



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 06:04 am BST

PDB ID : 6GJV
Title : apo-structure of IMPDH from Pseudomonas aeruginosa
Authors : Labesse, G.; Haouz, A.; Alexandre, T.; Munier-Lehmann, H.
Deposited on : 2018-05-17
Resolution : 2.11 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

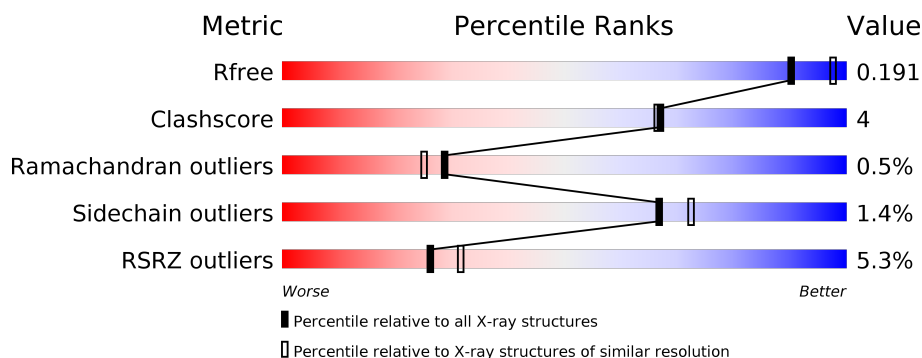
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	509	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>10%</div> <div>17%</div> </div> </div>
1	B	509	<div> <div>14%</div> <div> <div></div> <div>70%</div> <div>11%</div> <div>19%</div> </div> </div>
1	C	509	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>8%</div> <div>16%</div> </div> </div>
1	D	509	<div> <div>2%</div> <div> <div></div> <div>58%</div> <div>38%</div> </div> </div>
1	E	509	<div> <div>0%</div> <div> <div></div> <div>77%</div> <div>6%</div> <div>17%</div> </div> </div>
1	F	509	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>8%</div> <div>16%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	509	<div><div>%</div><div><div></div><div>77%</div><div>6%</div><div>17%</div></div></div>
1	H	509	<div><div>%</div><div><div></div><div>57%</div><div>5%</div><div>38%</div></div></div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 25755 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 Inosine-5'-monophosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	425	Total	C	N	O	S	0	1	0
			3150	1988	549	596	17			
1	B	414	Total	C	N	O	S	0	1	0
			3038	1916	531	574	17			
1	C	429	Total	C	N	O	S	0	0	0
			3182	2005	559	601	17			
1	D	318	Total	C	N	O	S	0	0	0
			2315	1453	413	436	13			
1	E	424	Total	C	N	O	S	1	6	0
			3183	2007	558	601	17			
1	F	426	Total	C	N	O	S	0	3	0
			3178	2004	558	599	17			
1	G	424	Total	C	N	O	S	0	3	0
			3172	1999	555	601	17			
1	H	317	Total	C	N	O	S	0	1	0
			2314	1451	414	436	13			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP Q9HXM5
A	-18	GLY	-	expression tag	UNP Q9HXM5
A	-17	SER	-	expression tag	UNP Q9HXM5
A	-16	SER	-	expression tag	UNP Q9HXM5
A	-15	HIS	-	expression tag	UNP Q9HXM5
A	-14	HIS	-	expression tag	UNP Q9HXM5
A	-13	HIS	-	expression tag	UNP Q9HXM5
A	-12	HIS	-	expression tag	UNP Q9HXM5
A	-11	HIS	-	expression tag	UNP Q9HXM5
A	-10	HIS	-	expression tag	UNP Q9HXM5
A	-9	SER	-	expression tag	UNP Q9HXM5
A	-8	SER	-	expression tag	UNP Q9HXM5
A	-7	GLY	-	expression tag	UNP Q9HXM5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	expression tag	UNP Q9HXM5
A	-5	VAL	-	expression tag	UNP Q9HXM5
A	-4	PRO	-	expression tag	UNP Q9HXM5
A	-3	ARG	-	expression tag	UNP Q9HXM5
A	-2	GLY	-	expression tag	UNP Q9HXM5
A	-1	SER	-	expression tag	UNP Q9HXM5
A	0	HIS	-	expression tag	UNP Q9HXM5
B	-19	MET	-	initiating methionine	UNP Q9HXM5
B	-18	GLY	-	expression tag	UNP Q9HXM5
B	-17	SER	-	expression tag	UNP Q9HXM5
B	-16	SER	-	expression tag	UNP Q9HXM5
B	-15	HIS	-	expression tag	UNP Q9HXM5
B	-14	HIS	-	expression tag	UNP Q9HXM5
B	-13	HIS	-	expression tag	UNP Q9HXM5
B	-12	HIS	-	expression tag	UNP Q9HXM5
B	-11	HIS	-	expression tag	UNP Q9HXM5
B	-10	HIS	-	expression tag	UNP Q9HXM5
B	-9	SER	-	expression tag	UNP Q9HXM5
B	-8	SER	-	expression tag	UNP Q9HXM5
B	-7	GLY	-	expression tag	UNP Q9HXM5
B	-6	LEU	-	expression tag	UNP Q9HXM5
B	-5	VAL	-	expression tag	UNP Q9HXM5
B	-4	PRO	-	expression tag	UNP Q9HXM5
B	-3	ARG	-	expression tag	UNP Q9HXM5
B	-2	GLY	-	expression tag	UNP Q9HXM5
B	-1	SER	-	expression tag	UNP Q9HXM5
B	0	HIS	-	expression tag	UNP Q9HXM5
C	-19	MET	-	initiating methionine	UNP Q9HXM5
C	-18	GLY	-	expression tag	UNP Q9HXM5
C	-17	SER	-	expression tag	UNP Q9HXM5
C	-16	SER	-	expression tag	UNP Q9HXM5
C	-15	HIS	-	expression tag	UNP Q9HXM5
C	-14	HIS	-	expression tag	UNP Q9HXM5
C	-13	HIS	-	expression tag	UNP Q9HXM5
C	-12	HIS	-	expression tag	UNP Q9HXM5
C	-11	HIS	-	expression tag	UNP Q9HXM5
C	-10	HIS	-	expression tag	UNP Q9HXM5
C	-9	SER	-	expression tag	UNP Q9HXM5
C	-8	SER	-	expression tag	UNP Q9HXM5
C	-7	GLY	-	expression tag	UNP Q9HXM5
C	-6	LEU	-	expression tag	UNP Q9HXM5
C	-5	VAL	-	expression tag	UNP Q9HXM5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	PRO	-	expression tag	UNP Q9HXM5
C	-3	ARG	-	expression tag	UNP Q9HXM5
C	-2	GLY	-	expression tag	UNP Q9HXM5
C	-1	SER	-	expression tag	UNP Q9HXM5
C	0	HIS	-	expression tag	UNP Q9HXM5
D	-19	MET	-	initiating methionine	UNP Q9HXM5
D	-18	GLY	-	expression tag	UNP Q9HXM5
D	-17	SER	-	expression tag	UNP Q9HXM5
D	-16	SER	-	expression tag	UNP Q9HXM5
D	-15	HIS	-	expression tag	UNP Q9HXM5
D	-14	HIS	-	expression tag	UNP Q9HXM5
D	-13	HIS	-	expression tag	UNP Q9HXM5
D	-12	HIS	-	expression tag	UNP Q9HXM5
D	-11	HIS	-	expression tag	UNP Q9HXM5
D	-10	HIS	-	expression tag	UNP Q9HXM5
D	-9	SER	-	expression tag	UNP Q9HXM5
D	-8	SER	-	expression tag	UNP Q9HXM5
D	-7	GLY	-	expression tag	UNP Q9HXM5
D	-6	LEU	-	expression tag	UNP Q9HXM5
D	-5	VAL	-	expression tag	UNP Q9HXM5
D	-4	PRO	-	expression tag	UNP Q9HXM5
D	-3	ARG	-	expression tag	UNP Q9HXM5
D	-2	GLY	-	expression tag	UNP Q9HXM5
D	-1	SER	-	expression tag	UNP Q9HXM5
D	0	HIS	-	expression tag	UNP Q9HXM5
E	-19	MET	-	initiating methionine	UNP Q9HXM5
E	-18	GLY	-	expression tag	UNP Q9HXM5
E	-17	SER	-	expression tag	UNP Q9HXM5
E	-16	SER	-	expression tag	UNP Q9HXM5
E	-15	HIS	-	expression tag	UNP Q9HXM5
E	-14	HIS	-	expression tag	UNP Q9HXM5
E	-13	HIS	-	expression tag	UNP Q9HXM5
E	-12	HIS	-	expression tag	UNP Q9HXM5
E	-11	HIS	-	expression tag	UNP Q9HXM5
E	-10	HIS	-	expression tag	UNP Q9HXM5
E	-9	SER	-	expression tag	UNP Q9HXM5
E	-8	SER	-	expression tag	UNP Q9HXM5
E	-7	GLY	-	expression tag	UNP Q9HXM5
E	-6	LEU	-	expression tag	UNP Q9HXM5
E	-5	VAL	-	expression tag	UNP Q9HXM5
E	-4	PRO	-	expression tag	UNP Q9HXM5
E	-3	ARG	-	expression tag	UNP Q9HXM5

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	expression tag	UNP Q9HXM5
E	-1	SER	-	expression tag	UNP Q9HXM5
E	0	HIS	-	expression tag	UNP Q9HXM5
F	-19	MET	-	initiating methionine	UNP Q9HXM5
F	-18	GLY	-	expression tag	UNP Q9HXM5
F	-17	SER	-	expression tag	UNP Q9HXM5
F	-16	SER	-	expression tag	UNP Q9HXM5
F	-15	HIS	-	expression tag	UNP Q9HXM5
F	-14	HIS	-	expression tag	UNP Q9HXM5
F	-13	HIS	-	expression tag	UNP Q9HXM5
F	-12	HIS	-	expression tag	UNP Q9HXM5
F	-11	HIS	-	expression tag	UNP Q9HXM5
F	-10	HIS	-	expression tag	UNP Q9HXM5
F	-9	SER	-	expression tag	UNP Q9HXM5
F	-8	SER	-	expression tag	UNP Q9HXM5
F	-7	GLY	-	expression tag	UNP Q9HXM5
F	-6	LEU	-	expression tag	UNP Q9HXM5
F	-5	VAL	-	expression tag	UNP Q9HXM5
F	-4	PRO	-	expression tag	UNP Q9HXM5
F	-3	ARG	-	expression tag	UNP Q9HXM5
F	-2	GLY	-	expression tag	UNP Q9HXM5
F	-1	SER	-	expression tag	UNP Q9HXM5
F	0	HIS	-	expression tag	UNP Q9HXM5
G	-19	MET	-	initiating methionine	UNP Q9HXM5
G	-18	GLY	-	expression tag	UNP Q9HXM5
G	-17	SER	-	expression tag	UNP Q9HXM5
G	-16	SER	-	expression tag	UNP Q9HXM5
G	-15	HIS	-	expression tag	UNP Q9HXM5
G	-14	HIS	-	expression tag	UNP Q9HXM5
G	-13	HIS	-	expression tag	UNP Q9HXM5
G	-12	HIS	-	expression tag	UNP Q9HXM5
G	-11	HIS	-	expression tag	UNP Q9HXM5
G	-10	HIS	-	expression tag	UNP Q9HXM5
G	-9	SER	-	expression tag	UNP Q9HXM5
G	-8	SER	-	expression tag	UNP Q9HXM5
G	-7	GLY	-	expression tag	UNP Q9HXM5
G	-6	LEU	-	expression tag	UNP Q9HXM5
G	-5	VAL	-	expression tag	UNP Q9HXM5
G	-4	PRO	-	expression tag	UNP Q9HXM5
G	-3	ARG	-	expression tag	UNP Q9HXM5
G	-2	GLY	-	expression tag	UNP Q9HXM5
G	-1	SER	-	expression tag	UNP Q9HXM5

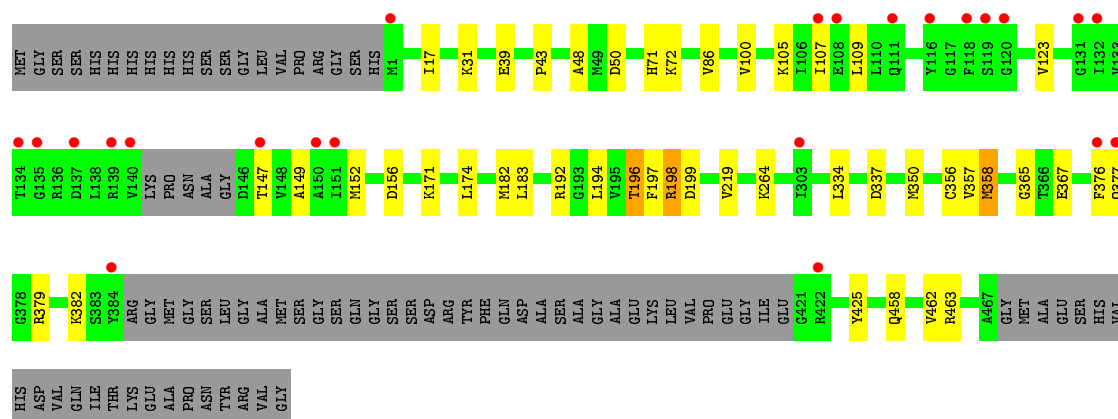
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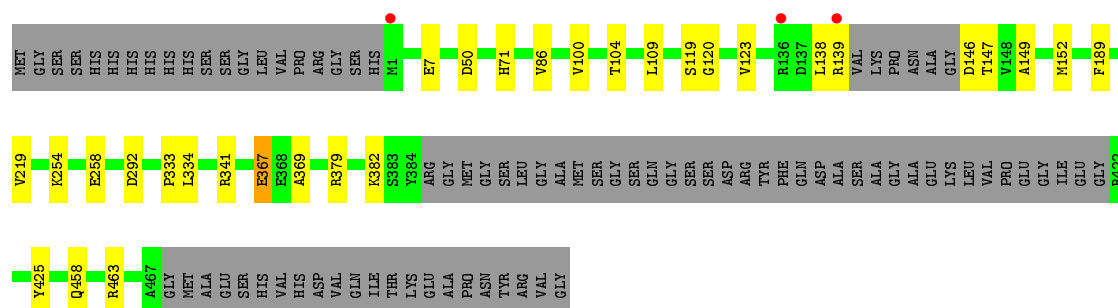
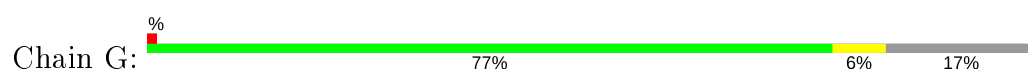
Chain	Residue	Modelled	Actual	Comment	Reference
G	0	HIS	-	expression tag	UNP Q9HXM5
H	-19	MET	-	initiating methionine	UNP Q9HXM5
H	-18	GLY	-	expression tag	UNP Q9HXM5
H	-17	SER	-	expression tag	UNP Q9HXM5
H	-16	SER	-	expression tag	UNP Q9HXM5
H	-15	HIS	-	expression tag	UNP Q9HXM5
H	-14	HIS	-	expression tag	UNP Q9HXM5
H	-13	HIS	-	expression tag	UNP Q9HXM5
H	-12	HIS	-	expression tag	UNP Q9HXM5
H	-11	HIS	-	expression tag	UNP Q9HXM5
H	-10	HIS	-	expression tag	UNP Q9HXM5
H	-9	SER	-	expression tag	UNP Q9HXM5
H	-8	SER	-	expression tag	UNP Q9HXM5
H	-7	GLY	-	expression tag	UNP Q9HXM5
H	-6	LEU	-	expression tag	UNP Q9HXM5
H	-5	VAL	-	expression tag	UNP Q9HXM5
H	-4	PRO	-	expression tag	UNP Q9HXM5
H	-3	ARG	-	expression tag	UNP Q9HXM5
H	-2	GLY	-	expression tag	UNP Q9HXM5
H	-1	SER	-	expression tag	UNP Q9HXM5
H	0	HIS	-	expression tag	UNP Q9HXM5

- Molecule 2 is water.

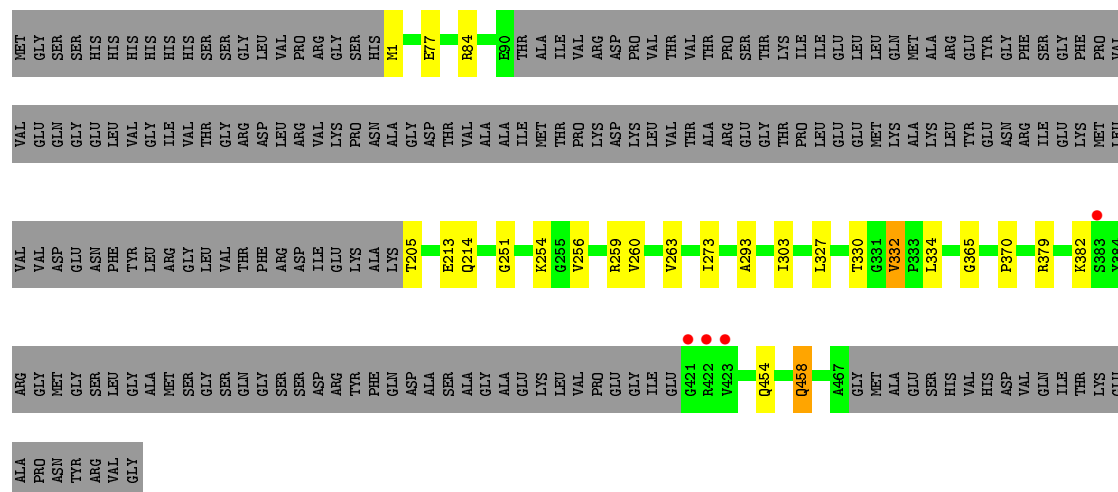
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	240	Total	O	0	0
			240	240		
2	B	256	Total	O	0	0
			256	256		
2	C	307	Total	O	0	0
			307	307		
2	D	211	Total	O	0	0
			211	211		
2	E	350	Total	O	0	0
			350	350		
2	F	289	Total	O	0	0
			289	289		
2	G	315	Total	O	0	0
			315	315		
2	H	255	Total	O	0	0
			255	255		



- Molecule 1: Inosine-5'-monophosphate dehydrogenase



- Molecule 1: Inosine-5'-monophosphate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	115.30Å 194.64Å 122.92Å 90.00° 117.75° 90.00°	Depositor
Resolution (Å)	44.63 – 2.11 49.35 – 2.11	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.63-2.11) 100.0 (49.35-2.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.160 , 0.191 0.159 , 0.191	Depositor DCC
R_{free} test set	13805 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	33.3	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.016 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	25755	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.38	0/3193	0.57	1/4319 (0.0%)
1	B	0.36	0/3078	0.56	1/4158 (0.0%)
1	C	0.38	1/3222 (0.0%)	0.57	0/4355
1	D	0.37	0/2343	0.55	0/3166
1	E	0.38	0/3246	0.57	0/4385
1	F	0.36	0/3227	0.55	1/4361 (0.0%)
1	G	0.40	0/3218	0.57	0/4348
1	H	0.41	0/2346	0.57	0/3169
All	All	0.38	1/23873 (0.0%)	0.57	3/32261 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	356	CYS	CB-SG	-5.09	1.73	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	359	MET	CA-CB-CG	5.56	122.76	113.30
1	B	102	PRO	N-CA-CB	5.52	109.93	103.30
1	F	358	MET	CG-SD-CE	5.48	108.96	100.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3150	0	3241	33	0
1	B	3038	0	3090	34	0
1	C	3182	0	3280	25	0
1	D	2315	0	2384	10	0
1	E	3183	0	3284	21	0
1	F	3178	0	3288	32	0
1	G	3172	0	3273	19	0
1	H	2314	0	2380	18	0
2	A	240	0	0	2	0
2	B	256	0	0	2	0
2	C	307	0	0	10	0
2	D	211	0	0	1	0
2	E	350	0	0	3	0
2	F	289	0	0	6	0
2	G	315	0	0	2	0
2	H	255	0	0	5	0
All	All	25755	0	24220	186	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 4.

All (186) 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:455:MET:HE3	1:A:459:PRO:HG2	1.48	0.92
1:B:7:GLU:OE2	1:B:463:ARG:NH2	2.09	0.86
1:F:350:MET:HE3	1:F:357:VAL:HG23	1.60	0.84
1:C:425:TYR:O	2:C:501:HOH:O	1.96	0.83
1:A:455:MET:CE	1:A:459:PRO:HG2	2.09	0.82
1:F:123:VAL:HG11	1:F:152:MET:HE2	1.65	0.78
1:B:198:ARG:HH11	1:B:202:LYS:HD3	1.49	0.78
1:A:90:GLU:HG3	1:A:95:ARG:HH21	1.49	0.78
1:F:350:MET:CE	1:F:357:VAL:HG23	2.15	0.76
1:F:196:THR:HG22	1:F:199:ASP:H	1.53	0.74
1:B:181:LYS:HB3	1:B:194:LEU:HD11	1.69	0.73
1:D:365:GLY:HA2	1:D:382:LYS:HG3	1.71	0.72
1:E:155:LYS:HA	1:E:158:LEU:HD22	1.72	0.72
1:C:426:LYS:HA	2:C:501:HOH:O	1.91	0.70
1:C:261:ARG:NH1	2:C:504:HOH:O	2.24	0.69
1:A:365:GLY:HA2	1:A:382:LYS:HG3	1.72	0.69
1:F:358:MET:HE3	2:F:580:HOH:O	1.93	0.68
1:F:337:ASP:OD2	1:F:358:MET:HE1	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:367:GLU:HG2	1:A:425:TYR:OH	1.93	0.68
1:D:367:GLU:HG2	1:D:425:TYR:OH	1.94	0.68
1:A:111:GLN:OE1	1:A:114:ARG:NH1	2.27	0.68
1:C:365:GLY:HA2	1:C:382:LYS:HG3	1.77	0.67
1:B:7:GLU:CD	1:B:463:ARG:HH22	1.98	0.67
1:B:108:GLU:O	1:B:112:MET:HG3	1.95	0.66
1:C:422:ARG:N	2:C:508:HOH:O	2.28	0.66
1:H:379:ARG:NH1	2:H:503:HOH:O	2.29	0.64
1:A:90:GLU:O	1:A:91:THR:OG1	2.16	0.64
1:A:1:MET:N	2:A:502:HOH:O	2.29	0.63
1:A:175:TYR:HA	1:B:118:PHE:HE2	1.62	0.63
1:B:379:ARG:NH2	1:B:381:TYR:OH	2.31	0.63
1:C:214:GLN:OE1	1:C:216:ARG:NH1	2.32	0.62
1:F:105:LYS:HD3	1:F:147:THR:HG22	1.80	0.61
1:F:196:THR:HG21	2:F:743:HOH:O	2.01	0.61
1:B:198:ARG:NH1	1:B:202:LYS:HD3	2.15	0.61
1:E:97:PRO:HG2	1:E:191:LEU:HD21	1.83	0.61
1:A:86:VAL:HG11	1:A:219:VAL:HB	1.83	0.60
1:A:285:ALA:O	1:A:288:GLU:HG3	2.01	0.60
1:C:78:GLN:HG3	2:C:737:HOH:O	2.00	0.60
1:E:216:ARG:NH2	2:E:504:HOH:O	2.30	0.59
2:C:501:HOH:O	1:E:380:SER:N	2.35	0.59
1:F:149:ALA:HA	1:F:152:MET:HE3	1.84	0.59
1:F:365:GLY:HA2	1:F:382:LYS:HG3	1.84	0.59
1:F:198:ARG:NH1	2:F:504:HOH:O	2.36	0.58
1:C:384:TYR:O	2:C:502:HOH:O	2.17	0.58
1:F:367:GLU:HG2	1:F:425:TYR:OH	2.04	0.58
1:A:340:ILE:HG23	1:A:345:ASP:HB2	1.86	0.58
1:E:123:VAL:HG11	1:E:152:MET:HE1	1.87	0.57
1:D:2:LEU:HD13	1:D:4:ILE:HD11	1.87	0.57
1:H:365:GLY:HA2	1:H:382:LYS:HG3	1.86	0.56
1:G:7:GLU:OE1	1:G:463:ARG:NH2	2.30	0.56
1:G:369:ALA:O	1:G:382:LYS:NZ	2.38	0.56
1:B:77:GLU:OE2	1:B:84:ARG:NH1	2.39	0.56
1:C:123:VAL:HG21	1:C:152:MET:SD	2.46	0.56
1:B:7:GLU:CD	1:B:463:ARG:NH2	2.59	0.55
1:H:327:LEU:HD13	1:H:332:VAL:HG22	1.88	0.54
1:F:337:ASP:CG	1:F:358:MET:HE1	2.28	0.54
1:G:367[A]:GLU:HG2	1:G:425:TYR:OH	2.07	0.54
1:E:365:GLY:HA2	1:E:382:LYS:HG3	1.90	0.54
1:H:77:GLU:OE1	1:H:84:ARG:NH2	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:374:GLU:OE1	1:C:383:SER:HB3	2.08	0.53
1:D:303:ILE:O	1:D:303:ILE:HG22	2.09	0.53
1:C:273:ILE:HG12	1:C:293:ALA:HB3	1.90	0.53
1:C:50:ASP:HA	1:C:71:HIS:CD2	2.44	0.53
1:C:374:GLU:OE2	1:C:422:ARG:NE	2.42	0.52
1:H:454:GLN:HG2	2:H:714:HOH:O	2.09	0.52
1:H:303:ILE:O	1:H:303:ILE:CD1	2.58	0.52
1:B:50:ASP:HA	1:B:71:HIS:CD2	2.45	0.51
1:E:50:ASP:HA	1:E:71:HIS:CD2	2.45	0.51
1:G:119:SER:N	1:G:120[A]:GLY:HA2	2.25	0.51
1:D:50:ASP:HA	1:D:71:HIS:CD2	2.46	0.51
1:A:351:VAL:HG22	1:A:455:MET:CE	2.41	0.51
1:G:254:LYS:HE3	1:G:258:GLU:OE2	2.11	0.51
1:C:0:HIS:N	2:C:512:HOH:O	2.43	0.51
1:B:86:VAL:HG11	1:B:219:VAL:HB	1.93	0.50
1:B:365:GLY:HA2	1:B:382:LYS:HG3	1.93	0.50
1:A:302:SER:HB3	1:A:303:ILE:HD12	1.92	0.50
1:F:50:ASP:HA	1:F:71:HIS:CD2	2.47	0.49
1:G:119:SER:N	1:G:120[B]:GLY:HA2	2.27	0.49
1:A:340:ILE:HD12	1:A:357:VAL:HG11	1.95	0.49
1:G:123:VAL:HG11	1:G:152:MET:HE1	1.93	0.49
1:C:379:ARG:HD3	1:E:379:ARG:HH12	1.78	0.49
1:F:48:ALA:HB1	1:F:71:HIS:HA	1.94	0.49
1:G:146:ASP:N	1:G:146:ASP:OD1	2.45	0.49
1:A:2:LEU:HD13	1:A:4:ILE:HD11	1.95	0.48
1:D:273:ILE:HG12	1:D:293:ALA:HB3	1.94	0.48
1:G:50:ASP:HA	1:G:71:HIS:CD2	2.48	0.48
1:A:351:VAL:HG22	1:A:455:MET:HE2	1.94	0.48
1:A:374:GLU:OE2	1:A:422:ARG:HD3	2.14	0.48
1:C:187:GLU:CD	1:C:187:GLU:H	2.16	0.48
1:B:166:PRO:HB2	1:B:169:GLU:HG3	1.95	0.48
1:C:377:GLN:NE2	1:C:381:TYR:OH	2.46	0.48
1:F:358:MET:HE2	2:F:694:HOH:O	2.13	0.48
1:G:109:LEU:HD22	1:G:138:LEU:HD21	1.95	0.48
1:G:86:VAL:HG11	1:G:219:VAL:HB	1.95	0.48
1:B:1:MET:N	2:B:507:HOH:O	2.42	0.48
1:F:100:VAL:HG11	1:F:109:LEU:HD23	1.96	0.48
1:H:370:PRO:O	1:H:382:LYS:HE3	2.14	0.48
1:F:86:VAL:HG11	1:F:219:VAL:HB	1.95	0.47
1:H:330:THR:OG1	1:H:332:VAL:HG13	2.13	0.47
1:A:50:ASP:HA	1:A:71:HIS:CD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:501:HOH:O	1:E:380:SER:OG	2.16	0.47
1:H:303:ILE:HD12	1:H:303:ILE:C	2.35	0.47
1:B:187:GLU:H	1:B:187:GLU:CD	2.18	0.47
1:B:2:LEU:HD13	1:B:4:ILE:HD11	1.97	0.47
1:F:17:ILE:HD11	1:F:462:VAL:HG13	1.97	0.47
1:B:132:ILE:HG22	1:B:158:LEU:HD21	1.95	0.46
1:A:100:VAL:HG23	1:A:123:VAL:HG22	1.97	0.46
1:A:90:GLU:HG3	1:A:95:ARG:NH2	2.24	0.46
1:F:31:LYS:HE2	1:F:39:GLU:OE1	2.15	0.46
1:F:379:ARG:HD3	1:G:379:ARG:NH2	2.30	0.46
1:B:128:GLU:N	2:B:512:HOH:O	2.48	0.46
1:D:86:VAL:HG11	1:D:219:VAL:HB	1.97	0.46
1:E:105:LYS:HD3	1:E:147:THR:HG22	1.98	0.46
1:F:264:LYS:HA	2:F:531:HOH:O	2.15	0.46
1:F:43:PRO:HB2	1:F:356:CYS:HA	1.96	0.46
1:A:156:ASP:N	1:A:156:ASP:OD1	2.46	0.46
1:F:183:LEU:HD23	1:F:194:LEU:HD13	1.98	0.45
1:C:304:CYS:HB2	2:C:717:HOH:O	2.15	0.45
1:C:375:LEU:CD2	1:E:432:ILE:HG12	2.47	0.45
1:B:155:LYS:HE2	1:B:189:PHE:CD2	2.52	0.45
1:E:86:VAL:HG11	1:E:219:VAL:HB	1.97	0.45
1:B:340:ILE:HG23	1:B:345:ASP:HB2	1.98	0.45
1:H:256:VAL:O	1:H:260:VAL:HG13	2.16	0.45
1:F:196:THR:CG2	1:F:199:ASP:H	2.24	0.45
1:B:284:LYS:O	1:B:288:GLU:HG3	2.17	0.44
1:B:132:ILE:HG23	1:B:153:THR:HB	1.99	0.44
1:A:384:TYR:O	2:A:501:HOH:O	2.20	0.44
1:B:91:THR:HG22	1:B:92:ALA:H	1.82	0.44
1:C:17:ILE:HD11	1:C:462:VAL:HG13	2.00	0.44
1:H:205:THR:N	2:H:517:HOH:O	2.51	0.44
1:E:107:ILE:HG12	2:E:688:HOH:O	2.17	0.44
1:H:273:ILE:HG12	1:H:293:ALA:HB3	1.99	0.44
1:B:306:THR:HG22	1:B:312:VAL:O	2.18	0.43
1:D:205:THR:HG22	1:D:207:PRO:HD3	2.00	0.43
1:E:139:ARG:NH1	2:E:512:HOH:O	2.47	0.43
1:H:251:GLY:HA2	1:H:256:VAL:HG21	2.00	0.43
1:E:22:GLU:HG3	1:H:254:LYS:HB2	2.00	0.43
1:A:110:LEU:HD11	1:A:138:LEU:HD22	2.00	0.43
1:B:370:PRO:O	1:B:382:LYS:HE3	2.19	0.43
1:C:109:LEU:HD22	1:C:138:LEU:HD21	1.99	0.43
1:F:376:PHE:CE2	1:F:377:GLN:HG3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:VAL:HG23	1:A:152:MET:HE1	2.00	0.43
1:C:340:ILE:HG23	1:C:345:ASP:HB2	1.99	0.43
1:F:105:LYS:HB3	1:F:107:ILE:HG22	2.01	0.43
1:E:273:ILE:HG12	1:E:293:ALA:HB3	2.01	0.43
1:G:341:ARG:HB2	1:G:341:ARG:HE	1.63	0.43
1:F:174:LEU:HD21	1:F:182:MET:HB2	2.01	0.42
1:A:131:GLY:HA2	1:A:158:LEU:HD11	2.00	0.42
1:A:192:ARG:HA	1:A:192:ARG:HD3	1.78	0.42
1:A:2:LEU:HD11	1:C:317:ILE:HG22	2.01	0.42
1:D:17:ILE:HD11	1:D:462:VAL:HG13	2.01	0.42
1:G:149:ALA:HA	1:G:152:MET:CE	2.50	0.42
1:B:259:ARG:O	1:B:263:VAL:HG23	2.20	0.42
1:C:174:LEU:HD21	1:C:182:MET:HB2	2.01	0.42
1:E:149:ALA:HA	1:E:152:MET:HE2	2.02	0.42
1:E:156:ASP:OD1	1:E:156:ASP:N	2.53	0.42
1:F:192:ARG:HA	1:F:192:ARG:HD3	1.83	0.42
1:A:340:ILE:HD12	1:A:357:VAL:CG1	2.50	0.41
1:C:131:GLY:HA2	1:C:158:LEU:HD11	2.02	0.41
1:H:254:LYS:NZ	2:H:520:HOH:O	2.53	0.41
1:A:303:ILE:HG22	1:A:303:ILE:O	2.20	0.41
1:A:165:THR:HG22	1:A:169:GLU:HB3	2.02	0.41
1:F:72:LYS:HE3	2:F:693:HOH:O	2.20	0.41
1:E:149:ALA:HA	1:E:152:MET:CE	2.50	0.41
1:G:292:ASP:O	1:G:333:PRO:HD2	2.21	0.41
1:F:171:LYS:HE3	1:F:197:PHE:CZ	2.56	0.41
1:G:100:VAL:HG22	1:G:104:THR:CB	2.50	0.41
1:H:458:GLN:HG3	2:H:714:HOH:O	2.21	0.41
1:B:264:LYS:HA	1:B:264:LYS:HD2	1.82	0.41
1:B:129:LEU:HB2	1:B:191:LEU:HB2	2.03	0.41
1:B:458:GLN:HB2	1:B:459:PRO:HD3	2.03	0.41
1:G:120[B]:GLY:HA2	2:G:681:HOH:O	2.20	0.41
1:E:346:LEU:HG	1:E:357:VAL:HG21	2.03	0.40
1:G:120[A]:GLY:HA2	2:G:681:HOH:O	2.20	0.40
1:G:146:ASP:HB2	1:G:147:THR:H	1.69	0.40
1:B:273:ILE:HG12	1:B:293:ALA:HB3	2.03	0.40
1:B:327:LEU:HD23	1:B:327:LEU:HA	1.86	0.40
1:D:23:VAL:HB	2:D:525:HOH:O	2.20	0.40
1:E:76:ILE:HD13	1:E:231:GLU:HG2	2.03	0.40
1:B:106:ILE:HD11	1:B:151:ILE:HD13	2.02	0.40
1:A:444:MET:HG2	1:A:455:MET:SD	2.62	0.40
1:B:198:ARG:O	1:B:202:LYS:HG2	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:149:ALA:HA	1:F:152:MET:CE	2.51	0.40
1:A:17:ILE:HD11	1:A:462:VAL:HG13	2.03	0.40
1:H:213:GLU:OE2	1:H:214:GLN:HG3	2.22	0.40
1:H:259:ARG:O	1:H:263:VAL:HG23	2.22	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	420/509 (82%)	404 (96%)	13 (3%)	3 (1%)	22	17
1	B	405/509 (80%)	386 (95%)	13 (3%)	6 (2%)	10	5
1	C	423/509 (83%)	410 (97%)	10 (2%)	3 (1%)	22	17
1	D	312/509 (61%)	304 (97%)	8 (3%)	0	100	100
1	E	424/509 (83%)	415 (98%)	9 (2%)	0	100	100
1	F	423/509 (83%)	411 (97%)	11 (3%)	1 (0%)	47	48
1	G	421/509 (83%)	410 (97%)	10 (2%)	1 (0%)	47	48
1	H	312/509 (61%)	304 (97%)	7 (2%)	1 (0%)	41	40
All	All	3140/4072 (77%)	3044 (97%)	81 (3%)	15 (0%)	29	25

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	THR
1	B	327	LEU
1	C	146	ASP
1	A	92	ALA
1	B	328	GLU
1	B	423	VAL

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Mol	Chain	Res	Type
1	B	89	HIS
1	B	103	SER
1	A	120	GLY
1	F	458	GLN
1	C	145	GLY
1	B	93	ILE
1	C	458	GLN
1	H	458	GLN
1	G	458	GLN

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	324/390 (83%)	320 (99%)	4 (1%)	71	77
1	B	305/390 (78%)	301 (99%)	4 (1%)	69	74
1	C	327/390 (84%)	322 (98%)	5 (2%)	65	70
1	D	233/390 (60%)	229 (98%)	4 (2%)	60	66
1	E	331/390 (85%)	326 (98%)	5 (2%)	65	70
1	F	329/390 (84%)	324 (98%)	5 (2%)	65	70
1	G	328/390 (84%)	323 (98%)	5 (2%)	65	70
1	H	233/390 (60%)	230 (99%)	3 (1%)	69	74
All	All	2410/3120 (77%)	2375 (98%)	35 (2%)	67	70

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	189	PHE
1	A	302	SER
1	A	334	LEU
1	A	341	ARG
1	B	2	LEU
1	B	189	PHE

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Mol	Chain	Res	Type
1	B	205	THR
1	B	334	LEU
1	C	1	MET
1	C	2	LEU
1	C	96	ASP
1	C	187	GLU
1	C	334	LEU
1	D	2	LEU
1	D	204	LYS
1	D	334	LEU
1	D	376	PHE
1	E	158	LEU
1	E	189	PHE
1	E	265[A]	GLN
1	E	265[B]	GLN
1	E	334	LEU
1	F	156	ASP
1	F	196	THR
1	F	198	ARG
1	F	334	LEU
1	F	463	ARG
1	G	139	ARG
1	G	189	PHE
1	G	334	LEU
1	G	367[A]	GLU
1	G	367[B]	GLU
1	H	1	MET
1	H	332	VAL
1	H	334	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	377	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 carbohydrates 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	425/509 (83%)	0.21	27 (6%) 19 23	23, 36, 98, 133	22 (5%)
1	B	414/509 (81%)	0.61	71 (17%) 1 1	23, 40, 90, 108	107 (25%)
1	C	429/509 (84%)	0.07	25 (5%) 23 28	24, 35, 78, 137	19 (4%)
1	D	318/509 (62%)	-0.19	11 (3%) 44 50	27, 38, 69, 113	19 (5%)
1	E	424/509 (83%)	-0.21	4 (0%) 84 86	21, 34, 61, 85	22 (5%)
1	F	426/509 (83%)	-0.03	23 (5%) 25 31	26, 41, 71, 102	24 (5%)
1	G	424/509 (83%)	-0.33	3 (0%) 87 89	24, 36, 66, 97	19 (4%)
1	H	317/509 (62%)	-0.07	4 (1%) 77 80	23, 33, 63, 97	20 (6%)
All	All	3177/4072 (78%)	0.02	168 (5%) 26 31	21, 37, 81, 137	252 (7%)

All (168) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	174	LEU	8.4
1	B	151	ILE	8.2
1	B	109	LEU	8.0
1	B	118	PHE	7.9
1	A	175	TYR	7.6
1	A	172	ALA	7.5
1	B	104	THR	6.7
1	B	102	PRO	6.6
1	B	129	LEU	6.4
1	B	189	PHE	6.4
1	B	150	ALA	6.3
1	B	119	SER	6.2
1	B	113	ALA	6.1
1	B	111	GLN	6.1
1	B	130	VAL	6.1
1	B	191	LEU	5.8

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Mol	Chain	Res	Type	RSRZ
1	B	105	LYS	5.6
1	C	145	GLY	5.6
1	C	144	ALA	5.5
1	B	103	SER	5.5
1	B	112	MET	5.4
1	C	140	VAL	5.4
1	B	121	PHE	5.4
1	A	178	ARG	5.4
1	B	99	THR	5.3
1	B	138	LEU	5.3
1	B	106	ILE	5.2
1	B	154	PRO	5.0
1	B	107	ILE	5.0
1	A	167	LEU	4.9
1	B	98	VAL	4.8
1	F	140	VAL	4.8
1	B	135	GLY	4.7
1	B	149	ALA	4.7
1	B	123	VAL	4.7
1	B	116	TYR	4.5
1	B	128	GLU	4.5
1	C	107	ILE	4.4
1	B	110	LEU	4.4
1	B	153	THR	4.4
1	H	421	GLY	4.3
1	B	108	GLU	4.2
1	B	188	ASN	4.2
1	B	132	ILE	4.1
1	F	107	ILE	4.1
1	C	149	ALA	4.0
1	B	137	ASP	4.0
1	B	97	PRO	4.0
1	A	176	GLU	4.0
1	A	423	VAL	3.8
1	C	111	GLN	3.8
1	C	139	ARG	3.8
1	A	197	PHE	3.8
1	B	136	ARG	3.8
1	E	303	ILE	3.7
1	F	120	GLY	3.7
1	B	156	ASP	3.6
1	B	122	PRO	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	148	VAL	3.6
1	C	150	ALA	3.6
1	A	422	ARG	3.5
1	B	152	MET	3.5
1	D	379	ARG	3.5
1	B	157	LYS	3.5
1	F	151	ILE	3.5
1	A	177	ASN	3.4
1	A	173	LYS	3.4
1	B	185	VAL	3.4
1	B	120	GLY	3.4
1	A	377	GLN	3.4
1	C	376	PHE	3.4
1	A	196	THR	3.3
1	C	138	LEU	3.3
1	F	111[A]	GLN	3.3
1	A	170	MET	3.3
1	C	106	ILE	3.2
1	B	134	THR	3.2
1	F	134	THR	3.2
1	C	116	TYR	3.2
1	C	141	LYS	3.2
1	B	94	VAL	3.2
1	A	376	PHE	3.2
1	C	146	ASP	3.2
1	B	192	ARG	3.1
1	B	158	LEU	3.1
1	C	108	GLU	3.1
1	D	384	TYR	3.1
1	B	92	ALA	3.0
1	A	104	THR	3.0
1	F	139	ARG	3.0
1	B	379	ARG	3.0
1	D	377	GLN	3.0
1	C	110	LEU	3.0
1	A	179	ILE	2.9
1	B	190	TYR	2.9
1	F	150	ALA	2.9
1	B	193	GLY	2.9
1	D	423	VAL	2.9
1	B	202	LYS	2.9
1	F	131	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	384	TYR	2.8
1	F	118	PHE	2.8
1	F	376	PHE	2.8
1	F	303	ILE	2.8
1	B	131	GLY	2.8
1	F	147	THR	2.7
1	D	303	ILE	2.7
1	G	1	MET	2.7
1	A	171	LYS	2.7
1	G	136	ARG	2.7
1	B	183	LEU	2.7
1	D	376	PHE	2.6
1	B	186	ASP	2.6
1	B	133	VAL	2.5
1	D	1	MET	2.5
1	B	383	SER	2.5
1	H	383	SER	2.5
1	F	116	TYR	2.5
1	B	161	ALA	2.5
1	A	159	VAL	2.5
1	C	98	VAL	2.5
1	D	422	ARG	2.5
1	B	165	THR	2.4
1	F	108	GLU	2.4
1	C	377	GLN	2.4
1	C	422	ARG	2.4
1	H	423	VAL	2.4
1	A	194	LEU	2.4
1	B	163	GLU	2.4
1	C	147	THR	2.4
1	H	422	ARG	2.4
1	B	194	LEU	2.4
1	G	139	ARG	2.4
1	E	265[A]	GLN	2.3
1	F	377	GLN	2.3
1	A	168	GLU	2.3
1	D	205	THR	2.3
1	C	151	ILE	2.3
1	B	375	LEU	2.3
1	A	137	ASP	2.3
1	F	132	ILE	2.3
1	C	120	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	135	GLY	2.3
1	B	376	PHE	2.3
1	A	375	LEU	2.3
1	A	125	GLU	2.3
1	B	377	GLN	2.3
1	C	115	GLU	2.2
1	C	137	ASP	2.2
1	F	1	MET	2.2
1	A	158	LEU	2.2
1	E	178	ARG	2.2
1	A	160	THR	2.2
1	B	93	ILE	2.2
1	B	184	VAL	2.1
1	B	384	TYR	2.1
1	F	137	ASP	2.1
1	A	120	GLY	2.1
1	B	117	GLY	2.1
1	D	90	GLU	2.1
1	D	91	THR	2.1
1	E	176	GLU	2.1
1	B	164	GLY	2.1
1	F	422	ARG	2.1
1	B	162	ARG	2.0
1	B	160	THR	2.0
1	B	114	ARG	2.0
1	F	119	SER	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.