



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 06:54 PM UTC

PDB ID : 8K2O / pdb_00008k2o
Title : Crystal structure of Fhb7-M10
Authors : Yang, J.; Lei, X.G.; Xiao, J.Y.
Deposited on : 2023-07-13
Resolution : 2.01 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

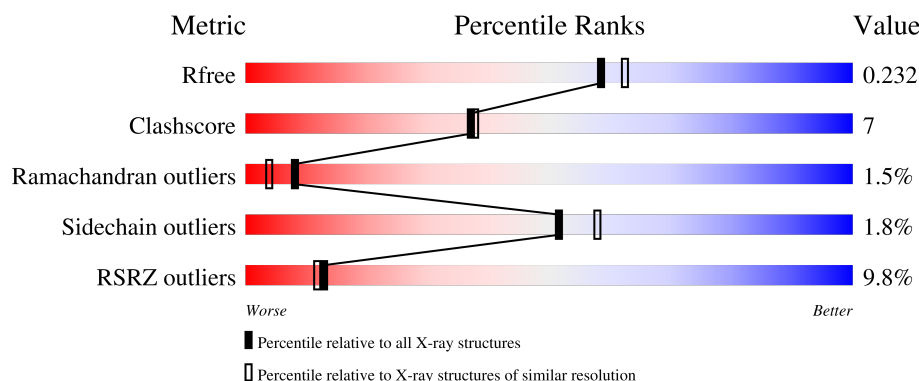
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	285	9% 87% 7% • •
1	B	285	8% 80% 14% • •
1	C	285	8% 80% 12% • 7%
1	D	285	12% 77% 15% • 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9031 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.

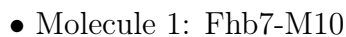
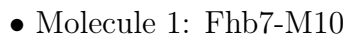
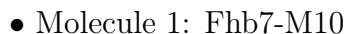
- Molecule 1 is a protein called Fhb7-M10.

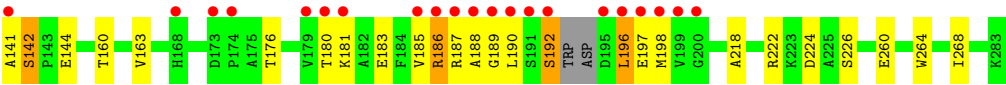
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2094	1338	358	386	12			
1	B	276	Total	C	N	O	S	0	0	0
			2128	1362	365	389	12			
1	C	266	Total	C	N	O	S	0	0	0
			2070	1326	353	379	12			
1	D	267	Total	C	N	O	S	0	0	0
			2051	1316	351	372	12			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	195	Total	O	0	0
			195	195		
2	B	167	Total	O	0	0
			167	167		
2	C	182	Total	O	0	0
			182	182		
2	D	144	Total	O	0	0
			144	144		

- Molecule 1: Fhb7-M10





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.51Å 112.16Å 135.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	51.80 – 2.01 51.80 – 2.01	Depositor EDS
% Data completeness (in resolution range)	99.9 (51.80-2.01) 99.9 (51.80-2.01)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.20.1	Depositor
R, R_{free}	0.200 , 0.232 0.200 , 0.232	Depositor DCC
R_{free} test set	2000 reflections (2.51%)	wwPDB-VP
Wilson B-factor (Å ²)	31.5	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9031	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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	0.36	0/2151	0.60	1/2931 (0.0%)
1	B	0.32	0/2186	0.59	2/2974 (0.1%)
1	C	0.35	0/2126	0.54	0/2890
1	D	0.34	0/2104	0.57	0/2860
All	All	0.34	0/8567	0.58	3/11655 (0.0%)

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	C	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	62	ILE	CA-C-N	-8.90	103.96	121.41
1	B	62	ILE	C-N-CA	-8.90	103.96	121.41
1	A	52	LEU	CA-CB-CG	6.17	137.90	116.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	19	ARG	Sidechain
1	C	190	LEU	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2094	0	2005	24	0
1	B	2128	0	2071	39	0
1	C	2070	0	2015	31	0
1	D	2051	0	1998	32	0
2	A	195	0	0	7	0
2	B	167	0	0	12	0
2	C	182	0	0	8	0
2	D	144	0	0	6	0
All	All	9031	0	8089	117	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 7.

All (117) 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:73:LYS:HE3	1:B:98:ASP:HB3	1.50	0.94
1:C:223:LYS:NZ	2:C:301:HOH:O	2.02	0.92
1:A:67:SER:O	2:A:301:HOH:O	1.91	0.87
1:A:54:ASP:HB2	1:A:180:THR:HG21	1.62	0.82
1:B:186:ARG:NH2	2:B:303:HOH:O	2.04	0.81
1:D:41:LYS:NZ	1:D:104:GLN:OE1	2.14	0.80
1:B:122:ASP:OD2	2:B:301:HOH:O	2.01	0.77
1:D:25:THR:O	2:D:301:HOH:O	2.03	0.77
1:B:54:ASP:O	2:B:302:HOH:O	2.02	0.76
1:C:72:LEU:N	2:C:303:HOH:O	2.19	0.76
1:D:139:ILE:N	2:D:303:HOH:O	2.18	0.75
1:B:54:ASP:HB2	1:B:180:THR:HG21	1.70	0.74
1:A:195:ASP:O	1:A:196:LEU:HB2	1.87	0.73
1:C:65:GLU:HG2	1:C:66:PRO:HD2	1.69	0.71
1:A:54:ASP:O	2:A:302:HOH:O	2.08	0.70
1:B:9:THR:O	2:B:304:HOH:O	2.10	0.69
1:D:18:GLN:HB2	1:D:25:THR:HG23	1.76	0.68
1:B:165:LEU:HD13	1:B:207:MET:HE2	1.76	0.68
1:A:185:VAL:O	1:A:186:ARG:HB3	1.94	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:GLN:NE2	2:B:308:HOH:O	2.21	0.67
1:C:54:ASP:HB2	1:C:180:THR:HG21	1.76	0.66
1:D:98:ASP:OD2	2:D:302:HOH:O	2.13	0.66
1:A:147:ASP:OD2	2:A:303:HOH:O	2.14	0.66
1:D:19:ARG:O	1:D:25:THR:HG21	1.95	0.66
1:B:185:VAL:O	1:B:186:ARG:HB3	1.96	0.64
1:C:185:VAL:HG13	1:C:190:LEU:O	1.98	0.64
1:A:222:ARG:HG2	1:A:268:ILE:HD11	1.80	0.63
1:A:52:LEU:HD23	2:A:340:HOH:O	2.00	0.62
1:D:188:ALA:O	1:D:190:LEU:N	2.28	0.62
1:C:127:ARG:HG3	2:C:407:HOH:O	1.99	0.62
1:D:54:ASP:HB2	1:D:180:THR:HG21	1.82	0.61
1:B:76:LYS:HG3	2:B:309:HOH:O	2.01	0.61
1:A:52:LEU:HD12	1:A:56:GLU:HG2	1.83	0.60
1:C:22:VAL:HG22	2:C:333:HOH:O	2.01	0.59
1:A:181:LYS:HG2	1:A:193:TRP:CD2	2.38	0.59
1:D:181:LYS:O	1:D:185:VAL:HG12	2.03	0.59
1:D:71:LEU:N	2:D:307:HOH:O	2.36	0.59
2:B:324:HOH:O	1:D:52:LEU:HD21	2.02	0.58
1:B:168:HIS:HB2	1:B:197:GLU:HG3	1.85	0.58
1:B:60:LYS:HD3	1:D:77:PRO:HA	1.84	0.57
1:A:167:VAL:HG21	1:B:129:MET:CE	2.34	0.57
1:D:57:ARG:O	1:D:61:GLU:HG3	2.04	0.57
1:B:39:ASN:OD1	1:B:283:LYS:HE3	2.05	0.57
1:B:56:GLU:HB2	1:D:74:GLU:HG3	1.87	0.57
1:D:188:ALA:C	1:D:190:LEU:H	2.11	0.56
1:C:54:ASP:O	2:C:302:HOH:O	2.17	0.56
1:B:72:LEU:N	2:B:309:HOH:O	2.22	0.56
1:C:52:LEU:HD23	1:C:56:GLU:HG2	1.88	0.56
1:B:185:VAL:HG23	1:B:190:LEU:HB3	1.87	0.56
1:D:142:SER:HB3	1:D:144:GLU:O	2.06	0.55
1:A:181:LYS:O	1:A:185:VAL:HG12	2.06	0.55
1:C:147:ASP:HB2	2:C:403:HOH:O	2.06	0.55
1:A:66:PRO:HB3	1:C:66:PRO:HB3	1.89	0.54
1:C:18:GLN:NE2	1:C:61:GLU:OE2	2.28	0.54
1:D:65:GLU:HG2	1:D:66:PRO:HD2	1.88	0.54
1:C:84:ILE:HG12	1:C:99:ILE:HG12	1.90	0.54
1:A:234:ARG:HD2	2:A:396:HOH:O	2.07	0.53
1:A:167:VAL:HG21	1:B:129:MET:HE3	1.91	0.53
1:A:131:GLN:NE2	2:A:311:HOH:O	2.37	0.53
1:B:250:ARG:CZ	1:B:258:TRP:HE1	2.22	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:ALA:O	2:B:305:HOH:O	2.19	0.52
1:C:176:THR:O	1:C:180:THR:HG23	2.10	0.52
1:A:77:PRO:HA	1:C:60:LYS:HD2	1.91	0.52
1:C:127:ARG:NH2	2:C:309:HOH:O	2.42	0.52
1:A:201:GLU:HG2	2:B:301:HOH:O	2.11	0.51
1:B:72:LEU:O	1:B:74:GLU:N	2.45	0.50
1:B:52:LEU:HD21	2:D:308:HOH:O	2.12	0.50
1:B:54:ASP:O	1:B:55:ILE:HB	2.11	0.50
1:B:52:LEU:HD13	1:B:56:GLU:HG2	1.93	0.49
1:B:256:SER:OG	2:B:306:HOH:O	2.19	0.49
1:D:183:GLU:OE1	1:D:187:ARG:HD2	2.13	0.49
1:C:22:VAL:C	1:C:24:GLU:H	2.20	0.49
1:C:23:ALA:HA	1:C:282:VAL:HG21	1.95	0.48
1:B:165:LEU:HD22	1:B:207:MET:HE2	1.95	0.48
1:D:18:GLN:HB2	1:D:25:THR:CG2	2.41	0.48
1:D:160:THR:O	1:D:163:VAL:HG22	2.13	0.48
1:B:55:ILE:HG12	1:B:184:PHE:CZ	2.49	0.47
1:B:283:LYS:OXT	2:B:307:HOH:O	2.20	0.47
1:D:25:THR:HA	2:D:337:HOH:O	2.14	0.47
1:D:176:THR:O	1:D:180:THR:HG23	2.15	0.46
1:A:190:LEU:HD12	1:A:196:LEU:HG	1.95	0.46
1:D:54:ASP:O	1:D:55:ILE:HB	2.14	0.46
1:C:207:MET:HE3	1:C:207:MET:HB3	1.61	0.46
1:D:222:ARG:HG2	1:D:268:ILE:HD11	1.97	0.46
1:D:260:GLU:HG3	1:D:264:TRP:NE1	2.31	0.46
1:B:195:ASP:O	1:B:196:LEU:HB2	2.15	0.46
1:D:197:GLU:O	1:D:198:MET:HG2	2.16	0.45
1:C:220:LEU:O	1:C:223:LYS:HG2	2.17	0.44
1:A:185:VAL:HG23	1:A:190:LEU:HB2	1.99	0.44
1:B:19:ARG:HA	1:B:19:ARG:HD2	1.79	0.44
1:C:50:VAL:HG12	1:C:52:LEU:HD12	1.99	0.44
1:D:218:ALA:O	1:D:222:ARG:HG3	2.18	0.43
1:B:73:LYS:HE3	1:B:98:ASP:CB	2.35	0.43
1:B:207:MET:HE3	1:B:257:GLU:CD	2.43	0.43
1:A:185:VAL:O	1:A:186:ARG:CB	2.65	0.43
1:C:181:LYS:HB3	1:C:193:TRP:HB2	2.01	0.42
1:C:238:ALA:HA	1:C:241:ILE:HD12	2.01	0.42
1:B:78:TYR:HB2	1:D:78:TYR:HB2	2.01	0.42
1:C:24:GLU:HG3	2:C:348:HOH:O	2.18	0.42
1:D:52:LEU:HD22	1:D:56:GLU:CD	2.45	0.42
1:D:224:ASP:OD1	1:D:226:SER:OG	2.33	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:167:VAL:HG21	1:B:129:MET:HE2	2.02	0.42
1:C:198:MET:HE3	1:C:198:MET:HB2	1.85	0.42
1:C:220:LEU:O	1:C:223:LYS:HE3	2.20	0.41
1:B:73:LYS:HE2	1:B:93:ILE:HG23	2.03	0.41
1:A:212:ASN:ND2	2:A:319:HOH:O	2.54	0.41
1:A:168:HIS:ND1	1:A:197:GLU:OE2	2.27	0.41
1:C:22:VAL:O	1:C:23:ALA:HB3	2.20	0.41
1:C:250:ARG:HD3	1:C:258:TRP:CE2	2.55	0.41
1:B:143:PRO:O	1:B:144:GLU:HB2	2.20	0.41
1:B:198:MET:HE2	1:B:198:MET:HB3	1.76	0.41
1:C:18:GLN:HG3	1:C:26:CYS:HA	2.03	0.41
1:D:192:SER:OG	1:D:196:LEU:HB2	2.21	0.41
1:B:182:ALA:HA	1:B:185:VAL:HG12	2.03	0.40
1:B:281:GLU:OE1	1:B:283:LYS:HE2	2.21	0.40
1:C:130:GLN:CG	1:D:63:GLY:HA2	2.51	0.40
1:C:207:MET:HE1	1:C:264:TRP:CH2	2.56	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	269/285 (94%)	253 (94%)	14 (5%)	2 (1%)	18	14
1	B	274/285 (96%)	248 (90%)	19 (7%)	7 (3%)	4	1
1	C	260/285 (91%)	247 (95%)	11 (4%)	2 (1%)	16	11
1	D	259/285 (91%)	244 (94%)	10 (4%)	5 (2%)	6	3
All	All	1062/1140 (93%)	992 (93%)	54 (5%)	16 (2%)	8	4

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	186	ARG
1	A	196	LEU
1	D	186	ARG
1	B	186	ARG
1	B	196	LEU
1	D	189	GLY
1	C	9	THR
1	D	141	ALA
1	B	71	LEU
1	D	95	ASP
1	B	95	ASP
1	B	173	ASP
1	C	22	VAL
1	D	142	SER
1	B	73	LYS
1	B	70	GLY

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	210/230 (91%)	208 (99%)	2 (1%)	68	75
1	B	215/230 (94%)	210 (98%)	5 (2%)	44	49
1	C	212/230 (92%)	209 (99%)	3 (1%)	59	66
1	D	207/230 (90%)	202 (98%)	5 (2%)	43	47
All	All	844/920 (92%)	829 (98%)	15 (2%)	51	58

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	ILE
1	A	195	ASP
1	B	9	THR
1	B	62	ILE
1	B	137	SER
1	B	144	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	217	LEU
1	C	22	VAL
1	C	65	GLU
1	C	142	SER
1	D	9	THR
1	D	71	LEU
1	D	186	ARG
1	D	192	SER
1	D	196	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	85	HIS
1	A	119	GLN
1	A	265	HIS
1	D	152	ASN
1	D	208	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	273/285 (95%)	0.23	26 (9%) 14 12	18, 28, 53, 71	0
1	B	276/285 (96%)	0.40	24 (8%) 16 15	23, 34, 59, 70	0
1	C	266/285 (93%)	0.28	22 (8%) 17 16	20, 29, 56, 67	0
1	D	267/285 (93%)	0.55	34 (12%) 8 7	21, 32, 61, 84	0
All	All	1082/1140 (94%)	0.36	106 (9%) 13 12	18, 31, 58, 84	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	72	LEU	6.4
1	D	141	ALA	6.1
1	D	197	GLU	6.0
1	C	190	LEU	5.4
1	A	141	ALA	5.3
1	D	189	GLY	5.1
1	C	142	SER	4.8
1	B	69	PHE	4.7
1	A	140	ARG	4.6
1	A	72	LEU	4.6
1	A	8	SER	4.6
1	D	195	ASP	4.4
1	B	68	ALA	4.4
1	D	71	LEU	4.4
1	D	196	LEU	4.3
1	C	143	PRO	4.1
1	C	135	PRO	4.1
1	B	71	LEU	3.9
1	D	185	VAL	3.9
1	D	74	GLU	3.9
1	C	8	SER	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	195	ASP	3.9
1	C	191	SER	3.8
1	C	189	GLY	3.7
1	B	189	GLY	3.6
1	D	22	VAL	3.6
1	D	188	ALA	3.5
1	D	168	HIS	3.5
1	D	139	ILE	3.5
1	A	71	LEU	3.4
1	C	22	VAL	3.4
1	D	21	PRO	3.4
1	D	191	SER	3.3
1	A	63	GLY	3.3
1	D	23	ALA	3.2
1	D	190	LEU	3.2
1	B	70	GLY	3.2
1	C	185	VAL	3.2
1	A	129	MET	3.2
1	B	75	GLY	3.1
1	B	174	PRO	3.1
1	C	62	ILE	3.1
1	B	188	ALA	3.1
1	D	52	LEU	3.0
1	C	23	ALA	2.9
1	B	172	LEU	2.9
1	D	192	SER	2.9
1	D	181	LYS	2.9
1	A	185	VAL	2.9
1	D	140	ARG	2.9
1	C	67	SER	2.9
1	B	196	LEU	2.8
1	A	142	SER	2.8
1	D	187	ARG	2.8
1	D	63	GLY	2.8
1	A	143	PRO	2.8
1	A	195	ASP	2.7
1	B	63	GLY	2.7
1	D	180	THR	2.7
1	B	67	SER	2.7
1	D	8	SER	2.7
1	D	200	GLY	2.7
1	A	74	GLU	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	66	PRO	2.7
1	B	185	VAL	2.7
1	A	189	GLY	2.7
1	C	63	GLY	2.7
1	B	190	LEU	2.6
1	A	67	SER	2.6
1	A	139	ILE	2.6
1	A	188	ALA	2.6
1	A	75	GLY	2.5
1	B	182	ALA	2.5
1	A	52	LEU	2.5
1	B	62	ILE	2.5
1	D	55	ILE	2.5
1	D	179	VAL	2.5
1	D	199	VAL	2.5
1	A	182	ALA	2.5
1	C	52	LEU	2.4
1	C	55	ILE	2.4
1	B	180	THR	2.4
1	C	180	THR	2.4
1	D	186	ARG	2.4
1	D	174	PRO	2.4
1	B	194	ASP	2.4
1	D	94	GLY	2.3
1	B	73	LYS	2.3
1	C	188	ALA	2.3
1	B	143	PRO	2.2
1	B	186	ARG	2.2
1	B	9	THR	2.2
1	D	9	THR	2.2
1	D	173	ASP	2.2
1	D	198	MET	2.2
1	A	55	ILE	2.1
1	A	186	ARG	2.1
1	A	196	LEU	2.1
1	C	283	LYS	2.1
1	A	272	LEU	2.1
1	C	128	ASP	2.0
1	C	186	ARG	2.0
1	C	9	THR	2.0
1	A	73	LYS	2.0
1	A	178	GLU	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	77	PRO	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.