



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:19 PM UTC

PDB ID : 1ZMO / pdb_00001zmo
Title : Apo structure of haloalcohol dehalogenase HheA of *Arthrobacter* sp. AD2
Authors : de Jong, R.M.; Kalk, K.H.; Tang, L.; Janssen, D.B.; Dijkstra, B.W.
Deposited on : 2005-05-10
Resolution : 2.00 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

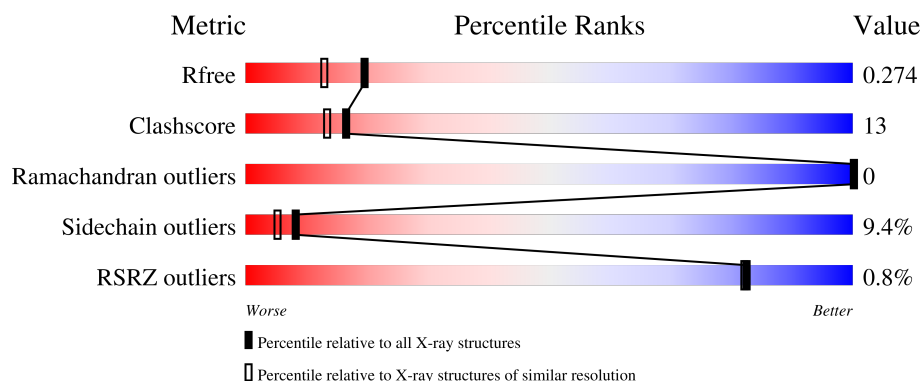
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	244	68% 26% 5%
1	B	244	68% 27% 5%
1	C	244	70% 25% .
1	D	244	% 65% 30% 5%
1	E	244	% 71% 23% 5%

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Mol	Chain	Length	Quality of chain
1	F	244	% 63% 31% 6%
1	G	244	% 67% 29%
1	H	244	% 69% 24% 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 15416 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 halohydrin dehalogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1823	1165	313	341	4			
1	B	243	Total	C	N	O	S	0	0	0
			1825	1167	316	338	4			
1	C	243	Total	C	N	O	S	0	0	0
			1818	1162	313	339	4			
1	D	243	Total	C	N	O	S	0	0	0
			1815	1160	315	336	4			
1	E	243	Total	C	N	O	S	0	0	0
			1822	1162	316	340	4			
1	F	243	Total	C	N	O	S	0	0	0
			1822	1165	315	338	4			
1	G	243	Total	C	N	O	S	0	0	0
			1811	1159	312	336	4			
1	H	243	Total	C	N	O	S	0	0	0
			1818	1163	315	336	4			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	115	Total	O	0	0
			115	115		
2	B	111	Total	O	0	0
			111	111		
2	C	102	Total	O	0	0
			102	102		
2	D	111	Total	O	0	0
			111	111		
2	E	93	Total	O	0	0
			93	93		
2	F	110	Total	O	0	0
			110	110		

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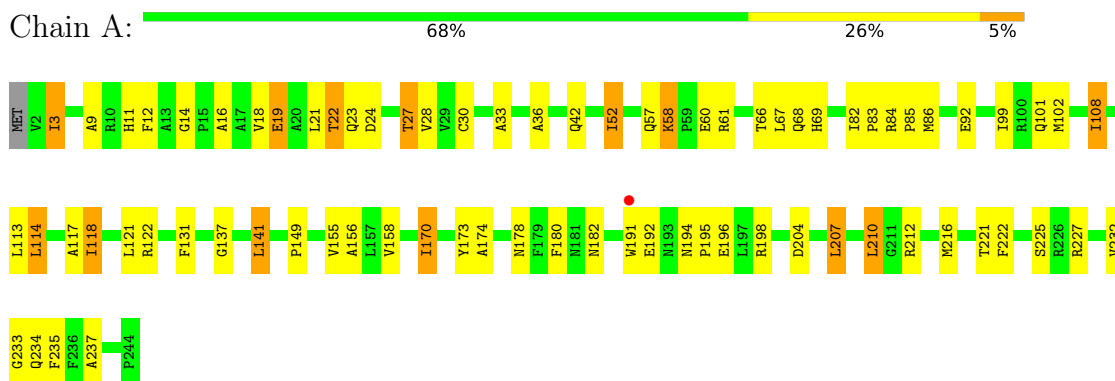
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	111	Total 111	O 111	0	0
2	H	109	Total 109	O 109	0	0

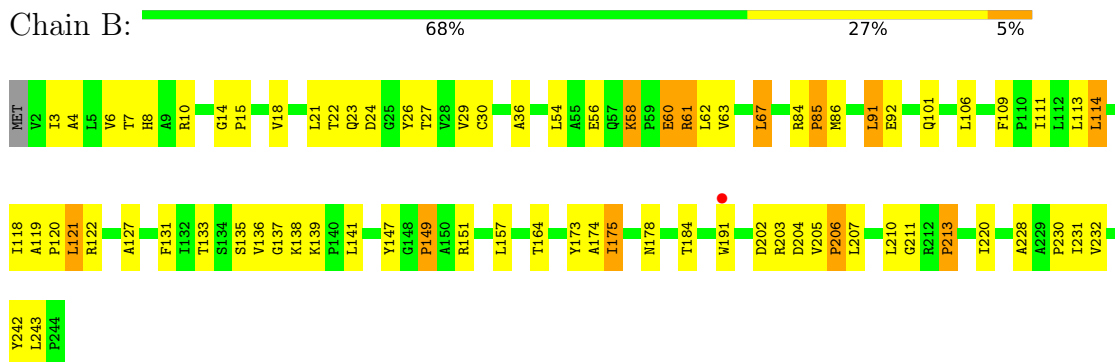
3 Residue-property plots [i](#)

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.

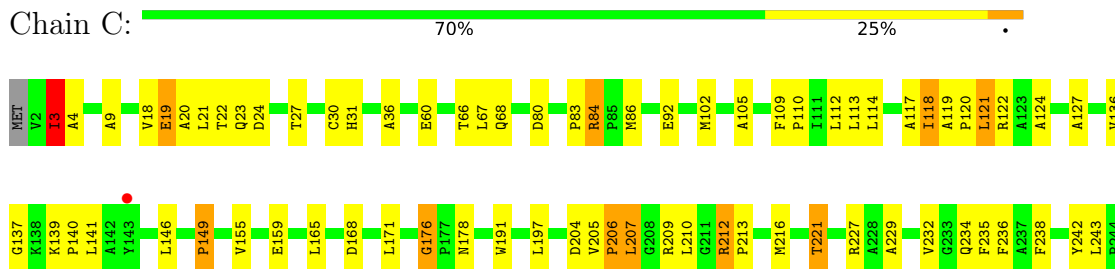
- Molecule 1: halohydrin dehalogenase



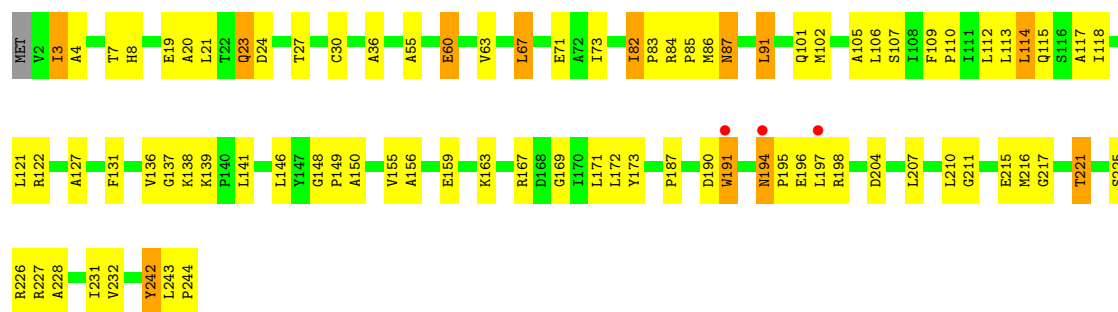
- Molecule 1: halohydrin dehalogenase



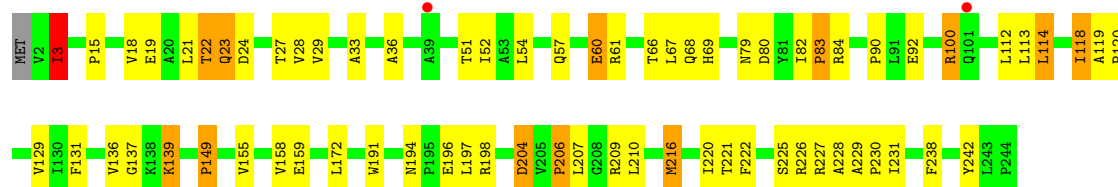
- Molecule 1: halohydrin dehalogenase



- Molecule 1: halohydrin dehalogenase



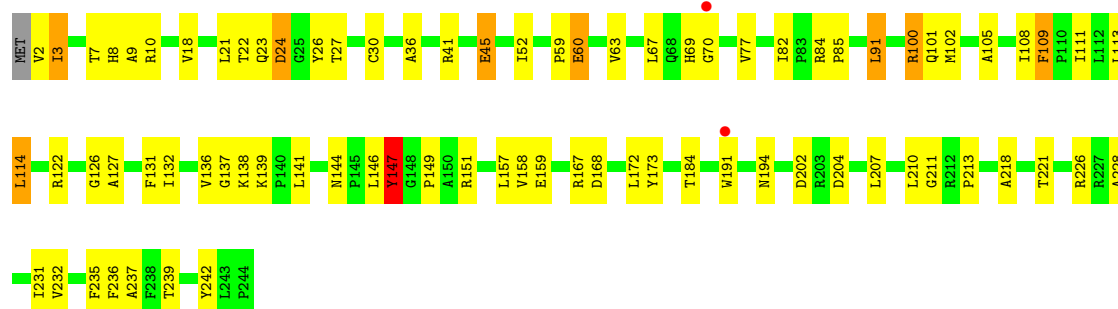
- Molecule 1: halohydrin dehalogenase



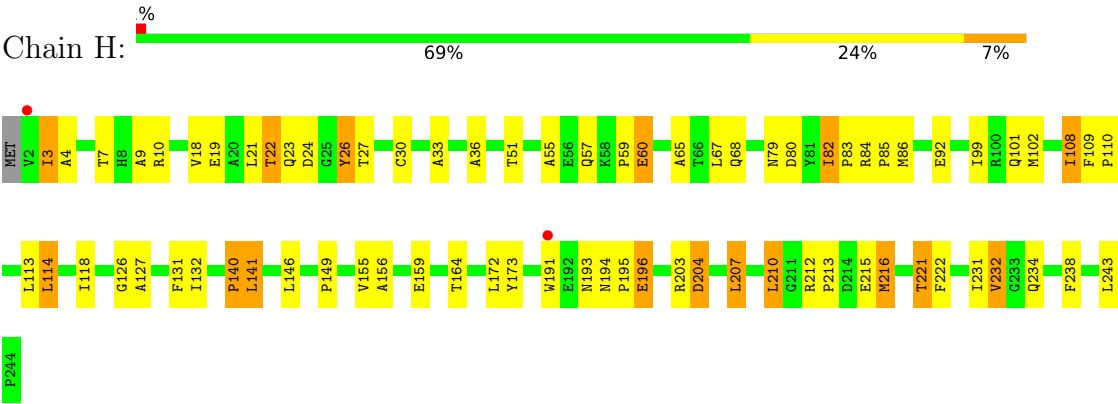
- Molecule 1: halohydrin dehalogenase



- Molecule 1: halohydrin dehalogenase



● Molecule 1: halohydrin dehalogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	64.98Å 77.72Å 111.42Å 97.29° 89.97° 112.98°	Depositor
Resolution (Å)	20.05 – 2.00 20.05 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.3 (20.05-2.00) 89.3 (20.05-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.72 (at 1.99Å)	Xtriage
Refinement program	CNS 1.1, CNX	Depositor
R, R_{free}	0.218 , 0.273 0.218 , 0.274	Depositor DCC
R_{free} test set	6020 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.521	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 56.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.267 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15416	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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	1.14	3/1867 (0.2%)	1.25	13/2552 (0.5%)
1	B	1.12	3/1869 (0.2%)	1.30	21/2553 (0.8%)
1	C	1.10	1/1862 (0.1%)	1.28	21/2546 (0.8%)
1	D	1.17	3/1859 (0.2%)	1.28	16/2541 (0.6%)
1	E	1.14	0/1865	1.29	17/2548 (0.7%)
1	F	1.18	4/1866 (0.2%)	1.31	23/2550 (0.9%)
1	G	1.11	3/1855 (0.2%)	1.28	18/2537 (0.7%)
1	H	1.15	3/1862 (0.2%)	1.29	17/2545 (0.7%)
All	All	1.14	20/14905 (0.1%)	1.29	146/20372 (0.7%)

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	G	0	1
1	H	0	1
All	All	0	2

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	216	MET	SD-CE	7.09	1.97	1.79
1	D	118	ILE	CA-CB	6.67	1.61	1.54
1	H	232	VAL	CA-CB	6.58	1.60	1.54
1	B	6	VAL	CA-CB	6.48	1.62	1.54
1	C	20	ALA	CA-CB	5.87	1.62	1.53
1	F	111	ILE	CA-CB	5.75	1.61	1.54
1	F	118	ILE	CA-CB	5.74	1.60	1.54
1	B	205	VAL	CA-CB	5.71	1.59	1.53
1	H	99	ILE	CA-CB	5.56	1.61	1.54
1	G	26	TYR	CA-C	-5.54	1.45	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	221	THR	CA-CB	5.49	1.61	1.53
1	B	133	THR	CA-CB	5.35	1.62	1.53
1	A	108	ILE	CA-CB	5.35	1.61	1.54
1	A	99	ILE	CA-CB	5.26	1.60	1.54
1	F	82	ILE	CA-CB	5.14	1.60	1.54
1	G	218	ALA	CA-CB	5.14	1.61	1.53
1	H	65	ALA	CA-CB	5.05	1.61	1.53
1	A	16	ALA	CA-CB	5.04	1.61	1.53
1	G	111	ILE	CA-CB	5.03	1.60	1.54
1	F	162	ALA	CA-CB	5.01	1.61	1.53

All (146) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	139	LYS	CA-C-N	8.80	128.82	119.76
1	E	139	LYS	C-N-CA	8.80	128.82	119.76
1	G	204	ASP	N-CA-C	8.59	123.94	113.38
1	D	84	ARG	N-CA-C	-8.58	98.94	109.83
1	G	30	CYS	N-CA-C	8.25	122.36	109.07
1	A	27	THR	N-CA-C	-8.09	96.09	108.96
1	G	137	GLY	N-CA-C	-7.82	104.70	114.16
1	C	118	ILE	N-CA-C	7.73	117.80	110.30
1	G	126	GLY	N-CA-C	7.70	119.89	112.04
1	F	137	GLY	N-CA-C	-7.67	105.14	114.66
1	B	205	VAL	N-CA-C	7.22	114.69	107.55
1	C	137	GLY	N-CA-C	-7.11	106.85	114.67
1	A	204	ASP	N-CA-C	7.08	122.92	113.72
1	E	118	ILE	N-CA-C	7.06	117.14	110.30
1	B	242	TYR	N-CA-C	-7.02	103.71	111.36
1	H	27	THR	N-CA-C	-6.99	97.84	108.96
1	C	27	THR	N-CA-C	-6.93	97.44	108.73
1	H	68	GLN	N-CA-C	-6.92	104.80	113.18
1	B	109	PHE	CA-C-N	-6.89	112.60	119.56
1	B	109	PHE	C-N-CA	-6.89	112.60	119.56
1	F	27	THR	N-CA-C	-6.88	97.34	108.41
1	B	30	CYS	N-CA-C	6.85	120.09	109.07
1	B	137	GLY	N-CA-C	-6.76	106.47	115.32
1	H	204	ASP	N-CA-C	6.74	122.33	113.30
1	F	84	ARG	N-CA-C	-6.72	101.30	109.83
1	G	27	THR	N-CA-C	-6.72	97.59	108.41
1	F	68	GLN	N-CA-C	-6.70	105.08	113.18
1	E	27	THR	N-CA-C	-6.69	98.35	109.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	137	GLY	N-CA-C	-6.68	106.08	114.16
1	D	27	THR	N-CA-C	-6.67	97.66	108.41
1	F	155	VAL	N-CA-C	-6.67	103.98	110.72
1	E	84	ARG	N-CA-C	-6.60	100.97	110.08
1	C	216	MET	N-CA-C	-6.59	104.18	111.36
1	E	80	ASP	N-CA-C	6.59	119.91	110.10
1	G	109	PHE	CA-C-N	-6.57	112.92	119.56
1	G	109	PHE	C-N-CA	-6.57	112.92	119.56
1	B	14	GLY	CA-C-N	6.56	125.79	118.97
1	B	14	GLY	C-N-CA	6.56	125.79	118.97
1	D	242	TYR	N-CA-C	-6.54	104.15	111.28
1	F	186	PHE	CA-C-N	-6.50	113.29	119.85
1	F	186	PHE	C-N-CA	-6.50	113.29	119.85
1	H	216	MET	N-CA-C	-6.45	104.33	111.36
1	E	204	ASP	N-CA-C	6.43	121.41	113.50
1	B	211	GLY	N-CA-C	-6.38	103.91	112.14
1	G	36	ALA	N-CA-C	-6.36	104.78	112.54
1	H	146	LEU	N-CA-C	-6.30	105.49	113.43
1	B	27	THR	N-CA-C	-6.19	97.62	108.20
1	B	3	ILE	N-CA-C	6.17	117.19	108.36
1	H	36	ALA	N-CA-C	-6.17	104.86	112.38
1	A	216	MET	N-CA-C	-6.15	104.66	111.36
1	F	147	TYR	N-CA-C	-6.09	104.72	111.36
1	G	157	LEU	N-CA-C	-6.04	104.78	111.36
1	G	102	MET	N-CA-C	-6.03	104.61	111.07
1	B	118	ILE	N-CA-C	6.02	116.49	110.23
1	E	137	GLY	N-CA-C	-6.02	108.05	114.67
1	A	52	ILE	N-CA-C	-5.99	99.49	108.12
1	F	118	ILE	N-CA-C	5.99	116.45	110.23
1	C	205	VAL	CA-C-N	5.97	126.12	119.32
1	C	205	VAL	C-N-CA	5.97	126.12	119.32
1	D	36	ALA	N-CA-C	-5.93	105.31	112.54
1	F	36	ALA	N-CA-C	-5.89	104.98	111.82
1	D	55	ALA	N-CA-C	-5.89	105.18	112.90
1	C	80	ASP	N-CA-C	5.88	118.78	109.96
1	D	204	ASP	N-CA-C	5.87	120.60	113.38
1	F	211	GLY	N-CA-C	-5.87	104.57	112.14
1	D	155	VAL	N-CA-C	-5.85	104.81	110.72
1	E	28	VAL	N-CA-C	5.85	116.02	107.37
1	B	147	TYR	N-CA-C	-5.83	104.92	111.28
1	F	204	ASP	N-CA-C	5.83	121.11	113.30
1	D	171	LEU	N-CA-C	5.80	118.11	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	151	ARG	N-CA-C	-5.78	104.58	111.69
1	D	169	GLY	N-CA-C	-5.77	106.19	115.08
1	D	30	CYS	N-CA-C	5.74	118.31	109.07
1	F	203	ARG	N-CA-C	5.68	117.47	111.28
1	A	68	GLN	N-CA-C	-5.62	105.68	112.54
1	F	30	CYS	N-CA-C	5.62	118.56	109.40
1	B	204	ASP	N-CA-C	5.61	120.19	113.23
1	A	137	GLY	N-CA-C	-5.61	108.50	114.67
1	B	36	ALA	N-CA-C	-5.59	105.56	112.38
1	D	159	GLU	N-CA-C	5.59	117.38	111.28
1	F	55	ALA	N-CA-C	-5.59	105.18	112.23
1	H	140	PRO	N-CA-C	5.57	120.07	111.21
1	C	242	TYR	N-CA-C	-5.56	105.13	111.14
1	G	184	THR	N-CA-C	5.56	120.07	111.56
1	F	176	GLY	CA-C-N	-5.54	115.04	120.52
1	F	176	GLY	C-N-CA	-5.54	115.04	120.52
1	C	178	ASN	N-CA-C	-5.53	100.03	108.76
1	C	212	ARG	N-CA-C	5.52	118.30	110.24
1	F	82	ILE	CA-C-N	-5.50	114.27	119.76
1	F	82	ILE	C-N-CA	-5.50	114.27	119.76
1	A	36	ALA	N-CA-C	-5.48	105.70	112.38
1	D	118	ILE	N-CA-C	5.48	115.92	110.23
1	G	9	ALA	N-CA-C	5.48	118.01	111.71
1	C	146	LEU	N-CA-C	-5.46	105.67	112.93
1	G	144	ASN	CA-C-N	-5.45	114.20	119.87
1	G	144	ASN	C-N-CA	-5.45	114.20	119.87
1	G	105	ALA	N-CA-C	5.45	117.30	111.36
1	E	83	PRO	N-CA-C	5.41	119.68	111.19
1	B	106	LEU	N-CA-C	5.40	120.02	113.38
1	C	238	PHE	N-CA-C	-5.37	99.63	108.49
1	F	242	TYR	N-CA-C	-5.37	105.51	111.36
1	F	171	LEU	N-CA-C	5.37	117.38	108.20
1	C	84	ARG	N-CA-C	-5.35	102.69	110.08
1	C	36	ALA	N-CA-C	-5.35	105.86	112.38
1	H	55	ALA	N-CA-C	-5.33	105.92	112.90
1	C	105	ALA	N-CA-C	5.31	117.15	111.36
1	B	184	THR	N-CA-C	5.30	119.67	111.56
1	E	36	ALA	N-CA-C	-5.28	105.94	112.38
1	A	141	LEU	N-CA-C	-5.28	100.43	109.24
1	H	80	ASP	N-CA-C	5.26	117.86	109.96
1	H	132	ILE	N-CA-C	-5.25	100.20	108.23
1	C	68	GLN	N-CA-C	-5.24	106.88	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	26	TYR	N-CA-C	5.23	117.87	110.50
1	C	3	ILE	N-CA-C	5.23	116.83	108.89
1	A	178	ASN	N-CA-C	-5.22	100.79	108.99
1	A	118	ILE	N-CA-C	5.22	115.36	110.30
1	G	147	TYR	N-CA-C	-5.21	105.78	111.82
1	F	106	LEU	N-CA-C	5.20	119.78	113.38
1	A	28	VAL	N-CA-C	5.19	115.05	107.37
1	G	211	GLY	N-CA-C	-5.19	104.48	111.54
1	B	111	ILE	N-CA-C	-5.18	105.49	110.72
1	A	212	ARG	N-CA-C	5.18	117.80	110.24
1	E	216	MET	N-CA-C	-5.17	105.53	111.07
1	H	193	ASN	N-CA-C	5.17	120.43	113.72
1	B	157	LEU	N-CA-C	-5.14	105.75	111.36
1	H	221	THR	N-CA-CB	-5.14	102.31	110.28
1	D	112	LEU	N-CA-C	-5.14	105.68	111.28
1	B	203	ARG	N-CA-C	5.13	116.69	111.14
1	E	51	THR	N-CA-C	5.13	117.96	109.85
1	H	82	ILE	CA-C-N	5.11	125.40	119.93
1	H	82	ILE	C-N-CA	5.11	125.40	119.93
1	C	176	GLY	CA-C-N	-5.11	115.30	120.31
1	C	176	GLY	C-N-CA	-5.11	115.30	120.31
1	C	243	LEU	N-CA-C	-5.10	103.72	110.40
1	E	242	TYR	N-CA-C	-5.10	105.63	111.14
1	E	68	GLN	N-CA-C	-5.10	106.57	112.89
1	D	105	ALA	N-CA-C	5.08	116.90	111.36
1	G	3	ILE	N-CA-C	5.08	115.22	108.11
1	A	155	VAL	N-CA-C	-5.07	105.60	110.72
1	D	211	GLY	N-CA-C	-5.07	106.10	111.93
1	F	99	ILE	CB-CA-C	-5.07	105.48	111.97
1	C	140	PRO	N-CA-C	5.04	119.23	111.21
1	H	51	THR	N-CA-C	5.04	117.97	109.95
1	E	3	ILE	N-CA-C	5.04	116.55	108.89
1	H	141	LEU	N-CA-C	-5.04	100.83	109.24
1	E	238	PHE	N-CA-C	-5.04	100.18	108.49

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	147	TYR	Sidechain
1	H	26	TYR	Sidechain

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	1823	0	1783	57	0
1	B	1825	0	1795	40	0
1	C	1818	0	1773	54	0
1	D	1815	0	1773	63	0
1	E	1822	0	1787	46	0
1	F	1822	0	1786	63	0
1	G	1811	0	1769	48	0
1	H	1818	0	1782	43	0
2	A	115	0	0	5	0
2	B	111	0	0	6	0
2	C	102	0	0	12	0
2	D	111	0	0	2	0
2	E	93	0	0	0	0
2	F	110	0	0	6	0
2	G	111	0	0	4	0
2	H	109	0	0	5	0
All	All	15416	0	14248	370	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 13.

All (370) 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:F:237:ALA:HB1	2:F:297:HOH:O	1.68	0.94
1:F:58:LYS:HB2	1:F:61:ARG:HG3	1.55	0.88
1:G:59:PRO:HG3	1:G:109:PHE:CD2	2.09	0.87
1:D:110:PRO:HA	2:D:343:HOH:O	1.74	0.87
1:D:3:ILE:HD11	1:D:73:ILE:HG12	1.55	0.86
1:F:3:ILE:HD12	1:F:3:ILE:C	2.01	0.85
1:E:194:ASN:HD21	1:E:196:GLU:HB3	1.42	0.84
1:C:234:GLN:HG3	2:C:309:HOH:O	1.77	0.84
1:B:206:PRO:O	1:D:167:ARG:HD3	1.80	0.82
1:D:194:ASN:C	1:D:194:ASN:HD22	1.85	0.82
1:A:82:ILE:HD12	1:A:83:PRO:HD2	1.62	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:207:LEU:HD22	1:C:209:ARG:HG2	1.63	0.79
1:H:203:ARG:HG2	1:H:204:ASP:OD1	1.84	0.78
1:D:19:GLU:OE1	1:D:23:GLN:NE2	2.14	0.78
1:C:118:ILE:HD12	2:C:341:HOH:O	1.83	0.77
1:H:18:VAL:O	1:H:22:THR:HB	1.85	0.77
1:C:229:ALA:O	1:C:232:VAL:HG13	1.85	0.76
1:E:194:ASN:C	1:E:194:ASN:HD22	1.93	0.76
1:G:18:VAL:O	1:G:22:THR:HG22	1.86	0.76
1:C:171:LEU:HD22	1:C:232:VAL:HG11	1.69	0.75
1:D:194:ASN:ND2	1:D:196:GLU:H	1.85	0.74
1:F:234:GLN:HE22	1:G:236:PHE:HA	1.53	0.74
1:A:114:LEU:HD13	1:B:91:LEU:HD11	1.69	0.74
1:D:87:ASN:H	1:D:87:ASN:HD22	1.36	0.73
1:A:58:LYS:HZ2	1:A:61:ARG:HD2	1.53	0.72
1:D:194:ASN:C	1:D:194:ASN:ND2	2.47	0.71
1:B:58:LYS:HB2	1:B:61:ARG:HG3	1.70	0.71
1:C:114:LEU:HD12	1:D:91:LEU:HD13	1.72	0.71
1:B:202:ASP:OD1	1:D:167:ARG:NH2	2.24	0.70
1:H:194:ASN:HD21	1:H:196:GLU:HB2	1.55	0.70
1:E:114:LEU:HD13	1:F:91:LEU:HD11	1.74	0.70
1:F:239:THR:HG23	2:F:297:HOH:O	1.91	0.69
1:B:18:VAL:O	1:B:22:THR:HG22	1.93	0.69
1:C:118:ILE:HB	2:C:341:HOH:O	1.93	0.68
1:H:194:ASN:C	1:H:194:ASN:HD22	2.00	0.68
1:B:61:ARG:NH1	2:B:331:HOH:O	2.27	0.68
1:B:114:LEU:HG	1:B:131:PHE:HZ	1.59	0.68
1:H:59:PRO:HD2	1:H:60:GLU:OE1	1.93	0.68
1:C:114:LEU:HD13	1:C:114:LEU:C	2.19	0.67
1:C:114:LEU:CD1	1:D:91:LEU:HD13	2.25	0.67
1:E:18:VAL:O	1:E:22:THR:HB	1.93	0.66
1:D:194:ASN:ND2	1:D:196:GLU:N	2.43	0.66
1:B:174:ALA:C	1:B:175:ILE:HD13	2.20	0.66
1:C:92:GLU:HG2	2:C:296:HOH:O	1.95	0.65
1:F:234:GLN:NE2	1:G:237:ALA:H	1.95	0.65
1:B:58:LYS:CB	1:B:61:ARG:HG3	2.26	0.65
1:H:194:ASN:HD21	1:H:196:GLU:CB	2.10	0.65
1:A:121:LEU:HB3	1:A:170:ILE:HD13	1.78	0.64
1:G:167:ARG:HG2	1:G:167:ARG:HH11	1.62	0.64
1:E:207:LEU:HD22	1:E:209:ARG:HG2	1.80	0.64
1:D:146:LEU:C	1:D:149:PRO:HD2	2.23	0.64
1:G:82:ILE:HD11	1:G:101:GLN:HG2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:19:GLU:O	1:H:23:GLN:HB2	1.98	0.63
1:H:86:MET:HE1	1:H:101:GLN:NE2	2.13	0.63
1:D:136:VAL:HA	1:D:139:LYS:O	1.98	0.63
1:G:114:LEU:HG	1:G:131:PHE:HZ	1.64	0.63
1:B:10:ARG:NE	2:B:298:HOH:O	2.32	0.63
1:G:91:LEU:HD11	1:H:114:LEU:HD13	1.80	0.63
1:G:2:VAL:N	2:G:320:HOH:O	2.31	0.63
1:D:24:ASP:OD1	1:D:221:THR:HG23	1.99	0.62
1:F:136:VAL:HA	1:F:139:LYS:O	1.99	0.62
1:D:19:GLU:CD	1:D:23:GLN:HE22	2.07	0.62
1:G:149:PRO:HA	1:H:156:ALA:HB3	1.79	0.62
1:D:87:ASN:HD22	1:D:87:ASN:N	1.97	0.62
1:E:194:ASN:ND2	1:E:196:GLU:H	1.97	0.62
1:A:122:ARG:HH12	1:B:92:GLU:CD	2.08	0.62
1:D:194:ASN:HD21	1:D:196:GLU:CB	2.13	0.62
1:A:117:ALA:O	1:A:121:LEU:HB2	2.00	0.61
1:H:92:GLU:HG2	2:H:289:HOH:O	2.00	0.61
1:C:212:ARG:NH2	2:C:256:HOH:O	2.32	0.61
1:D:82:ILE:HG12	1:D:83:PRO:HD2	1.81	0.61
1:A:11:HIS:HE1	2:A:344:HOH:O	1.84	0.61
1:G:77:VAL:CG1	1:G:132:ILE:HD11	2.31	0.61
1:H:3:ILE:HD12	1:H:4:ALA:N	2.15	0.61
1:A:156:ALA:HB3	1:B:149:PRO:HA	1.83	0.60
1:H:194:ASN:C	1:H:194:ASN:ND2	2.59	0.60
1:E:114:LEU:CD1	1:F:91:LEU:HD11	2.30	0.60
1:A:82:ILE:CD1	1:A:83:PRO:HD2	2.31	0.60
1:D:60:GLU:H	1:D:60:GLU:CD	2.10	0.60
1:C:121:LEU:HG	1:C:127:ALA:HB3	1.84	0.59
1:E:114:LEU:HD13	1:F:91:LEU:CD1	2.32	0.59
1:F:83:PRO:O	1:F:86:MET:HB2	2.01	0.59
1:D:7:THR:O	1:D:8:HIS:C	2.46	0.58
1:C:221:THR:HG21	2:C:285:HOH:O	2.04	0.58
1:A:114:LEU:HD13	1:B:91:LEU:CD1	2.32	0.58
1:E:194:ASN:ND2	1:E:196:GLU:HB3	2.17	0.58
1:B:175:ILE:HD13	1:B:175:ILE:N	2.20	0.57
1:E:194:ASN:HD21	1:E:196:GLU:CB	2.15	0.57
1:B:206:PRO:O	1:D:167:ARG:CD	2.52	0.57
1:D:20:ALA:HB2	1:D:217:GLY:HA3	1.86	0.57
1:F:3:ILE:C	1:F:3:ILE:CD1	2.71	0.57
1:F:99:ILE:HD12	1:F:146:LEU:HD22	1.85	0.57
1:E:118:ILE:HD13	1:F:92:GLU:HG3	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:ASN:C	1:A:194:ASN:HD22	2.13	0.56
1:D:194:ASN:ND2	1:D:197:LEU:H	2.03	0.56
1:F:200:ARG:HA	1:F:203:ARG:NH1	2.19	0.56
1:A:237:ALA:H	1:C:234:GLN:NE2	2.02	0.56
1:B:60:GLU:H	1:B:60:GLU:CD	2.13	0.56
1:D:187:PRO:O	1:D:190:ASP:HB2	2.05	0.56
1:F:106:LEU:HB2	1:F:150:ALA:HB1	1.87	0.56
1:D:173:TYR:CE1	1:D:232:VAL:HA	2.39	0.56
1:B:56:GLU:OE2	1:B:58:LYS:HB2	2.06	0.56
1:F:3:ILE:HD12	1:F:3:ILE:O	2.04	0.56
1:E:29:VAL:CG1	1:E:54:LEU:HD11	2.35	0.56
1:G:41:ARG:O	1:G:45:GLU:HG3	2.06	0.56
1:A:121:LEU:HB3	1:A:170:ILE:CD1	2.35	0.55
1:H:173:TYR:CE1	1:H:232:VAL:HA	2.42	0.55
1:B:230:PRO:HD3	1:D:215:GLU:OE1	2.05	0.55
1:C:204:ASP:O	1:C:206:PRO:HD3	2.07	0.55
1:H:114:LEU:HG	1:H:131:PHE:HZ	1.72	0.55
1:A:192:GLU:HA	1:A:198:ARG:HH11	1.71	0.55
1:C:114:LEU:HD12	1:D:91:LEU:CD1	2.37	0.55
1:G:91:LEU:CD1	1:H:114:LEU:HD13	2.37	0.55
1:G:194:ASN:HD22	1:G:194:ASN:C	2.14	0.54
1:C:227:ARG:NH1	2:C:306:HOH:O	2.41	0.54
1:D:242:TYR:CD1	1:D:243:LEU:HD23	2.42	0.54
1:A:19:GLU:O	1:A:23:GLN:HB2	2.08	0.54
1:A:122:ARG:NH1	2:A:347:HOH:O	2.27	0.54
1:H:59:PRO:HG3	1:H:109:PHE:CD2	2.43	0.54
1:C:171:LEU:HB3	1:C:232:VAL:HG12	1.89	0.54
1:F:20:ALA:HB2	1:F:217:GLY:HA3	1.90	0.54
1:E:194:ASN:ND2	1:E:196:GLU:N	2.56	0.54
1:F:3:ILE:HD11	1:F:73:ILE:HG23	1.90	0.54
1:A:18:VAL:O	1:A:22:THR:HB	2.08	0.54
1:A:222:PHE:O	1:A:225:SER:OG	2.20	0.53
1:A:225:SER:HB2	1:A:227:ARG:HD3	1.90	0.53
1:D:3:ILE:C	1:D:3:ILE:HD12	2.33	0.53
1:E:3:ILE:HD11	1:E:66:THR:CG2	2.38	0.53
1:F:19:GLU:OE2	1:F:23:GLN:OE1	2.26	0.53
1:F:117:ALA:HB1	1:F:121:LEU:HD22	1.89	0.53
1:E:194:ASN:C	1:E:194:ASN:ND2	2.60	0.53
1:E:3:ILE:HD11	1:E:66:THR:HG21	1.90	0.53
1:G:122:ARG:HD3	1:G:168:ASP:CG	2.34	0.53
1:A:9:ALA:HB2	1:A:30:CYS:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:24:ASP:OD1	1:G:221:THR:HG23	2.08	0.53
1:A:173:TYR:CE1	1:A:232:VAL:HA	2.44	0.53
1:B:29:VAL:CG1	1:B:54:LEU:HD11	2.39	0.53
1:G:122:ARG:HH12	1:H:92:GLU:CD	2.17	0.53
1:A:3:ILE:HD11	1:A:66:THR:HG22	1.91	0.53
1:B:228:ALA:O	1:B:231:ILE:HG12	2.09	0.53
1:A:114:LEU:CD1	1:B:91:LEU:HD11	2.39	0.52
1:D:121:LEU:HG	1:D:127:ALA:CB	2.39	0.52
1:F:37:ASP:OD1	1:F:39:ALA:HB3	2.10	0.52
1:F:18:VAL:O	1:F:22:THR:HG22	2.10	0.52
1:G:167:ARG:HG2	1:G:167:ARG:NH1	2.24	0.52
1:B:84:ARG:HG3	1:B:85:PRO:HA	1.91	0.52
1:G:59:PRO:HG3	1:G:109:PHE:CE2	2.44	0.52
1:A:33:ALA:HA	1:A:57:GLN:OE1	2.09	0.52
1:A:84:ARG:HG3	1:A:85:PRO:HA	1.92	0.52
1:E:52:ILE:HD13	1:E:69:HIS:CD2	2.45	0.52
1:A:122:ARG:NH1	1:B:92:GLU:OE2	2.43	0.51
1:F:234:GLN:NE2	1:G:236:PHE:HA	2.23	0.51
1:G:63:VAL:O	1:G:67:LEU:HD13	2.10	0.51
1:E:204:ASP:O	1:E:206:PRO:HD3	2.11	0.51
1:D:195:PRO:HA	1:D:198:ARG:NH2	2.25	0.51
1:A:180:PHE:O	1:A:182:ASN:ND2	2.43	0.51
1:D:146:LEU:O	1:D:149:PRO:HD2	2.11	0.51
1:D:243:LEU:HB3	1:D:244:PRO:HA	1.92	0.51
1:F:173:TYR:CE1	1:F:232:VAL:HA	2.44	0.50
1:A:210:LEU:HB2	2:A:279:HOH:O	2.12	0.50
1:C:213:PRO:HB3	2:C:304:HOH:O	2.11	0.50
1:D:121:LEU:HG	1:D:127:ALA:HB3	1.93	0.50
1:F:3:ILE:HD11	1:F:73:ILE:HG12	1.94	0.50
1:F:135:SER:HB2	1:F:243:LEU:HD11	1.93	0.50
1:H:109:PHE:HB3	1:H:110:PRO:CD	2.41	0.50
1:A:234:GLN:HG2	1:A:235:PHE:N	2.26	0.50
1:F:99:ILE:HD12	1:F:146:LEU:CD2	2.41	0.50
1:E:114:LEU:HG	1:E:131:PHE:HZ	1.76	0.50
1:F:234:GLN:HG2	2:F:318:HOH:O	2.10	0.50
1:A:194:ASN:HD21	1:A:196:GLU:CB	2.25	0.50
1:F:121:LEU:HG	1:F:127:ALA:HB3	1.94	0.50
1:A:194:ASN:C	1:A:194:ASN:ND2	2.69	0.49
1:F:82:ILE:HD13	1:F:105:ALA:HB3	1.94	0.49
1:D:86:MET:O	1:D:102:MET:HE1	2.12	0.49
1:D:194:ASN:HD22	1:D:196:GLU:H	1.55	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:194:ASN:HD22	1:E:196:GLU:H	1.60	0.49
1:F:21:LEU:N	1:F:21:LEU:CD1	2.75	0.49
1:E:225:SER:HB2	1:E:227:ARG:HD3	1.93	0.49
1:B:173:TYR:CE1	1:B:232:VAL:HA	2.48	0.49
1:E:228:ALA:O	1:E:231:ILE:HG12	2.13	0.49
1:F:202:ASP:O	1:G:167:ARG:NH2	2.46	0.49
1:C:92:GLU:CD	1:D:122:ARG:HH12	2.21	0.48
1:D:194:ASN:HD22	1:D:195:PRO:N	2.10	0.48
1:F:60:GLU:CD	1:F:60:GLU:H	2.20	0.48
1:G:138:LYS:HG2	1:G:235:PHE:CD2	2.49	0.48
1:H:86:MET:O	1:H:102:MET:HE1	2.13	0.48
1:G:173:TYR:CE1	1:G:232:VAL:HA	2.49	0.48
1:E:82:ILE:HD12	1:E:83:PRO:HD2	1.95	0.48
1:F:238:PHE:N	2:F:297:HOH:O	2.46	0.48
1:C:136:VAL:HA	1:C:139:LYS:O	2.14	0.48
1:F:138:LYS:HD3	1:F:235:PHE:CG	2.49	0.48
1:D:106:LEU:HB2	1:D:150:ALA:HB1	1.95	0.48
1:E:136:VAL:HA	1:E:139:LYS:O	2.15	0.47
1:B:10:ARG:CD	2:B:298:HOH:O	2.62	0.47
1:E:229:ALA:HB3	1:E:230:PRO:HD3	1.96	0.47
1:D:83:PRO:O	1:D:86:MET:HB2	2.14	0.47
1:B:119:ALA:HB3	1:B:120:PRO:CD	2.45	0.47
1:H:10:ARG:NH2	2:H:290:HOH:O	2.47	0.47
1:D:136:VAL:HG23	1:D:148:GLY:HA2	1.97	0.47
1:F:187:PRO:O	1:F:190:ASP:HB2	2.14	0.47
1:G:7:THR:O	1:G:8:HIS:C	2.58	0.47
1:A:194:ASN:ND2	1:A:196:GLU:H	2.12	0.47
1:F:56:GLU:OE2	1:F:58:LYS:CG	2.62	0.47
1:G:213:PRO:HG2	2:G:305:HOH:O	2.14	0.47
1:H:213:PRO:HG2	2:H:285:HOH:O	2.13	0.47
1:G:194:ASN:C	1:G:194:ASN:ND2	2.72	0.47
1:A:23:GLN:NE2	2:A:324:HOH:O	2.47	0.47
1:C:207:LEU:O	1:C:207:LEU:HD23	2.15	0.47
1:G:100:ARG:HG3	2:G:285:HOH:O	2.15	0.47
1:D:114:LEU:HG	1:D:131:PHE:HZ	1.79	0.46
1:B:4:ALA:HB2	1:B:26:TYR:CD2	2.50	0.46
1:G:127:ALA:HA	1:G:226:ARG:NH2	2.30	0.46
1:H:7:THR:O	1:H:79:ASN:HB3	2.14	0.46
1:D:109:PHE:HB3	1:D:110:PRO:CD	2.45	0.46
1:G:45:GLU:HB3	2:G:306:HOH:O	2.14	0.46
1:H:33:ALA:HA	1:H:57:GLN:OE1	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:GLN:NE2	1:G:10:ARG:HH21	2.12	0.46
1:A:52:ILE:HG21	1:A:69:HIS:CE1	2.49	0.46
1:A:194:ASN:HA	1:A:195:PRO:HD2	1.83	0.46
1:F:114:LEU:O	1:F:115:GLN:C	2.56	0.46
1:F:237:ALA:CB	2:F:297:HOH:O	2.44	0.46
1:A:86:MET:HE1	1:A:101:GLN:NE2	2.31	0.46
1:E:92:GLU:OE2	1:F:122:ARG:NH1	2.46	0.46
1:F:21:LEU:HD11	1:F:220:ILE:HG22	1.98	0.46
1:G:138:LYS:HG2	1:G:235:PHE:CE2	2.51	0.46
1:A:14:GLY:O	1:A:18:VAL:HG23	2.15	0.46
1:C:109:PHE:N	1:C:110:PRO:HD2	2.30	0.46
1:E:191:TRP:CH2	1:E:198:ARG:HG3	2.50	0.46
2:B:348:HOH:O	1:D:163:LYS:HE3	2.16	0.46
1:C:117:ALA:HB1	1:C:121:LEU:HD22	1.97	0.46
1:C:118:ILE:CB	2:C:341:HOH:O	2.55	0.46
1:H:155:VAL:O	1:H:159:GLU:HG3	2.15	0.45
1:D:225:SER:O	1:D:226:ARG:HB2	2.15	0.45
1:E:155:VAL:O	1:E:159:GLU:HG3	2.17	0.45
1:G:158:VAL:HG13	1:G:172:LEU:HB3	1.98	0.45
1:E:92:GLU:CD	1:F:122:ARG:HH12	2.24	0.45
1:H:194:ASN:ND2	1:H:196:GLU:H	2.15	0.45
1:D:114:LEU:O	1:D:115:GLN:C	2.57	0.45
1:G:239:THR:HB	1:G:242:TYR:HB3	1.98	0.45
1:A:158:VAL:HG21	1:A:174:ALA:HB2	1.99	0.45
1:D:141:LEU:HD23	1:D:141:LEU:HA	1.86	0.45
1:E:149:PRO:HA	1:F:156:ALA:HB3	1.99	0.45
1:A:3:ILE:HD11	1:A:66:THR:CG2	2.47	0.45
1:C:18:VAL:O	1:C:22:THR:HG22	2.17	0.45
1:C:114:LEU:HD13	1:C:114:LEU:O	2.16	0.45
1:C:149:PRO:HA	1:D:156:ALA:HB3	1.99	0.45
1:B:86:MET:HE1	1:B:101:GLN:NE2	2.32	0.44
1:B:202:ASP:O	1:D:167:ARG:NH1	2.47	0.44
1:H:210:LEU:HB2	2:H:256:HOH:O	2.15	0.44
1:A:141:LEU:HD23	1:A:141:LEU:HA	1.79	0.44
1:E:52:ILE:HG21	1:E:69:HIS:CD2	2.52	0.44
1:E:79:ASN:CG	1:E:79:ASN:O	2.60	0.44
1:F:102:MET:HE2	1:F:102:MET:HB2	1.84	0.44
1:B:63:VAL:O	1:B:67:LEU:HD13	2.17	0.44
1:C:119:ALA:HB3	1:C:120:PRO:CD	2.47	0.44
1:E:60:GLU:H	1:E:60:GLU:CD	2.24	0.44
1:F:63:VAL:O	1:F:67:LEU:HD22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:243:LEU:HD23	1:H:243:LEU:HA	1.69	0.44
1:B:141:LEU:HD23	1:B:141:LEU:HA	1.91	0.44
1:D:67:LEU:HD12	1:D:67:LEU:HA	1.69	0.44
1:E:119:ALA:HB3	1:E:120:PRO:CD	2.47	0.44
1:E:216:MET:O	1:E:220:ILE:HG12	2.17	0.44
1:F:56:GLU:OE2	1:F:58:LYS:HG2	2.18	0.44
1:C:109:PHE:HB3	1:C:110:PRO:CD	2.48	0.44
1:H:108:ILE:N	1:H:108:ILE:CD1	2.81	0.44
1:H:164:THR:HG23	2:H:253:HOH:O	2.17	0.44
1:C:118:ILE:HD13	1:C:165:LEU:HD21	1.99	0.44
1:D:194:ASN:HD21	1:D:196:GLU:N	2.14	0.44
1:C:141:LEU:HD23	1:C:141:LEU:HA	1.88	0.43
1:D:110:PRO:CA	2:D:343:HOH:O	2.46	0.43
1:C:3:ILE:HD12	1:C:4:ALA:O	2.17	0.43
1:G:147:TYR:O	1:G:151:ARG:HG2	2.18	0.43
1:B:7:THR:O	1:B:8:HIS:C	2.61	0.43
1:C:155:VAL:O	1:C:159:GLU:HG3	2.17	0.43
1:E:225:SER:O	1:E:226:ARG:HB2	2.19	0.43
1:G:141:LEU:HD23	1:G:141:LEU:HA	1.84	0.43
1:H:126:GLY:O	1:H:127:ALA:HB2	2.19	0.43
1:H:194:ASN:HA	1:H:195:PRO:HD2	1.83	0.43
1:A:207:LEU:HD23	1:A:207:LEU:HA	1.65	0.43
1:A:82:ILE:HD12	1:A:82:ILE:HA	1.82	0.43
1:A:92:GLU:CD	1:B:122:ARG:HH12	2.27	0.43
1:C:102:MET:HE2	1:C:102:MET:HB2	1.88	0.43
1:E:19:GLU:OE2	1:E:23:GLN:NE2	2.52	0.43
1:F:35:PHE:HB3	1:F:41:ARG:HG2	2.00	0.43
1:F:191:TRP:CE2	1:F:198:ARG:HD2	2.54	0.43
1:A:234:GLN:HE22	1:C:236:PHE:HA	1.83	0.43
1:F:3:ILE:HD11	1:F:73:ILE:HA	2.01	0.43
1:G:59:PRO:HD2	1:G:60:GLU:OE1	2.17	0.43
1:G:159:GLU:HB3	1:H:140:PRO:HG3	2.01	0.43
1:B:164:THR:HG23	2:B:266:HOH:O	2.19	0.43
1:F:206:PRO:O	1:G:167:ARG:NH1	2.51	0.43
1:H:141:LEU:HD23	1:H:141:LEU:HA	1.88	0.43
1:C:3:ILE:HD11	1:C:66:THR:HG21	2.00	0.43
1:D:71:GLU:OE2	1:D:71:GLU:HA	2.18	0.43
1:F:237:ALA:C	2:F:297:HOH:O	2.62	0.43
1:G:3:ILE:HD12	1:G:70:GLY:HA3	2.00	0.43
1:G:228:ALA:O	1:G:231:ILE:HG12	2.19	0.43
1:C:121:LEU:HG	1:C:127:ALA:CB	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:63:VAL:HG21	1:D:117:ALA:HB2	2.01	0.42
1:C:92:GLU:OE2	1:D:122:ARG:NH1	2.48	0.42
1:C:118:ILE:CD1	2:C:341:HOH:O	2.56	0.42
1:E:129:VAL:HB	1:E:172:LEU:HD12	2.01	0.42
1:G:122:ARG:HD3	1:G:168:ASP:OD2	2.19	0.42
1:A:122:ARG:NE	2:A:347:HOH:O	2.42	0.42
1:A:194:ASN:HD21	1:A:196:GLU:HB3	1.84	0.42
1:E:191:TRP:HA	1:E:197:LEU:HD23	2.00	0.42
1:E:222:PHE:HB2	1:H:222:PHE:CE1	2.54	0.42
1:H:84:ARG:HG3	1:H:85:PRO:HA	2.01	0.42
1:C:118:ILE:CG1	2:C:341:HOH:O	2.66	0.42
1:C:191:TRP:HA	1:C:197:LEU:HD23	2.02	0.42
1:F:35:PHE:CD1	1:F:53:ALA:HB1	2.53	0.42
1:D:3:ILE:HG13	1:D:73:ILE:HA	2.00	0.42
1:F:82:ILE:CD1	1:F:105:ALA:HB3	2.49	0.42
1:A:232:VAL:HG22	1:A:233:GLY:N	2.34	0.42
1:C:19:GLU:O	1:C:22:THR:HG22	2.19	0.42
1:D:187:PRO:HG2	1:D:190:ASP:HB2	2.02	0.42
1:D:228:ALA:O	1:D:231:ILE:HG12	2.20	0.42
1:E:158:VAL:HG13	1:E:172:LEU:HB3	2.01	0.42
1:B:230:PRO:HD3	1:D:215:GLU:CD	2.45	0.42
1:F:141:LEU:HD23	1:F:141:LEU:HA	1.91	0.42
1:F:167:ARG:NH2	1:G:202:ASP:OD2	2.53	0.42
1:C:9:ALA:HB2	1:C:30:CYS:HB3	2.01	0.41
1:E:33:ALA:HA	1:E:57:GLN:OE1	2.19	0.41
1:F:109:PHE:HB3	1:F:110:PRO:CD	2.50	0.41
1:B:135:SER:HB3	1:B:178:ASN:HB2	2.02	0.41
1:A:3:ILE:CD1	1:A:66:THR:HG22	2.51	0.41
1:D:191:TRP:CE2	1:D:198:ARG:HG2	2.55	0.41
1:F:205:VAL:O	1:F:206:PRO:C	2.63	0.41
1:H:207:LEU:HD23	1:H:207:LEU:HA	1.73	0.41
1:H:216:MET:HE2	1:H:238:PHE:CD2	2.55	0.41
1:A:108:ILE:HD12	1:A:108:ILE:N	2.33	0.41
1:H:82:ILE:HD12	1:H:83:PRO:HD2	2.03	0.41
1:C:207:LEU:HD22	1:C:209:ARG:CG	2.43	0.41
1:F:7:THR:O	1:F:8:HIS:C	2.63	0.41
1:G:136:VAL:HA	1:G:139:LYS:O	2.21	0.41
1:D:3:ILE:HD12	1:D:4:ALA:N	2.35	0.41
1:F:22:THR:HG23	1:F:23:GLN:N	2.36	0.41
1:A:237:ALA:H	1:C:234:GLN:HE22	1.66	0.41
1:C:112:LEU:HA	1:C:112:LEU:HD23	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:ALA:HB3	1:C:120:PRO:HD3	2.01	0.41
1:C:124:ALA:HB3	2:C:300:HOH:O	2.21	0.41
1:H:231:ILE:HB	1:H:234:GLN:NE2	2.36	0.41
1:A:58:LYS:NZ	1:A:61:ARG:HD2	2.30	0.41
1:B:58:LYS:O	1:B:62:LEU:HG	2.21	0.41
1:D:82:ILE:HD11	1:D:101:GLN:HG2	2.03	0.41
1:F:117:ALA:O	1:F:121:LEU:HB2	2.21	0.41
1:F:234:GLN:HE22	1:G:236:PHE:CA	2.29	0.41
1:G:146:LEU:C	1:G:149:PRO:HD2	2.45	0.41
1:H:109:PHE:HB3	1:H:110:PRO:HD3	2.03	0.41
1:B:121:LEU:HG	1:B:127:ALA:HB3	2.03	0.41
1:C:176:GLY:HA3	1:C:235:PHE:CZ	2.56	0.41
1:E:82:ILE:HD12	1:E:82:ILE:HA	1.93	0.40
1:E:112:LEU:HA	1:E:112:LEU:HD23	1.84	0.40
1:G:52:ILE:HG21	1:G:69:HIS:CD2	2.56	0.40
1:H:212:ARG:HB2	1:H:215:GLU:HG3	2.03	0.40
1:E:207:LEU:HD23	1:E:207:LEU:O	2.21	0.40
1:H:9:ALA:HB2	1:H:30:CYS:HB3	2.02	0.40
1:A:12:PHE:O	1:A:182:ASN:HB2	2.22	0.40
1:C:122:ARG:HD3	1:C:168:ASP:CG	2.46	0.40
1:C:3:ILE:HD12	1:C:4:ALA:N	2.37	0.40
1:C:83:PRO:O	1:C:86:MET:HB2	2.21	0.40
1:A:3:ILE:HA	1:A:27:THR:O	2.21	0.40
1:A:86:MET:O	1:A:102:MET:HE1	2.21	0.40
1:A:114:LEU:O	1:A:118:ILE:HG13	2.21	0.40
1:A:114:LEU:HG	1:A:131:PHE:HZ	1.86	0.40
1:B:136:VAL:HA	1:B:139:LYS:O	2.22	0.40
1:B:213:PRO:HG3	2:B:294:HOH:O	2.22	0.40
1:C:171:LEU:HB3	1:C:232:VAL:CG1	2.52	0.40
1:E:100:ARG:NH2	1:F:104:GLU:OE1	2.53	0.40
1:G:108:ILE:HD13	1:G:108:ILE:HA	1.86	0.40

There are no symmetry-related clashes.

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	241/244 (99%)	232 (96%)	9 (4%)	0	100	100
1	B	241/244 (99%)	226 (94%)	15 (6%)	0	100	100
1	C	241/244 (99%)	228 (95%)	13 (5%)	0	100	100
1	D	241/244 (99%)	227 (94%)	14 (6%)	0	100	100
1	E	241/244 (99%)	228 (95%)	13 (5%)	0	100	100
1	F	241/244 (99%)	229 (95%)	12 (5%)	0	100	100
1	G	241/244 (99%)	227 (94%)	14 (6%)	0	100	100
1	H	241/244 (99%)	231 (96%)	10 (4%)	0	100	100
All	All	1928/1952 (99%)	1828 (95%)	100 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	184/193 (95%)	168 (91%)	16 (9%)	9	6
1	B	184/193 (95%)	161 (88%)	23 (12%)	4	2
1	C	182/193 (94%)	166 (91%)	16 (9%)	9	6
1	D	181/193 (94%)	162 (90%)	19 (10%)	6	4
1	E	184/193 (95%)	167 (91%)	17 (9%)	8	5
1	F	183/193 (95%)	167 (91%)	16 (9%)	9	6
1	G	181/193 (94%)	167 (92%)	14 (8%)	12	8
1	H	182/193 (94%)	165 (91%)	17 (9%)	8	5
All	All	1461/1544 (95%)	1323 (91%)	138 (9%)	8	5

All (138) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ILE
1	A	19	GLU
1	A	21	LEU
1	A	22	THR
1	A	24	ASP
1	A	58	LYS
1	A	60	GLU
1	A	67	LEU
1	A	113	LEU
1	A	114	LEU
1	A	149	PRO
1	A	170	ILE
1	A	191	TRP
1	A	207	LEU
1	A	210	LEU
1	A	221	THR
1	B	15	PRO
1	B	21	LEU
1	B	23	GLN
1	B	24	ASP
1	B	58	LYS
1	B	60	GLU
1	B	61	ARG
1	B	67	LEU
1	B	85	PRO
1	B	91	LEU
1	B	113	LEU
1	B	114	LEU
1	B	121	LEU
1	B	138	LYS
1	B	149	PRO
1	B	175	ILE
1	B	191	TRP
1	B	206	PRO
1	B	207	LEU
1	B	210	LEU
1	B	213	PRO
1	B	220	ILE
1	B	243	LEU
1	C	3	ILE
1	C	19	GLU
1	C	21	LEU
1	C	23	GLN

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Mol	Chain	Res	Type
1	C	24	ASP
1	C	31	HIS
1	C	60	GLU
1	C	67	LEU
1	C	84	ARG
1	C	113	LEU
1	C	121	LEU
1	C	149	PRO
1	C	206	PRO
1	C	207	LEU
1	C	210	LEU
1	C	221	THR
1	D	3	ILE
1	D	21	LEU
1	D	23	GLN
1	D	60	GLU
1	D	67	LEU
1	D	82	ILE
1	D	85	PRO
1	D	87	ASN
1	D	91	LEU
1	D	107	SER
1	D	113	LEU
1	D	114	LEU
1	D	138	LYS
1	D	172	LEU
1	D	191	TRP
1	D	194	ASN
1	D	207	LEU
1	D	210	LEU
1	D	227	ARG
1	E	3	ILE
1	E	15	PRO
1	E	21	LEU
1	E	22	THR
1	E	23	GLN
1	E	24	ASP
1	E	60	GLU
1	E	61	ARG
1	E	67	LEU
1	E	90	PRO
1	E	100	ARG

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Mol	Chain	Res	Type
1	E	113	LEU
1	E	114	LEU
1	E	149	PRO
1	E	206	PRO
1	E	210	LEU
1	E	221	THR
1	F	3	ILE
1	F	23	GLN
1	F	58	LYS
1	F	60	GLU
1	F	61	ARG
1	F	67	LEU
1	F	71	GLU
1	F	85	PRO
1	F	91	LEU
1	F	113	LEU
1	F	114	LEU
1	F	121	LEU
1	F	138	LYS
1	F	198	ARG
1	F	207	LEU
1	F	210	LEU
1	G	21	LEU
1	G	23	GLN
1	G	24	ASP
1	G	45	GLU
1	G	60	GLU
1	G	84	ARG
1	G	85	PRO
1	G	91	LEU
1	G	100	ARG
1	G	113	LEU
1	G	114	LEU
1	G	191	TRP
1	G	207	LEU
1	G	210	LEU
1	H	3	ILE
1	H	21	LEU
1	H	22	THR
1	H	24	ASP
1	H	60	GLU
1	H	67	LEU

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Mol	Chain	Res	Type
1	H	108	ILE
1	H	113	LEU
1	H	114	LEU
1	H	118	ILE
1	H	149	PRO
1	H	172	LEU
1	H	191	TRP
1	H	196	GLU
1	H	207	LEU
1	H	210	LEU
1	H	221	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	23	GLN
1	A	42	GLN
1	A	101	GLN
1	A	193	ASN
1	A	194	ASN
1	B	48	ASN
1	B	101	GLN
1	B	193	ASN
1	C	42	GLN
1	C	69	HIS
1	C	234	GLN
1	D	23	GLN
1	D	48	ASN
1	D	87	ASN
1	D	181	ASN
1	D	193	ASN
1	D	194	ASN
1	E	23	GLN
1	E	42	GLN
1	E	69	HIS
1	E	101	GLN
1	E	193	ASN
1	E	194	ASN
1	F	23	GLN
1	F	42	GLN
1	F	181	ASN

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Mol	Chain	Res	Type
1	F	193	ASN
1	F	234	GLN
1	G	23	GLN
1	G	48	ASN
1	G	69	HIS
1	G	144	ASN
1	G	181	ASN
1	G	194	ASN
1	H	42	GLN
1	H	48	ASN
1	H	101	GLN
1	H	194	ASN

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	243/244 (99%)	0.04	1 (0%) 88 88	9, 18, 29, 44	0
1	B	243/244 (99%)	0.02	1 (0%) 88 88	8, 18, 31, 41	0
1	C	243/244 (99%)	0.09	1 (0%) 88 88	9, 19, 33, 42	0
1	D	243/244 (99%)	0.15	3 (1%) 76 76	10, 18, 33, 45	0
1	E	243/244 (99%)	0.12	2 (0%) 82 82	9, 18, 34, 46	0
1	F	243/244 (99%)	0.14	3 (1%) 76 76	9, 18, 32, 47	0
1	G	243/244 (99%)	0.07	2 (0%) 82 82	9, 18, 32, 38	0
1	H	243/244 (99%)	0.06	2 (0%) 82 82	9, 18, 29, 40	0
All	All	1944/1952 (99%)	0.09	15 (0%) 82 82	8, 18, 33, 47	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	191	TRP	3.8
1	H	191	TRP	3.8
1	D	191	TRP	3.7
1	B	191	TRP	3.6
1	D	194	ASN	3.4
1	G	191	TRP	3.3
1	C	143	TYR	2.5
1	F	197	LEU	2.3
1	E	39	ALA	2.2
1	D	197	LEU	2.2
1	F	22	THR	2.2
1	F	195	PRO	2.1
1	E	101	GLN	2.1
1	H	2	VAL	2.1
1	G	70	GLY	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.