



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 09:14 pm BST

PDB ID : 6C43
Title : 2.9 Angstrom Resolution Crystal Structure of Gamma-Aminobutyraldehyde Dehydrogenase from Salmonella typhimurium.
Authors : Minasov, G.; Shuvalova, L.; Dubrovskaya, I.; Winsor, J.; Tekleab, H.; Kwon, K.; Anderson, W.F.; Satchell, K.J.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2018-01-11
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

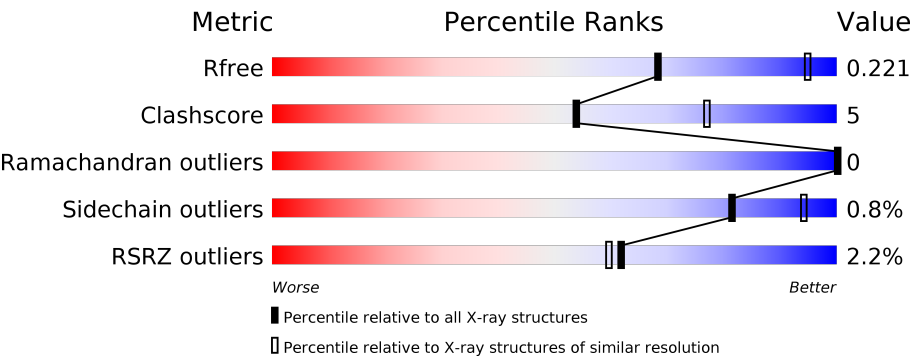
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 on 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	477	<div><div>2%</div><div>87%</div><div>12%</div><div>.</div></div>
1	B	477	<div><div>3%</div><div>86%</div><div>13%</div><div>..</div></div>
1	C	477	<div><div>%</div><div>85%</div><div>13%</div><div>..</div></div>
1	D	477	<div><div>2%</div><div>85%</div><div>14%</div><div>..</div></div>
1	E	477	<div><div>3%</div><div>86%</div><div>13%</div><div>..</div></div>
1	F	477	<div><div>2%</div><div>86%</div><div>13%</div><div>.</div></div>

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Mol	Chain	Length	Quality of chain
1	G	477	<div><div></div><div>4%</div><div>86%</div><div>13%</div><div></div></div>
1	H	477	<div><div></div><div>2%</div><div>88%</div><div>11%</div><div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 29036 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 Gamma-aminobutyraldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	B	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	C	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	D	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	E	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	F	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	G	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			
1	H	473	Total	C	N	O	S	0	0	0
			3591	2269	626	678	18			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP A0A0W3SL32
A	-1	ASN	-	expression tag	UNP A0A0W3SL32
A	0	ALA	-	expression tag	UNP A0A0W3SL32
B	-2	SER	-	expression tag	UNP A0A0W3SL32
B	-1	ASN	-	expression tag	UNP A0A0W3SL32
B	0	ALA	-	expression tag	UNP A0A0W3SL32
C	-2	SER	-	expression tag	UNP A0A0W3SL32
C	-1	ASN	-	expression tag	UNP A0A0W3SL32
C	0	ALA	-	expression tag	UNP A0A0W3SL32
D	-2	SER	-	expression tag	UNP A0A0W3SL32
D	-1	ASN	-	expression tag	UNP A0A0W3SL32
D	0	ALA	-	expression tag	UNP A0A0W3SL32
E	-2	SER	-	expression tag	UNP A0A0W3SL32

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-1	ASN	-	expression tag	UNP A0A0W3SL32
E	0	ALA	-	expression tag	UNP A0A0W3SL32
F	-2	SER	-	expression tag	UNP A0A0W3SL32
F	-1	ASN	-	expression tag	UNP A0A0W3SL32
F	0	ALA	-	expression tag	UNP A0A0W3SL32
G	-2	SER	-	expression tag	UNP A0A0W3SL32
G	-1	ASN	-	expression tag	UNP A0A0W3SL32
G	0	ALA	-	expression tag	UNP A0A0W3SL32
H	-2	SER	-	expression tag	UNP A0A0W3SL32
H	-1	ASN	-	expression tag	UNP A0A0W3SL32
H	0	ALA	-	expression tag	UNP A0A0W3SL32

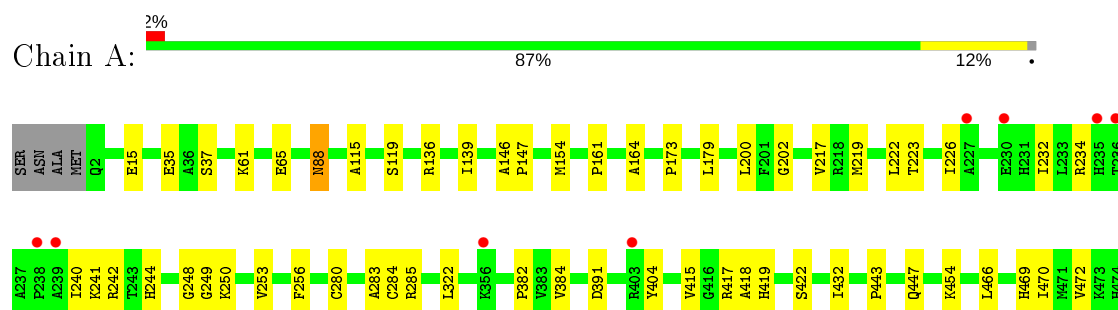
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	58	Total O 59 59	0	1
2	B	26	Total O 26 26	0	0
2	C	64	Total O 64 64	0	0
2	D	25	Total O 25 25	0	0
2	E	34	Total O 34 34	0	0
2	F	35	Total O 35 35	0	0
2	G	23	Total O 23 23	0	0
2	H	42	Total O 42 42	0	0

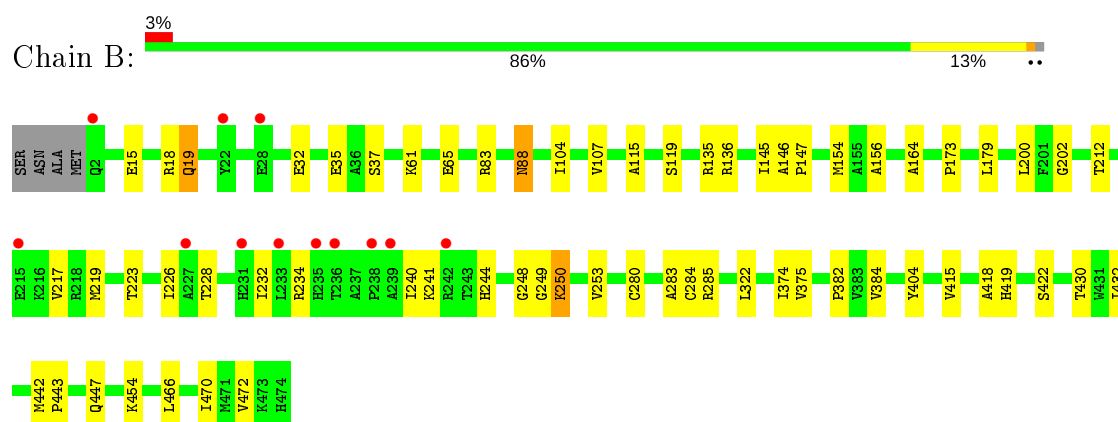
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

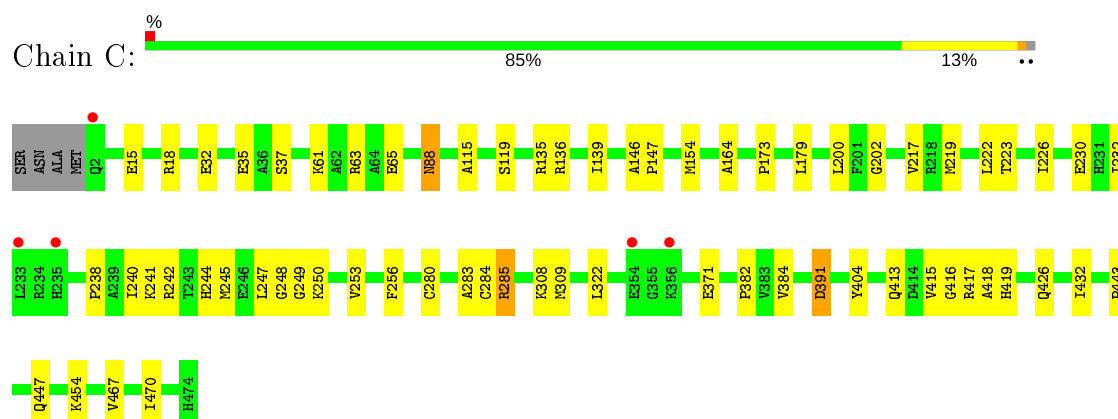
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase



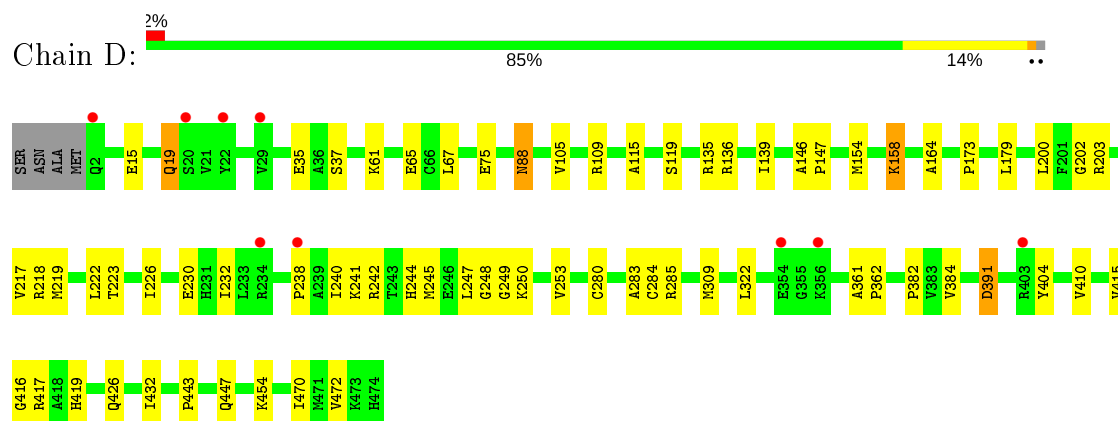
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase



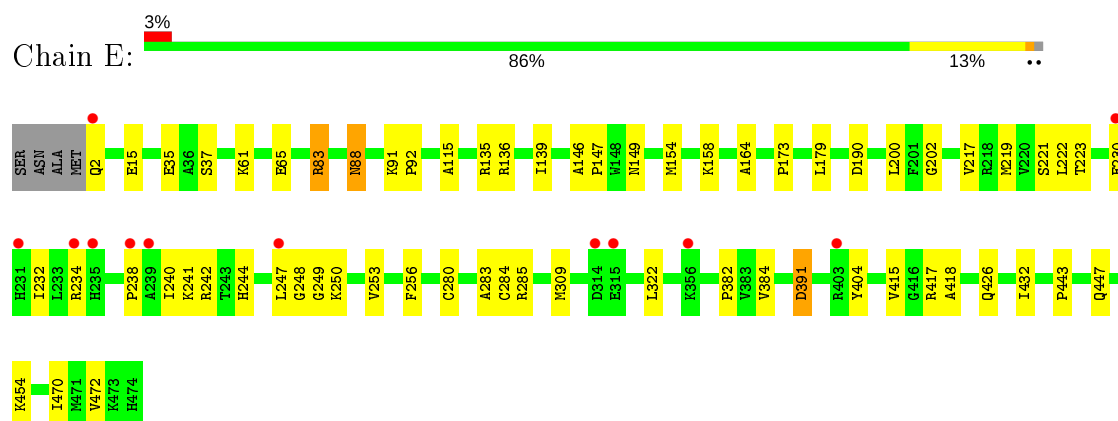
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase



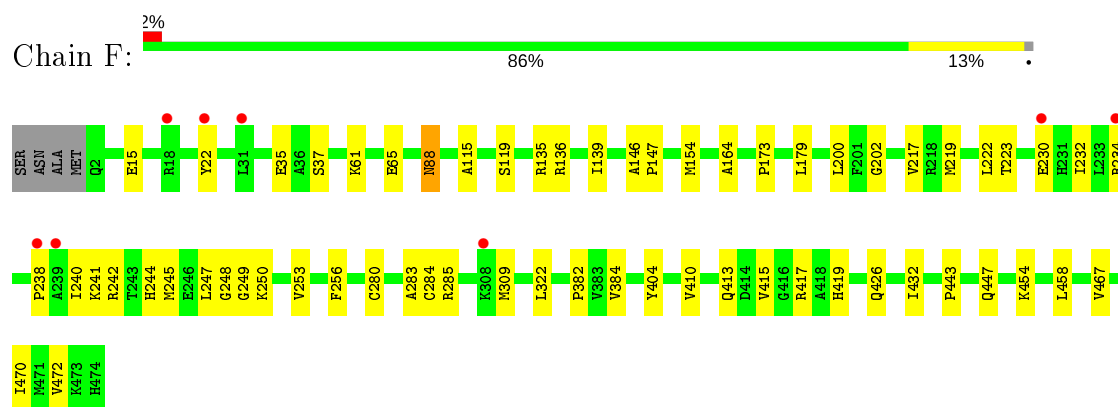
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase



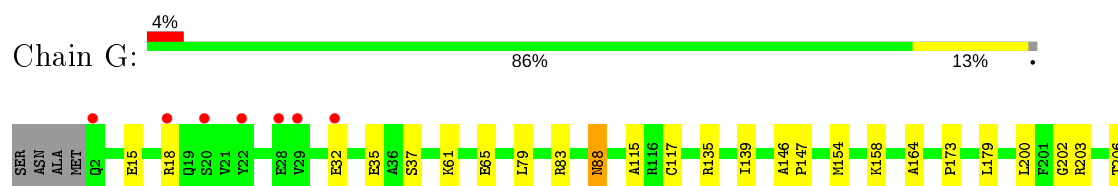
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

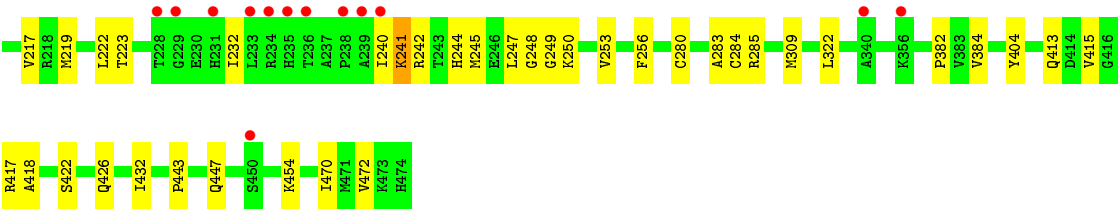


- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

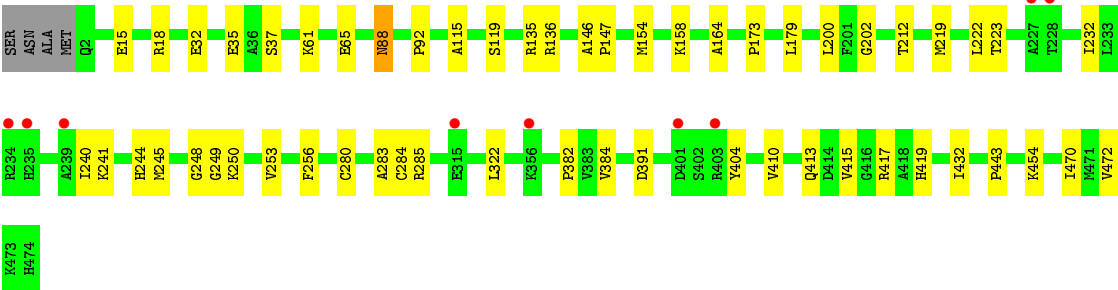
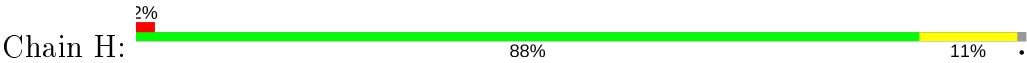


- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase





● Molecule 1: Gamma-aminobutyraldehyde dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	81.89 Å 87.80 Å 145.07 Å 82.76° 77.91° 72.86°	Depositor
Resolution (Å)	29.73 – 2.90 29.72 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.4 (29.73-2.90) 98.4 (29.72-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.90 Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.189 , 0.219 0.190 , 0.221	Depositor DCC
R_{free} test set	4155 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	56.7	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	29036	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% 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

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	0.40	0/3663	0.71	1/4975 (0.0%)
1	B	0.40	0/3663	0.71	3/4975 (0.1%)
1	C	0.43	2/3663 (0.1%)	0.83	4/4975 (0.1%)
1	D	0.40	0/3663	0.72	4/4975 (0.1%)
1	E	0.41	0/3663	0.71	2/4975 (0.0%)
1	F	0.40	0/3663	0.70	0/4975
1	G	0.41	0/3663	0.74	4/4975 (0.1%)
1	H	0.39	0/3663	0.69	2/4975 (0.0%)
All	All	0.41	2/29304 (0.0%)	0.73	20/39800 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	285	ARG	CZ-NH1	-6.90	1.24	1.33
1	C	371	GLU	CD-OE1	-5.86	1.19	1.25

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	285	ARG	NE-CZ-NH1	-20.12	110.24	120.30
1	C	285	ARG	NE-CZ-NH2	16.13	128.36	120.30
1	C	63	ARG	NE-CZ-NH2	11.71	126.16	120.30
1	C	63	ARG	NE-CZ-NH1	-10.75	114.92	120.30
1	G	203	ARG	NE-CZ-NH1	-10.50	115.05	120.30
1	G	241	LYS	CA-CB-CG	9.17	133.57	113.40
1	D	158	LYS	CD-CE-NZ	-8.02	93.26	111.70
1	B	250	LYS	CB-CG-CD	7.44	130.95	111.60
1	D	158	LYS	CB-CG-CD	6.41	128.27	111.60
1	G	203	ARG	CG-CD-NE	-6.33	98.50	111.80
1	E	83	ARG	CB-CG-CD	6.15	127.58	111.60
1	D	158	LYS	CA-CB-CG	6.05	126.72	113.40
1	G	158	LYS	CD-CE-NZ	5.98	125.45	111.70
1	A	391	ASP	CB-CG-OD1	5.52	123.27	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	158	LYS	CD-CE-NZ	5.45	124.23	111.70
1	D	218	ARG	CG-CD-NE	5.42	123.17	111.80
1	E	136	ARG	NE-CZ-NH2	5.19	122.89	120.30
1	H	391	ASP	CB-CG-OD1	5.11	122.90	118.30
1	B	83	ARG	CG-CD-NE	5.08	122.47	111.80
1	B	250	LYS	CD-CE-NZ	5.01	123.22	111.70

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	3591	0	3571	40	0
1	B	3591	0	3571	51	0
1	C	3591	0	3571	51	1
1	D	3591	0	3571	59	0
1	E	3591	0	3571	44	1
1	F	3591	0	3571	45	1
1	G	3591	0	3571	39	1
1	H	3591	0	3571	37	0
2	A	59	0	0	1	0
2	B	26	0	0	0	0
2	C	64	0	0	0	0
2	D	25	0	0	0	0
2	E	34	0	0	0	0
2	F	35	0	0	0	0
2	G	23	0	0	0	0
2	H	42	0	0	1	0
All	All	29036	0	28568	306	2

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 5.

All (306) 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:D:19:GLN:HE21	1:D:19:GLN:HA	1.28	0.96
1:B:250:LYS:NZ	1:B:374:ILE:O	2.07	0.86
1:B:104:ILE:O	1:B:107:VAL:HG12	1.77	0.84
1:C:217:VAL:O	1:C:241:LYS:NZ	2.12	0.82
1:B:250:LYS:HG2	1:B:285:ARG:HD2	1.61	0.81
1:B:19:GLN:HE21	1:B:19:GLN:HA	1.45	0.81
1:E:91:LYS:NZ	1:E:149:ASN:O	2.18	0.77
1:C:240:ILE:HD13	1:D:247:LEU:CD1	2.14	0.76
1:C:238:PRO:HA	1:D:230:GLU:HG3	1.68	0.75
1:C:247:LEU:CD1	1:D:240:ILE:HD13	2.17	0.75
1:E:234:ARG:HG3	1:F:234:ARG:HG3	1.69	0.74
1:B:250:LYS:HD3	1:B:375:VAL:O	1.87	0.73
1:B:415:VAL:HG22	1:C:415:VAL:HA	1.70	0.73
1:E:432:ILE:HD12	1:F:472:VAL:HG22	1.71	0.72
1:E:247:LEU:CD1	1:F:240:ILE:HD13	2.21	0.70
1:C:432:ILE:HD12	1:D:472:VAL:HG22	1.72	0.70
1:D:309:MET:HE1	1:D:361:ALA:HA	1.76	0.68
1:C:240:ILE:HD13	1:D:247:LEU:HD13	1.76	0.68
1:C:230:GLU:HG3	1:D:238:PRO:HA	1.75	0.67
1:E:418:ALA:HB1	1:E:432:ILE:HD13	1.78	0.66
1:C:418:ALA:HB1	1:C:432:ILE:HD13	1.77	0.66
1:B:418:ALA:HB1	1:B:432:ILE:HD13	1.78	0.65
1:D:309:MET:CE	1:D:361:ALA:HA	2.25	0.65
1:G:418:ALA:HB1	1:G:432:ILE:HD13	1.78	0.65
1:C:247:LEU:HD13	1:D:240:ILE:HD13	1.77	0.65
1:A:418:ALA:HB1	1:A:432:ILE:HD13	1.78	0.65
1:A:472:VAL:HG22	1:B:432:ILE:HD12	1.79	0.64
1:A:470:ILE:HD12	1:B:422:SER:HB2	1.79	0.62
1:H:222:LEU:HB3	1:H:245:MET:HE2	1.82	0.61
1:C:240:ILE:HG21	1:D:247:LEU:HD12	1.82	0.60
1:A:432:ILE:HD12	1:B:472:VAL:HG22	1.83	0.60
1:E:247:LEU:HD13	1:F:240:ILE:HD13	1.82	0.60
1:D:19:GLN:HA	1:D:19:GLN:NE2	2.09	0.59
1:F:415:VAL:HG23	1:H:413:GLN:O	2.03	0.59
1:D:309:MET:HE3	1:D:362:PRO:HD3	1.86	0.58
1:F:322:LEU:HD21	1:F:382:PRO:HD3	1.85	0.58
1:G:245:MET:HE2	1:G:247:LEU:HD11	1.86	0.57
1:F:222:LEU:HB3	1:F:245:MET:HE2	1.86	0.57
1:E:322:LEU:HD21	1:E:382:PRO:HD3	1.87	0.57
1:A:322:LEU:HD21	1:A:382:PRO:HD3	1.87	0.56
1:D:322:LEU:HD21	1:D:382:PRO:HD3	1.87	0.56
1:A:249:GLY:HA2	1:A:404:TYR:CD1	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:249:GLY:HA2	1:D:404:TYR:CD1	2.40	0.56
1:G:322:LEU:HD21	1:G:382:PRO:HD3	1.86	0.56
1:B:249:GLY:HA2	1:B:404:TYR:CD1	2.40	0.56
1:H:249:GLY:HA2	1:H:404:TYR:CD1	2.41	0.56
1:H:322:LEU:HD21	1:H:382:PRO:HD3	1.87	0.56
1:E:249:GLY:HA2	1:E:404:TYR:CD1	2.40	0.56
1:B:322:LEU:HD21	1:B:382:PRO:HD3	1.86	0.55
1:C:322:LEU:HD21	1:C:382:PRO:HD3	1.86	0.55
1:G:222:LEU:HD22	1:G:232:ILE:HD12	1.88	0.55
1:F:249:GLY:HA2	1:F:404:TYR:CD1	2.41	0.55
1:F:415:VAL:HG22	1:H:415:VAL:HA	1.88	0.55
1:B:19:GLN:HE21	1:B:19:GLN:CA	2.13	0.55
1:C:222:LEU:HD22	1:C:232:ILE:HD12	1.88	0.55
1:C:222:LEU:HB3	1:C:245:MET:HE2	1.89	0.55
1:G:249:GLY:HA2	1:G:404:TYR:CD1	2.41	0.55
1:F:415:VAL:HG23	1:H:413:GLN:C	2.28	0.55
1:F:222:LEU:HD22	1:F:232:ILE:HD12	1.88	0.54
1:F:415:VAL:HA	1:H:415:VAL:HG22	1.88	0.54
1:A:222:LEU:HD22	1:A:232:ILE:HD12	1.89	0.54
1:C:249:GLY:HA2	1:C:404:TYR:CD1	2.41	0.54
1:D:222:LEU:HD22	1:D:232:ILE:HD12	1.88	0.54
1:G:248:GLY:HA2	1:G:280:CYS:SG	2.48	0.54
1:F:248:GLY:HA2	1:F:280:CYS:SG	2.48	0.54
1:E:222:LEU:HD22	1:E:232:ILE:HD12	1.89	0.53
1:G:432:ILE:HD12	1:H:472:VAL:HG22	1.88	0.53
1:C:247:LEU:HD12	1:D:240:ILE:HD13	1.91	0.53
1:B:19:GLN:NE2	1:B:19:GLN:HA	2.21	0.53
1:B:145:ILE:HD11	1:B:212:THR:CG2	2.38	0.53
1:A:248:GLY:HA2	1:A:280:CYS:SG	2.49	0.53
1:E:415:VAL:HG23	1:G:413:GLN:O	2.09	0.53
1:A:415:VAL:HG22	1:D:415:VAL:HA	1.90	0.53
1:B:107:VAL:HG11	1:B:156:ALA:HB1	1.90	0.53
1:B:248:GLY:HA2	1:B:280:CYS:SG	2.49	0.52
1:C:247:LEU:HD12	1:D:240:ILE:HG21	1.90	0.52
1:H:222:LEU:HD22	1:H:232:ILE:HD12	1.92	0.52
1:C:61:LYS:HE2	1:C:65:GLU:OE2	2.09	0.52
1:H:248:GLY:HA2	1:H:280:CYS:SG	2.50	0.52
1:D:248:GLY:HA2	1:D:280:CYS:SG	2.50	0.52
1:D:61:LYS:HE2	1:D:65:GLU:OE2	2.10	0.52
1:E:230:GLU:HG3	1:F:238:PRO:HA	1.91	0.52
1:C:240:ILE:HD13	1:D:247:LEU:HD12	1.89	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:248:GLY:HA2	1:E:280:CYS:SG	2.50	0.52
1:C:248:GLY:HA2	1:C:280:CYS:SG	2.49	0.52
1:G:61:LYS:HE2	1:G:65:GLU:OE2	2.10	0.51
1:A:61:LYS:HE2	1:A:65:GLU:OE2	2.10	0.51
1:C:240:ILE:CD1	1:D:226:ILE:HG23	2.39	0.51
1:F:61:LYS:HE2	1:F:65:GLU:OE2	2.10	0.51
1:A:470:ILE:CD1	1:B:430:THR:HB	2.41	0.51
1:B:88:ASN:HB2	1:B:179:LEU:HD12	1.93	0.51
1:E:88:ASN:HB2	1:E:179:LEU:HD12	1.93	0.51
1:G:115:ALA:HB2	1:G:164:ALA:HB1	1.93	0.51
1:H:61:LYS:HE2	1:H:65:GLU:OE2	2.11	0.51
1:C:115:ALA:HB2	1:C:164:ALA:HB1	1.93	0.50
1:E:61:LYS:HE2	1:E:65:GLU:OE2	2.11	0.50
1:C:88:ASN:HB2	1:C:179:LEU:HD12	1.93	0.50
1:H:88:ASN:HB2	1:H:179:LEU:HD12	1.93	0.50
1:F:88:ASN:HB2	1:F:179:LEU:HD12	1.93	0.50
1:G:88:ASN:HB2	1:G:179:LEU:HD12	1.93	0.50
1:E:391:ASP:OD1	1:E:391:ASP:N	2.45	0.50
1:H:115:ALA:HB2	1:H:164:ALA:HB1	1.93	0.50
1:D:391:ASP:N	1:D:391:ASP:OD1	2.45	0.50
1:E:415:VAL:HA	1:G:415:VAL:HG22	1.94	0.50
1:B:115:ALA:HB2	1:B:164:ALA:HB1	1.94	0.50
1:D:88:ASN:HB2	1:D:179:LEU:HD12	1.93	0.50
1:C:240:ILE:HD11	1:D:226:ILE:HG23	1.93	0.50
1:F:115:ALA:HB2	1:F:164:ALA:HB1	1.93	0.50
1:A:88:ASN:HB2	1:A:179:LEU:HD12	1.93	0.50
1:A:115:ALA:HB2	1:A:164:ALA:HB1	1.93	0.49
1:C:250:LYS:HD2	1:C:285:ARG:HD3	1.94	0.49
1:D:105:VAL:HG12	1:D:109:ARG:HE	1.75	0.49
1:E:238:PRO:HA	1:F:230:GLU:HG3	1.94	0.49
1:D:105:VAL:CG1	1:D:109:ARG:HE	2.25	0.49
1:B:61:LYS:HE2	1:B:65:GLU:OE2	2.12	0.49
1:D:410:VAL:HB	1:D:432:ILE:HD13	1.94	0.49
1:G:245:MET:CE	1:G:247:LEU:HD11	2.42	0.49
1:C:147:PRO:HD3	1:C:223:THR:HB	1.95	0.49
1:H:410:VAL:HB	1:H:432:ILE:HD13	1.95	0.49
1:C:391:ASP:OD1	1:C:391:ASP:N	2.45	0.49
1:D:115:ALA:HB2	1:D:164:ALA:HB1	1.93	0.49
1:C:226:ILE:HG23	1:D:240:ILE:CD1	2.43	0.49
1:F:147:PRO:HD3	1:F:223:THR:HB	1.95	0.49
1:E:240:ILE:HD13	1:F:247:LEU:CD1	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:135:ARG:HD2	1:C:470:ILE:HD11	1.94	0.48
1:D:245:MET:HE2	1:D:247:LEU:HD11	1.95	0.48
1:A:147:PRO:HD3	1:A:223:THR:HB	1.95	0.48
1:A:470:ILE:HD13	1:B:430:THR:HB	1.95	0.48
1:E:115:ALA:HB2	1:E:164:ALA:HB1	1.94	0.48
1:E:147:PRO:HD3	1:E:223:THR:HB	1.95	0.48
1:F:419:HIS:CD2	1:G:472:VAL:HG11	2.49	0.48
1:F:410:VAL:HB	1:F:432:ILE:HD13	1.94	0.48
1:D:245:MET:CE	1:D:247:LEU:HD11	2.44	0.48
1:A:35:GLU:HG2	1:A:202:GLY:HA2	1.96	0.48
1:B:228:THR:O	1:B:232:ILE:HG12	2.12	0.48
1:E:158:LYS:HD2	1:E:221:SER:OG	2.14	0.48
1:B:147:PRO:HD3	1:B:223:THR:HB	1.96	0.48
1:F:413:GLN:O	1:H:415:VAL:HG23	2.13	0.48
1:C:173:PRO:HD2	1:C:200:LEU:O	2.14	0.47
1:D:75:GLU:OE2	1:D:109:ARG:NH1	2.47	0.47
1:H:147:PRO:HD3	1:H:223:THR:HB	1.96	0.47
1:D:147:PRO:HD3	1:D:223:THR:HB	1.95	0.47
1:G:422:SER:HB2	1:H:470:ILE:HD13	1.95	0.47
1:B:145:ILE:HD11	1:B:212:THR:HG21	1.96	0.47
1:F:173:PRO:HD2	1:F:200:LEU:O	2.14	0.47
1:F:250:LYS:HG3	1:F:285:ARG:HD2	1.97	0.47
1:G:173:PRO:HD2	1:G:200:LEU:O	2.14	0.47
1:G:35:GLU:HG2	1:G:202:GLY:HA2	1.97	0.47
1:B:472:VAL:HG11	1:D:419:HIS:CD2	2.50	0.47
1:E:135:ARG:HD2	1:E:470:ILE:HD11	1.97	0.46
1:F:426:GLN:HB3	1:F:447:GLN:OE1	2.16	0.46
1:G:147:PRO:HD3	1:G:223:THR:HB	1.96	0.46
1:H:35:GLU:HG2	1:H:202:GLY:HA2	1.97	0.46
1:A:173:PRO:HD2	1:A:200:LEU:O	2.15	0.46
1:B:173:PRO:HD2	1:B:200:LEU:O	2.15	0.46
1:B:443:PRO:HB2	1:B:454:LYS:HD3	1.98	0.46
1:E:454:LYS:NZ	1:F:467:VAL:O	2.36	0.46
1:H:173:PRO:HD2	1:H:200:LEU:O	2.15	0.46
1:D:173:PRO:HD2	1:D:200:LEU:O	2.15	0.46
1:E:472:VAL:HG11	1:H:419:HIS:CD2	2.50	0.46
1:D:426:GLN:HB3	1:D:447:GLN:OE1	2.16	0.46
1:F:35:GLU:HG2	1:F:202:GLY:HA2	1.98	0.46
1:E:173:PRO:HD2	1:E:200:LEU:O	2.15	0.46
1:H:250:LYS:HG3	1:H:285:ARG:HD2	1.97	0.46
1:G:253:VAL:HG21	1:G:283:ALA:HB1	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:253:VAL:HG21	1:E:283:ALA:HB1	1.98	0.45
1:D:253:VAL:HG21	1:D:283:ALA:HB1	1.98	0.45
1:E:250:LYS:HG3	1:E:285:ARG:HD2	1.96	0.45
1:E:426:GLN:HB3	1:E:447:GLN:OE1	2.16	0.45
1:C:35:GLU:HG2	1:C:202:GLY:HA2	1.98	0.45
1:D:35:GLU:HG2	1:D:202:GLY:HA2	1.97	0.45
1:D:250:LYS:HG3	1:D:285:ARG:HD2	1.98	0.45
1:D:443:PRO:HB2	1:D:454:LYS:HD3	1.99	0.45
1:A:250:LYS:HG3	1:A:285:ARG:HD2	1.98	0.45
1:C:222:LEU:HB3	1:C:245:MET:CE	2.47	0.45
1:E:443:PRO:HB2	1:E:454:LYS:HD3	1.98	0.45
1:F:135:ARG:HD2	1:F:470:ILE:HD11	1.98	0.45
1:C:226:ILE:HG23	1:D:240:ILE:HD11	1.99	0.45
1:H:135:ARG:HD2	1:H:470:ILE:HD11	1.98	0.45
1:F:119:SER:O	1:F:136:ARG:HD2	2.17	0.45
1:C:253:VAL:HG21	1:C:283:ALA:HB1	1.98	0.45
1:H:253:VAL:HG21	1:H:283:ALA:HB1	1.99	0.45
1:A:253:VAL:HG21	1:A:283:ALA:HB1	1.99	0.44
1:G:426:GLN:HB3	1:G:447:GLN:OE1	2.16	0.44
1:C:309:MET:HE2	1:C:309:MET:HB2	1.86	0.44
1:C:426:GLN:HB3	1:C:447:GLN:OE1	2.17	0.44
1:E:35:GLU:HG2	1:E:202:GLY:HA2	1.99	0.44
1:B:253:VAL:HG21	1:B:283:ALA:HB1	1.98	0.44
1:D:119:SER:O	1:D:136:ARG:HD2	2.17	0.44
1:D:219:MET:CE	1:D:244:HIS:HB2	2.48	0.44
1:H:222:LEU:HB3	1:H:245:MET:CE	2.48	0.44
1:H:443:PRO:HB2	1:H:454:LYS:HD3	1.99	0.44
1:A:161:PRO:HG3	2:A:517:HOH:O	2.16	0.44
1:F:222:LEU:HB3	1:F:245:MET:CE	2.47	0.44
1:B:119:SER:O	1:B:136:ARG:HD2	2.17	0.44
1:C:119:SER:O	1:C:136:ARG:HD2	2.17	0.44
1:A:119:SER:O	1:A:136:ARG:HD2	2.18	0.44
1:F:253:VAL:HG21	1:F:283:ALA:HB1	1.98	0.44
1:H:119:SER:O	1:H:136:ARG:HD2	2.18	0.44
1:A:466:LEU:HD13	1:B:447:GLN:CD	2.38	0.44
1:B:35:GLU:HG2	1:B:202:GLY:HA2	1.99	0.44
1:A:472:VAL:HG11	1:C:419:HIS:CD2	2.53	0.44
1:E:2:GLN:HG2	1:E:2:GLN:O	2.17	0.44
1:G:250:LYS:HG3	1:G:285:ARG:HD2	1.99	0.44
1:B:419:HIS:NE2	1:C:419:HIS:NE2	2.63	0.44
1:F:147:PRO:HG3	1:F:223:THR:CG2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:443:PRO:HB2	1:F:454:LYS:HD3	1.99	0.44
1:G:219:MET:CE	1:G:244:HIS:HB2	2.48	0.44
1:A:219:MET:CE	1:A:244:HIS:HB2	2.48	0.43
1:C:219:MET:CE	1:C:244:HIS:HB2	2.48	0.43
1:G:443:PRO:HB2	1:G:454:LYS:HD3	1.99	0.43
1:B:18:ARG:HB3	1:B:32:GLU:HG2	2.00	0.43
1:C:443:PRO:HB2	1:C:454:LYS:HD3	1.99	0.43
1:D:147:PRO:HG3	1:D:223:THR:CG2	2.49	0.43
1:H:284:CYS:HB3	2:H:510:HOH:O	2.17	0.43
1:D:146:ALA:HB3	1:D:173:PRO:HA	2.01	0.43
1:A:443:PRO:HB2	1:A:454:LYS:HD3	1.99	0.43
1:F:146:ALA:HB3	1:F:173:PRO:HA	2.00	0.43
1:G:79:LEU:HD21	1:G:83:ARG:HH11	1.83	0.43
1:A:147:PRO:HG3	1:A:223:THR:CG2	2.48	0.43
1:B:146:ALA:HB3	1:B:173:PRO:HA	2.01	0.43
1:E:146:ALA:HB3	1:E:173:PRO:HA	2.01	0.43
1:F:219:MET:CE	1:F:244:HIS:HB2	2.49	0.43
1:G:147:PRO:HG3	1:G:223:THR:CG2	2.48	0.43
1:C:147:PRO:HG3	1:C:223:THR:CG2	2.48	0.43
1:D:135:ARG:HD2	1:D:470:ILE:HD11	1.99	0.43
1:B:147:PRO:HG3	1:B:223:THR:CG2	2.49	0.43
1:B:219:MET:CE	1:B:244:HIS:HB2	2.49	0.43
1:B:135:ARG:HD2	1:B:470:ILE:HD11	2.00	0.43
1:C:18:ARG:HB3	1:C:32:GLU:HG2	2.00	0.43
1:E:147:PRO:HG3	1:E:223:THR:CG2	2.49	0.43
1:H:18:ARG:HB3	1:H:32:GLU:HG2	2.01	0.43
1:A:447:GLN:CD	1:B:466:LEU:HD13	2.39	0.43
1:C:146:ALA:HB3	1:C:173:PRO:HA	2.01	0.43
1:A:419:HIS:NE2	1:D:419:HIS:NE2	2.66	0.43
1:E:284:CYS:HA	1:E:384:VAL:HG22	2.01	0.42
1:G:309:MET:HE2	1:G:309:MET:HB2	1.91	0.42
1:G:135:ARG:HD2	1:G:470:ILE:HD11	2.01	0.42
1:H:219:MET:CE	1:H:244:HIS:HB2	2.49	0.42
1:A:146:ALA:HB3	1:A:173:PRO:HA	2.01	0.42
1:E:219:MET:CE	1:E:244:HIS:HB2	2.49	0.42
1:A:234:ARG:HG3	1:B:234:ARG:HG3	2.01	0.42
1:E:139:ILE:HD13	1:E:242:ARG:HD2	2.02	0.42
1:G:146:ALA:HB3	1:G:173:PRO:HA	2.01	0.42
1:H:147:PRO:HG3	1:H:223:THR:CG2	2.49	0.42
1:C:284:CYS:HA	1:C:384:VAL:HG22	2.02	0.42
1:D:309:MET:HE3	1:D:361:ALA:HA	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:247:LEU:HD12	1:F:240:ILE:HG21	2.01	0.42
1:F:256:PHE:CZ	1:F:417:ARG:HD3	2.55	0.42
1:F:15:GLU:HB2	1:F:37:SER:HB3	2.02	0.42
1:C:256:PHE:CZ	1:C:417:ARG:HD3	2.55	0.42
1:C:467:VAL:O	1:D:454:LYS:NZ	2.36	0.42
1:D:284:CYS:HA	1:D:384:VAL:HG22	2.01	0.42
1:E:309:MET:HE2	1:E:309:MET:HB2	1.90	0.42
1:G:15:GLU:HB2	1:G:37:SER:HB3	2.01	0.42
1:E:415:VAL:HG22	1:G:415:VAL:HA	2.02	0.42
1:D:15:GLU:HB2	1:D:37:SER:HB3	2.01	0.42
1:E:15:GLU:HB2	1:E:37:SER:HB3	2.01	0.42
1:F:284:CYS:HA	1:F:384:VAL:HG22	2.02	0.42
1:A:15:GLU:HB2	1:A:37:SER:HB3	2.01	0.42
1:E:415:VAL:HG23	1:G:413:GLN:C	2.40	0.42
1:A:284:CYS:HA	1:A:384:VAL:HG22	2.02	0.41
1:H:146:ALA:HB3	1:H:173:PRO:HA	2.01	0.41
1:B:284:CYS:HA	1:B:384:VAL:HG22	2.01	0.41
1:H:284:CYS:HA	1:H:384:VAL:HG22	2.02	0.41
1:B:217:VAL:O	1:B:241:LYS:HE3	2.20	0.41
1:G:284:CYS:HA	1:G:384:VAL:HG22	2.02	0.41
1:A:256:PHE:CZ	1:A:417:ARG:HD3	2.56	0.41
1:F:139:ILE:HD13	1:F:242:ARG:HD2	2.02	0.41
1:F:413:GLN:C	1:H:415:VAL:HG23	2.41	0.41
1:A:240:ILE:HD11	1:B:226:ILE:HG23	2.01	0.41
1:B:19:GLN:NE2	1:B:19:GLN:CA	2.82	0.41
1:C:15:GLU:HB2	1:C:37:SER:HB3	2.02	0.41
1:C:415:VAL:HG13	1:C:416:GLY:N	2.35	0.41
1:H:256:PHE:CZ	1:H:417:ARG:HD3	2.55	0.41
1:D:19:GLN:OE1	1:D:203:ARG:NE	2.54	0.41
1:D:217:VAL:O	1:D:241:LYS:HE3	2.21	0.41
1:E:256:PHE:CZ	1:E:417:ARG:HD3	2.55	0.41
1:G:147:PRO:HG2	1:G:154:MET:HG3	2.01	0.41
1:A:217:VAL:O	1:A:241:LYS:HE3	2.21	0.41
1:B:15:GLU:HB2	1:B:37:SER:HB3	2.02	0.41
1:B:415:VAL:HG23	1:C:413:GLN:O	2.21	0.41
1:A:469:HIS:NE2	1:B:442:MET:HG3	2.36	0.41
1:D:67:LEU:HD11	1:D:164:ALA:HB2	2.03	0.41
1:H:240:ILE:HG22	1:H:240:ILE:O	2.21	0.41
1:G:240:ILE:O	1:G:240:ILE:HG22	2.21	0.41
1:G:18:ARG:HB3	1:G:32:GLU:HG2	2.03	0.41
1:G:256:PHE:CZ	1:G:417:ARG:HD3	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:139:ILE:HD13	1:D:242:ARG:HD2	2.03	0.41
1:F:458:LEU:HD11	1:G:117:CYS:HB3	2.03	0.41
1:H:15:GLU:HB2	1:H:37:SER:HB3	2.02	0.41
1:H:212:THR:O	1:H:241:LYS:HE3	2.21	0.41
1:A:226:ILE:HG23	1:B:240:ILE:HD11	2.03	0.40
1:D:240:ILE:O	1:D:240:ILE:HG22	2.21	0.40
1:A:139:ILE:HD13	1:A:242:ARG:HD2	2.03	0.40
1:B:240:ILE:HG22	1:B:240:ILE:O	2.21	0.40
1:E:240:ILE:O	1:E:240:ILE:HG22	2.21	0.40
1:F:309:MET:HE2	1:F:309:MET:HB2	1.95	0.40
1:G:139:ILE:HD13	1:G:242:ARG:HD2	2.03	0.40
1:A:240:ILE:CD1	1:B:226:ILE:HG23	2.52	0.40
1:C:139:ILE:HD13	1:C:242:ARG:HD2	2.04	0.40
1:F:217:VAL:O	1:F:241:LYS:HE3	2.21	0.40
1:G:217:VAL:O	1:G:241:LYS:HE2	2.21	0.40
1:A:422:SER:HB2	1:B:470:ILE:HD13	2.04	0.40
1:D:415:VAL:HG13	1:D:416:GLY:N	2.36	0.40
1:E:217:VAL:O	1:E:241:LYS:HE3	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:22:TYR:OH	1:G:206:THR:CG2[1_455]	1.98	0.22
1:C:308:LYS:CB	1:E:190:ASP:OD1[1_664]	2.18	0.02

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 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	471/477 (99%)	454 (96%)	17 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
1	C	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
1	D	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
1	E	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
1	F	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
1	G	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
1	H	471/477 (99%)	454 (96%)	17 (4%)	0	100	100
All	All	3768/3816 (99%)	3632 (96%)	136 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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	373/376 (99%)	371 (100%)	2 (0%)	88	96
1	B	373/376 (99%)	370 (99%)	3 (1%)	81	94
1	C	373/376 (99%)	370 (99%)	3 (1%)	81	94
1	D	373/376 (99%)	367 (98%)	6 (2%)	62	86
1	E	373/376 (99%)	368 (99%)	5 (1%)	69	90
1	F	373/376 (99%)	371 (100%)	2 (0%)	88	96
1	G	373/376 (99%)	372 (100%)	1 (0%)	92	98
1	H	373/376 (99%)	370 (99%)	3 (1%)	81	94
All	All	2984/3008 (99%)	2959 (99%)	25 (1%)	81	94

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	88	ASN
1	A	154	MET

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Mol	Chain	Res	Type
1	B	19	GLN
1	B	88	ASN
1	B	154	MET
1	C	88	ASN
1	C	154	MET
1	C	391	ASP
1	D	19	GLN
1	D	88	ASN
1	D	154	MET
1	D	158	LYS
1	D	391	ASP
1	D	417	ARG
1	E	83	ARG
1	E	88	ASN
1	E	92	PRO
1	E	154	MET
1	E	391	ASP
1	F	88	ASN
1	F	154	MET
1	G	88	ASN
1	H	88	ASN
1	H	92	PRO
1	H	154	MET

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

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 carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	473/477 (99%)	-0.38	8 (1%) 70 69	33, 51, 92, 145	0
1	B	473/477 (99%)	-0.21	12 (2%) 57 55	34, 65, 115, 168	0
1	C	473/477 (99%)	-0.41	5 (1%) 80 80	30, 49, 82, 117	0
1	D	473/477 (99%)	-0.20	9 (1%) 66 65	36, 68, 104, 139	0
1	E	473/477 (99%)	-0.32	12 (2%) 57 55	35, 55, 92, 130	0
1	F	473/477 (99%)	-0.24	8 (1%) 70 69	37, 68, 105, 143	0
1	G	473/477 (99%)	-0.19	20 (4%) 36 32	38, 67, 108, 164	0
1	H	473/477 (99%)	-0.34	9 (1%) 66 65	32, 54, 97, 160	0
All	All	3784/3816 (99%)	-0.29	83 (2%) 62 59	30, 59, 103, 168	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	235	HIS	6.3
1	G	238	PRO	5.7
1	G	239	ALA	5.0
1	G	233	LEU	5.0
1	B	239	ALA	4.9
1	B	235	HIS	4.3
1	D	2	GLN	4.2
1	G	231	HIS	4.1
1	F	238	PRO	4.1
1	A	238	PRO	4.0
1	G	236	THR	3.9
1	H	227	ALA	3.9
1	C	2	GLN	3.9
1	B	238	PRO	3.6
1	A	239	ALA	3.6
1	B	236	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	D	22	TYR	3.5
1	B	227	ALA	3.4
1	B	22	TYR	3.4
1	E	314	ASP	3.3
1	H	239	ALA	3.2
1	A	227	ALA	3.2
1	D	403	ARG	3.1
1	A	235	HIS	3.1
1	H	234	ARG	3.1
1	A	403	ARG	3.1
1	B	2	GLN	3.0
1	G	229	GLY	3.0
1	E	356	LYS	3.0
1	G	2	GLN	2.9
1	C	356	LYS	2.9
1	B	215	GLU	2.9
1	A	236	THR	2.8
1	H	403	ARG	2.8
1	E	315	GLU	2.8
1	A	356	LYS	2.8
1	H	315	GLU	2.7
1	G	28	GLU	2.7
1	H	235	HIS	2.7
1	G	22	TYR	2.6
1	G	356	LYS	2.6
1	G	20	SER	2.5
1	C	235	HIS	2.5
1	E	235	HIS	2.5
1	F	234	ARG	2.4
1	E	247	LEU	2.4
1	B	233	LEU	2.4
1	H	228	THR	2.4
1	A	230	GLU	2.4
1	D	20	SER	2.4
1	G	228	THR	2.3
1	E	403	ARG	2.3
1	E	239	ALA	2.3
1	G	234	ARG	2.3
1	G	29	VAL	2.3
1	E	230	GLU	2.3
1	F	239	ALA	2.3
1	C	354	GLU	2.3

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Mol	Chain	Res	Type	RSRZ
1	E	234	ARG	2.3
1	B	242	ARG	2.2
1	E	238	PRO	2.2
1	F	22	TYR	2.2
1	B	28	GLU	2.2
1	F	308	LYS	2.2
1	B	231	HIS	2.2
1	D	238	PRO	2.2
1	D	234	ARG	2.2
1	D	356	LYS	2.2
1	F	230	GLU	2.1
1	G	32	GLU	2.1
1	H	401	ASP	2.1
1	G	450	SER	2.1
1	C	233	LEU	2.1
1	G	340	ALA	2.1
1	H	356	LYS	2.1
1	G	18	ARG	2.1
1	D	29	VAL	2.1
1	E	2	GLN	2.1
1	G	240	ILE	2.1
1	D	354	GLU	2.1
1	E	231	HIS	2.1
1	F	18	ARG	2.0
1	F	31	LEU	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.