 (4%)	24	33

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LEU
1	A	154	LEU
1	A	263	ASN
1	A	273	LYS
1	A	276	VAL
1	A	285	PRO

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Mol	Chain	Res	Type
1	A	324	LEU
1	A	339	LEU
1	A	349	LEU
1	A	372	GLU
1	A	399	LEU
1	A	403	LEU
1	A	476	LEU
1	A	488	LEU
1	A	514	GLU
1	A	557	LEU
1	A	612	GLU
1	A	628	ILE
1	A	638	VAL
1	A	683	ARG
1	A	696	GLU
1	A	706	LEU
1	A	718	LEU
1	A	728	MSE
1	B	46	VAL
1	B	128	LYS
1	B	132	ASP
1	B	137	LYS
1	B	139	LYS
1	B	154	LEU
1	B	245	LEU
1	B	250	LEU
1	B	285	PRO
1	B	312	PHE
1	B	339	LEU
1	B	387	GLU
1	B	395	GLN
1	B	398	ARG
1	B	417	GLU
1	B	476	LEU
1	B	488	LEU
1	B	515	ILE
1	B	516	THR
1	B	545	ASP
1	B	557	LEU
1	B	638	VAL
1	B	687	GLU
1	B	696	GLU

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Mol	Chain	Res	Type
1	B	709	ARG
1	B	718	LEU
1	B	724	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	168	ASN
1	A	191	ASN
1	A	263	ASN
1	A	560	ASN
1	A	601	ASN
1	A	608	ASN
1	A	701	ASN
1	B	39	ASN
1	B	111	HIS
1	B	199	ASN
1	B	263	ASN
1	B	414	GLN
1	B	477	ASN
1	B	625	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 ⓘ

Of 8 ligands modelled in this entry, 2 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	SO4	A	1048	-	4,4,4	0.39	0	6,6,6	0.08	0
4	HCS	A	1050	-	6,7,7	0.94	0	5,8,8	0.74	0
4	HCS	A	1049	-	6,7,7	0.95	0	5,8,8	0.57	0
2	SO4	B	1047	-	4,4,4	0.29	0	6,6,6	0.19	0
5	MRY	A	1053	-	7,7,7	0.31	0	8,8,8	0.42	0
5	MRY	B	1054	-	7,7,7	0.33	0	8,8,8	0.47	0

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	HCS	A	1050	-	-	1/7/7/7	-
5	MRY	B	1054	-	-	1/8/8/8	-
4	HCS	A	1049	-	-	1/7/7/7	-
5	MRY	A	1053	-	-	1/8/8/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1053	MRY	OAD-CAC-CAE-CAG
5	B	1054	MRY	OAD-CAC-CAE-CAG
4	A	1049	HCS	O-C-CA-N
4	A	1050	HCS	OXT-C-CA-N

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1054	MRY	1	0

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 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	705/766 (92%)	-1.62	0 100 100	9, 25, 42, 58	0
1	B	695/766 (90%)	-1.63	0 100 100	7, 25, 39, 56	0
All	All	1400/1532 (91%)	-1.62	0 100 100	7, 25, 41, 58	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MRY	A	1053	8/8	0.99	0.04	25,26,27,28	0
5	MRY	B	1054	8/8	0.99	0.07	27,33,36,39	0
3	ZN	A	1051	1/1	1.00	0.01	27,27,27,27	0
3	ZN	B	1052	1/1	1.00	0.02	30,30,30,30	0
4	HCS	A	1049	8/8	1.00	0.02	11,15,21,25	0
4	HCS	A	1050	8/8	1.00	0.02	15,21,25,28	0
2	SO4	A	1048	5/5	1.00	0.01	24,26,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	B	1047	5/5	1.00	0.02	25,26,27,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.