



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 23, 2026 – 04:56 AM UTC

PDB ID : 6C43 / pdb_00006c43
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 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
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

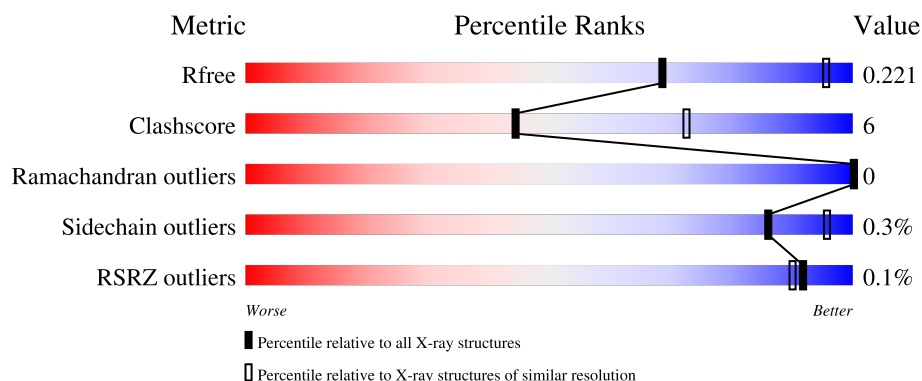
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

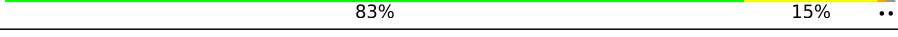
The reported resolution of this entry is 2.90 Å.

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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (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 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	477	
1	B	477	
1	C	477	
1	D	477	
1	E	477	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain	
1	F	477		.
1	G	477		..
1	H	477		..

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

Continued on next page...

Continued from previous page...

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

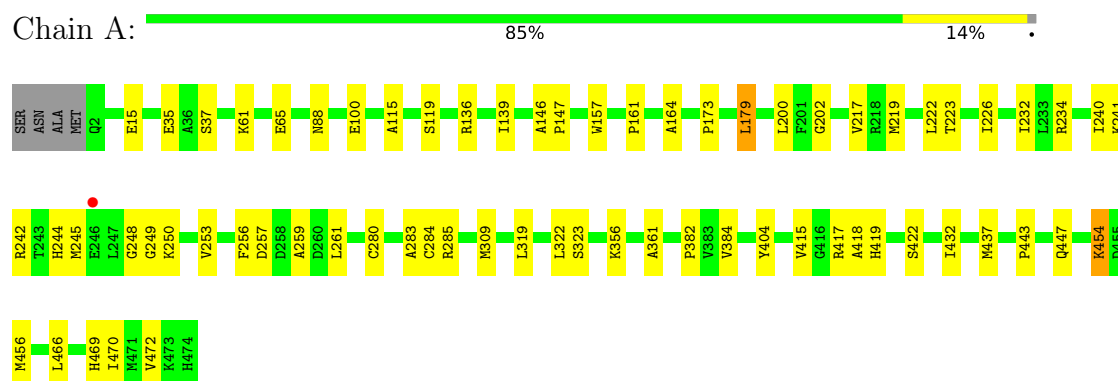
- Molecule 2 is water.

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

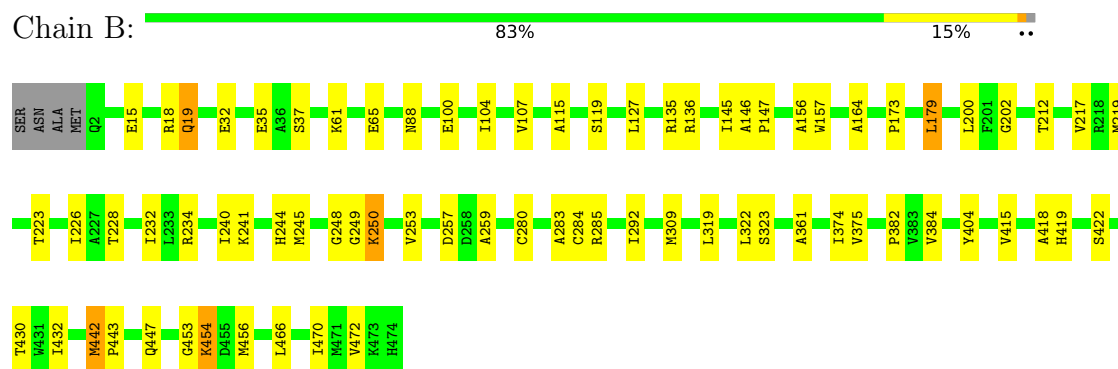
3 Residue-property plots

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.

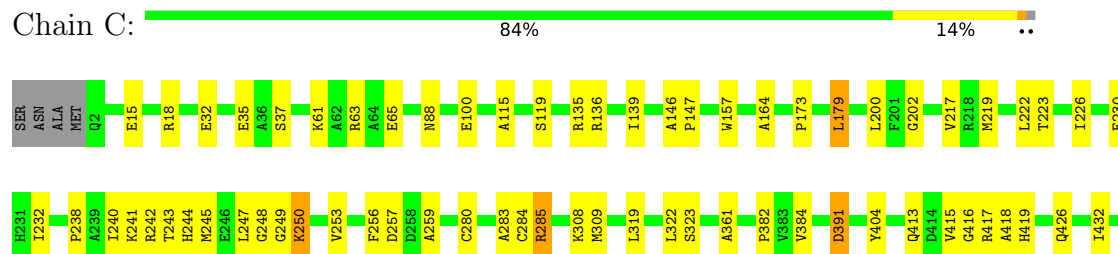
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

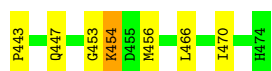


- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase



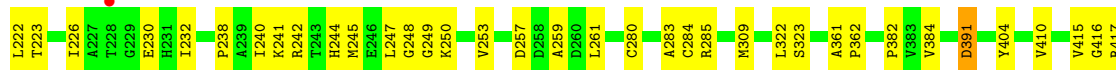
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase





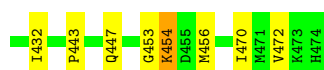
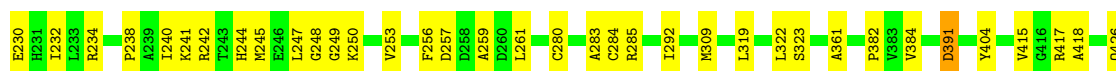
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

Chain D: 83% 15% ..



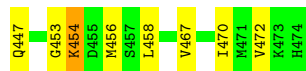
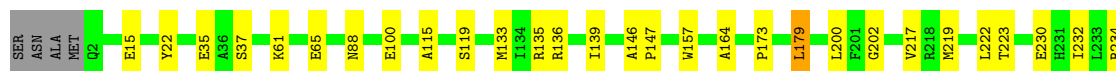
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

Chain E: 84% 15% ..



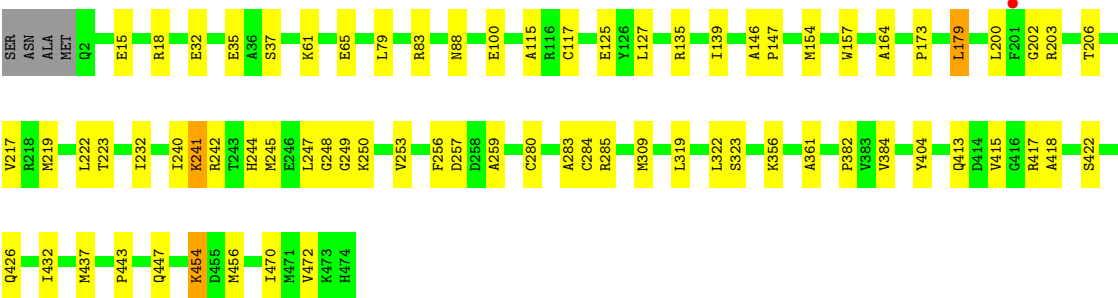
- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

Chain F: 84% 15% .

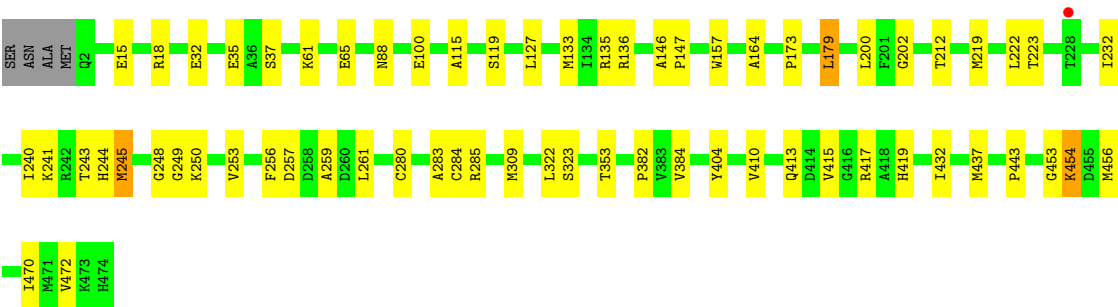
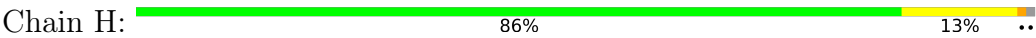


- Molecule 1: Gamma-aminobutyraldehyde dehydrogenase

Chain G: 84% 15% ..



● 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	81.89Å 87.80Å 145.07Å 82.76° 77.91° 72.86°	Depositor
Resolution (Å)	29.73 – 2.90 29.73 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.4 (29.73-2.90) 98.3 (29.73-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.90Å)	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 (4.98%)	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 , 33.9	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	0.52	0/3663	1.02	10/4975 (0.2%)
1	B	0.51	0/3663	1.02	12/4975 (0.2%)
1	C	0.54	1/3663 (0.0%)	1.05	12/4975 (0.2%)
1	D	0.52	0/3663	1.03	12/4975 (0.2%)
1	E	0.53	0/3663	1.02	11/4975 (0.2%)
1	F	0.52	0/3663	1.01	8/4975 (0.2%)
1	G	0.53	0/3663	1.03	11/4975 (0.2%)
1	H	0.51	0/3663	1.01	10/4975 (0.2%)
All	All	0.52	1/29304 (0.0%)	1.02	86/39800 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	285	ARG	CZ-NH1	-6.20	1.24	1.32

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	285	ARG	NE-CZ-NH1	-11.26	110.24	121.50
1	C	285	ARG	NE-CZ-NH2	10.18	128.36	119.20
1	G	241	LYS	CA-CB-CG	9.73	133.57	114.10
1	B	250	LYS	CB-CG-CD	8.54	130.95	111.30
1	H	100	GLU	N-CA-C	8.24	119.96	110.97
1	F	100	GLU	N-CA-C	8.19	119.90	110.97
1	G	100	GLU	N-CA-C	8.14	119.84	110.97
1	A	100	GLU	N-CA-C	8.10	119.80	110.97
1	D	100	GLU	N-CA-C	7.95	119.63	110.97
1	E	100	GLU	N-CA-C	7.90	119.58	110.97
1	C	100	GLU	N-CA-C	7.84	119.52	110.97
1	C	63	ARG	NE-CZ-NH2	7.73	126.16	119.20
1	B	100	GLU	N-CA-C	7.59	119.25	110.97
1	D	158	LYS	CB-CG-CD	7.38	128.27	111.30
1	E	83	ARG	CB-CG-CD	7.08	127.58	111.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	63	ARG	NE-CZ-NH1	-6.58	114.92	121.50
1	H	323	SER	N-CA-C	6.56	119.27	111.33
1	C	323	SER	N-CA-C	6.55	119.26	111.33
1	B	323	SER	N-CA-C	6.50	119.19	111.33
1	F	323	SER	N-CA-C	6.49	119.18	111.33
1	D	323	SER	N-CA-C	6.48	119.17	111.33
1	A	323	SER	N-CA-C	6.47	119.16	111.33
1	G	203	ARG	NE-CZ-NH1	-6.45	115.05	121.50
1	E	323	SER	N-CA-C	6.38	119.05	111.33
1	G	323	SER	N-CA-C	6.35	119.01	111.33
1	D	158	LYS	CA-CB-CG	6.31	126.72	114.10
1	H	259	ALA	N-CA-C	6.17	118.67	110.53
1	A	259	ALA	N-CA-C	6.15	118.65	110.53
1	E	259	ALA	N-CA-C	6.15	118.65	110.53
1	G	203	ARG	CG-CD-NE	-6.14	98.50	112.00
1	C	259	ALA	N-CA-C	6.05	118.52	110.53
1	F	259	ALA	N-CA-C	6.00	118.46	110.53
1	D	259	ALA	N-CA-C	6.00	118.46	110.53
1	G	259	ALA	N-CA-C	5.93	118.36	110.53
1	B	259	ALA	N-CA-C	5.92	118.34	110.53
1	D	158	LYS	CD-CE-NZ	-5.83	93.26	111.90
1	H	453	GLY	N-CA-C	-5.54	105.78	112.48
1	F	179	LEU	N-CA-C	5.51	117.74	111.02
1	D	179	LEU	N-CA-C	5.48	117.71	111.02
1	E	245	MET	N-CA-C	5.47	117.88	109.07
1	B	245	MET	N-CA-C	5.47	117.88	109.07
1	C	454	LYS	N-CA-C	5.45	118.45	109.72
1	H	179	LEU	N-CA-C	5.45	117.67	111.02
1	B	454	LYS	N-CA-C	5.44	118.42	109.72
1	G	179	LEU	N-CA-C	5.44	117.66	111.02
1	D	454	LYS	N-CA-C	5.43	118.41	109.72
1	E	261	LEU	N-CA-C	5.42	117.19	111.28
1	H	454	LYS	N-CA-C	5.42	118.40	109.72
1	A	454	LYS	N-CA-C	5.41	118.37	109.72
1	F	454	LYS	N-CA-C	5.40	118.36	109.72
1	B	179	LEU	N-CA-C	5.40	117.61	111.02
1	E	454	LYS	N-CA-C	5.37	118.31	109.72
1	F	453	GLY	N-CA-C	-5.35	106.00	112.48
1	A	245	MET	N-CA-C	5.34	117.67	109.07
1	A	179	LEU	N-CA-C	5.32	117.52	111.02
1	G	454	LYS	N-CA-C	5.31	118.22	109.72
1	C	179	LEU	N-CA-C	5.28	117.46	111.02

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	179	LEU	N-CA-C	5.27	117.45	111.02
1	E	257	ASP	CA-C-N	5.27	129.03	120.60
1	E	257	ASP	C-N-CA	5.27	129.03	120.60
1	C	453	GLY	N-CA-C	-5.26	106.11	112.48
1	C	257	ASP	CA-C-N	5.25	128.99	120.60
1	C	257	ASP	C-N-CA	5.25	128.99	120.60
1	D	257	ASP	CA-C-N	5.23	128.97	120.60
1	D	257	ASP	C-N-CA	5.23	128.97	120.60
1	F	257	ASP	CA-C-N	5.22	128.95	120.60
1	F	257	ASP	C-N-CA	5.22	128.95	120.60
1	B	257	ASP	CA-C-N	5.19	128.91	120.60
1	B	257	ASP	C-N-CA	5.19	128.91	120.60
1	A	257	ASP	CA-C-N	5.16	128.85	120.60
1	A	257	ASP	C-N-CA	5.16	128.85	120.60
1	G	257	ASP	CA-C-N	5.16	128.85	120.60
1	G	257	ASP	C-N-CA	5.16	128.85	120.60
1	B	453	GLY	N-CA-C	-5.14	106.25	112.48
1	A	356	LYS	N-CA-C	5.12	116.54	111.07
1	H	257	ASP	CA-C-N	5.11	128.77	120.60
1	H	257	ASP	C-N-CA	5.11	128.77	120.60
1	H	261	LEU	N-CA-C	5.10	116.83	111.28
1	D	218	ARG	CG-CD-NE	5.08	123.17	112.00
1	H	245	MET	N-CA-C	5.05	117.20	109.07
1	G	356	LYS	N-CA-C	5.04	116.47	111.07
1	B	442	MET	CA-C-N	5.04	124.94	119.85
1	B	442	MET	C-N-CA	5.04	124.94	119.85
1	E	453	GLY	N-CA-C	-5.01	106.42	112.48
1	D	261	LEU	N-CA-C	5.01	116.74	111.28
1	A	261	LEU	N-CA-C	5.01	116.74	111.28

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	45	0
1	B	3591	0	3571	58	0
1	C	3591	0	3571	56	1
1	D	3591	0	3571	60	0
1	E	3591	0	3571	49	1
1	F	3591	0	3571	52	1
1	G	3591	0	3571	44	1
1	H	3591	0	3571	45	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	343	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 6.

All (343) 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:G:245:MET:HE3	1:G:247:LEU:HD11	1.54	0.89
1:B:250:LYS:NZ	1:B:374:ILE:O	2.07	0.86
1:D:245:MET:HE3	1:D:247:LEU:HD11	1.56	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: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:D:309:MET:HE1	1:D:361:ALA:HA	1.71	0.72
1:E:91:LYS:NZ	1:E:149:ASN:O	2.18	0.71
1:E:247:LEU:CD1	1:F:240:ILE:HD13	2.21	0.70

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:432:ILE:HD12	1:D:472:VAL:HG22	1.72	0.70
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:D:219:MET:HE2	1:D:244:HIS:HB2	1.80	0.64
1:G:219:MET:HE2	1:G:244:HIS:HB2	1.80	0.63
1:C:219:MET:HE2	1:C:244:HIS:HB2	1.80	0.63
1:A:470:ILE:HD12	1:B:422:SER:HB2	1.79	0.62
1:H:309:MET:HE1	1:H:353:THR:HG21	1.81	0.62
1:H:219:MET:HE2	1:H:244:HIS:HB2	1.81	0.62
1:F:219:MET:HE2	1:F:244:HIS:HB2	1.81	0.61
1:E:219:MET:HE2	1:E:244:HIS:HB2	1.82	0.61
1:A:219:MET:HE2	1:A:244:HIS:HB2	1.81	0.61
1:C:240:ILE:HG21	1:D:247:LEU:HD12	1.82	0.60
1:B:219:MET:HE2	1:B:244:HIS:HB2	1.81	0.60
1:D:309:MET:HE3	1:D:362:PRO:HD3	1.83	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:C	2.28	0.59
1:F:415:VAL:HG23	1:H:413:GLN:O	2.03	0.59
1:A:309:MET:HE1	1:A:361:ALA:HA	1.86	0.58
1:F:322:LEU:HD21	1:F:382:PRO:HD3	1.85	0.58
1:B:309:MET:HE1	1:B:361:ALA:HA	1.85	0.57
1:C:309:MET:HE1	1:C:361:ALA:HA	1.85	0.57
1:G:309:MET:HE1	1:G:361:ALA:HA	1.87	0.57
1:E:309:MET:HE1	1:E:361:ALA:HA	1.86	0.57
1:E:322:LEU:HD21	1:E:382:PRO:HD3	1.87	0.57
1:A:249:GLY:HA2	1:A:404:TYR:CD1	2.40	0.56
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: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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:309:MET:HE1	1:F:361:ALA:HA	1.86	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:G:249:GLY:HA2	1:G:404:TYR:CD1	2.41	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:B:157:TRP:NE1	1:B:456:MET:HE3	2.23	0.53
1:A:157:TRP:NE1	1:A:456:MET:HE3	2.24	0.53
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:F:157:TRP:NE1	1:F:456:MET:HE3	2.24	0.53
1:G:157:TRP:NE1	1:G:456:MET:HE3	2.23	0.53
1:C:157:TRP:NE1	1:C:456:MET:HE3	2.24	0.53
1:C:247:LEU:HD12	1:D:240:ILE:HD13	1.91	0.53
1:B:19:GLN:HA	1:B:19:GLN:NE2	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:H:157:TRP:NE1	1:H:456:MET:HE3	2.23	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:61:LYS:HE2	1:D:65:GLU:OE2	2.10	0.52
1:D:248:GLY:HA2	1:D:280:CYS:SG	2.50	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:E:157:TRP:NE1	1:E:456:MET:HE3	2.25	0.52
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:D:157:TRP:NE1	1:D:456:MET:HE3	2.24	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:H:222:LEU:HB3	1:H:245:MET:HE1	1.94	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:OD1	1:D:391:ASP:N	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:C:240:ILE:HD11	1:D:226:ILE:HG23	1.93	0.50
1:D:88:ASN:HB2	1:D:179:LEU:HD12	1.93	0.50
1:F:115:ALA:HB2	1:F:164:ALA:HB1	1.93	0.50
1:F:222:LEU:HB3	1:F:245:MET:HE1	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:C:222:LEU:HB3	1:C:245:MET:HE1	1.94	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:C:147:PRO:HD3	1:C:223:THR:HB	1.95	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:226:ILE:HG23	1:D:240:ILE:CD1	2.43	0.49
1:D:115:ALA:HB2	1:D:164:ALA:HB1	1.93	0.49
1:E:240:ILE:HD13	1:F:247:LEU:CD1	2.42	0.49
1:F:147:PRO:HD3	1:F:223:THR:HB	1.95	0.49
1:C:135:ARG:HD2	1:C:470:ILE:HD11	1.94	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:410:VAL:HB	1:F:432:ILE:HD13	1.94	0.48
1:F:419:HIS:CD2	1:G:472:VAL:HG11	2.49	0.48
1:A:35:GLU:HG2	1:A:202:GLY:HA2	1.96	0.48
1:A:437:MET:HE2	1:B:127:LEU:HD21	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:A:466:LEU:HD13	1:B:447:GLN:CD	2.38	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:A:447:GLN:CD	1:B:466:LEU:HD13	2.39	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:35:GLU:HG2	1:G:202:GLY:HA2	1.97	0.47
1:G:173:PRO:HD2	1:G:200:LEU:O	2.14	0.47
1:B:472:VAL:HG11	1:D:419:HIS:CD2	2.50	0.47
1:E:415:VAL:HG23	1:G:413:GLN:C	2.40	0.46
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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:413:GLN:C	1:H:415:VAL:HG23	2.41	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
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:G:309:MET:HE2	1:G:319:LEU:HD23	1.98	0.45
1:C:243:THR:HB	1:C:245:MET:HE3	1.99	0.45
1:E:309:MET:HE2	1:E:319:LEU:HD23	1.98	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:F:243:THR:HB	1:F:245:MET:HE3	1.99	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:A:309:MET:HE2	1:A:319:LEU:HD23	1.99	0.44
1:F:222:LEU:HB3	1:F:245:MET:CE	2.47	0.44
1:F:309:MET:HE2	1:F:319:LEU:HD23	1.99	0.44
1:G:426:GLN:HB3	1:G:447:GLN:OE1	2.16	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:G:437:MET:HE2	1:H:127:LEU:HD21	2.00	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

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:127:LEU:HD21	1:H:437:MET:HE2	1.98	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:B:309:MET:HE2	1:B:319:LEU:HD23	1.99	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:C:309:MET:HE2	1:C:319:LEU:HD23	1.99	0.44
1:H:243:THR:HB	1:H:245:MET:HE3	1.99	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:472:VAL:HG11	1:C:419:HIS:CD2	2.53	0.44
1:B:35:GLU:HG2	1:B:202:GLY:HA2	1.99	0.44
1:E:2:GLN:O	1:E:2:GLN:HG2	2.17	0.44
1:G:250:LYS:HG3	1:G:285:ARG:HD2	1.99	0.44
1:B:309:MET:HE3	1:B:309:MET:HB2	1.92	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
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: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:309:MET:HE3	1:A:309:MET:HB2	1.93	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:H:244:HIS:O	1:H:245:MET:HE2	2.18	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:135:ARG:HD2	1:B:470:ILE:HD11	2.00	0.43
1:B:147:PRO:HG3	1:B:223:THR:CG2	2.49	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:ARG:HB3	1:C:32:GLU:HG2	2.00	0.43
1:C:466:LEU:HD13	1:D:447:GLN:CD	2.44	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:419:HIS:NE2	1:D:419:HIS:NE2	2.66	0.43
1:C:146:ALA:HB3	1:C:173:PRO:HA	2.01	0.43
1:E:284:CYS:HA	1:E:384:VAL:HG22	2.01	0.42
1:G:135:ARG:HD2	1:G:470:ILE:HD11	2.01	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:F:244:HIS:O	1:F:245:MET:HE2	2.19	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:E:247:LEU:HD12	1:F:240:ILE:HG21	2.01	0.42
1:F:15:GLU:HB2	1:F:37:SER:HB3	2.02	0.42
1:F:256:PHE:CZ	1:F:417:ARG:HD3	2.55	0.42
1:C:244:HIS:O	1:C:245:MET:HE2	2.20	0.42
1:C:256:PHE:CZ	1:C:417:ARG:HD3	2.55	0.42
1:D:284:CYS:HA	1:D:384:VAL:HG22	2.01	0.42
1:E:415:VAL:HG22	1:G:415:VAL:HA	2.02	0.42
1:G:15:GLU:HB2	1:G:37:SER:HB3	2.01	0.42
1:B:219:MET:CE	1:B:244:HIS:HB2	2.49	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:F:133:MET:HE1	1:H:133:MET:HE1	2.02	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:A:240:ILE:HD11	1:B:226:ILE:HG23	2.01	0.41
1:B:19:GLN:CA	1:B:19:GLN:NE2	2.82	0.41
1:C:15:GLU:HB2	1:C:37:SER:HB3	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:E:135:ARG:HH22	1:G:125:GLU:CD	2.28	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:A:469:HIS:NE2	1:B:442:MET:HG3	2.36	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:D:67:LEU:HD11	1:D:164:ALA:HB2	2.03	0.41
1:G:18:ARG:HB3	1:G:32:GLU:HG2	2.03	0.41
1:G:240:ILE:HG22	1:G:240:ILE:O	2.21	0.41
1:G:256:PHE:CZ	1:G:417:ARG:HD3	2.56	0.41
1:H:240:ILE:O	1:H:240:ILE:HG22	2.21	0.41
1:D:139:ILE:HD13	1:D:242:ARG:HD2	2.03	0.41
1:F:133:MET:CE	1:H:133:MET:HE1	2.51	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:B:292:ILE:HD12	1:B:292:ILE:C	2.47	0.40
1:C:309:MET:HE3	1:C:309:MET:HB2	1.91	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:O	1:B:240:ILE:HG22	2.21	0.40
1:E:240:ILE:O	1:E:240:ILE:HG22	2.21	0.40
1:E:292:ILE:HD12	1:E:292:ILE:C	2.47	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:D:309:MET:HE1	1:D:361:ALA:CA	2.46	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
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 [i](#)

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%)	373 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	373/376 (99%)	372 (100%)	1 (0%)	86	96
1	C	373/376 (99%)	371 (100%)	2 (0%)	81	93
1	D	373/376 (99%)	369 (99%)	4 (1%)	65	88
1	E	373/376 (99%)	370 (99%)	3 (1%)	73	90
1	F	373/376 (99%)	373 (100%)	0	100	100
1	G	373/376 (99%)	373 (100%)	0	100	100
1	H	373/376 (99%)	373 (100%)	0	100	100
All	All	2984/3008 (99%)	2974 (100%)	10 (0%)	86	96

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	19	GLN
1	C	250	LYS
1	C	391	ASP
1	D	19	GLN
1	D	158	LYS
1	D	391	ASP
1	D	417	ARG
1	E	32	GLU
1	E	83	ARG
1	E	391	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	352	GLN
1	A	413	GLN
1	B	231	HIS
1	B	352	GLN
1	B	413	GLN
1	C	57	GLN
1	C	352	GLN
1	C	413	GLN
1	D	8	ASN
1	D	19	GLN
1	D	205	GLN
1	D	352	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	19	GLN
1	E	352	GLN
1	F	8	ASN
1	F	19	GLN
1	F	352	GLN
1	G	19	GLN
1	G	352	GLN
1	G	413	GLN
1	H	8	ASN
1	H	19	GLN
1	H	352	GLN
1	H	413	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	473/477 (99%)	-0.42	1 (0%) 91 89	33, 51, 92, 145	0
1	B	473/477 (99%)	-0.22	0 100 100	34, 65, 115, 168	0
1	C	473/477 (99%)	-0.49	0 100 100	30, 49, 82, 117	0
1	D	473/477 (99%)	-0.27	1 (0%) 91 89	36, 68, 104, 139	0
1	E	473/477 (99%)	-0.39	0 100 100	35, 55, 92, 130	0
1	F	473/477 (99%)	-0.26	1 (0%) 91 89	37, 68, 105, 143	0
1	G	473/477 (99%)	-0.26	1 (0%) 91 89	38, 67, 108, 164	0
1	H	473/477 (99%)	-0.35	1 (0%) 91 89	32, 54, 97, 160	0
All	All	3784/3816 (99%)	-0.33	5 (0%) 92 90	30, 59, 103, 168	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	446	GLY	2.5
1	G	201	PHE	2.4
1	A	246	GLU	2.4
1	H	228	THR	2.4
1	D	228	THR	2.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.