



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:04 AM UTC

PDB ID : 3PU2 / pdb_00003pu2
Title : Crystal Structure of the Q3J4M4_RHOS4 protein from Rhodobacter sphaeroides. Northeast Structural Genomics Consortium Target RhR263.
Authors : Vorobiev, S.; Chen, Y.; Seetharaman, J.; Sahdev, S.; Xiao, R.; Ciccocanti, C.; Wang, D.; Everett, J.K.; Nair, R.; Acton, T.B.; Rost, B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2010-12-03
Resolution : 2.61 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

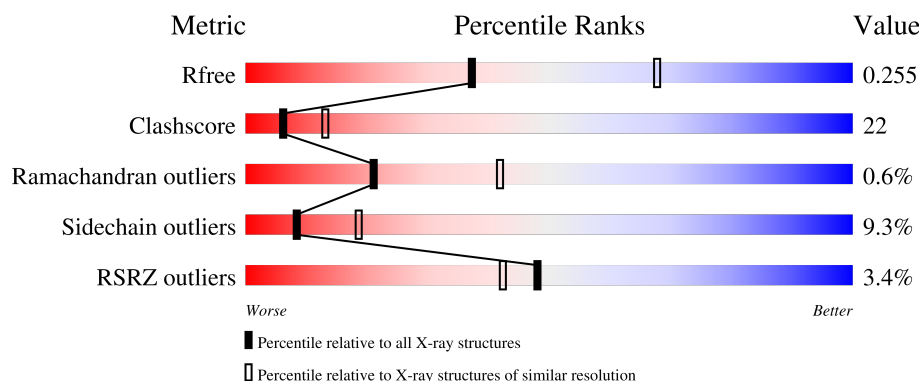
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.61 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	164	
1	B	164	
1	C	164	
1	D	164	

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Mol	Chain	Length	Quality of chain
1	E	164	4% 52% 33% 7% 8%
1	F	164	3% 52% 32% 5% 10%
1	G	164	2% 59% 27% 5% 9%
1	H	164	3% 54% 31% 5% 10%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 9371 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 uncharacterized protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	Se	0	1	0
			1167	744	203	213	1	6			
1	B	153	Total	C	N	O	S	Se	0	1	0
			1181	751	205	218	1	6			
1	C	152	Total	C	N	O	S	Se	0	1	0
			1182	752	204	219	1	6			
1	D	149	Total	C	N	O	S	Se	0	1	0
			1157	739	201	210	1	6			
1	E	151	Total	C	N	O	S	Se	0	0	0
			1169	744	203	216	1	5			
1	F	148	Total	C	N	O	S	Se	0	0	0
			1142	730	197	209	1	5			
1	G	149	Total	C	N	O	S	Se	0	0	0
			1153	734	203	210	1	5			
1	H	148	Total	C	N	O	S	Se	0	0	0
			1148	733	200	209	1	5			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	LEU	-	expression tag	UNP Q3J4M4
A	170	GLU	-	expression tag	UNP Q3J4M4
A	171	HIS	-	expression tag	UNP Q3J4M4
A	172	HIS	-	expression tag	UNP Q3J4M4
A	173	HIS	-	expression tag	UNP Q3J4M4
A	174	HIS	-	expression tag	UNP Q3J4M4
A	175	HIS	-	expression tag	UNP Q3J4M4
A	176	HIS	-	expression tag	UNP Q3J4M4
B	169	LEU	-	expression tag	UNP Q3J4M4
B	170	GLU	-	expression tag	UNP Q3J4M4
B	171	HIS	-	expression tag	UNP Q3J4M4
B	172	HIS	-	expression tag	UNP Q3J4M4
B	173	HIS	-	expression tag	UNP Q3J4M4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	174	HIS	-	expression tag	UNP Q3J4M4
B	175	HIS	-	expression tag	UNP Q3J4M4
B	176	HIS	-	expression tag	UNP Q3J4M4
C	169	LEU	-	expression tag	UNP Q3J4M4
C	170	GLU	-	expression tag	UNP Q3J4M4
C	171	HIS	-	expression tag	UNP Q3J4M4
C	172	HIS	-	expression tag	UNP Q3J4M4
C	173	HIS	-	expression tag	UNP Q3J4M4
C	174	HIS	-	expression tag	UNP Q3J4M4
C	175	HIS	-	expression tag	UNP Q3J4M4
C	176	HIS	-	expression tag	UNP Q3J4M4
D	169	LEU	-	expression tag	UNP Q3J4M4
D	170	GLU	-	expression tag	UNP Q3J4M4
D	171	HIS	-	expression tag	UNP Q3J4M4
D	172	HIS	-	expression tag	UNP Q3J4M4
D	173	HIS	-	expression tag	UNP Q3J4M4
D	174	HIS	-	expression tag	UNP Q3J4M4
D	175	HIS	-	expression tag	UNP Q3J4M4
D	176	HIS	-	expression tag	UNP Q3J4M4
E	169	LEU	-	expression tag	UNP Q3J4M4
E	170	GLU	-	expression tag	UNP Q3J4M4
E	171	HIS	-	expression tag	UNP Q3J4M4
E	172	HIS	-	expression tag	UNP Q3J4M4
E	173	HIS	-	expression tag	UNP Q3J4M4
E	174	HIS	-	expression tag	UNP Q3J4M4
E	175	HIS	-	expression tag	UNP Q3J4M4
E	176	HIS	-	expression tag	UNP Q3J4M4
F	169	LEU	-	expression tag	UNP Q3J4M4
F	170	GLU	-	expression tag	UNP Q3J4M4
F	171	HIS	-	expression tag	UNP Q3J4M4
F	172	HIS	-	expression tag	UNP Q3J4M4
F	173	HIS	-	expression tag	UNP Q3J4M4
F	174	HIS	-	expression tag	UNP Q3J4M4
F	175	HIS	-	expression tag	UNP Q3J4M4
F	176	HIS	-	expression tag	UNP Q3J4M4
G	169	LEU	-	expression tag	UNP Q3J4M4
G	170	GLU	-	expression tag	UNP Q3J4M4
G	171	HIS	-	expression tag	UNP Q3J4M4
G	172	HIS	-	expression tag	UNP Q3J4M4
G	173	HIS	-	expression tag	UNP Q3J4M4
G	174	HIS	-	expression tag	UNP Q3J4M4
G	175	HIS	-	expression tag	UNP Q3J4M4

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Chain	Residue	Modelled	Actual	Comment	Reference
G	176	HIS	-	expression tag	UNP Q3J4M4
H	169	LEU	-	expression tag	UNP Q3J4M4
H	170	GLU	-	expression tag	UNP Q3J4M4
H	171	HIS	-	expression tag	UNP Q3J4M4
H	172	HIS	-	expression tag	UNP Q3J4M4
H	173	HIS	-	expression tag	UNP Q3J4M4
H	174	HIS	-	expression tag	UNP Q3J4M4
H	175	HIS	-	expression tag	UNP Q3J4M4
H	176	HIS	-	expression tag	UNP Q3J4M4

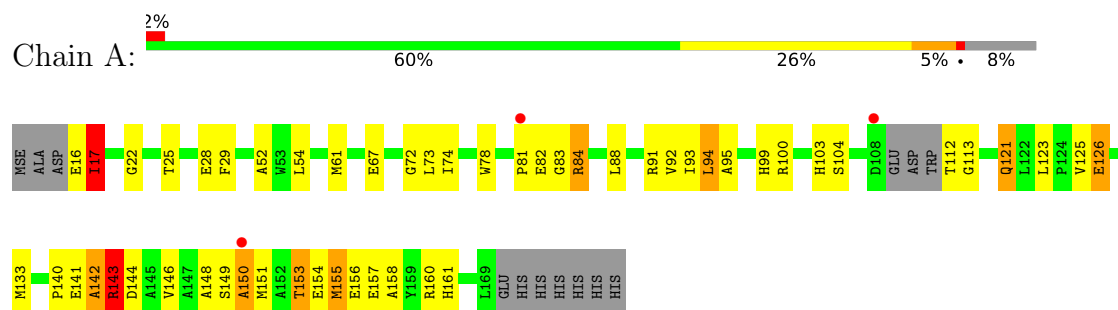
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	14	Total O 14 14	0	0
2	B	14	Total O 14 14	0	0
2	C	10	Total O 10 10	0	0
2	D	8	Total O 8 8	0	0
2	E	7	Total O 7 7	0	0
2	F	10	Total O 10 10	0	0
2	G	4	Total O 4 4	0	0
2	H	5	Total O 5 5	0	0

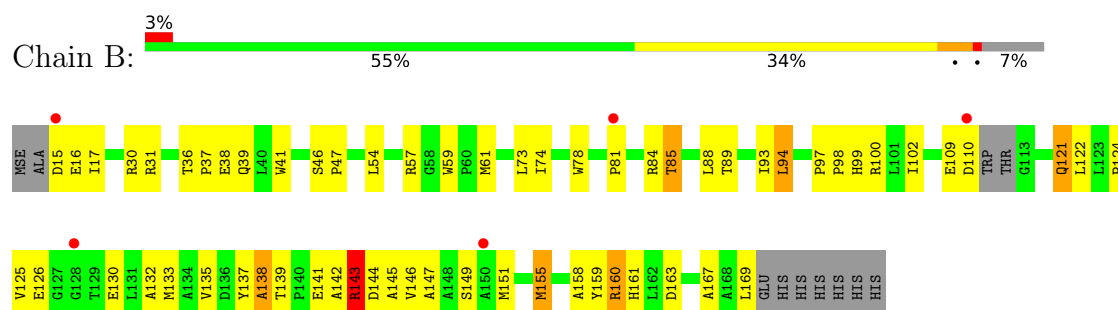
3 Residue-property plots [i](#)

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.

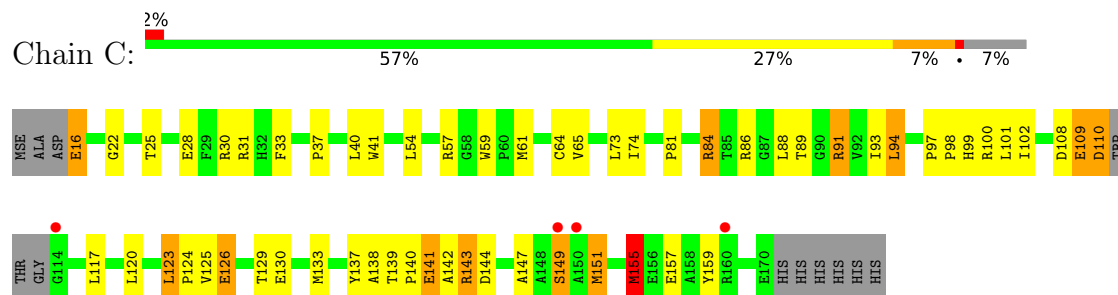
- Molecule 1: uncharacterized protein



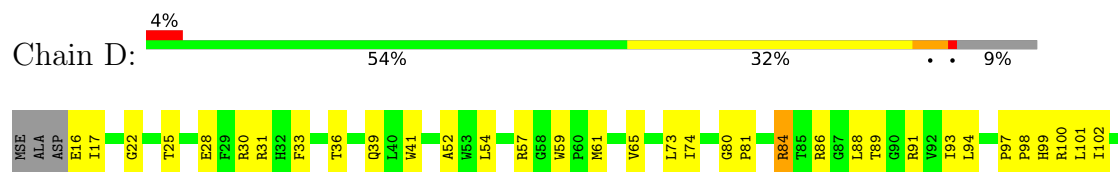
- Molecule 1: uncharacterized protein



- Molecule 1: uncharacterized protein

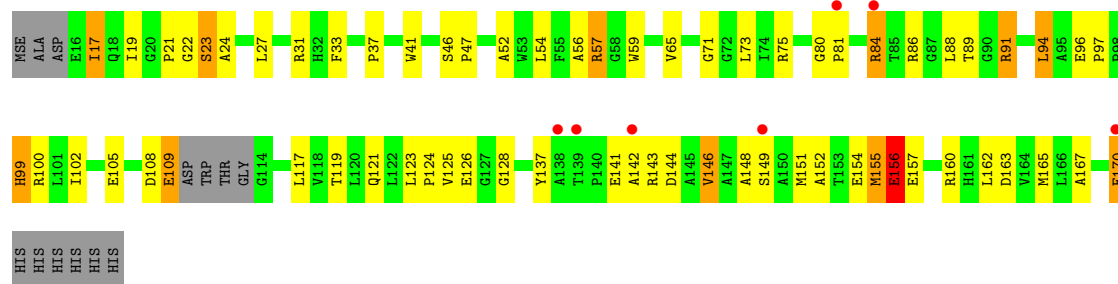


- Molecule 1: uncharacterized protein

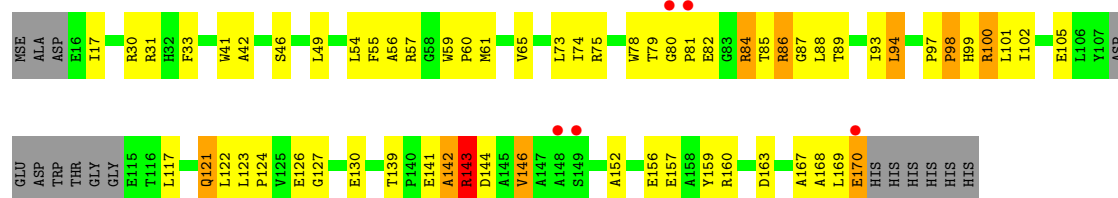




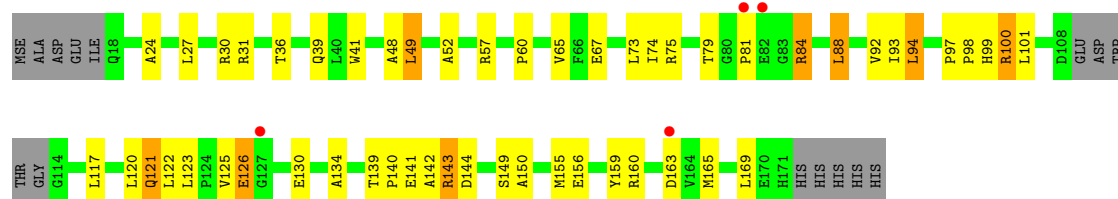
- Molecule 1: uncharacterized protein



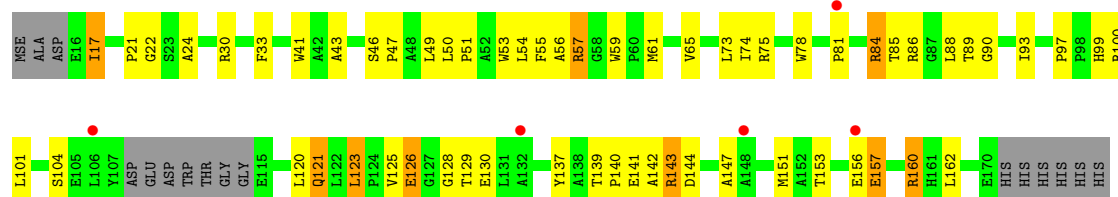
- Molecule 1: uncharacterized protein



- Molecule 1: uncharacterized protein



- Molecule 1: uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	75.21Å 35.86Å 230.47Å 90.00° 98.98° 90.00°	Depositor
Resolution (Å)	29.10 – 2.61 29.10 – 2.61	Depositor EDS
% Data completeness (in resolution range)	93.3 (29.10-2.61) 95.6 (29.10-2.61)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.80 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.6 _289	Depositor
R, R_{free}	0.230 , 0.263 0.225 , 0.255	Depositor DCC
R_{free} test set	1833 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.679	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9371	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 70.39 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.1464e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.67	2/1194 (0.2%)	0.99	4/1619 (0.2%)
1	B	0.64	4/1208 (0.3%)	1.01	3/1638 (0.2%)
1	C	0.72	4/1209 (0.3%)	0.99	4/1639 (0.2%)
1	D	0.66	2/1184 (0.2%)	1.01	6/1606 (0.4%)
1	E	0.50	0/1193	1.08	8/1618 (0.5%)
1	F	0.49	0/1166	0.97	7/1583 (0.4%)
1	G	0.47	0/1178	0.97	3/1598 (0.2%)
1	H	0.56	0/1172	1.07	7/1590 (0.4%)
All	All	0.60	12/9504 (0.1%)	1.01	42/12891 (0.3%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	155[A]	MSE	SE-CE	-6.48	1.76	1.95
1	A	155[B]	MSE	SE-CE	-6.48	1.76	1.95
1	B	155[A]	MSE	SE-CE	-5.78	1.78	1.95
1	B	155[B]	MSE	SE-CE	-5.78	1.78	1.95
1	D	155[A]	MSE	SE-CE	-5.53	1.78	1.95
1	D	155[B]	MSE	SE-CE	-5.53	1.78	1.95
1	C	155[A]	MSE	SE-CE	-5.22	1.79	1.95
1	C	155[B]	MSE	SE-CE	-5.22	1.79	1.95
1	B	155[A]	MSE	CG-SE	-5.18	1.79	1.95
1	B	155[B]	MSE	CG-SE	-5.18	1.79	1.95
1	C	155[A]	MSE	CG-SE	-5.16	1.79	1.95
1	C	155[B]	MSE	CG-SE	-5.16	1.79	1.95

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	99	HIS	N-CA-C	11.57	123.97	111.36
1	D	108	ASP	N-CA-C	9.74	124.88	110.17
1	E	146	VAL	N-CA-C	8.96	120.77	111.00
1	E	146	VAL	CB-CA-C	-7.34	102.41	112.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	142	ALA	N-CA-C	-7.11	103.27	112.23
1	C	142	ALA	N-CA-C	-6.98	103.69	111.71
1	A	17	ILE	N-CA-C	6.64	117.60	108.84
1	A	143	ARG	N-CA-C	-6.59	104.50	112.54
1	H	49	LEU	N-CA-C	6.54	118.49	111.36
1	D	80	GLY	CA-C-N	6.46	126.49	119.90
1	D	80	GLY	C-N-CA	6.46	126.49	119.90
1	H	160	ARG	N-CA-C	-6.42	104.28	111.28
1	F	143	ARG	N-CA-C	-6.39	104.97	112.89
1	D	107	TYR	N-CA-C	-6.29	97.40	110.80
1	G	149	SER	N-CA-C	6.24	117.77	110.41
1	E	156	GLU	N-CA-C	-6.22	104.58	111.36
1	F	127	GLY	N-CA-C	-6.21	98.46	113.18
1	E	167	ALA	N-CA-C	-6.15	104.58	111.28
1	A	142	ALA	N-CA-C	-6.12	104.52	112.23
1	B	85	THR	N-CA-C	5.89	118.43	109.95
1	C	65	VAL	N-CA-C	-5.87	99.17	107.75
1	F	87	GLY	N-CA-C	5.73	119.47	111.10
1	G	142	ALA	N-CA-C	-5.71	105.20	111.82
1	B	142	ALA	N-CA-C	-5.66	105.10	112.23
1	H	151	MSE	N-CA-C	-5.66	105.49	112.90
1	B	143	ARG	N-CA-C	-5.62	106.38	113.18
1	H	147	ALA	N-CA-C	5.61	120.57	111.37
1	H	142	ALA	N-CA-C	-5.56	105.23	112.23
1	E	17	ILE	N-CA-C	5.49	116.13	108.89
1	A	83	GLY	N-CA-C	5.48	121.21	114.69
1	F	100	ARG	N-CA-C	5.35	117.89	108.75
1	F	167	ALA	N-CA-C	-5.32	105.13	111.03
1	G	100	ARG	N-CA-C	5.22	117.67	108.75
1	D	65	VAL	N-CA-C	-5.20	100.88	108.17
1	F	142	ALA	N-CA-C	-5.19	105.80	111.82
1	E	128	GLY	N-CA-C	-5.10	101.09	113.18
1	D	100	ARG	N-CA-C	5.09	117.46	108.75
1	F	146	VAL	N-CA-C	5.08	116.97	111.58
1	C	100	ARG	N-CA-C	5.07	117.42	108.75
1	E	100	ARG	N-CA-C	5.01	116.54	108.32
1	C	149	SER	N-CA-C	5.01	121.46	110.80
1	H	85	THR	N-CA-C	5.00	118.11	110.36

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	1167	0	1157	50	0
1	B	1181	0	1160	67	0
1	C	1182	0	1163	55	0
1	D	1157	0	1146	52	0
1	E	1169	0	1150	56	0
1	F	1142	0	1126	58	0
1	G	1153	0	1134	44	0
1	H	1148	0	1137	49	0
2	A	14	0	0	3	0
2	B	14	0	0	0	0
2	C	10	0	0	2	0
2	D	8	0	0	0	0
2	E	7	0	0	2	0
2	F	10	0	0	1	0
2	G	4	0	0	0	0
2	H	5	0	0	0	0
All	All	9371	0	9173	414	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 22.

All (414) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:MSE:HE2	1:B:155[B]:MSE:HE1	1.28	1.13
1:B:133:MSE:CE	1:B:155[B]:MSE:CE	2.29	1.10
1:B:133:MSE:HE2	1:B:155[B]:MSE:CE	1.82	1.08
1:F:17:ILE:HG22	1:F:31:ARG:HB2	1.37	1.03
1:B:155[A]:MSE:HE3	1:B:159:TYR:CZ	1.94	1.01
1:B:133:MSE:CE	1:B:155[B]:MSE:HE2	1.93	0.97
1:D:150:ALA:O	1:D:154:GLU:HG2	1.66	0.96
1:D:146:VAL:CG1	1:D:151:MSE:HE3	1.96	0.95
1:D:143:ARG:HD2	1:D:143:ARG:O	1.71	0.90
1:E:57:ARG:HD2	1:E:154:GLU:OE2	1.72	0.90
1:B:133:MSE:HE1	1:B:155[B]:MSE:CE	2.03	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:57:ARG:HG3	1:E:57:ARG:HH11	1.40	0.87
1:A:94:LEU:HD12	1:A:94:LEU:C	2.00	0.87
1:D:150:ALA:O	1:D:154:GLU:CG	2.23	0.85
1:D:151:MSE:SE	1:D:155[B]:MSE:CE	2.75	0.84
1:F:82:GLU:HB2	1:F:84:ARG:HD3	1.59	0.84
1:C:133:MSE:HE1	1:C:155[B]:MSE:HE2	1.60	0.83
1:B:143:ARG:HD2	1:B:143:ARG:O	1.79	0.82
1:A:84:ARG:HH11	1:A:84:ARG:HB2	1.46	0.80
1:D:146:VAL:HG11	1:D:151:MSE:HE3	1.63	0.78
1:E:59:TRP:CZ2	1:E:81:PRO:HG3	2.19	0.78
1:B:133:MSE:HE1	1:B:155[B]:MSE:HE2	1.62	0.78
1:E:57:ARG:HH11	1:E:57:ARG:CG	1.96	0.77
1:D:81:PRO:O	1:D:84:ARG:HG2	1.85	0.77
1:E:81:PRO:O	1:E:84:ARG:HG2	1.83	0.77
1:E:17:ILE:HG22	1:E:156:GLU:HG2	1.67	0.76
1:B:155[A]:MSE:HE3	1:B:159:TYR:OH	1.86	0.75
1:F:60:PRO:HD2	1:F:79:THR:O	1.88	0.74
1:H:84:ARG:HB2	1:H:84:ARG:NH1	2.02	0.74
1:E:151:MSE:O	1:E:155:MSE:HB2	1.89	0.73
1:H:84:ARG:HB2	1:H:84:ARG:HH11	1.53	0.73
1:C:84:ARG:HB2	1:C:84:ARG:NH1	2.04	0.72
1:E:137:TYR:CE1	1:E:146:VAL:HG21	2.25	0.72
1:C:157:GLU:HG3	2:C:225:HOH:O	1.91	0.71
1:G:30:ARG:HD2	1:G:130:GLU:OE2	1.92	0.70
1:D:81:PRO:O	1:D:84:ARG:CG	2.40	0.69
1:E:137:TYR:OH	1:E:151:MSE:HE1	1.92	0.69
1:C:81:PRO:O	1:C:84:ARG:HG2	1.93	0.69
1:C:84:ARG:HH11	1:C:84:ARG:CB	2.05	0.69
1:F:56:ALA:HB1	1:F:157:GLU:OE1	1.92	0.69
1:C:147:ALA:C	1:C:149:SER:H	1.98	0.69
1:B:155[A]:MSE:CE	1:B:159:TYR:OH	2.41	0.68
1:E:143:ARG:O	1:E:143:ARG:NH1	2.26	0.68
1:D:149:SER:C	1:D:151:MSE:H	1.99	0.68
1:H:143:ARG:HD2	1:H:143:ARG:O	1.93	0.68
1:B:125:VAL:HG12	1:B:126:GLU:N	2.09	0.68
1:E:27:LEU:HD21	1:E:143:ARG:CZ	2.23	0.67
1:A:94:LEU:HD21	1:A:100:ARG:NH2	2.08	0.67
1:B:73:LEU:HD12	1:B:74:ILE:H	1.60	0.67
1:C:84:ARG:HB2	1:C:84:ARG:HH11	1.60	0.66
1:A:146:VAL:HG12	1:A:151:MSE:HE3	1.77	0.66
1:F:54:LEU:O	1:F:61:MSE:HE3	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:141:GLU:O	1:F:144:ASP:HB2	1.96	0.66
1:D:151:MSE:SE	1:D:155[B]:MSE:HE1	2.45	0.66
1:A:149:SER:C	1:A:151:MSE:H	2.04	0.66
1:H:73:LEU:HD12	1:H:74:ILE:H	1.59	0.66
1:E:73:LEU:HB2	1:E:91:ARG:NH1	2.11	0.65
1:H:73:LEU:HD12	1:H:74:ILE:N	2.12	0.65
1:A:22:GLY:HA2	1:A:28:GLU:OE2	1.97	0.65
1:A:16:GLU:HG2	1:A:17:ILE:H	1.61	0.65
1:A:146:VAL:CG1	1:A:151:MSE:HE3	2.27	0.65
1:H:139:THR:OG1	1:H:141:GLU:HG2	1.97	0.64
1:D:156:GLU:HG3	1:D:160:ARG:NH2	2.13	0.64
1:E:108:ASP:C	1:E:109:GLU:HG2	2.23	0.64
1:A:84:ARG:HH11	1:A:84:ARG:CB	2.11	0.63
1:B:94:LEU:C	1:B:94:LEU:HD12	2.22	0.63
1:F:94:LEU:HD12	1:F:94:LEU:C	2.24	0.63
1:C:73:LEU:HD12	1:C:74:ILE:H	1.63	0.63
1:D:149:SER:O	1:D:151:MSE:N	2.25	0.63
1:C:141:GLU:N	1:C:141:GLU:OE1	2.32	0.63
1:H:81:PRO:O	1:H:84:ARG:HG2	1.98	0.63
1:B:36:THR:OG1	1:B:39:GLN:HG3	1.98	0.62
1:E:102:ILE:HG12	1:E:119:THR:HG23	1.81	0.62
1:A:73:LEU:HD12	1:A:74:ILE:H	1.63	0.62
1:C:30:ARG:HD2	1:C:130:GLU:OE2	2.00	0.62
1:G:88:LEU:C	1:G:88:LEU:HD12	2.23	0.62
1:C:151:MSE:SE	1:C:155[B]:MSE:HE1	2.49	0.62
1:H:33:PHE:CE1	1:H:162:LEU:HD23	2.35	0.62
1:G:73:LEU:HD12	1:G:74:ILE:H	1.65	0.62
1:F:88:LEU:HD22	1:F:105:GLU:HG2	1.80	0.61
1:C:37:PRO:HG3	1:C:124:PRO:HG3	1.82	0.61
1:C:151:MSE:SE	1:C:155[B]:MSE:CE	2.98	0.61
1:C:123:LEU:N	1:C:123:LEU:HD23	2.15	0.61
1:B:15:ASP:CG	1:B:16:GLU:H	2.09	0.61
1:D:143:ARG:HD2	1:D:143:ARG:C	2.25	0.61
1:C:37:PRO:HD3	1:C:129:THR:OG1	2.00	0.61
1:F:86:ARG:NH2	2:F:230:HOH:O	2.32	0.61
1:C:139:THR:OG1	1:C:141:GLU:HG2	2.01	0.60
1:A:94:LEU:C	1:A:94:LEU:CD1	2.72	0.60
1:D:41:TRP:CE3	1:D:98:PRO:O	2.54	0.60
1:B:155[A]:MSE:HE3	1:B:159:TYR:CE1	2.36	0.60
1:D:146:VAL:HG11	1:D:151:MSE:CE	2.31	0.60
1:E:146:VAL:CG1	1:E:151:MSE:HE3	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:30:ARG:HD2	1:F:130:GLU:OE2	2.01	0.60
1:C:125:VAL:HG12	1:C:126:GLU:N	2.16	0.60
1:C:25:THR:HG22	1:C:140:PRO:HA	1.84	0.59
1:C:147:ALA:C	1:C:149:SER:N	2.60	0.59
1:G:84:ARG:HH11	1:G:84:ARG:HB2	1.68	0.59
1:B:41:TRP:CE3	1:B:98:PRO:O	2.56	0.58
1:C:31:ARG:CZ	2:C:205:HOH:O	2.51	0.58
1:C:101:LEU:HD23	1:C:120:LEU:HD12	1.85	0.58
1:D:149:SER:C	1:D:151:MSE:N	2.60	0.58
1:B:138:ALA:HB1	1:E:96:GLU:OE1	2.03	0.58
1:D:30:ARG:HD2	1:D:130:GLU:OE2	2.03	0.58
1:E:71:GLY:O	1:E:91:ARG:HD2	2.03	0.58
1:F:65:VAL:HB	1:F:75:ARG:HB3	1.85	0.58
1:G:156:GLU:HG3	1:G:160:ARG:NH2	2.19	0.58
1:H:84:ARG:HH11	1:H:84:ARG:CB	2.15	0.58
1:F:82:GLU:OE2	1:F:82:GLU:HA	2.03	0.58
1:B:88:LEU:C	1:B:88:LEU:HD12	2.28	0.58
1:G:84:ARG:HB2	1:G:84:ARG:NH1	2.18	0.58
1:A:88:LEU:HD12	1:A:88:LEU:C	2.29	0.58
1:C:16:GLU:HB2	1:C:31:ARG:NH1	2.19	0.58
1:C:73:LEU:HD12	1:C:74:ILE:N	2.18	0.58
1:H:54:LEU:O	1:H:61:MSE:HE3	2.03	0.58
1:D:94:LEU:C	1:D:94:LEU:HD12	2.29	0.57
1:F:81:PRO:O	1:F:82:GLU:HB2	2.03	0.57
1:A:81:PRO:O	1:A:82:GLU:HB2	2.02	0.57
1:D:140:PRO:HD2	1:D:141:GLU:OE1	2.05	0.57
1:G:100:ARG:HD3	1:G:121:GLN:NE2	2.19	0.57
1:G:41:TRP:CE3	1:G:98:PRO:O	2.57	0.57
1:E:65:VAL:HB	1:E:75:ARG:HB3	1.86	0.57
1:F:73:LEU:HD12	1:F:74:ILE:H	1.70	0.57
1:C:94:LEU:HD12	1:C:94:LEU:C	2.30	0.57
1:A:52:ALA:HB2	1:D:52:ALA:HB2	1.87	0.57
1:F:31:ARG:NH2	1:F:163:ASP:OD2	2.38	0.56
1:A:29:PHE:CD2	1:A:155[B]:MSE:HE1	2.40	0.56
1:E:73:LEU:HB2	1:E:91:ARG:HH12	1.70	0.56
1:E:31:ARG:NH2	1:E:163:ASP:OD2	2.38	0.56
1:D:88:LEU:C	1:D:88:LEU:HD12	2.31	0.56
1:G:73:LEU:HD12	1:G:74:ILE:N	2.21	0.56
1:H:137:TYR:CD2	1:H:143:ARG:HG2	2.40	0.56
1:H:157:GLU:HA	1:H:160:ARG:NH1	2.20	0.56
1:B:93:ILE:HG21	1:E:117:LEU:HD13	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:23:SER:OG	1:E:24:ALA:N	2.39	0.56
1:F:55:PHE:HB2	1:F:59:TRP:O	2.06	0.56
1:A:112:THR:HG22	1:A:113:GLY:N	2.21	0.56
1:H:129:THR:HG22	1:H:130:GLU:N	2.21	0.56
1:B:41:TRP:HE3	1:B:98:PRO:O	1.89	0.55
1:F:86:ARG:O	1:F:86:ARG:HG3	2.01	0.55
1:G:49:LEU:CD1	1:G:49:LEU:N	2.69	0.55
1:A:150:ALA:O	1:A:154:GLU:HG3	2.05	0.55
1:E:27:LEU:HD21	1:E:143:ARG:NH2	2.20	0.55
1:E:88:LEU:C	1:E:88:LEU:HD12	2.31	0.55
1:B:125:VAL:CG1	1:B:126:GLU:N	2.69	0.55
1:E:57:ARG:CG	1:E:57:ARG:NH1	2.64	0.55
1:D:135:VAL:O	1:D:137:TYR:CE2	2.59	0.55
1:B:59:TRP:CZ2	1:B:81:PRO:HD2	2.42	0.55
1:A:29:PHE:CE2	1:A:155[B]:MSE:HE1	2.42	0.55
1:A:157:GLU:HA	1:A:160:ARG:NH1	2.21	0.55
1:E:109:GLU:N	2:E:243:HOH:O	2.40	0.54
1:D:54:LEU:O	1:D:61:MSE:HE3	2.08	0.54
1:H:88:LEU:HD12	1:H:88:LEU:C	2.32	0.54
1:E:94:LEU:C	1:E:94:LEU:HD12	2.33	0.54
1:F:54:LEU:HD23	1:F:61:MSE:HE1	1.89	0.54
1:G:84:ARG:HH11	1:G:84:ARG:CB	2.20	0.54
1:C:125:VAL:CG1	1:C:126:GLU:N	2.70	0.54
1:D:16:GLU:OE1	1:D:16:GLU:N	2.40	0.54
1:B:100:ARG:NH2	1:E:137:TYR:O	2.41	0.54
1:D:150:ALA:O	1:D:154:GLU:HG3	2.07	0.54
1:F:88:LEU:HD12	1:F:88:LEU:C	2.32	0.54
1:F:156:GLU:HG3	1:F:160:ARG:NH2	2.22	0.54
1:D:102:ILE:HG12	1:D:119:THR:HG23	1.88	0.54
1:F:86:ARG:O	1:F:86:ARG:CG	2.56	0.54
1:F:142:ALA:O	1:F:146:VAL:HG23	2.07	0.54
1:G:100:ARG:HD3	1:G:121:GLN:HE22	1.71	0.54
1:H:100:ARG:HD3	1:H:121:GLN:HE22	1.73	0.54
1:E:59:TRP:CD2	1:E:80:GLY:HA2	2.43	0.53
1:D:93:ILE:HD12	1:G:93:ILE:HG23	1.89	0.53
1:E:170:GLU:C	1:E:170:GLU:OE1	2.52	0.53
1:H:24:ALA:O	1:H:140:PRO:HA	2.08	0.53
1:E:57:ARG:CD	1:E:154:GLU:OE2	2.50	0.53
1:C:41:TRP:CE3	1:C:98:PRO:O	2.62	0.53
1:C:93:ILE:HD12	1:F:93:ILE:HG23	1.89	0.53
1:G:94:LEU:C	1:G:94:LEU:HD12	2.33	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:TYR:N	1:D:137:TYR:CD2	2.76	0.53
1:F:82:GLU:HB2	1:F:84:ARG:CD	2.35	0.53
1:C:57:ARG:HG2	1:C:57:ARG:HH11	1.73	0.53
1:F:57:ARG:O	1:F:59:TRP:HD1	1.92	0.53
1:F:79:THR:OG1	1:F:85:THR:HG22	2.09	0.53
1:H:17:ILE:N	1:H:156:GLU:OE1	2.37	0.53
1:H:57:ARG:NH1	1:H:57:ARG:HG3	2.24	0.53
1:D:57:ARG:HH11	1:D:57:ARG:HG2	1.74	0.53
1:H:57:ARG:HG3	1:H:57:ARG:HH11	1.74	0.52
1:F:80:GLY:O	1:F:81:PRO:C	2.52	0.52
1:H:90:GLY:HA3	1:H:104:SER:O	2.09	0.52
1:A:148:ALA:O	1:F:124:PRO:HG2	2.09	0.52
1:A:149:SER:O	1:A:151:MSE:N	2.41	0.52
1:A:84:ARG:HB2	1:A:84:ARG:NH1	2.20	0.52
1:A:93:ILE:HD12	1:H:93:ILE:HG23	1.92	0.52
1:B:109:GLU:O	1:B:110:ASP:CB	2.58	0.52
1:B:138:ALA:CB	1:E:96:GLU:OE1	2.57	0.52
1:C:33:PHE:HZ	1:C:159:TYR:CD1	2.28	0.52
1:F:139:THR:OG1	1:F:141:GLU:HG2	2.10	0.52
1:G:60:PRO:HD2	1:G:79:THR:O	2.09	0.52
1:A:141:GLU:O	1:A:144:ASP:HB2	2.09	0.52
1:A:157:GLU:HG3	2:A:217:HOH:O	2.10	0.52
1:A:153:THR:HG22	1:A:154:GLU:N	2.24	0.51
1:B:16:GLU:HG3	1:B:17:ILE:H	1.75	0.51
1:G:31:ARG:NH2	1:G:163:ASP:OD2	2.37	0.51
1:A:73:LEU:HD12	1:A:74:ILE:N	2.24	0.51
1:G:41:TRP:HE3	1:G:98:PRO:O	1.94	0.51
1:G:165:MSE:HE3	1:G:169:LEU:HD11	1.92	0.51
1:B:16:GLU:HG3	1:B:17:ILE:N	2.25	0.51
1:D:36:THR:OG1	1:D:39:GLN:HG3	2.10	0.51
1:D:142:ALA:O	1:D:146:VAL:HG23	2.09	0.51
1:B:30:ARG:HD2	1:B:130:GLU:OE2	2.10	0.51
1:B:139:THR:HG22	1:E:99:HIS:CD2	2.45	0.51
1:C:88:LEU:HD12	1:C:88:LEU:C	2.36	0.51
1:H:65:VAL:HB	1:H:75:ARG:HB3	1.91	0.51
1:E:94:LEU:HD13	1:E:96:GLU:OE2	2.11	0.51
1:B:100:ARG:HD3	1:B:121:GLN:NE2	2.25	0.51
1:D:31:ARG:NH2	1:D:163:ASP:OD2	2.39	0.51
1:E:125:VAL:HG12	1:E:126:GLU:N	2.25	0.51
1:G:156:GLU:HG3	1:G:160:ARG:HH22	1.74	0.51
1:C:151:MSE:SE	1:C:155[B]:MSE:HE3	2.61	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:17:ILE:CG2	1:E:156:GLU:HG2	2.37	0.51
1:G:81:PRO:O	1:G:84:ARG:HG2	2.10	0.51
1:F:17:ILE:HG23	1:F:156:GLU:OE1	2.12	0.50
1:C:137:TYR:CD2	1:C:143:ARG:HG2	2.46	0.50
1:E:157:GLU:HA	1:E:160:ARG:NH1	2.26	0.50
1:F:73:LEU:HD12	1:F:74:ILE:N	2.26	0.50
1:G:143:ARG:HD2	1:G:143:ARG:O	2.10	0.50
1:H:156:GLU:OE2	1:H:156:GLU:O	2.30	0.50
1:H:50:LEU:HB2	1:H:51:PRO:HD3	1.92	0.50
1:E:57:ARG:HB2	1:E:59:TRP:HD1	1.77	0.50
1:E:108:ASP:O	1:E:109:GLU:HG2	2.11	0.50
1:E:33:PHE:CZ	1:E:162:LEU:HD23	2.47	0.50
1:F:49:LEU:N	1:F:49:LEU:HD12	2.27	0.50
1:C:54:LEU:O	1:C:61:MSE:HE3	2.11	0.49
1:B:61:MSE:HE2	1:B:78:TRP:CZ2	2.47	0.49
1:D:106:LEU:HG	1:D:107:TYR:O	2.11	0.49
1:G:141:GLU:O	1:G:144:ASP:HB2	2.12	0.49
1:A:133:MSE:CE	1:A:155[B]:MSE:HE2	2.42	0.49
1:A:158:ALA:O	1:A:161:HIS:HB2	2.11	0.49
1:C:133:MSE:CE	1:C:155[B]:MSE:HE2	2.36	0.49
1:D:143:ARG:O	1:D:143:ARG:CD	2.53	0.49
1:G:24:ALA:O	1:G:140:PRO:HA	2.13	0.49
1:A:112:THR:HG22	1:A:113:GLY:H	1.77	0.49
1:B:38:GLU:OE1	1:B:38:GLU:N	2.46	0.49
1:B:160:ARG:NH2	1:B:161:HIS:CD2	2.81	0.49
1:D:151:MSE:SE	1:D:155[B]:MSE:HE3	2.61	0.49
1:H:53:TRP:CD2	1:H:162:LEU:HD13	2.48	0.49
1:B:141:GLU:O	1:B:144:ASP:HB2	2.12	0.49
1:E:37:PRO:HG3	1:E:124:PRO:HG3	1.95	0.49
1:A:133:MSE:HE2	1:A:155[B]:MSE:HE2	1.95	0.49
1:E:146:VAL:HG11	1:E:151:MSE:CE	2.43	0.49
1:A:25:THR:CG2	1:A:140:PRO:HG3	2.44	0.48
1:E:146:VAL:HG11	1:E:151:MSE:HE3	1.95	0.48
1:F:33:PHE:HZ	1:F:159:TYR:CD1	2.32	0.48
1:H:30:ARG:HD2	1:H:130:GLU:OE2	2.12	0.48
1:A:142:ALA:O	1:A:146:VAL:HG23	2.13	0.48
1:C:84:ARG:NH1	1:C:84:ARG:CB	2.71	0.48
1:D:125:VAL:HG12	1:D:126:GLU:N	2.29	0.48
1:A:143:ARG:NH1	2:A:248:HOH:O	2.45	0.48
1:B:15:ASP:CG	1:B:16:GLU:N	2.72	0.48
1:B:94:LEU:CD2	1:B:102:ILE:HD12	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:TYR:OH	1:B:146:VAL:HG11	2.13	0.48
1:G:101:LEU:HD23	1:G:120:LEU:HD12	1.95	0.48
1:H:125:VAL:HG12	1:H:126:GLU:N	2.27	0.48
1:G:65:VAL:HB	1:G:75:ARG:HB3	1.96	0.48
1:A:125:VAL:HG12	1:A:126:GLU:N	2.28	0.48
1:B:143:ARG:HD2	1:B:143:ARG:C	2.38	0.48
1:A:156:GLU:HG3	1:A:160:ARG:NH2	2.29	0.47
1:D:25:THR:HG22	1:D:140:PRO:HA	1.95	0.47
1:D:30:ARG:HA	1:D:131:LEU:O	2.14	0.47
1:A:149:SER:C	1:A:151:MSE:N	2.69	0.47
1:B:31:ARG:NH2	1:B:163:ASP:OD2	2.48	0.47
1:C:59:TRP:CE2	1:C:86:ARG:NH2	2.82	0.47
1:G:49:LEU:N	1:G:49:LEU:HD12	2.29	0.47
1:B:121:GLN:HB3	1:B:132:ALA:HB3	1.95	0.47
1:B:73:LEU:HD12	1:B:74:ILE:N	2.27	0.47
1:D:135:VAL:O	1:D:137:TYR:HE2	1.97	0.47
1:D:22:GLY:HA2	1:D:28:GLU:OE2	2.14	0.47
1:A:67:GLU:O	1:A:92:VAL:HG21	2.15	0.47
1:C:137:TYR:OH	1:C:151:MSE:HE1	2.14	0.47
1:A:94:LEU:HD12	1:A:94:LEU:O	2.13	0.47
1:G:141:GLU:HA	1:G:144:ASP:OD1	2.14	0.47
1:H:78:TRP:HE3	1:H:86:ARG:O	1.98	0.47
1:H:141:GLU:O	1:H:144:ASP:HB2	2.15	0.47
1:F:143:ARG:HD2	1:F:143:ARG:O	2.15	0.47
1:C:108:ASP:C	1:C:109:GLU:HG3	2.39	0.46
1:E:102:ILE:CG1	1:E:119:THR:HG23	2.44	0.46
1:H:123:LEU:H	1:H:123:LEU:HD23	1.80	0.46
1:B:94:LEU:HG	1:B:102:ILE:HD12	1.97	0.46
1:H:33:PHE:CZ	1:H:162:LEU:HD23	2.50	0.46
1:H:143:ARG:HD2	1:H:143:ARG:C	2.41	0.46
1:C:25:THR:HG22	1:C:140:PRO:CA	2.45	0.46
1:F:81:PRO:O	1:F:82:GLU:CB	2.63	0.46
1:E:86:ARG:NH2	2:E:207:HOH:O	2.37	0.46
1:F:100:ARG:HD3	1:F:121:GLN:NE2	2.30	0.46
1:B:143:ARG:O	1:B:147:ALA:HB2	2.16	0.46
1:G:57:ARG:HH11	1:G:57:ARG:HG2	1.81	0.46
1:G:139:THR:OG1	1:G:141:GLU:HG2	2.16	0.46
1:B:97:PRO:HA	1:B:98:PRO:HA	1.36	0.46
1:B:121:GLN:C	1:B:122:LEU:HD12	2.40	0.46
1:C:16:GLU:HB2	1:C:31:ARG:HH11	1.81	0.46
1:C:22:GLY:HA2	1:C:28:GLU:OE2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:88:LEU:HD12	1:G:88:LEU:O	2.16	0.45
1:A:94:LEU:HD12	1:A:95:ALA:N	2.31	0.45
1:B:141:GLU:HA	1:B:144:ASP:OD1	2.16	0.45
1:F:143:ARG:HD2	1:F:143:ARG:C	2.41	0.45
1:F:31:ARG:HE	1:F:159:TYR:HB3	1.82	0.45
1:D:73:LEU:HB2	1:D:91:ARG:NH1	2.31	0.45
1:G:97:PRO:HA	1:G:98:PRO:HA	1.47	0.45
1:B:109:GLU:O	1:B:110:ASP:CG	2.59	0.45
1:C:97:PRO:HA	1:C:98:PRO:HA	1.55	0.45
1:C:102:ILE:HD13	1:F:117:LEU:HD22	1.99	0.45
1:H:21:PRO:HG2	1:H:22:GLY:H	1.81	0.45
1:F:17:ILE:HD11	1:F:152:ALA:O	2.17	0.44
1:B:84:ARG:HG2	1:B:85:THR:N	2.31	0.44
1:G:67:GLU:O	1:G:92:VAL:HG21	2.18	0.44
1:F:41:TRP:CE3	1:F:98:PRO:O	2.71	0.44
1:B:94:LEU:HD21	1:B:100:ARG:NH2	2.32	0.44
1:C:117:LEU:HD22	1:F:102:ILE:HD13	1.99	0.44
1:D:17:ILE:HG23	1:D:17:ILE:O	2.18	0.44
1:H:43:ALA:HB1	1:H:162:LEU:HD11	2.00	0.44
1:B:167:ALA:C	1:B:169:LEU:N	2.75	0.44
1:H:100:ARG:HD3	1:H:121:GLN:NE2	2.32	0.44
1:C:73:LEU:HB2	1:C:91:ARG:NH1	2.32	0.44
1:G:155:MSE:HG3	1:G:159:TYR:CE2	2.52	0.44
1:A:143:ARG:HD2	1:A:143:ARG:O	2.17	0.44
1:B:167:ALA:C	1:B:169:LEU:H	2.26	0.44
1:C:141:GLU:HA	1:C:144:ASP:HB2	2.00	0.44
1:G:143:ARG:O	1:G:143:ARG:CD	2.65	0.44
1:D:146:VAL:HG12	1:D:151:MSE:HE3	1.93	0.43
1:E:160:ARG:O	1:E:163:ASP:HB2	2.18	0.43
1:G:36:THR:OG1	1:G:39:GLN:HG3	2.19	0.43
1:E:52:ALA:HB3	1:E:165:MSE:SE	2.69	0.43
1:E:56:ALA:O	1:E:57:ARG:HB2	2.19	0.43
1:F:41:TRP:CE2	1:F:97:PRO:HB3	2.54	0.43
1:F:78:TRP:HE3	1:F:86:ARG:O	2.02	0.43
1:A:93:ILE:HG23	1:H:93:ILE:HD12	1.99	0.43
1:C:138:ALA:O	1:F:121:GLN:NE2	2.51	0.43
1:H:56:ALA:O	1:H:57:ARG:C	2.60	0.43
1:H:123:LEU:HD23	1:H:123:LEU:N	2.34	0.43
1:E:141:GLU:HA	1:E:144:ASP:CG	2.44	0.43
1:F:49:LEU:N	1:F:49:LEU:CD1	2.82	0.43
1:B:125:VAL:CG1	1:B:126:GLU:H	2.30	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:MSE:HE3	1:B:135:VAL:CG2	2.48	0.43
1:A:61:MSE:HE2	1:A:78:TRP:CZ2	2.54	0.43
1:A:81:PRO:O	1:A:82:GLU:CB	2.67	0.43
1:D:137:TYR:OH	1:D:151:MSE:HE1	2.19	0.43
1:F:61:MSE:HE2	1:F:78:TRP:CZ2	2.53	0.43
1:H:41:TRP:CZ2	1:H:97:PRO:HG3	2.54	0.43
1:A:72:GLY:O	1:A:91:ARG:HD2	2.19	0.42
1:G:143:ARG:HD2	1:G:143:ARG:C	2.44	0.42
1:B:59:TRP:CE2	1:B:81:PRO:HD2	2.55	0.42
1:C:31:ARG:HE	1:C:159:TYR:HB3	1.84	0.42
1:F:94:LEU:HG	1:F:102:ILE:HD12	2.00	0.42
1:F:100:ARG:HD3	1:F:121:GLN:HE22	1.84	0.42
1:H:157:GLU:OE2	1:H:160:ARG:NH1	2.52	0.42
1:B:57:ARG:HH11	1:B:57:ARG:HG2	1.85	0.42
1:B:139:THR:OG1	1:B:141:GLU:HG2	2.19	0.42
1:B:151:MSE:O	1:B:155[B]:MSE:HB2	2.19	0.42
1:D:73:LEU:HD12	1:D:74:ILE:H	1.84	0.42
1:D:160:ARG:O	1:D:163:ASP:HB2	2.20	0.42
1:F:84:ARG:O	1:F:84:ARG:HG2	2.19	0.42
1:H:101:LEU:HD23	1:H:120:LEU:HD12	2.01	0.42
1:H:129:THR:CG2	1:H:130:GLU:N	2.81	0.42
1:C:61:MSE:SE	1:C:64:CYS:HB2	2.70	0.42
1:F:168:ALA:O	1:F:170:GLU:N	2.51	0.42
1:H:46:SER:HA	1:H:47:PRO:HD3	1.94	0.42
1:B:37:PRO:HG3	1:B:124:PRO:HG3	2.01	0.42
1:D:59:TRP:CE2	1:D:86:ARG:NH2	2.88	0.42
1:E:17:ILE:HG23	1:E:17:ILE:O	2.19	0.42
1:E:21:PRO:HG2	1:E:22:GLY:H	1.84	0.42
1:A:54:LEU:O	1:A:61:MSE:HE3	2.19	0.42
1:B:158:ALA:O	1:B:161:HIS:HB2	2.19	0.42
1:B:160:ARG:HH22	1:B:161:HIS:CD2	2.38	0.42
1:E:59:TRP:CE3	1:E:80:GLY:HA2	2.55	0.42
1:G:88:LEU:C	1:G:88:LEU:CD1	2.93	0.42
1:H:55:PHE:CD1	1:H:55:PHE:C	2.97	0.42
1:H:100:ARG:HA	1:H:120:LEU:O	2.20	0.42
1:B:109:GLU:O	1:B:110:ASP:HB3	2.19	0.42
1:E:46:SER:HA	1:E:47:PRO:HD3	1.87	0.42
1:F:79:THR:HG22	1:F:80:GLY:N	2.35	0.42
1:A:100:ARG:HD3	1:A:121:GLN:NE2	2.34	0.41
1:B:54:LEU:O	1:B:61:MSE:HE3	2.20	0.41
1:D:97:PRO:HA	1:D:98:PRO:HA	1.47	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:102:ILE:HD13	1:G:117:LEU:HD22	2.02	0.41
1:B:145:ALA:O	1:B:149:SER:OG	2.37	0.41
1:D:57:ARG:HG2	1:D:57:ARG:NH1	2.35	0.41
1:F:97:PRO:HA	1:F:98:PRO:HA	1.57	0.41
1:B:46:SER:HA	1:B:47:PRO:HD3	1.85	0.41
1:F:82:GLU:OE2	1:F:82:GLU:CA	2.69	0.41
1:A:16:GLU:N	2:A:206:HOH:O	2.53	0.41
1:B:125:VAL:HG12	1:B:126:GLU:H	1.82	0.41
1:C:117:LEU:HD21	1:F:117:LEU:HD21	2.02	0.41
1:H:143:ARG:O	1:H:143:ARG:CD	2.65	0.41
1:C:40:LEU:HD12	1:C:40:LEU:HA	1.89	0.41
1:E:19:ILE:HD11	1:E:152:ALA:CB	2.51	0.41
1:G:160:ARG:O	1:G:163:ASP:HB2	2.20	0.41
1:G:125:VAL:HG12	1:G:126:GLU:N	2.35	0.41
1:B:100:ARG:HD3	1:B:121:GLN:HE22	1.85	0.41
1:C:109:GLU:O	1:C:110:ASP:C	2.64	0.41
1:C:141:GLU:CD	1:C:141:GLU:H	2.28	0.41
1:D:33:PHE:HZ	1:D:159:TYR:CD1	2.39	0.41
1:E:41:TRP:CE2	1:E:97:PRO:HB3	2.56	0.41
1:F:122:LEU:N	1:F:122:LEU:HD12	2.36	0.41
1:G:48:ALA:C	1:G:49:LEU:HD12	2.46	0.41
1:G:52:ALA:HB3	1:G:165:MSE:SE	2.71	0.41
1:H:55:PHE:HB2	1:H:59:TRP:O	2.21	0.41
1:A:156:GLU:HG3	1:A:160:ARG:HH22	1.86	0.41
1:G:27:LEU:O	1:G:134:ALA:HA	2.21	0.41
1:C:93:ILE:CD1	1:F:93:ILE:HG23	2.52	0.40
1:C:94:LEU:HG	1:C:102:ILE:HD12	2.03	0.40
1:G:122:LEU:HD12	1:G:122:LEU:N	2.36	0.40
1:H:125:VAL:O	1:H:128:GLY:O	2.39	0.40
1:E:141:GLU:HA	1:E:144:ASP:OD1	2.21	0.40
1:F:42:ALA:O	1:F:46:SER:HB3	2.21	0.40
1:H:125:VAL:CG1	1:H:126:GLU:N	2.84	0.40
1:D:146:VAL:O	1:D:149:SER:HB2	2.21	0.40
1:A:103:HIS:HD2	1:A:104:SER:O	2.05	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	148/164 (90%)	136 (92%)	11 (7%)	1 (1%)	18	38
1	B	150/164 (92%)	141 (94%)	8 (5%)	1 (1%)	18	38
1	C	149/164 (91%)	141 (95%)	8 (5%)	0	100	100
1	D	146/164 (89%)	136 (93%)	9 (6%)	1 (1%)	18	38
1	E	147/164 (90%)	140 (95%)	6 (4%)	1 (1%)	18	38
1	F	144/164 (88%)	130 (90%)	13 (9%)	1 (1%)	18	38
1	G	145/164 (88%)	133 (92%)	11 (8%)	1 (1%)	18	38
1	H	144/164 (88%)	130 (90%)	13 (9%)	1 (1%)	18	38
All	All	1173/1312 (89%)	1087 (93%)	79 (7%)	7 (1%)	21	42

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	148	ALA
1	F	169	LEU
1	A	150	ALA
1	D	150	ALA
1	G	150	ALA
1	H	57	ARG
1	B	138	ALA

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	118/123 (96%)	109 (92%)	9 (8%)	12	27
1	B	119/123 (97%)	113 (95%)	6 (5%)	22	46
1	C	120/123 (98%)	105 (88%)	15 (12%)	4	9
1	D	117/123 (95%)	104 (89%)	13 (11%)	6	12
1	E	118/123 (96%)	102 (86%)	16 (14%)	3	7
1	F	115/123 (94%)	103 (90%)	12 (10%)	7	14
1	G	116/123 (94%)	107 (92%)	9 (8%)	11	26
1	H	116/123 (94%)	107 (92%)	9 (8%)	11	26
All	All	939/984 (95%)	850 (90%)	89 (10%)	8	18

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ILE
1	A	84	ARG
1	A	94	LEU
1	A	99	HIS
1	A	121	GLN
1	A	123	LEU
1	A	126	GLU
1	A	143	ARG
1	A	153	THR
1	B	89	THR
1	B	94	LEU
1	B	99	HIS
1	B	121	GLN
1	B	143	ARG
1	B	160	ARG
1	C	16	GLU
1	C	84	ARG
1	C	89	THR
1	C	91	ARG
1	C	94	LEU
1	C	99	HIS
1	C	109	GLU
1	C	110	ASP
1	C	123	LEU
1	C	126	GLU
1	C	141	GLU
1	C	143	ARG

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Mol	Chain	Res	Type
1	C	151	MSE
1	C	155[A]	MSE
1	C	155[B]	MSE
1	D	84	ARG
1	D	89	THR
1	D	99	HIS
1	D	101	LEU
1	D	119	THR
1	D	121	GLN
1	D	123	LEU
1	D	143	ARG
1	D	151	MSE
1	D	153	THR
1	D	154	GLU
1	D	155[A]	MSE
1	D	155[B]	MSE
1	E	23	SER
1	E	54	LEU
1	E	57	ARG
1	E	84	ARG
1	E	89	THR
1	E	91	ARG
1	E	94	LEU
1	E	99	HIS
1	E	105	GLU
1	E	109	GLU
1	E	121	GLN
1	E	123	LEU
1	E	149	SER
1	E	155	MSE
1	E	156	GLU
1	E	170	GLU
1	F	84	ARG
1	F	86	ARG
1	F	89	THR
1	F	94	LEU
1	F	98	PRO
1	F	99	HIS
1	F	101	LEU
1	F	121	GLN
1	F	123	LEU
1	F	126	GLU

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Mol	Chain	Res	Type
1	F	143	ARG
1	F	170	GLU
1	G	49	LEU
1	G	84	ARG
1	G	88	LEU
1	G	94	LEU
1	G	99	HIS
1	G	121	GLN
1	G	123	LEU
1	G	126	GLU
1	G	143	ARG
1	H	17	ILE
1	H	84	ARG
1	H	89	THR
1	H	121	GLN
1	H	123	LEU
1	H	126	GLU
1	H	143	ARG
1	H	153	THR
1	H	157	GLU

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	18	GLN
1	A	103	HIS
1	B	18	GLN
1	B	99	HIS
1	B	103	HIS
1	B	121	GLN
1	C	18	GLN
1	C	103	HIS
1	C	121	GLN
1	D	18	GLN
1	D	103	HIS
1	D	121	GLN
1	E	18	GLN
1	E	76	GLN
1	E	99	HIS
1	F	18	GLN
1	F	32	HIS
1	F	121	GLN

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Mol	Chain	Res	Type
1	F	161	HIS
1	G	18	GLN
1	H	18	GLN
1	H	32	HIS
1	H	103	HIS
1	H	121	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 ⓘ

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	146/164 (89%)	0.10	3 (2%) 63 58	22, 57, 95, 118	0
1	B	148/164 (90%)	0.08	5 (3%) 48 42	30, 55, 112, 161	0
1	C	147/164 (89%)	-0.12	4 (2%) 56 50	22, 50, 100, 140	0
1	D	144/164 (87%)	0.31	6 (4%) 40 35	28, 64, 125, 188	0
1	E	146/164 (89%)	0.30	7 (4%) 35 30	42, 63, 98, 138	0
1	F	143/164 (87%)	0.32	5 (3%) 47 41	35, 64, 120, 156	0
1	G	144/164 (87%)	0.35	4 (2%) 55 49	34, 71, 138, 180	0
1	H	143/164 (87%)	0.43	5 (3%) 47 41	34, 70, 114, 164	0
All	All	1161/1312 (88%)	0.22	39 (3%) 48 42	22, 62, 117, 188	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	145	ALA	4.3
1	C	149	SER	4.2
1	D	147	ALA	4.0
1	F	80	GLY	3.9
1	E	81	PRO	3.9
1	B	150	ALA	3.9
1	B	81	PRO	3.8
1	D	148	ALA	3.8
1	D	108	ASP	3.5
1	H	156	GLU	3.1
1	H	81	PRO	3.0
1	D	150	ALA	2.9
1	A	150	ALA	2.8
1	G	81	PRO	2.8
1	H	132	ALA	2.7
1	C	114	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	E	138	ALA	2.6
1	A	108	ASP	2.5
1	B	128	GLY	2.5
1	C	150	ALA	2.5
1	E	139	THR	2.4
1	E	170	GLU	2.4
1	H	148	ALA	2.3
1	G	82	GLU	2.3
1	E	84	ARG	2.3
1	E	149	SER	2.3
1	F	81	PRO	2.3
1	G	127	GLY	2.3
1	H	106	LEU	2.3
1	F	170	GLU	2.2
1	E	142	ALA	2.2
1	A	81	PRO	2.2
1	B	110	ASP	2.2
1	D	149	SER	2.1
1	F	148	ALA	2.1
1	B	15	ASP	2.1
1	F	149	SER	2.1
1	G	163	ASP	2.0
1	C	160	ARG	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.