



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:23 AM UTC

PDB ID : 8E1E / pdb_00008e1e
Title : Scaffolding protein functional sites using deep learning
Authors : Bera, A.K.; Gerben, S.; Baker, D.
Deposited on : 2022-08-10
Resolution : 4.27 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

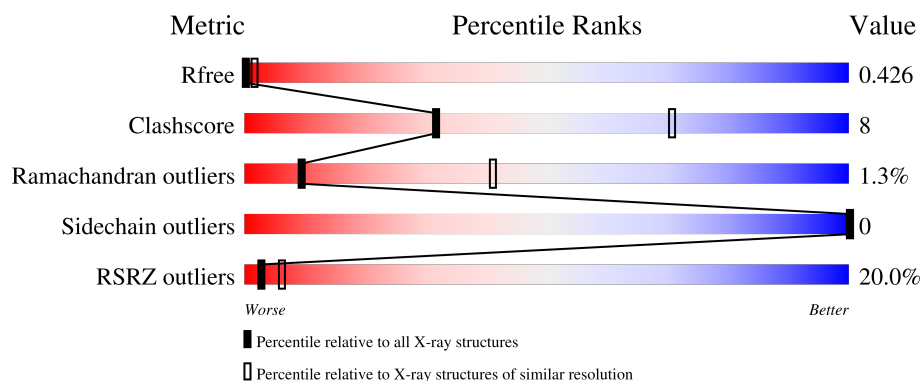
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.27 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	1029 (4.62-3.90)
Clashscore	190562	1074 (4.62-3.90)
Ramachandran outliers	187476	1001 (4.62-3.88)
Sidechain outliers	187428	1021 (4.66-3.86)
RSRZ outliers	180081	1026 (4.62-3.90)

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	243	
1	B	243	
1	C	243	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5622 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.

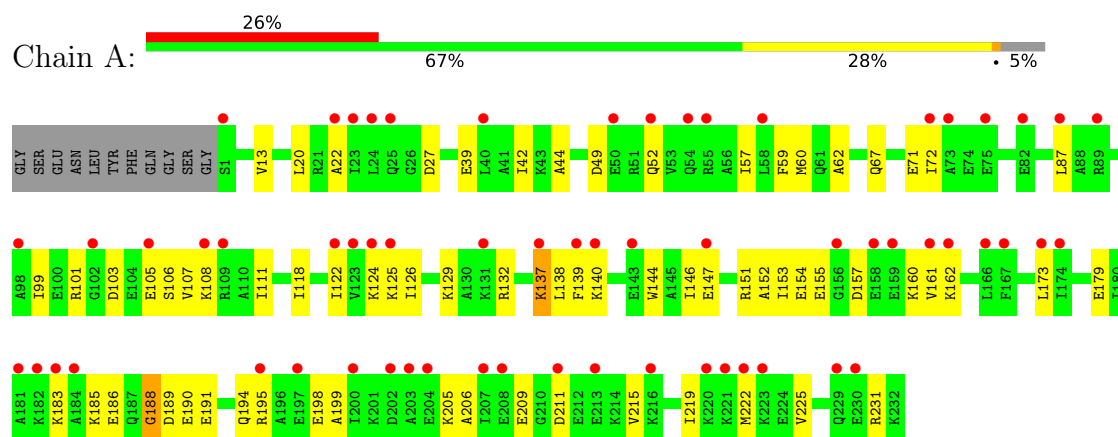
- Molecule 1 is a protein called SG122_C3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	232	Total	C	N	O	S	0	0	0
			1874	1197	320	353	4			
1	B	232	Total	C	N	O	S	0	0	0
			1874	1197	320	353	4			
1	C	232	Total	C	N	O	S	0	0	0
			1874	1197	320	353	4			

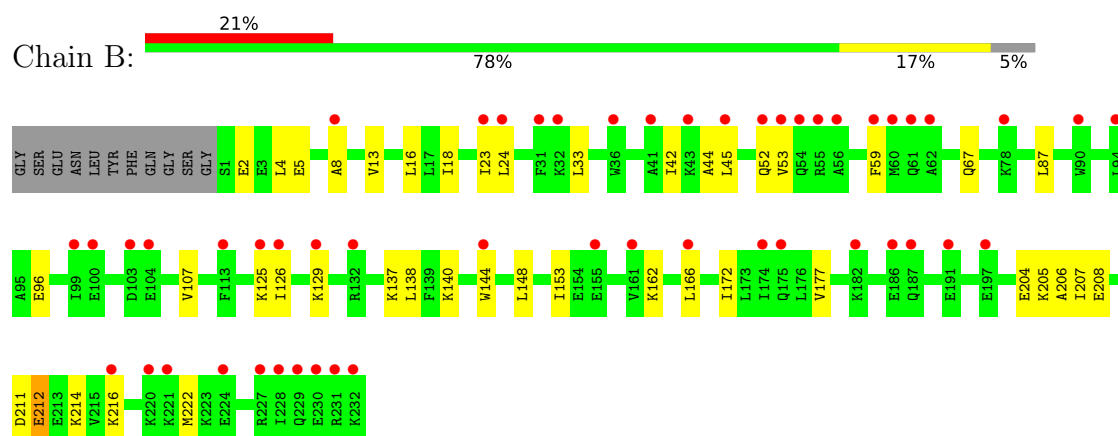
3 Residue-property plots [i](#)

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.

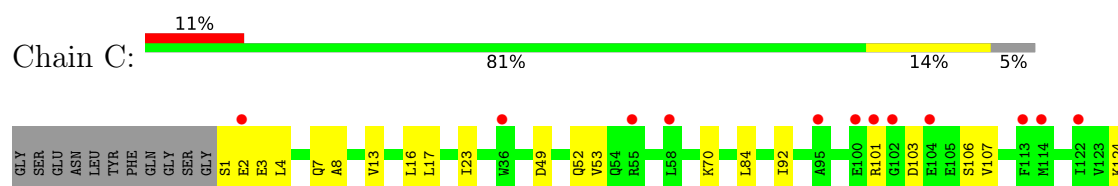
• Molecule 1: SG122_C3

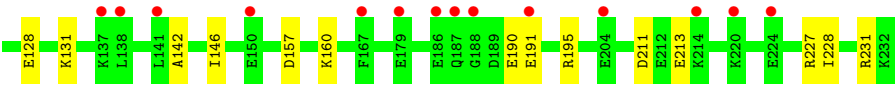


• Molecule 1: SG122_C3



• Molecule 1: SG122_C3





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	234.09Å 234.09Å 234.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	95.57 – 4.27 95.57 – 4.27	Depositor EDS
% Data completeness (in resolution range)	99.7 (95.57-4.27) 99.9 (95.57-4.27)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 4.30Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.405 , 0.427 0.405 , 0.426	Depositor DCC
R_{free} test set	741 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	282.1	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 256.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.147 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	5622	wwPDB-VP
Average B, all atoms (Å ²)	249.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.11% 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	0.16	0/1887	0.42	0/2517
1	B	0.14	0/1887	0.40	0/2517
1	C	0.13	0/1887	0.37	0/2517
All	All	0.15	0/5661	0.40	0/7551

There are no bond length outliers.

There are no bond angle outliers.

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	1874	0	1984	50	0
1	B	1874	0	1984	29	0
1	C	1874	0	1984	23	0
All	All	5622	0	5952	94	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 8.

All (94) 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:2:GLU:OE2	1:B:4:LEU:HG	1.79	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:ASP:HB3	1:A:52:GLN:HG2	1.61	0.82
1:C:227:ARG:HH12	1:C:231:ARG:HB3	1.47	0.80
1:A:138:LEU:HD12	1:A:139:PHE:H	1.49	0.77
1:A:231:ARG:O	1:A:231:ARG:NH1	2.17	0.75
1:A:129:LYS:HE2	1:A:132:ARG:HH11	1.55	0.70
1:B:2:GLU:OE1	1:B:5:GLU:OE1	2.11	0.69
1:C:227:ARG:NH1	1:C:231:ARG:HB3	2.08	0.68
1:A:206:ALA:HB1	1:A:215:VAL:HG22	1.75	0.68
1:A:160:LYS:HG2	1:A:161:VAL:H	1.58	0.68
1:B:24:LEU:HD11	1:C:23:ILE:HG12	1.75	0.67
1:A:39:GLU:HG2	1:C:17:LEU:HD13	1.76	0.67
1:A:179:GLU:O	1:A:183:LYS:NZ	2.29	0.65
1:A:42:ILE:HD11	1:C:13:VAL:HG12	1.82	0.62
1:C:228:ILE:O	1:C:231:ARG:HG2	2.01	0.61
1:C:190:GLU:OE1	1:C:190:GLU:N	2.25	0.60
1:A:147:GLU:O	1:A:151:ARG:HG2	2.03	0.59
1:B:5:GLU:HB2	1:B:45:LEU:HD22	1.85	0.58
1:B:24:LEU:HD21	1:C:23:ILE:HG23	1.85	0.57
1:A:124:LYS:HG3	1:A:146:ILE:HD13	1.86	0.57
1:A:101:ARG:NH2	1:A:106:SER:OG	2.37	0.57
1:B:33:LEU:HD11	1:B:96:GLU:HG2	1.86	0.57
1:A:129:LYS:HE2	1:A:132:ARG:NH1	2.20	0.56
1:A:137:LYS:HG3	1:A:138:LEU:H	1.70	0.56
1:A:151:ARG:HA	1:A:154:GLU:HG2	1.86	0.56
1:A:122:ILE:O	1:A:126:ILE:HD12	2.06	0.56
1:A:194:GLN:O	1:A:198:GLU:HG2	2.05	0.56
1:A:211:ASP:HB3	1:A:215:VAL:HG23	1.87	0.56
1:C:157:ASP:HB3	1:C:160:LYS:HB2	1.88	0.55
1:C:8:ALA:HB2	1:C:53:VAL:HG13	1.86	0.55
1:C:191:GLU:OE2	1:C:195:ARG:NH2	2.40	0.54
1:B:205:LYS:C	1:B:207:ILE:H	2.15	0.53
1:C:49:ASP:OD2	1:C:52:GLN:NE2	2.41	0.53
1:C:128:GLU:HA	1:C:131:LYS:HG2	1.90	0.53
1:A:67:GLN:NE2	1:A:71:GLU:OE2	2.36	0.52
1:B:177:VAL:HG21	1:B:222:MET:HG2	1.90	0.52
1:A:87:LEU:HD22	1:A:153:ILE:HD13	1.92	0.52
1:B:2:GLU:CD	1:B:4:LEU:HG	2.35	0.51
1:B:13:VAL:HG13	1:C:16:LEU:HD11	1.91	0.51
1:A:188:GLY:O	1:A:190:GLU:N	2.44	0.51
1:B:125:LYS:HE2	1:B:129:LYS:HE3	1.93	0.51
1:B:204:GLU:O	1:B:207:ILE:HG23	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ASP:O	1:A:107:VAL:HG23	2.11	0.51
1:A:13:VAL:HG13	1:B:16:LEU:HD11	1.93	0.50
1:A:103:ASP:HB3	1:A:106:SER:HB2	1.93	0.50
1:A:199:ALA:HB1	1:A:222:MET:HG2	1.92	0.50
1:A:219:ILE:HA	1:A:222:MET:HE3	1.92	0.50
1:B:211:ASP:OD1	1:B:214:LYS:HG2	2.12	0.50
1:A:138:LEU:CD1	1:A:139:PHE:H	2.22	0.49
1:A:138:LEU:O	1:A:140:LYS:N	2.43	0.48
1:A:155:GLU:HG3	1:A:157:ASP:HB2	1.94	0.48
1:B:166:LEU:HD21	1:B:216:LYS:HD3	1.95	0.47
1:A:195:ARG:HG3	1:A:225:VAL:HG22	1.97	0.47
1:A:13:VAL:HG12	1:B:42:ILE:HD11	1.96	0.47
1:C:211:ASP:OD1	1:C:213:GLU:HG2	2.15	0.47
1:B:18:ILE:HG12	1:B:67:GLN:HG3	1.97	0.46
1:C:103:ASP:O	1:C:107:VAL:HG23	2.16	0.46
1:C:1:SER:C	1:C:3:GLU:H	2.24	0.46
1:A:44:ALA:HA	1:A:52:GLN:NE2	2.31	0.46
1:C:4:LEU:HA	1:C:7:GLN:OE1	2.16	0.45
1:A:188:GLY:O	1:A:191:GLU:N	2.44	0.45
1:B:59:PHE:CE1	1:B:107:VAL:HG21	2.51	0.45
1:A:140:LYS:O	1:A:144:TRP:N	2.43	0.45
1:A:22:ALA:HB1	1:A:27:ASP:O	2.17	0.45
1:A:160:LYS:HB3	1:A:160:LYS:HE2	1.56	0.44
1:C:142:ALA:O	1:C:146:ILE:HG12	2.17	0.44
1:A:152:ALA:HB1	1:A:161:VAL:HG22	1.98	0.44
1:A:125:LYS:O	1:A:129:LYS:N	2.45	0.44
1:B:211:ASP:C	1:B:212:GLU:HG2	2.42	0.44
1:A:59:PHE:HE1	1:A:99:ILE:HA	1.83	0.44
1:A:185:LYS:NZ	1:A:186:GLU:HB2	2.33	0.44
1:C:101:ARG:HD3	1:C:103:ASP:HB2	2.00	0.44
1:C:103:ASP:HB3	1:C:106:SER:HB2	2.00	0.43
1:A:160:LYS:O	1:A:162:LYS:N	2.46	0.43
1:B:144:TRP:CZ2	1:B:148:LEU:HD22	2.52	0.43
1:A:152:ALA:HB1	1:A:160:LYS:HE3	2.00	0.43
1:A:62:ALA:HB2	1:A:111:ILE:HD11	2.00	0.43
1:B:137:LYS:O	1:B:140:LYS:N	2.52	0.42
1:A:105:GLU:O	1:A:108:LYS:HG2	2.19	0.42
1:B:126:ILE:HD12	1:B:172:ILE:HG12	2.02	0.42
1:B:8:ALA:HB2	1:B:53:VAL:HG13	2.02	0.42
1:B:44:ALA:HB1	1:B:52:GLN:HB3	2.02	0.42
1:B:205:LYS:O	1:B:207:ILE:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:LEU:HB2	1:C:124:LYS:NZ	2.35	0.42
1:A:185:LYS:HZ2	1:A:186:GLU:HB2	1.85	0.41
1:A:20:LEU:HB3	1:B:23:ILE:HD11	2.01	0.41
1:A:57:ILE:HA	1:A:60:MET:HE3	2.02	0.41
1:B:87:LEU:HD22	1:B:153:ILE:HG22	2.03	0.41
1:C:70:LYS:HG2	1:C:92:ILE:HD13	2.02	0.41
1:A:173:LEU:HD11	1:A:219:ILE:HG23	2.02	0.41
1:B:166:LEU:HD21	1:B:216:LYS:CD	2.51	0.41
1:A:72:ILE:HD12	1:A:118:ILE:HG12	2.03	0.40
1:B:205:LYS:C	1:B:207:ILE:N	2.80	0.40
1:A:205:LYS:O	1:A:209:GLU:HG3	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	230/243 (95%)	216 (94%)	11 (5%)	3 (1%)	9	41
1	B	230/243 (95%)	220 (96%)	5 (2%)	5 (2%)	5	29
1	C	230/243 (95%)	226 (98%)	3 (1%)	1 (0%)	30	66
All	All	690/729 (95%)	662 (96%)	19 (3%)	9 (1%)	9	41

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	ASP
1	B	138	LEU
1	A	188	GLY
1	B	162	LYS
1	B	206	ALA
1	A	137	LYS

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Mol	Chain	Res	Type
1	B	212	GLU
1	C	2	GLU
1	B	208	GLU

5.3.2 Protein sidechains [i](#)

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	191/199 (96%)	191 (100%)	0	100	100
1	B	191/199 (96%)	191 (100%)	0	100	100
1	C	191/199 (96%)	191 (100%)	0	100	100
All	All	573/597 (96%)	573 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
1	C	46	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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	232/243 (95%)	1.30	62 (26%)	1 4	232, 252, 262, 266	0
1	B	232/243 (95%)	1.21	51 (21%)	2 5	241, 253, 261, 264	0
1	C	232/243 (95%)	0.71	26 (11%)	10 14	240, 245, 252, 258	0
All	All	696/729 (95%)	1.07	139 (19%)	3 6	232, 249, 260, 266	0

All (139) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	59	PHE	7.8
1	B	103	ASP	5.3
1	A	23	ILE	4.6
1	B	52	GLN	4.4
1	B	229	GLN	4.2
1	B	55	ARG	4.1
1	B	191	GLU	4.1
1	B	24	LEU	4.0
1	A	124	LYS	4.0
1	B	224	GLU	4.0
1	B	155	GLU	4.0
1	A	123	VAL	3.7
1	C	214	LYS	3.7
1	A	197	GLU	3.6
1	A	105	GLU	3.6
1	C	114	MET	3.6
1	B	104	GLU	3.5
1	B	187	GLN	3.5
1	C	187	GLN	3.5
1	B	53	VAL	3.4
1	B	216	LYS	3.4
1	C	204	GLU	3.4
1	A	195	ARG	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	2	GLU	3.3
1	C	102	GLY	3.3
1	B	161	VAL	3.3
1	B	228	ILE	3.3
1	A	183	LYS	3.3
1	B	113	PHE	3.2
1	A	139	PHE	3.2
1	A	50	GLU	3.2
1	A	167	PHE	3.2
1	A	1	SER	3.1
1	A	220	LYS	3.1
1	A	221	LYS	3.1
1	A	102	GLY	3.1
1	A	55	ARG	3.1
1	B	126	ILE	3.0
1	A	52	GLN	3.0
1	A	184	ALA	3.0
1	A	54	GLN	3.0
1	A	131	LYS	2.9
1	B	125	LYS	2.9
1	B	61	GLN	2.9
1	A	207	ILE	2.9
1	A	40	LEU	2.9
1	A	204	GLU	2.9
1	B	182	LYS	2.9
1	C	188	GLY	2.9
1	C	55	ARG	2.8
1	B	54	GLN	2.8
1	A	122	ILE	2.8
1	A	25	GLN	2.8
1	C	58	LEU	2.8
1	B	31	PHE	2.8
1	B	227	ARG	2.8
1	B	36	TRP	2.8
1	A	108	LYS	2.7
1	B	60	MET	2.7
1	B	129	LYS	2.7
1	A	24	LEU	2.7
1	A	203	ALA	2.7
1	A	229	GLN	2.7
1	B	41	ALA	2.6
1	B	62	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	45	LEU	2.6
1	C	141	LEU	2.6
1	B	186	GLU	2.6
1	C	104	GLU	2.6
1	A	182	LYS	2.6
1	A	158	GLU	2.6
1	A	159	GLU	2.6
1	B	221	LYS	2.6
1	A	202	ASP	2.5
1	A	181	ALA	2.5
1	A	230	GLU	2.5
1	A	223	LYS	2.5
1	C	95	ALA	2.5
1	A	166	LEU	2.5
1	B	32	LYS	2.5
1	B	78	LYS	2.5
1	A	109	ARG	2.5
1	A	162	LYS	2.4
1	A	213	GLU	2.4
1	C	191	GLU	2.4
1	A	222	MET	2.4
1	C	113	PHE	2.4
1	C	179	GLU	2.4
1	A	98	ALA	2.4
1	A	72	ILE	2.4
1	B	232	LYS	2.4
1	B	8	ALA	2.4
1	B	94	LEU	2.4
1	C	150	GLU	2.3
1	B	144	TRP	2.3
1	A	22	ALA	2.3
1	C	138	LEU	2.3
1	A	156	GLY	2.3
1	B	230	GLU	2.3
1	A	137	LYS	2.3
1	B	99	ILE	2.3
1	A	87	LEU	2.3
1	A	147	GLU	2.3
1	C	220	LYS	2.2
1	B	56	ALA	2.2
1	A	82	GLU	2.2
1	A	208	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	100	GLU	2.2
1	B	220	LYS	2.2
1	A	174	ILE	2.2
1	C	36	TRP	2.2
1	A	73	ALA	2.2
1	A	200	ILE	2.2
1	A	161	VAL	2.2
1	A	140	LYS	2.2
1	A	211	ASP	2.2
1	B	100	GLU	2.2
1	C	224	GLU	2.2
1	A	58	LEU	2.1
1	B	132	ARG	2.1
1	B	166	LEU	2.1
1	A	75	GLU	2.1
1	B	23	ILE	2.1
1	B	90	TRP	2.1
1	C	167	PHE	2.1
1	C	101	ARG	2.1
1	A	125	LYS	2.1
1	A	216	LYS	2.1
1	B	175	GLN	2.1
1	B	174	ILE	2.1
1	A	173	LEU	2.1
1	C	137	LYS	2.1
1	B	197	GLU	2.1
1	B	231	ARG	2.1
1	B	43	LYS	2.0
1	A	89	ARG	2.0
1	A	143	GLU	2.0
1	C	122	ILE	2.0
1	C	186	GLU	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.