



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:11 AM UTC

PDB ID : 4XTK / pdb_00004xtk
Title : Structure of TM1797, a CAS1 protein from *Thermotoga maritima*
Authors : Petit, P.; Beloglazova, N.; Skarina, T.; Chang, C.; Edwards, A.; Joachimiak, A.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2015-01-23
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

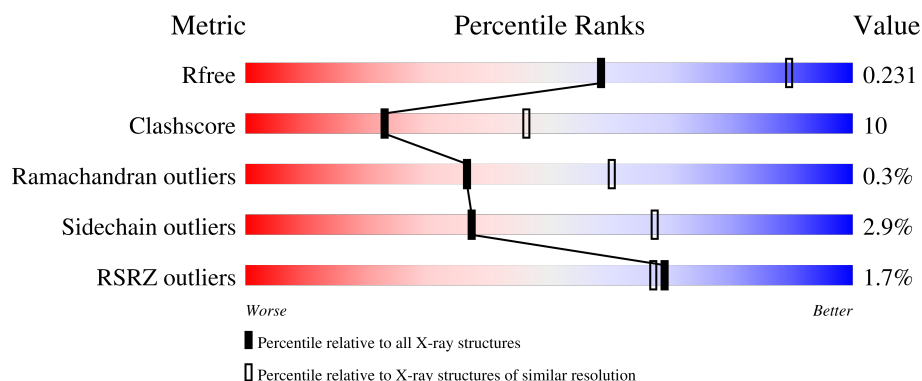
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	326	
1	B	326	
1	C	326	
1	D	326	
1	E	326	

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Mol	Chain	Length	Quality of chain
1	F	326	2% 74% 21%
1	G	326	2% 71% 21% 7%
1	H	326	2% 67% 22% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 20014 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 CRISPR-associated endonuclease Cas1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2476	1583	426	457	10			
1	B	309	Total	C	N	O	S	0	0	0
			2459	1575	415	459	10			
1	C	298	Total	C	N	O	S	0	0	0
			2382	1532	398	443	9			
1	D	311	Total	C	N	O	S	0	0	0
			2487	1595	423	460	9			
1	E	319	Total	C	N	O	S	0	0	0
			2609	1675	452	472	10			
1	F	314	Total	C	N	O	S	0	1	0
			2562	1651	443	458	10			
1	G	304	Total	C	N	O	S	0	1	0
			2465	1581	428	447	9			
1	H	300	Total	C	N	O	S	0	1	0
			2441	1569	421	442	9			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	ARG	-	expression tag	UNP Q9X2B7
A	-5	GLU	-	expression tag	UNP Q9X2B7
A	-4	LEU	-	expression tag	UNP Q9X2B7
A	-3	TYR	-	expression tag	UNP Q9X2B7
A	-2	PHE	-	expression tag	UNP Q9X2B7
A	-1	GLN	-	expression tag	UNP Q9X2B7
A	0	GLY	-	expression tag	UNP Q9X2B7
B	-6	ARG	-	expression tag	UNP Q9X2B7
B	-5	GLU	-	expression tag	UNP Q9X2B7
B	-4	LEU	-	expression tag	UNP Q9X2B7
B	-3	TYR	-	expression tag	UNP Q9X2B7
B	-2	PHE	-	expression tag	UNP Q9X2B7
B	-1	GLN	-	expression tag	UNP Q9X2B7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP Q9X2B7
C	-6	ARG	-	expression tag	UNP Q9X2B7
C	-5	GLU	-	expression tag	UNP Q9X2B7
C	-4	LEU	-	expression tag	UNP Q9X2B7
C	-3	TYR	-	expression tag	UNP Q9X2B7
C	-2	PHE	-	expression tag	UNP Q9X2B7
C	-1	GLN	-	expression tag	UNP Q9X2B7
C	0	GLY	-	expression tag	UNP Q9X2B7
D	-6	ARG	-	expression tag	UNP Q9X2B7
D	-5	GLU	-	expression tag	UNP Q9X2B7
D	-4	LEU	-	expression tag	UNP Q9X2B7
D	-3	TYR	-	expression tag	UNP Q9X2B7
D	-2	PHE	-	expression tag	UNP Q9X2B7
D	-1	GLN	-	expression tag	UNP Q9X2B7
D	0	GLY	-	expression tag	UNP Q9X2B7
E	-6	ARG	-	expression tag	UNP Q9X2B7
E	-5	GLU	-	expression tag	UNP Q9X2B7
E	-4	LEU	-	expression tag	UNP Q9X2B7
E	-3	TYR	-	expression tag	UNP Q9X2B7
E	-2	PHE	-	expression tag	UNP Q9X2B7
E	-1	GLN	-	expression tag	UNP Q9X2B7
E	0	GLY	-	expression tag	UNP Q9X2B7
F	-6	ARG	-	expression tag	UNP Q9X2B7
F	-5	GLU	-	expression tag	UNP Q9X2B7
F	-4	LEU	-	expression tag	UNP Q9X2B7
F	-3	TYR	-	expression tag	UNP Q9X2B7
F	-2	PHE	-	expression tag	UNP Q9X2B7
F	-1	GLN	-	expression tag	UNP Q9X2B7
F	0	GLY	-	expression tag	UNP Q9X2B7
G	-6	ARG	-	expression tag	UNP Q9X2B7
G	-5	GLU	-	expression tag	UNP Q9X2B7
G	-4	LEU	-	expression tag	UNP Q9X2B7
G	-3	TYR	-	expression tag	UNP Q9X2B7
G	-2	PHE	-	expression tag	UNP Q9X2B7
G	-1	GLN	-	expression tag	UNP Q9X2B7
G	0	GLY	-	expression tag	UNP Q9X2B7
H	-6	ARG	-	expression tag	UNP Q9X2B7
H	-5	GLU	-	expression tag	UNP Q9X2B7
H	-4	LEU	-	expression tag	UNP Q9X2B7
H	-3	TYR	-	expression tag	UNP Q9X2B7
H	-2	PHE	-	expression tag	UNP Q9X2B7
H	-1	GLN	-	expression tag	UNP Q9X2B7

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Chain	Residue	Modelled	Actual	Comment	Reference
H	0	GLY	-	expression tag	UNP Q9X2B7

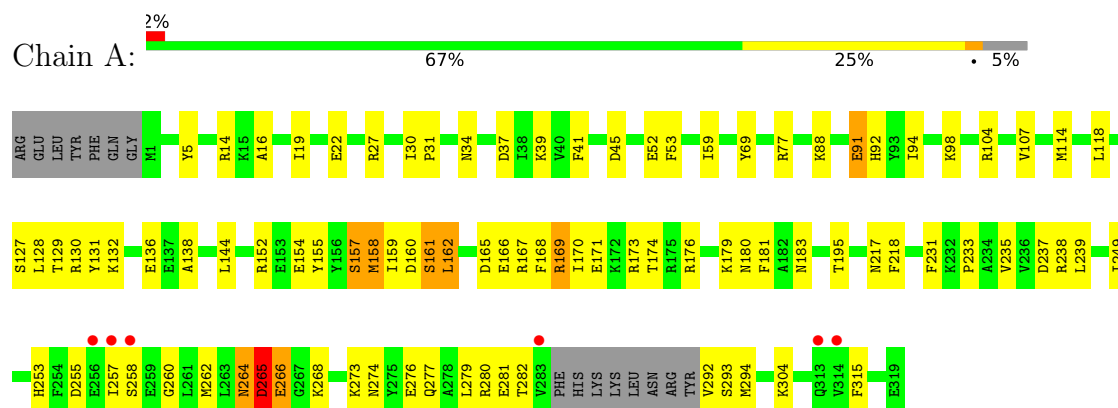
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	11	Total O 11 11	0	0
2	B	8	Total O 8 8	0	0
2	C	9	Total O 9 9	0	0
2	D	18	Total O 18 18	0	0
2	E	30	Total O 30 30	0	0
2	F	23	Total O 23 23	0	0
2	G	18	Total O 18 18	0	0
2	H	16	Total O 16 16	0	0

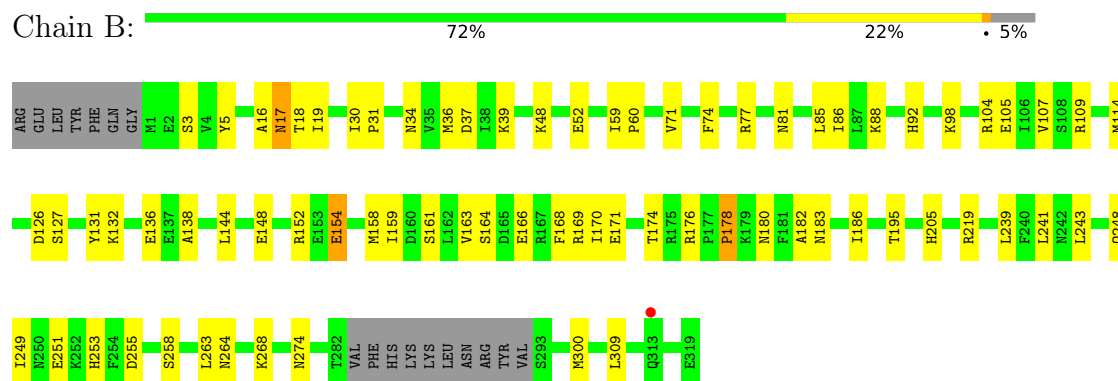
3 Residue-property plots

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.

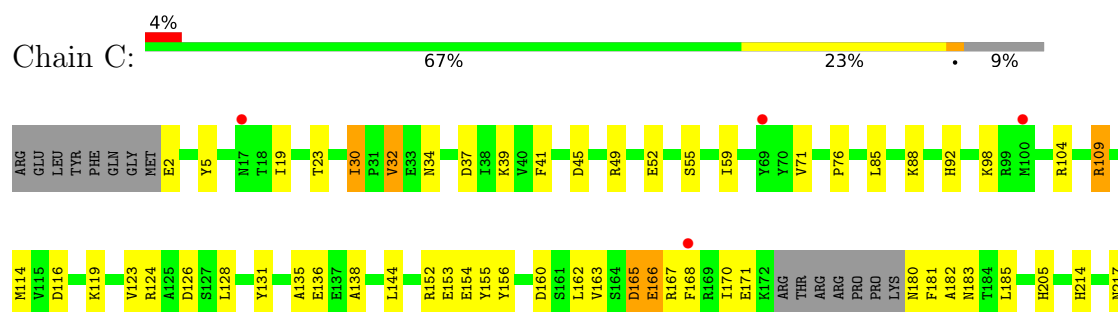
• Molecule 1: CRISPR-associated endonuclease Cas1

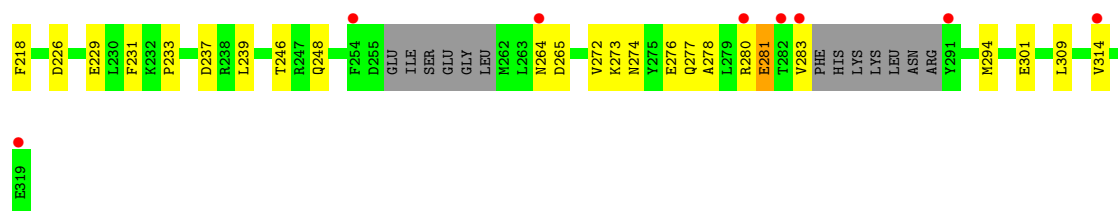


• Molecule 1: CRISPR-associated endonuclease Cas1

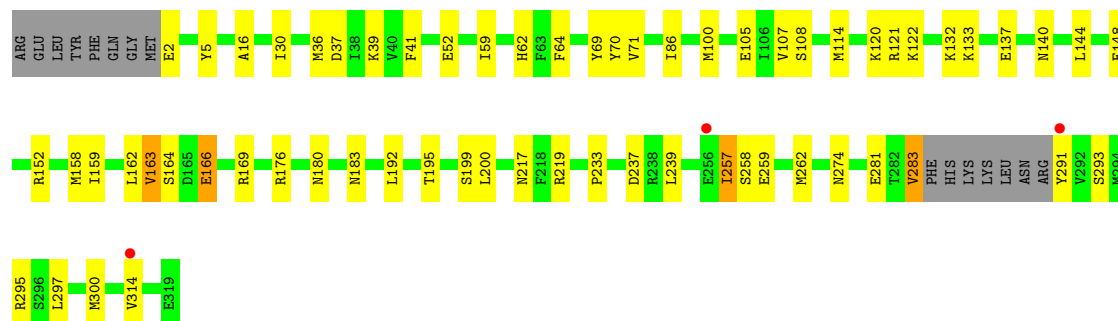
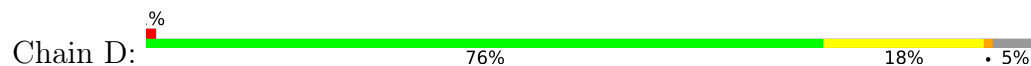


• Molecule 1: CRISPR-associated endonuclease Cas1

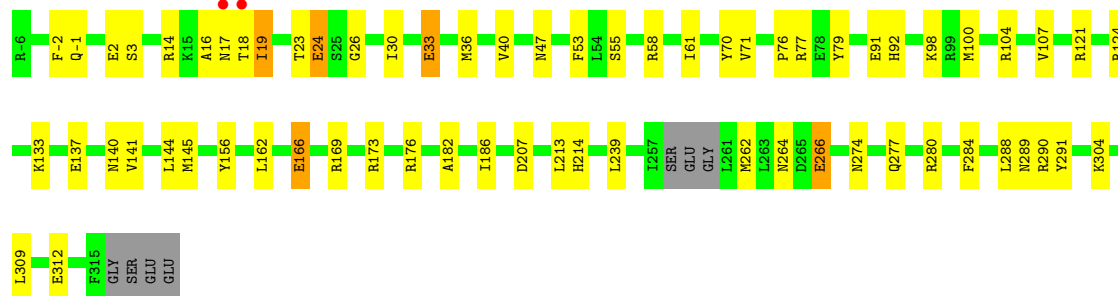
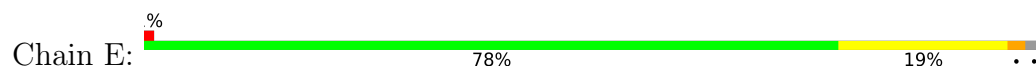




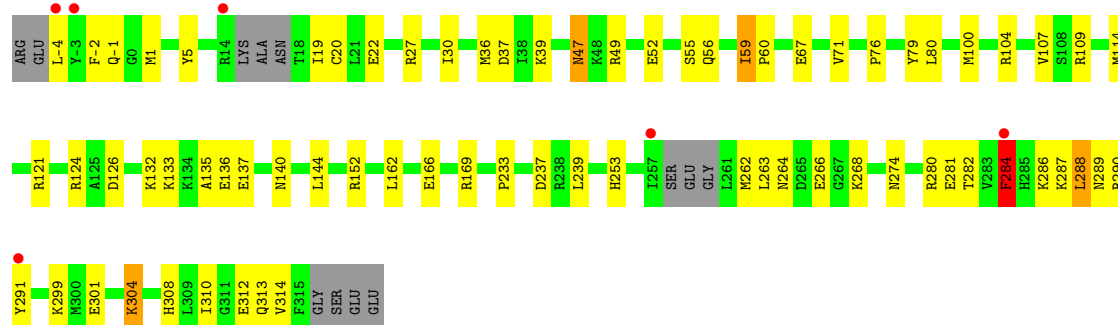
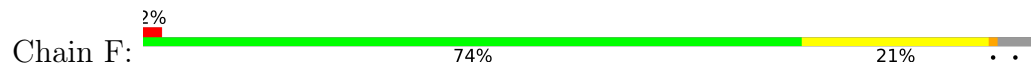
• Molecule 1: CRISPR-associated endonuclease Cas1



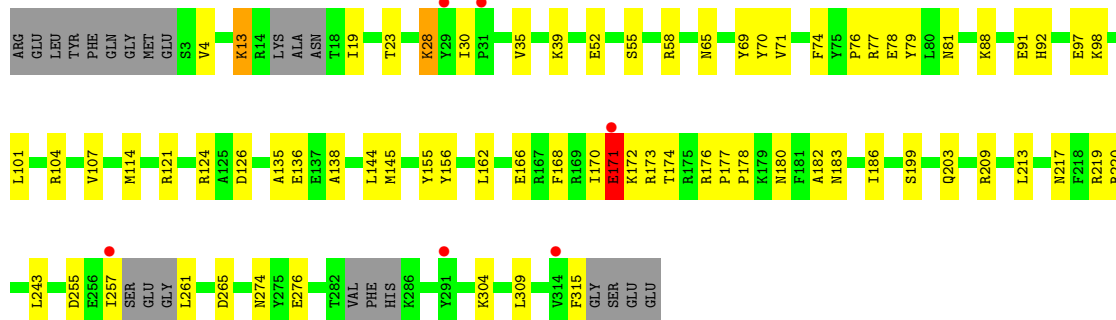
• Molecule 1: CRISPR-associated endonuclease Cas1



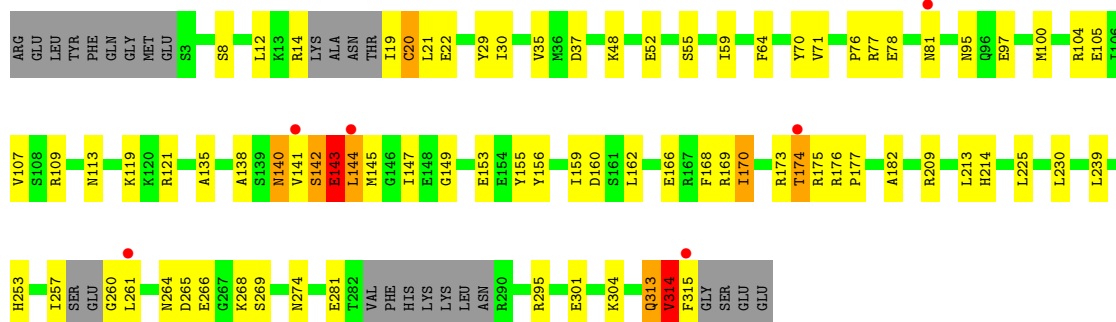
• Molecule 1: CRISPR-associated endonuclease Cas1



• Molecule 1: CRISPR-associated endonuclease Cas1



• Molecule 1: CRISPR-associated endonuclease Cas1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	92.28Å 94.56Å 106.39Å 93.03° 115.06° 102.97°	Depositor
Resolution (Å)	19.85 – 2.70 19.85 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.6 (19.85-2.70) 97.3 (19.85-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.09 (at 2.71Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1702)	Depositor
R, R_{free}	0.183 , 0.230 0.188 , 0.231	Depositor DCC
R_{free} test set	4202 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.346	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 67.8	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	0.95	EDS
Total number of atoms	20014	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.74% 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	0.45	0/2522	0.91	8/3402 (0.2%)
1	B	0.41	0/2505	0.88	6/3380 (0.2%)
1	C	0.41	0/2425	0.92	5/3269 (0.2%)
1	D	0.42	0/2534	0.84	3/3417 (0.1%)
1	E	0.37	0/2660	0.78	6/3581 (0.2%)
1	F	0.40	0/2616	0.88	6/3523 (0.2%)
1	G	0.39	0/2513	0.89	6/3385 (0.2%)
1	H	0.41	0/2490	0.90	10/3351 (0.3%)
All	All	0.41	0/20265	0.87	50/27308 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
1	C	0	2
1	D	0	1
1	G	0	1
1	H	0	2
All	All	0	9

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	176	ARG	CA-C-N	-9.84	110.24	120.38
1	B	176	ARG	C-N-CA	-9.84	110.24	120.38
1	F	288	LEU	N-CA-C	-9.37	95.09	109.24
1	D	176	ARG	CA-C-N	-9.30	110.80	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	176	ARG	C-N-CA	-9.30	110.80	120.38
1	A	217	ASN	CA-C-N	-8.47	110.25	123.12
1	A	217	ASN	C-N-CA	-8.47	110.25	123.12
1	F	288	LEU	CA-C-N	-8.22	110.72	123.13
1	F	288	LEU	C-N-CA	-8.22	110.72	123.13
1	A	176	ARG	CA-C-N	-8.14	112.00	120.38
1	A	176	ARG	C-N-CA	-8.14	112.00	120.38
1	C	217	ASN	CA-C-N	-7.97	111.17	122.41
1	C	217	ASN	C-N-CA	-7.97	111.17	122.41
1	G	171	GLU	CB-CA-C	-7.50	107.03	117.23
1	A	217	ASN	N-CA-C	-7.10	98.63	109.85
1	C	217	ASN	N-CA-C	-6.84	98.77	109.72
1	E	207	ASP	CA-C-N	6.67	126.30	119.56
1	E	207	ASP	C-N-CA	6.67	126.30	119.56
1	C	23	THR	N-CA-C	-6.64	100.96	110.59
1	G	91	GLU	CA-C-N	-6.52	110.58	122.38
1	G	91	GLU	C-N-CA	-6.52	110.58	122.38
1	H	174	THR	N-CA-C	6.35	121.17	113.17
1	F	284	PHE	CA-C-N	-6.33	112.64	122.59
1	F	284	PHE	C-N-CA	-6.33	112.64	122.59
1	H	140	ASN	N-CA-C	-6.26	97.47	110.80
1	C	278	ALA	N-CA-C	-6.03	104.70	111.28
1	H	177	PRO	CA-C-N	5.94	125.91	120.03
1	H	177	PRO	C-N-CA	5.94	125.91	120.03
1	H	20	CYS	N-CA-C	5.66	117.67	109.07
1	E	24	GLU	N-CA-C	-5.65	103.46	110.41
1	A	264	ASN	N-CA-C	5.64	117.35	110.41
1	A	260	GLY	N-CA-C	-5.64	103.19	110.69
1	H	143	GLU	CA-CB-CG	5.62	125.35	114.10
1	B	34	ASN	N-CA-C	5.56	119.76	112.92
1	F	59	ILE	N-CA-C	5.41	113.90	107.73
1	A	169	ARG	N-CA-C	5.33	118.37	110.48
1	H	35	VAL	N-CA-C	5.29	116.36	108.54
1	D	16	ALA	CB-CA-C	-5.25	110.50	116.54
1	H	314	VAL	N-CA-C	5.24	120.23	109.34
1	B	178	PRO	CA-C-O	-5.23	115.45	121.36
1	E	91	GLU	CA-C-N	-5.19	112.41	122.06
1	E	91	GLU	C-N-CA	-5.19	112.41	122.06
1	G	171	GLU	CB-CG-CD	5.17	121.39	112.60
1	G	28	LYS	CA-C-N	-5.14	115.94	122.77
1	G	28	LYS	C-N-CA	-5.14	115.94	122.77
1	B	258	SER	CB-CA-C	-5.12	110.65	116.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	26	GLY	N-CA-C	5.10	120.31	112.51
1	B	180	ASN	N-CA-C	5.07	114.84	108.45
1	H	143	GLU	CA-C-N	5.04	131.16	121.54
1	H	143	GLU	C-N-CA	5.04	131.16	121.54

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	16	ALA	Peptide
1	A	265	ASP	Peptide
1	B	16	ALA	Peptide
1	C	126	ASP	Peptide
1	C	165	ASP	Peptide
1	D	257	ILE	Peptide
1	G	65	ASN	Peptide
1	H	142	SER	Peptide
1	H	313	GLN	Peptide

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	2476	0	2405	71	0
1	B	2459	0	2386	49	0
1	C	2382	0	2310	56	0
1	D	2487	0	2426	44	0
1	E	2609	0	2567	47	0
1	F	2562	0	2528	50	1
1	G	2465	0	2421	53	1
1	H	2441	0	2404	71	0
2	A	11	0	0	0	0
2	B	8	0	0	0	0
2	C	9	0	0	1	0
2	D	18	0	0	0	0
2	E	30	0	0	1	0
2	F	23	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	18	0	0	0	0
2	H	16	0	0	3	0
All	All	20014	0	19447	409	1

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 10.

All (409) 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:154:GLU:OE2	1:A:158:MET:HE1	1.37	1.24
1:A:255:ASP:OD1	1:A:264:ASN:ND2	1.84	1.09
1:A:154:GLU:OE2	1:A:158:MET:CE	2.00	1.08
1:G:92:HIS:ND1	1:G:98:LYS:HD3	1.75	1.01
1:E:-1:GLN:NE2	1:F:36:MET:SD	2.50	0.85
1:H:143:GLU:CG	1:H:144:LEU:H	1.91	0.84
1:F:282:THR:HG21	1:F:291:TYR:HD2	1.45	0.82
1:G:92:HIS:HD2	1:G:309:LEU:HD23	1.42	0.81
1:D:239:LEU:HD11	1:D:274:ASN:HB3	1.62	0.81
1:A:255:ASP:OD1	1:A:264:ASN:CG	2.24	0.80
1:E:92:HIS:CD2	1:E:309:LEU:HD23	2.17	0.79
1:C:109:ARG:NH2	1:C:301:GLU:OE2	2.15	0.79
1:D:100:MET:HE1	1:D:140:ASN:HA	1.65	0.79
1:C:273:LYS:HA	1:C:276:GLU:HB2	1.64	0.79
1:G:104:ARG:NH1	1:G:135:ALA:O	2.16	0.78
1:E:166:GLU:HA	1:E:169:ARG:HG3	1.65	0.77
1:H:140:ASN:OD1	1:H:143:GLU:HB3	1.85	0.76
1:C:104:ARG:NH1	1:C:135:ALA:O	2.20	0.75
1:C:153:GLU:HA	1:C:156:TYR:HD2	1.52	0.74
1:B:52:GLU:HG2	1:G:71:VAL:HG11	1.70	0.74
1:D:163:VAL:O	1:D:169:ARG:NH2	2.22	0.72
1:E:288:LEU:HB3	1:E:290:ARG:HD2	1.71	0.72
1:A:157:SER:O	1:A:159:ILE:N	2.22	0.71
1:C:170:ILE:HG23	1:C:183:ASN:HD21	1.54	0.71
1:B:169:ARG:NH2	1:B:171:GLU:OE2	2.23	0.71
1:C:170:ILE:O	1:C:171:GLU:HG2	1.91	0.70
1:A:170:ILE:HG23	1:A:183:ASN:HD21	1.54	0.70
1:H:107:VAL:HG21	1:H:144:LEU:HG	1.73	0.70
1:D:257:ILE:HG22	1:D:258:SER:HB3	1.74	0.69
1:F:22:GLU:HG2	1:F:27:ARG:HG2	1.75	0.69
1:B:105:GLU:OE2	1:B:109:ARG:NH1	2.24	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:282:THR:HG21	1:F:291:TYR:CD2	2.26	0.69
1:G:92:HIS:CD2	1:G:309:LEU:HD23	2.26	0.69
1:C:45:ASP:HA	1:F:47:ASN:HB3	1.75	0.68
1:C:264:ASN:OD1	1:C:265:ASP:N	2.27	0.68
1:C:92:HIS:CD2	1:C:309:LEU:HD23	2.29	0.68
1:H:20:CYS:HB3	1:H:29:TYR:CD1	2.28	0.68
1:C:239:LEU:HD11	1:C:274:ASN:HB3	1.74	0.68
1:F:239:LEU:HD11	1:F:274:ASN:HB3	1.76	0.67
1:H:143:GLU:HG2	1:H:144:LEU:H	1.58	0.67
1:D:258:SER:OG	1:D:259:GLU:N	2.28	0.67
1:A:257:ILE:O	1:A:258:SER:OG	2.13	0.66
1:H:143:GLU:CD	1:H:144:LEU:H	2.02	0.66
1:G:92:HIS:HD2	1:G:309:LEU:CD2	2.07	0.66
1:A:255:ASP:OD1	1:A:264:ASN:OD1	2.13	0.66
1:G:39:LYS:NZ	1:G:203:GLN:O	2.28	0.66
1:H:265:ASP:OD2	2:H:410:HOH:O	2.13	0.66
1:A:104:ARG:NH2	1:A:138:ALA:O	2.25	0.66
1:A:19:ILE:HB	1:A:30:ILE:HG23	1.78	0.65
1:D:105:GLU:OE1	1:E:176:ARG:NH1	2.28	0.65
1:E:3:SER:HB2	1:F:-1:GLN:HE21	1.62	0.65
1:B:174:THR:HG22	1:B:183:ASN:HD22	1.60	0.64
1:D:257:ILE:HD11	1:D:262:MET:SD	2.37	0.64
1:A:154:GLU:O	1:A:157:SER:HB2	1.98	0.64
1:H:239:LEU:HD11	1:H:274:ASN:HB3	1.80	0.64
1:C:277:GLN:HA	1:C:280:ARG:H	1.61	0.64
1:E:55:SER:HA	1:E:76:PRO:HB3	1.80	0.63
1:H:14:ARG:HA	1:H:19:ILE:HG22	1.79	0.63
1:C:180:ASN:OD1	1:C:181:PHE:N	2.32	0.63
1:D:107:VAL:HG21	1:D:144:LEU:HG	1.80	0.63
1:C:128:LEU:HA	1:C:131:TYR:HD2	1.64	0.63
1:E:92:HIS:HD2	1:E:309:LEU:HD23	1.64	0.62
1:A:127:SER:OG	1:A:158:MET:HE3	1.99	0.62
1:D:52:GLU:HG2	1:H:71:VAL:HG11	1.78	0.62
1:A:239:LEU:HD11	1:A:274:ASN:HB3	1.82	0.62
1:B:255:ASP:OD2	1:B:264:ASN:ND2	2.32	0.62
1:G:55:SER:HA	1:G:76:PRO:HB3	1.82	0.62
1:G:180:ASN:HA	1:G:261:LEU:HD23	1.81	0.62
1:F:289:ASN:O	1:F:290:ARG:HG2	2.00	0.61
1:H:314:VAL:HG13	1:H:315:PHE:N	2.15	0.61
1:B:92:HIS:CD2	1:B:309:LEU:HD23	2.35	0.60
1:C:231:PHE:HD1	1:C:294:MET:HE1	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:124:ARG:NH1	1:F:126:ASP:OD2	2.33	0.60
1:E:239:LEU:HD11	1:E:274:ASN:HB3	1.84	0.60
1:G:243:LEU:HD21	1:G:274:ASN:ND2	2.16	0.60
1:A:159:ILE:O	1:A:160:ASP:C	2.43	0.60
1:A:304:LYS:HE3	1:A:315:PHE:HE1	1.66	0.59
1:C:49:ARG:HG2	1:C:49:ARG:HH11	1.67	0.59
1:G:304:LYS:HE3	1:G:315:PHE:CE1	2.37	0.59
1:D:257:ILE:HG13	1:D:262:MET:HB3	1.82	0.59
1:G:170:ILE:C	1:G:171:GLU:CG	2.75	0.59
1:C:152:ARG:HG3	1:C:156:TYR:CE2	2.38	0.59
1:F:286:LYS:O	1:F:287:LYS:HB2	2.02	0.59
1:H:295:ARG:NH1	2:H:406:HOH:O	2.29	0.59
1:G:58:ARG:NH2	1:G:78:GLU:OE2	2.35	0.59
1:C:277:GLN:O	1:C:281:GLU:HB2	2.02	0.59
1:B:107:VAL:HG21	1:B:144:LEU:HG	1.85	0.58
1:E:145:MET:HE2	1:E:213:LEU:HG	1.85	0.58
1:H:174:THR:HG23	1:H:176:ARG:H	1.67	0.58
1:E:33:GLU:OE1	1:F:304:LYS:NZ	2.35	0.58
1:B:239:LEU:HD11	1:B:274:ASN:HB3	1.84	0.58
1:H:55:SER:HA	1:H:76:PRO:HB3	1.85	0.58
1:C:218:PHE:N	2:C:403:HOH:O	2.36	0.58
1:H:313:GLN:CD	1:H:314:VAL:H	2.12	0.58
1:B:159:ILE:HG23	1:B:241:LEU:HD21	1.84	0.58
1:D:180:ASN:OD1	1:D:183:ASN:HB2	2.04	0.58
1:A:107:VAL:HG21	1:A:144:LEU:HG	1.86	0.58
1:F:67:GLU:OE1	1:F:67:GLU:N	2.36	0.58
1:H:104:ARG:NH2	1:H:138:ALA:O	2.37	0.58
1:E:2:GLU:CG	1:F:-4:LEU:CD2	2.53	0.58
1:G:171:GLU:OE2	1:G:180:ASN:ND2	2.36	0.58
1:G:173:ARG:NH1	1:G:183:ASN:OD1	2.35	0.57
1:A:159:ILE:O	1:A:161:SER:N	2.38	0.57
1:D:2:GLU:HB2	1:D:295:ARG:HH12	1.69	0.57
1:D:71:VAL:HG11	1:H:52:GLU:HG2	1.86	0.57
1:C:52:GLU:HG2	1:F:71:VAL:HG11	1.87	0.57
1:H:143:GLU:OE2	1:H:144:LEU:HB2	2.05	0.57
1:G:243:LEU:HD21	1:G:274:ASN:HD21	1.69	0.57
1:C:19:ILE:HG13	1:C:30:ILE:HB	1.86	0.56
1:A:165:ASP:HB3	1:A:168:PHE:HB2	1.87	0.56
1:A:257:ILE:HG22	1:A:258:SER:N	2.19	0.56
1:B:30:ILE:HD12	1:B:31:PRO:HD2	1.88	0.56
1:F:37:ASP:OD2	1:F:39:LYS:NZ	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:THR:N	1:A:183:ASN:OD1	2.38	0.56
1:G:104:ARG:NH2	1:G:138:ALA:O	2.38	0.56
1:C:123:VAL:HG21	1:C:162:LEU:HD11	1.87	0.56
1:A:233:PRO:HA	1:A:237:ASP:HB2	1.88	0.56
1:A:37:ASP:OD2	1:A:39:LYS:NZ	2.38	0.55
1:C:160:ASP:HA	1:C:163:VAL:HG23	1.88	0.55
1:A:30:ILE:HD12	1:A:31:PRO:HD2	1.89	0.55
1:C:165:ASP:O	1:C:167:ARG:N	2.40	0.55
1:B:85:LEU:HD22	1:B:205:HIS:HB2	1.89	0.55
1:D:120:LYS:HE2	1:D:121:ARG:NH1	2.22	0.55
1:G:170:ILE:C	1:G:171:GLU:HG3	2.31	0.55
1:G:255:ASP:O	1:G:261:LEU:HD12	2.07	0.55
1:H:105:GLU:HB2	1:H:315:PHE:HB2	1.88	0.55
1:C:32:VAL:HG23	1:C:59:ILE:HD11	1.87	0.55
1:H:143:GLU:CD	1:H:144:LEU:HB2	2.31	0.55
1:A:114:MET:HE2	1:A:152:ARG:CZ	2.37	0.55
1:G:171:GLU:OE1	1:G:178:PRO:HG2	2.07	0.54
1:A:5:TYR:OH	1:A:276:GLU:OE2	2.24	0.54
1:A:255:ASP:CG	1:A:264:ASN:OD1	2.51	0.54
1:C:231:PHE:CD1	1:C:294:MET:HE1	2.43	0.54
1:F:263:LEU:O	1:F:268:LYS:NZ	2.36	0.54
1:F:166:GLU:HA	1:F:169:ARG:HG3	1.90	0.54
1:C:2:GLU:HG2	1:C:34:ASN:O	2.08	0.54
1:B:219:ARG:NH1	1:G:81:ASN:HD21	2.07	0.53
1:H:100:MET:HE1	1:H:140:ASN:O	2.08	0.53
1:C:165:ASP:CB	1:C:168:PHE:HB2	2.39	0.53
1:E:3:SER:HB2	1:F:-1:GLN:NE2	2.22	0.53
1:E:264:ASN:OD1	1:E:266:GLU:HG3	2.09	0.53
1:C:153:GLU:HA	1:C:156:TYR:CD2	2.39	0.53
1:D:166:GLU:HA	1:D:169:ARG:HG2	1.89	0.53
1:H:141:VAL:C	1:H:143:GLU:OE1	2.52	0.53
1:E:23:THR:HG23	1:E:24:GLU:O	2.09	0.53
1:G:174:THR:HG23	1:G:176:ARG:H	1.73	0.53
1:A:282:THR:HG23	1:A:292:VAL:N	2.24	0.53
1:D:257:ILE:HG22	1:D:258:SER:CB	2.39	0.53
1:G:217:ASN:OD1	1:G:219:ARG:HB3	2.08	0.53
1:H:78:GLU:OE2	2:H:415:HOH:O	2.18	0.52
1:H:140:ASN:ND2	1:H:142:SER:OG	2.42	0.52
1:D:62:HIS:CG	1:D:195:THR:HG23	2.43	0.52
1:H:166:GLU:HG2	1:H:169:ARG:HD2	1.92	0.52
1:E:2:GLU:HG2	1:F:-4:LEU:CD2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:166:GLU:C	1:H:168:PHE:H	2.18	0.52
1:A:257:ILE:CG2	1:A:258:SER:N	2.73	0.52
1:B:114:MET:HE2	1:B:152:ARG:CZ	2.38	0.52
1:C:152:ARG:O	1:C:155:TYR:HB3	2.09	0.52
1:C:214:HIS:ND1	1:C:226:ASP:OD1	2.29	0.52
1:F:114:MET:HE2	1:F:152:ARG:CZ	2.40	0.52
1:D:120:LYS:HE2	1:D:121:ARG:HH12	1.73	0.52
1:F:19:ILE:HG13	1:F:30:ILE:HG13	1.90	0.52
1:G:58:ARG:HB3	1:G:79:TYR:HA	1.92	0.52
1:H:105:GLU:OE2	1:H:109:ARG:NH2	2.42	0.52
1:H:145:MET:HE2	1:H:213:LEU:HG	1.92	0.52
1:A:277:GLN:HA	1:A:280:ARG:HG2	1.92	0.51
1:B:3:SER:OG	1:B:37:ASP:OD1	2.19	0.51
1:G:114:MET:HE3	1:G:155:TYR:HB2	1.91	0.51
1:H:143:GLU:CG	1:H:144:LEU:N	2.67	0.51
1:B:71:VAL:HG11	1:G:52:GLU:HG2	1.92	0.51
1:B:158:MET:O	1:B:161:SER:OG	2.26	0.51
1:H:253:HIS:HA	1:H:264:ASN:HD21	1.76	0.51
1:A:255:ASP:CG	1:A:264:ASN:HD21	2.09	0.51
1:B:131:TYR:OH	1:B:154:GLU:OE1	2.25	0.51
1:A:157:SER:C	1:A:159:ILE:H	2.18	0.51
1:C:37:ASP:OD2	1:C:39:LYS:NZ	2.40	0.51
1:A:14:ARG:HD3	1:A:53:PHE:CD1	2.46	0.51
1:B:37:ASP:HA	1:B:59:ILE:HG23	1.93	0.50
1:F:140:ASN:ND2	2:F:416:HOH:O	2.43	0.50
1:F:107:VAL:HG21	1:F:144:LEU:HG	1.92	0.50
1:D:36:MET:HE1	1:D:200:LEU:HD23	1.93	0.50
1:G:171:GLU:OE2	1:G:180:ASN:OD1	2.30	0.50
1:E:58:ARG:HB3	1:E:79:TYR:HA	1.92	0.50
1:F:133:LYS:O	1:F:137:GLU:HG2	2.12	0.50
1:A:238:ARG:NH2	1:A:281:GLU:OE1	2.44	0.50
1:H:143:GLU:HG2	1:H:144:LEU:N	2.24	0.50
1:C:92:HIS:ND1	1:C:98:LYS:HD3	2.27	0.50
1:C:71:VAL:HG11	1:F:52:GLU:HG2	1.93	0.50
1:C:85:LEU:HD22	1:C:205:HIS:HB2	1.93	0.50
1:A:231:PHE:HD1	1:A:294:MET:HE1	1.75	0.50
1:C:116:ASP:HA	1:C:119:LYS:HB2	1.93	0.50
1:G:121:ARG:HD3	1:G:162:LEU:HD13	1.92	0.50
1:H:104:ARG:NH1	1:H:135:ALA:O	2.45	0.50
1:A:127:SER:C	1:A:129:THR:H	2.20	0.49
1:F:310:ILE:HD11	1:F:312:GLU:OE2	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:ILE:C	1:A:161:SER:N	2.69	0.49
1:F:308:HIS:HA	1:F:313:GLN:HB3	1.93	0.49
1:G:156:TYR:CE2	1:G:173:ARG:HD3	2.47	0.49
1:G:124:ARG:HH12	1:H:169:ARG:HH12	1.61	0.49
1:H:301:GLU:OE2	1:H:304:LYS:NZ	2.45	0.49
1:D:2:GLU:HB2	1:D:295:ARG:NH1	2.27	0.49
1:A:52:GLU:HG2	1:E:71:VAL:HG11	1.95	0.48
1:H:155:TYR:O	1:H:159:ILE:HG12	2.12	0.48
1:B:170:ILE:HG23	1:B:183:ASN:OD1	2.13	0.48
1:G:170:ILE:O	1:G:170:ILE:HG22	2.14	0.48
1:B:263:LEU:O	1:B:268:LYS:HE3	2.13	0.48
1:D:100:MET:CE	1:D:140:ASN:HA	2.39	0.48
1:G:92:HIS:CD2	1:G:309:LEU:CD2	2.91	0.48
1:G:171:GLU:CD	1:G:180:ASN:HD21	2.20	0.48
1:A:157:SER:C	1:A:159:ILE:N	2.72	0.48
1:A:165:ASP:O	1:A:167:ARG:N	2.46	0.48
1:C:170:ILE:HG12	1:C:182:ALA:HB3	1.96	0.48
1:H:20:CYS:HB3	1:H:29:TYR:CE1	2.48	0.48
1:A:218:PHE:HB3	2:E:407:HOH:O	2.13	0.48
1:C:233:PRO:HA	1:C:237:ASP:HB2	1.96	0.48
1:A:128:LEU:HA	1:A:131:TYR:HD2	1.79	0.48
1:F:284:PHE:HD2	1:F:284:PHE:N	2.11	0.48
1:F:284:PHE:N	1:F:284:PHE:CD2	2.81	0.48
1:H:301:GLU:CD	1:H:304:LYS:NZ	2.72	0.48
1:D:41:PHE:HZ	1:D:192:LEU:HD21	1.79	0.48
1:H:301:GLU:CD	1:H:304:LYS:HZ3	2.22	0.48
1:C:123:VAL:HG12	1:C:124:ARG:H	1.79	0.47
1:H:121:ARG:HD3	1:H:162:LEU:HD13	1.95	0.47
1:H:266:GLU:O	1:H:269:SER:OG	2.32	0.47
1:H:12:LEU:HA	1:H:20:CYS:O	2.14	0.47
1:A:5:TYR:CE2	1:A:39:LYS:HD2	2.49	0.47
1:D:69:TYR:OH	1:H:77:ARG:HD3	2.13	0.47
1:B:81:ASN:HD22	1:G:220:ARG:HH21	1.63	0.47
1:H:253:HIS:HA	1:H:264:ASN:ND2	2.28	0.47
1:C:170:ILE:HG23	1:C:183:ASN:ND2	2.24	0.47
1:G:276:GLU:N	1:G:276:GLU:OE1	2.47	0.47
1:H:257:ILE:O	1:H:260:GLY:N	2.47	0.47
1:E:133:LYS:O	1:E:137:GLU:HG3	2.14	0.47
1:E:284:PHE:HB2	1:E:291:TYR:CE1	2.50	0.47
1:F:55:SER:HA	1:F:76:PRO:HB3	1.96	0.47
1:H:155:TYR:CE1	1:H:159:ILE:HD11	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:ARG:NH1	1:E:70:TYR:O	2.42	0.47
1:C:5:TYR:CE2	1:C:39:LYS:HD2	2.50	0.47
1:F:52:GLU:O	1:F:56:GLN:HG3	2.15	0.47
1:H:20:CYS:HB3	1:H:29:TYR:HD1	1.80	0.47
1:H:105:GLU:OE2	1:H:109:ARG:NE	2.47	0.47
1:C:114:MET:HE2	1:C:152:ARG:CZ	2.44	0.47
1:A:5:TYR:HD2	1:A:41:PHE:HE2	1.63	0.46
1:E:107:VAL:HG21	1:E:144:LEU:HG	1.97	0.46
1:G:19:ILE:HG13	1:G:30:ILE:O	2.15	0.46
1:A:92:HIS:CE1	1:A:98:LYS:HD3	2.50	0.46
1:D:217:ASN:HB2	1:D:219:ARG:H	1.80	0.46
1:B:239:LEU:O	1:B:243:LEU:HB2	2.15	0.46
1:B:253:HIS:O	1:B:263:LEU:HD12	2.15	0.46
1:G:69:TYR:HD1	1:G:199:SER:HB3	1.80	0.46
1:B:178:PRO:HB3	1:B:183:ASN:O	2.15	0.46
1:D:114:MET:HE2	1:D:152:ARG:CZ	2.45	0.46
1:H:145:MET:CE	1:H:213:LEU:HG	2.46	0.46
1:B:39:LYS:HE3	1:B:195:THR:OG1	2.16	0.46
1:E:100:MET:HE2	1:E:104:ARG:NH2	2.31	0.46
1:G:145:MET:HE2	1:G:213:LEU:HG	1.97	0.46
1:D:5:TYR:CZ	1:D:39:LYS:HD2	2.51	0.46
1:A:281:GLU:O	1:A:293:SER:HA	2.16	0.45
1:C:104:ARG:NH2	1:C:138:ALA:O	2.47	0.45
1:H:8:SER:O	1:H:21:LEU:HD21	2.15	0.45
1:H:264:ASN:O	1:H:268:LYS:HG3	2.16	0.45
1:E:145:MET:HE1	1:E:214:HIS:CE1	2.52	0.45
1:H:170:ILE:HG12	1:H:182:ALA:HB1	1.97	0.45
1:D:64:PHE:CE2	1:D:70:TYR:HB2	2.51	0.45
1:F:132:LYS:O	1:F:136:GLU:HG3	2.16	0.45
1:D:100:MET:HB2	1:D:100:MET:HE2	1.61	0.45
1:E:277:GLN:OE1	1:E:280:ARG:NH2	2.44	0.45
1:B:168:PHE:O	1:B:182:ALA:HB2	2.17	0.45
1:B:174:THR:H	1:B:183:ASN:ND2	2.15	0.45
1:B:219:ARG:CZ	1:G:81:ASN:HD21	2.30	0.45
1:E:2:GLU:HG2	1:F:-4:LEU:HD23	1.99	0.45
1:F:109:ARG:NH1	1:F:301:GLU:OE1	2.50	0.45
1:A:173:ARG:H	1:A:173:ARG:HG3	1.32	0.45
1:E:14:ARG:HA	1:E:19:ILE:HG22	1.97	0.45
1:E:16:ALA:HB3	1:E:18:THR:HG22	1.97	0.45
1:E:140:ASN:OD1	1:E:141:VAL:N	2.50	0.45
1:F:121:ARG:HD3	1:F:162:LEU:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ARG:HD2	1:A:130:ARG:HA	1.61	0.45
1:B:92:HIS:CE1	1:B:98:LYS:HD3	2.52	0.45
1:D:37:ASP:HA	1:D:59:ILE:HG23	1.97	0.45
1:H:156:TYR:CE2	1:H:173:ARG:HD3	2.51	0.45
1:A:22:GLU:HG2	1:A:27:ARG:HG2	1.99	0.45
1:C:109:ARG:HH22	1:C:301:GLU:CD	2.22	0.45
1:C:165:ASP:O	1:C:168:PHE:N	2.50	0.45
1:C:182:ALA:O	1:C:185:LEU:N	2.50	0.45
1:F:286:LYS:C	1:F:288:LEU:H	2.23	0.45
1:A:132:LYS:O	1:A:136:GLU:HG3	2.17	0.44
1:A:69:TYR:OH	1:E:77:ARG:HD3	2.16	0.44
1:G:107:VAL:HG21	1:G:144:LEU:HG	2.00	0.44
1:B:178:PRO:HB3	1:B:183:ASN:C	2.42	0.44
1:A:118:LEU:HD11	1:A:155:TYR:HD1	1.81	0.44
1:E:92:HIS:CD2	1:E:309:LEU:CD2	2.96	0.44
1:C:88:LYS:HA	1:C:88:LYS:HD3	1.80	0.44
1:D:5:TYR:CE2	1:D:39:LYS:HD2	2.52	0.44
1:F:104:ARG:NH1	1:F:135:ALA:O	2.50	0.44
1:F:5:TYR:OH	1:F:299:LYS:NZ	2.50	0.44
1:C:49:ARG:HG2	1:C:49:ARG:NH1	2.31	0.44
1:C:152:ARG:NH2	1:C:229:GLU:O	2.44	0.44
1:B:17:ASN:OD1	1:B:18:THR:HG22	2.17	0.44
1:F:109:ARG:NH2	2:F:420:HOH:O	2.51	0.44
1:H:140:ASN:O	1:H:143:GLU:CD	2.61	0.44
1:A:45:ASP:OD1	1:E:47:ASN:HB3	2.18	0.43
1:C:5:TYR:HD2	1:C:41:PHE:HE2	1.66	0.43
1:E:-2:PHE:HB3	1:F:-2:PHE:HB3	1.99	0.43
1:G:97:GLU:O	1:G:101:LEU:HD13	2.18	0.43
1:A:157:SER:O	1:A:160:ASP:N	2.51	0.43
1:A:180:ASN:OD1	1:A:181:PHE:N	2.51	0.43
1:D:62:HIS:CD2	1:D:195:THR:HG23	2.53	0.43
1:H:37:ASP:HA	1:H:59:ILE:HG23	2.00	0.43
1:A:235:VAL:O	1:A:239:LEU:HB2	2.19	0.43
1:B:248:GLN:O	1:B:253:HIS:HE1	2.00	0.43
1:D:2:GLU:O	1:D:36:MET:N	2.41	0.43
1:F:253:HIS:O	1:F:263:LEU:HD12	2.18	0.43
1:B:126:ASP:O	1:B:127:SER:OG	2.27	0.43
1:D:36:MET:HE2	1:D:199:SER:HB3	2.01	0.43
1:A:265:ASP:O	1:A:268:LYS:HB2	2.18	0.43
1:A:276:GLU:O	1:A:279:LEU:N	2.42	0.43
1:B:300:MET:HE3	1:B:300:MET:HB2	1.86	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:17:ASN:CB	1:F:287:LYS:HE2	2.49	0.43
1:E:156:TYR:CE2	1:E:173:ARG:HD3	2.54	0.43
1:A:174:THR:HG21	1:A:179:LYS:N	2.34	0.43
1:D:107:VAL:HG22	1:D:148:GLU:HG3	2.01	0.43
1:H:48:LYS:HE3	1:H:52:GLU:OE2	2.19	0.43
1:B:5:TYR:CZ	1:B:39:LYS:HD2	2.54	0.43
1:D:133:LYS:O	1:D:137:GLU:HG3	2.19	0.43
1:E:2:GLU:HG3	1:F:-2:PHE:CD1	2.54	0.43
1:H:100:MET:HB2	1:H:100:MET:HE2	1.83	0.43
1:E:100:MET:HE2	1:E:104:ARG:HH21	1.84	0.42
1:B:48:LYS:HE3	1:B:52:GLU:OE2	2.19	0.42
1:B:132:LYS:O	1:B:136:GLU:HG3	2.19	0.42
1:F:264:ASN:OD1	1:F:266:GLU:HG2	2.19	0.42
1:H:95:ASN:OD1	1:H:97:GLU:HG2	2.18	0.42
1:H:149:GLY:O	1:H:153:GLU:HG3	2.19	0.42
1:A:255:ASP:HB2	1:A:262:MET:HG2	2.02	0.42
1:B:92:HIS:ND1	1:B:98:LYS:HD3	2.34	0.42
1:D:233:PRO:HA	1:D:237:ASP:HB2	2.00	0.42
1:E:289:ASN:C	1:E:290:ARG:HG3	2.45	0.42
1:H:141:VAL:HG12	1:H:141:VAL:O	2.20	0.42
1:G:13:LYS:HD2	1:G:13:LYS:HA	1.46	0.42
1:A:88:LYS:HA	1:A:88:LYS:HD3	1.79	0.42
1:E:145:MET:CE	1:E:213:LEU:HG	2.48	0.42
1:A:154:GLU:OE2	1:A:158:MET:HE2	2.10	0.42
1:B:107:VAL:HG22	1:B:148:GLU:HG3	2.02	0.42
1:E:19:ILE:HG23	1:E:53:PHE:CE2	2.55	0.42
1:C:55:SER:HA	1:C:76:PRO:HB3	2.01	0.42
1:E:40:VAL:HG21	1:E:61:ILE:HD11	2.01	0.42
1:B:186:ILE:HD13	1:B:186:ILE:HA	1.80	0.42
1:C:183:ASN:HD22	1:C:183:ASN:H	1.68	0.42
1:F:233:PRO:HA	1:F:237:ASP:HB2	2.01	0.42
1:G:168:PHE:O	1:G:182:ALA:HB2	2.19	0.42
1:H:30:ILE:O	1:H:30:ILE:HG13	2.20	0.42
1:E:92:HIS:ND1	1:E:98:LYS:HD3	2.35	0.42
1:F:1:MET:HE1	1:F:80:LEU:HD13	2.01	0.42
1:F:59:ILE:HA	1:F:60:PRO:HD3	1.92	0.42
1:H:113:ASN:ND2	1:H:230:LEU:O	2.43	0.42
1:B:104:ARG:NH2	1:B:138:ALA:O	2.53	0.42
1:D:108:SER:OG	1:D:132:LYS:NZ	2.53	0.42
1:C:123:VAL:HG12	1:C:124:ARG:N	2.33	0.41
1:D:283:VAL:O	1:D:291:TYR:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:121:ARG:HD3	1:E:162:LEU:HD13	2.02	0.41
1:A:304:LYS:HE3	1:A:315:PHE:CE1	2.51	0.41
1:C:246:THR:HG23	1:C:248:GLN:HG3	2.02	0.41
1:G:4:VAL:HG11	1:G:30:ILE:HD13	2.02	0.41
1:H:144:LEU:HA	1:H:147:ILE:HB	2.02	0.41
1:A:166:GLU:O	1:A:169:ARG:HG3	2.21	0.41
1:A:264:ASN:HB3	1:A:266:GLU:HG3	2.02	0.41
1:B:88:LYS:HD3	1:B:88:LYS:HA	1.81	0.41
1:G:171:GLU:HB2	1:G:172:LYS:H	1.30	0.41
1:D:281:GLU:O	1:D:293:SER:HA	2.20	0.41
1:E:36:MET:HE3	1:E:36:MET:HB2	1.91	0.41
1:G:145:MET:CE	1:G:213:LEU:HG	2.50	0.41
1:A:37:ASP:HA	1:A:59:ILE:HG23	2.02	0.41
1:D:219:ARG:NE	1:H:81:ASN:OD1	2.43	0.41
1:F:144:LEU:HD12	1:F:144:LEU:HA	1.92	0.41
1:A:39:LYS:HE3	1:A:195:THR:OG1	2.20	0.41
1:D:162:LEU:HD23	1:D:162:LEU:HA	1.78	0.41
1:E:19:ILE:H	1:E:19:ILE:HG13	1.75	0.41
1:H:214:HIS:CE1	1:H:225:LEU:HB3	2.56	0.41
1:H:313:GLN:NE2	1:H:314:VAL:HB	2.35	0.41
1:B:77:ARG:HD3	1:G:70:TYR:O	2.21	0.41
1:D:300:MET:HE3	1:D:300:MET:HB2	1.93	0.41
1:G:19:ILE:HD11	1:G:30:ILE:HB	2.03	0.41
1:G:243:LEU:HD23	1:G:243:LEU:HA	1.83	0.41
1:H:174:THR:O	1:H:175:ARG:HB2	2.21	0.41
1:B:36:MET:O	1:B:60:PRO:HD2	2.21	0.41
1:B:74:PHE:HA	1:G:74:PHE:HA	2.02	0.41
1:B:86:ILE:HG13	1:G:209:ARG:CZ	2.51	0.41
1:E:182:ALA:O	1:E:186:ILE:HG12	2.21	0.41
1:H:64:PHE:CE2	1:H:70:TYR:HB2	2.55	0.41
1:H:314:VAL:CG1	1:H:315:PHE:N	2.84	0.41
1:A:91:GLU:HA	1:A:94:ILE:HG12	2.03	0.41
1:F:100:MET:HE2	1:F:100:MET:HB2	1.94	0.41
1:G:182:ALA:O	1:G:186:ILE:HG12	2.21	0.41
1:C:32:VAL:HG23	1:C:59:ILE:CD1	2.50	0.40
1:F:76:PRO:HD2	1:F:79:TYR:HB3	2.02	0.40
1:H:153:GLU:HG2	1:H:173:ARG:O	2.21	0.40
1:H:160:ASP:HB3	1:H:169:ARG:NH2	2.35	0.40
1:A:127:SER:O	1:A:128:LEU:HB2	2.20	0.40
1:B:249:ILE:HA	1:B:253:HIS:HD1	1.86	0.40
1:C:104:ARG:NH1	1:C:136:GLU:HA	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:177:PRO:HA	1:G:178:PRO:HD2	1.81	0.40
1:D:86:ILE:HG13	1:H:209:ARG:NE	2.37	0.40
1:D:297:LEU:HD23	1:D:297:LEU:HA	1.90	0.40
1:E:14:ARG:HE	1:E:53:PHE:HD1	1.70	0.40
1:A:249:ILE:HA	1:A:253:HIS:ND1	2.36	0.40
1:B:104:ARG:HA	1:B:144:LEU:HD21	2.03	0.40
1:B:163:VAL:HG12	1:B:164:SER:N	2.36	0.40
1:A:161:SER:O	1:A:162:LEU:CB	2.70	0.40
1:C:104:ARG:HG2	1:C:144:LEU:HD11	2.02	0.40

All (1) 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:280:ARG:NH1	1:G:23:THR:O[1_455]	2.13	0.07

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	307/326 (94%)	284 (92%)	18 (6%)	5 (2%)	7	20
1	B	305/326 (94%)	294 (96%)	10 (3%)	1 (0%)	36	60
1	C	290/326 (89%)	275 (95%)	14 (5%)	1 (0%)	36	60
1	D	307/326 (94%)	298 (97%)	9 (3%)	0	100	100
1	E	315/326 (97%)	305 (97%)	10 (3%)	0	100	100
1	F	309/326 (95%)	298 (96%)	11 (4%)	0	100	100
1	G	297/326 (91%)	287 (97%)	10 (3%)	0	100	100
1	H	293/326 (90%)	280 (96%)	12 (4%)	1 (0%)	36	60
All	All	2423/2608 (93%)	2321 (96%)	94 (4%)	8 (0%)	36	60

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	158	MET
1	C	166	GLU
1	A	157	SER
1	A	162	LEU
1	A	265	ASP
1	A	266	GLU
1	B	17	ASN
1	H	143	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	261/300 (87%)	256 (98%)	5 (2%)	50	77
1	B	261/300 (87%)	257 (98%)	4 (2%)	57	81
1	C	253/300 (84%)	244 (96%)	9 (4%)	31	60
1	D	264/300 (88%)	255 (97%)	9 (3%)	32	62
1	E	277/300 (92%)	268 (97%)	9 (3%)	34	64
1	F	273/300 (91%)	265 (97%)	8 (3%)	37	67
1	G	263/300 (88%)	252 (96%)	11 (4%)	26	55
1	H	261/300 (87%)	254 (97%)	7 (3%)	39	69
All	All	2113/2400 (88%)	2051 (97%)	62 (3%)	37	67

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	91	GLU
1	A	161	SER
1	A	171	GLU
1	A	273	LYS
1	B	19	ILE

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Mol	Chain	Res	Type
1	B	154	GLU
1	B	166	GLU
1	B	251	GLU
1	C	30	ILE
1	C	32	VAL
1	C	109	ARG
1	C	154	GLU
1	C	166	GLU
1	C	272	VAL
1	C	281	GLU
1	C	283	VAL
1	C	314	VAL
1	D	30	ILE
1	D	122	LYS
1	D	158	MET
1	D	159	ILE
1	D	163	VAL
1	D	164	SER
1	D	166	GLU
1	D	283	VAL
1	D	314	VAL
1	E	19	ILE
1	E	30	ILE
1	E	33	GLU
1	E	124	ARG
1	E	166	GLU
1	E	262	MET
1	E	266	GLU
1	E	304	LYS
1	E	312	GLU
1	F	20	CYS
1	F	47	ASN
1	F	49	ARG
1	F	262	MET
1	F	281	GLU
1	F	284	PHE
1	F	304	LYS
1	F	314	VAL
1	G	13	LYS
1	G	28	LYS
1	G	35	VAL
1	G	77	ARG

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Mol	Chain	Res	Type
1	G	88	LYS
1	G	126	ASP
1	G	136	GLU
1	G	166	GLU
1	G	171	GLU
1	G	257	ILE
1	G	265	ASP
1	H	22	GLU
1	H	119	LYS
1	H	144	LEU
1	H	170	ILE
1	H	261	LEU
1	H	281	GLU
1	H	314	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	HIS
1	A	112	GLN
1	A	248	GLN
1	B	183	ASN
1	C	183	ASN
1	D	92	HIS
1	E	-1	GLN
1	E	248	GLN
1	F	-1	GLN
1	F	92	HIS
1	G	81	ASN
1	G	92	HIS
1	G	274	ASN
1	H	96	GLN
1	H	112	GLN
1	H	248	GLN
1	H	313	GLN

5.3.3 RNA ⓘ

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 ⓘ

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	311/326 (95%)	-0.07	6 (1%) 66 63	32, 81, 127, 154	0
1	B	309/326 (94%)	-0.09	1 (0%) 90 89	37, 81, 126, 142	0
1	C	298/326 (91%)	0.02	12 (4%) 42 38	35, 85, 133, 162	0
1	D	311/326 (95%)	-0.11	3 (0%) 79 78	36, 79, 126, 141	0
1	E	319/326 (97%)	-0.31	2 (0%) 85 85	32, 59, 100, 124	0
1	F	314/326 (96%)	-0.22	6 (1%) 66 63	30, 66, 108, 135	1 (0%)
1	G	304/326 (93%)	-0.13	6 (1%) 65 62	36, 72, 108, 135	1 (0%)
1	H	300/326 (92%)	-0.12	6 (2%) 65 62	36, 76, 109, 134	1 (0%)
All	All	2466/2608 (94%)	-0.13	42 (1%) 69 67	30, 74, 122, 162	3 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	283	VAL	4.7
1	H	81	ASN	4.1
1	F	-4	LEU	3.6
1	C	283	VAL	3.6
1	C	319	GLU	3.4
1	C	254	PHE	3.3
1	G	291	TYR	3.2
1	A	314	VAL	3.1
1	H	315	PHE	3.0
1	C	282	THR	2.9
1	A	258	SER	2.8
1	F	257	ILE	2.7
1	C	168	PHE	2.7
1	H	141	VAL	2.7
1	G	29	TYR	2.6
1	D	314	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	313	GLN	2.6
1	G	257	ILE	2.5
1	H	174	THR	2.5
1	F	291	TYR	2.5
1	D	256	GLU	2.4
1	C	280	ARG	2.4
1	H	144	LEU	2.3
1	C	100	MET	2.3
1	G	314	VAL	2.3
1	C	291	TYR	2.3
1	E	18	THR	2.3
1	G	171	GLU	2.2
1	F	-3	TYR	2.2
1	G	31	PRO	2.2
1	A	257	ILE	2.2
1	F	14	ARG	2.1
1	C	264	ASN	2.1
1	D	291	TYR	2.1
1	C	314	VAL	2.1
1	F	284	PHE	2.1
1	A	256	GLU	2.0
1	A	313	GLN	2.0
1	C	17	ASN	2.0
1	E	17	ASN	2.0
1	H	261	LEU	2.0
1	C	69	TYR	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 oligosaccharides 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.