



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:06 AM UTC

PDB ID : 9DSC / pdb_00009dsc
Title : Crystal structure of Apo-241_2F04-A95a mutant Fab
Authors : Lin, T.H.; Wilson, I.A.
Deposited on : 2024-09-26
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

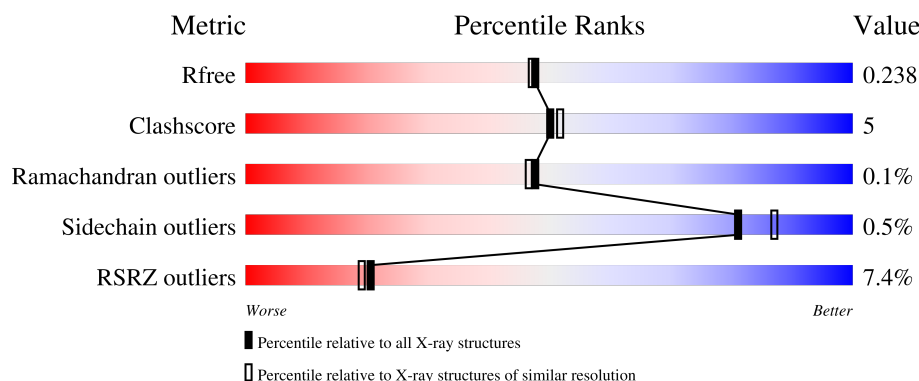
1 Overall quality at a glance

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	227	12% 85% 15%
1	B	227	10% 85% 14%
2	C	215	2% 95% .
2	F	215	5% 89% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7341 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 241_2F04-A95a, Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	S	0	0	0
			1693	1071	282	332	8			
1	B	227	Total	C	N	O	S	0	0	0
			1693	1071	282	332	8			

- Molecule 2 is a protein called 241_2F04-A95a, Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	214	Total	C	N	O	S	0	0	0
			1643	1026	281	331	5			
2	C	214	Total	C	N	O	S	0	0	0
			1643	1026	281	331	5			

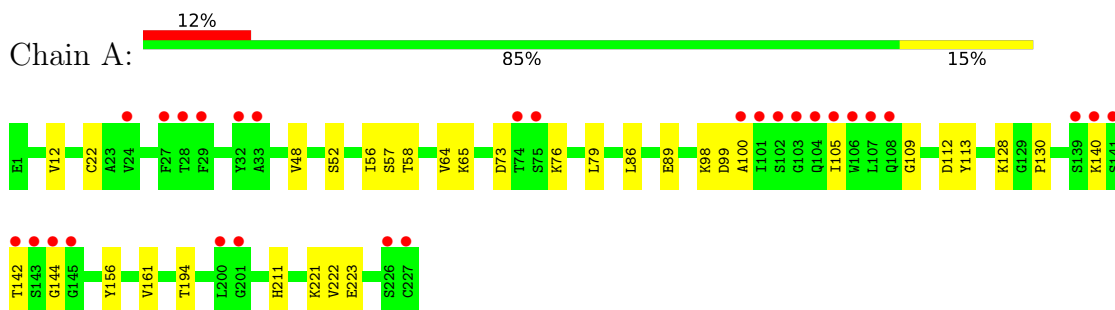
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	168	Total	O	0	0
			168	168		
3	F	138	Total	O	0	0
			138	138		
3	B	163	Total	O	0	0
			163	163		
3	C	200	Total	O	0	0
			200	200		

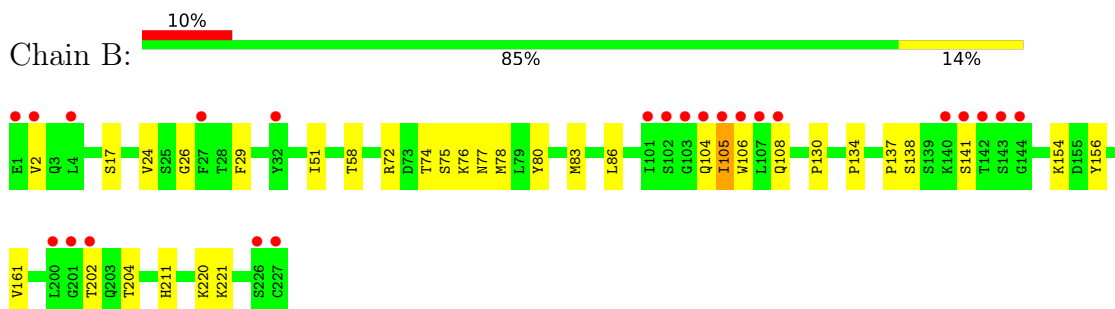
3 Residue-property plots

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.

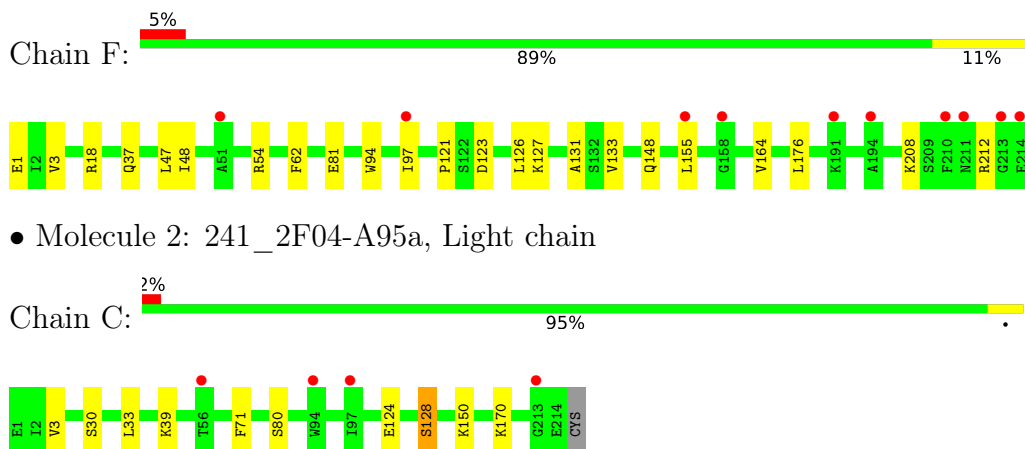
- Molecule 1: 241_2F04-A95a, Heavy chain



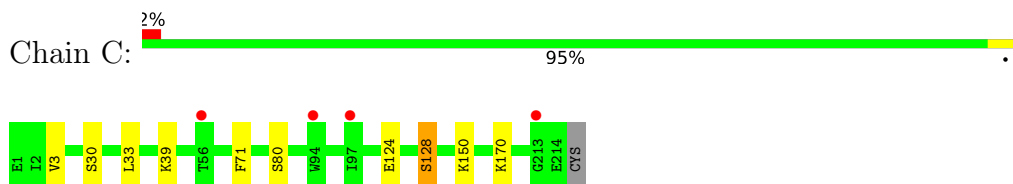
- Molecule 1: 241_2F04-A95a, Heavy chain



- Molecule 2: 241_2F04-A95a, Light chain



- Molecule 2: 241_2F04-A95a, Light chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.25Å 113.66Å 160.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.41 – 2.01 46.41 – 2.01	Depositor EDS
% Data completeness (in resolution range)	96.0 (46.41-2.01) 99.9 (46.41-2.01)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.00Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.200 , 0.236 0.202 , 0.238	Depositor DCC
R_{free} test set	4100 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7341	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.95% 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	0.11	0/1733	0.34	0/2357
1	B	0.11	0/1733	0.35	0/2357
2	C	0.10	0/1679	0.33	1/2282 (0.0%)
2	F	0.11	0/1679	0.33	0/2282
All	All	0.11	0/6824	0.34	1/9278 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	30	SER	CB-CA-C	-5.24	110.10	117.23

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	1693	0	1669	22	0
1	B	1693	0	1669	23	0
2	C	1643	0	1596	8	0
2	F	1643	0	1596	21	0
3	A	168	0	0	6	0
3	B	163	0	0	6	1
3	C	200	0	0	4	1
3	F	138	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7341	0	6530	70	1

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 5.

All (70) 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:154:LYS:NZ	3:B:303:HOH:O	2.06	0.89
1:B:137:PRO:O	3:B:301:HOH:O	1.92	0.86
1:B:141:SER:OG	3:B:302:HOH:O	1.93	0.85
1:B:141:SER:HB2	3:B:304:HOH:O	1.85	0.77
1:A:89:GLU:OE2	3:A:302:HOH:O	2.05	0.75
1:A:194:THR:O	3:A:301:HOH:O	2.04	0.74
2:F:3:VAL:O	3:F:301:HOH:O	2.05	0.74
2:F:148:GLN:HG2	2:F:155:LEU:HD11	1.73	0.71
1:B:104:GLN:HA	1:B:108:GLN:HG3	1.76	0.68
2:C:128:SER:HB2	3:C:381:HOH:O	1.97	0.65
1:B:138:SER:O	3:B:304:HOH:O	2.14	0.64
1:A:222:VAL:O	3:A:304:HOH:O	2.14	0.64
2:F:1:GLU:N	3:F:302:HOH:O	2.09	0.62
2:C:124:GLU:OE1	2:C:124:GLU:N	2.25	0.62
1:B:134:PRO:HD3	1:B:220:LYS:HE2	1.83	0.60
2:F:18:ARG:NH1	2:C:170:LYS:O	2.34	0.60
1:B:17:SER:OG	3:B:305:HOH:O	2.17	0.56
1:B:75:SER:OG	1:B:76:LYS:NZ	2.34	0.56
1:B:2:VAL:HA	1:B:26:GLY:HA3	1.88	0.56
1:A:221:LYS:HD3	1:A:223:GLU:OE1	2.07	0.55
1:B:104:GLN:HG3	1:B:105:ILE:H	1.71	0.55
1:B:72:ARG:HE	1:B:74:THR:HG22	1.70	0.55
1:B:130:PRO:HB3	1:B:156:TYR:HB3	1.90	0.54
1:B:105:ILE:HG23	1:B:106:TRP:H	1.71	0.54
2:C:150:LYS:HG3	3:C:426:HOH:O	2.08	0.54
1:A:65:LYS:NZ	3:A:306:HOH:O	2.25	0.51
2:F:212:ARG:NH2	3:F:308:HOH:O	2.43	0.51
2:F:127:LYS:HD2	2:F:127:LYS:C	2.36	0.50
2:F:54:ARG:NH1	3:F:309:HOH:O	2.43	0.49
2:F:121:PRO:HD3	2:F:133:VAL:HG22	1.94	0.49
2:F:54:ARG:HH11	2:F:54:ARG:HG3	1.77	0.49
1:B:104:GLN:CG	1:B:105:ILE:H	2.25	0.49
1:A:109:GLY:HA2	2:F:97:ILE:HD11	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:37:GLN:HB2	2:F:47:LEU:HD11	1.96	0.48
1:B:51:ILE:HG13	1:B:58:THR:HG22	1.96	0.47
2:F:123:ASP:HA	2:F:126:LEU:HB2	1.97	0.47
1:A:22:CYS:HB3	1:A:79:LEU:HB3	1.97	0.46
1:B:24:VAL:HG12	1:B:77:ASN:HB3	1.97	0.46
1:A:73:ASP:OD1	1:A:76:LYS:HG3	2.15	0.46
2:F:54:ARG:HD3	2:F:62:PHE:O	2.15	0.46
2:F:123:ASP:N	3:F:315:HOH:O	2.48	0.46
1:A:128:LYS:NZ	3:A:313:HOH:O	2.48	0.45
1:B:161:VAL:HG12	1:B:211:HIS:CD2	2.51	0.45
2:C:39:LYS:HD3	3:C:338:HOH:O	2.15	0.45
1:B:105:ILE:HG23	1:B:106:TRP:N	2.32	0.45
2:F:121:PRO:HG3	2:F:131:ALA:HB1	1.97	0.45
1:A:105:ILE:HG23	2:F:94:TRP:CE3	2.52	0.44
1:A:58:THR:OG1	3:A:303:HOH:O	2.11	0.44
1:A:140:LYS:HG2	1:A:140:LYS:O	2.17	0.44
2:F:47:LEU:O	2:F:48:ILE:HD13	2.18	0.43
1:B:24:VAL:HG11	1:B:29:PHE:CD1	2.54	0.43
1:A:161:VAL:HG12	1:A:211:HIS:CD2	2.54	0.43
1:A:100:ALA:HA	1:A:112:ASP:OD2	2.18	0.43
1:B:83:MET:HB3	1:B:86:LEU:HD21	2.00	0.43
1:A:12:VAL:HG11	1:A:86:LEU:HD13	2.00	0.42
1:A:142:THR:C	1:A:144:GLY:H	2.27	0.42
2:C:3:VAL:O	3:C:301:HOH:O	2.21	0.42
2:F:47:LEU:C	2:F:48:ILE:HD13	2.45	0.42
1:A:140:LYS:HD3	1:A:140:LYS:H	1.83	0.42
1:A:130:PRO:HB3	1:A:156:TYR:HB3	2.02	0.42
1:B:204:THR:HG22	1:B:221:LYS:HE3	2.01	0.42
2:C:33:LEU:HD13	2:C:71:PHE:CD2	2.55	0.42
1:A:56:ILE:H	1:A:56:ILE:HG12	1.64	0.41
1:A:52:SER:HB3	1:A:57:SER:HB2	2.03	0.41
2:F:18:ARG:HB3	2:C:80:SER:OG	2.20	0.41
1:A:98:LYS:HE3	1:A:113:TYR:CD2	2.55	0.41
1:A:48:VAL:HG13	1:A:64:VAL:HG21	2.03	0.41
1:B:78:MET:HB3	1:B:80:TYR:CE2	2.56	0.40
2:F:164:VAL:HG22	2:F:176:LEU:HD12	2.03	0.40
2:F:208:LYS:NZ	3:F:322:HOH:O	2.55	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:427:HOH:O	3:C:421:HOH:O[4_444]	2.13	0.07

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	225/227 (99%)	217 (96%)	8 (4%)	0	100	100
1	B	225/227 (99%)	221 (98%)	3 (1%)	1 (0%)	30	27
2	C	212/215 (99%)	207 (98%)	5 (2%)	0	100	100
2	F	212/215 (99%)	205 (97%)	7 (3%)	0	100	100
All	All	874/884 (99%)	850 (97%)	23 (3%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	105	ILE

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	190/190 (100%)	189 (100%)	1 (0%)	81	87
1	B	190/190 (100%)	189 (100%)	1 (0%)	81	87
2	C	185/186 (100%)	184 (100%)	1 (0%)	81	87
2	F	185/186 (100%)	184 (100%)	1 (0%)	81	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	750/752 (100%)	746 (100%)	4 (0%)	81	87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	ASP
2	F	81	GLU
1	B	202	THR
2	C	128	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	104	GLN
2	F	38	GLN
2	F	153	ASN
1	B	50	HIS
1	B	108	GLN
1	B	116	GLN
2	C	32	ASN
2	C	92	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

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	227/227 (100%)	0.59	28 (12%) 8 7	26, 44, 111, 159	0
1	B	227/227 (100%)	0.46	23 (10%) 12 11	23, 41, 88, 279	0
2	C	214/215 (99%)	0.03	4 (1%) 66 66	23, 37, 64, 88	0
2	F	214/215 (99%)	0.38	10 (4%) 36 35	32, 48, 72, 104	0
All	All	882/884 (99%)	0.37	65 (7%) 20 19	23, 43, 83, 279	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	107	LEU	7.2
1	B	106	TRP	6.6
1	B	101	ILE	6.0
1	A	106	TRP	5.8
1	A	107	LEU	5.5
1	A	101	ILE	5.3
1	B	103	GLY	5.1
1	B	105	ILE	4.9
1	A	105	ILE	4.8
1	A	102	SER	4.7
1	B	104	GLN	4.6
1	A	143	SER	4.2
1	A	103	GLY	4.0
1	B	2	VAL	4.0
1	B	102	SER	4.0
1	A	200	LEU	3.9
1	A	142	THR	3.8
1	A	29	PHE	3.6
1	B	202	THR	3.0
1	A	140	LYS	3.0
1	A	104	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	142	THR	2.9
1	A	145	GLY	2.9
1	A	32	TYR	2.9
1	B	226	SER	2.9
1	A	28	THR	2.9
1	B	227	CYS	2.8
1	B	141	SER	2.8
2	C	94	TRP	2.7
2	F	214	GLU	2.6
1	A	27	PHE	2.6
2	F	155	LEU	2.6
1	B	1	GLU	2.5
1	A	108	GLN	2.4
1	B	27	PHE	2.4
1	A	24	VAL	2.4
1	A	139	SER	2.4
2	F	97	ILE	2.4
1	A	75	SER	2.4
1	B	140	LYS	2.4
1	A	227	CYS	2.4
1	A	33	ALA	2.3
2	F	191	LYS	2.3
1	A	226	SER	2.3
2	F	213	GLY	2.3
2	C	213	GLY	2.3
1	A	144	GLY	2.3
1	A	201	GLY	2.3
1	B	4	LEU	2.3
2	F	194	ALA	2.3
2	C	97	ILE	2.2
1	B	143	SER	2.2
1	B	144	GLY	2.2
1	B	201	GLY	2.2
2	F	51	ALA	2.2
1	B	32	TYR	2.2
1	B	108	GLN	2.2
2	C	56	THR	2.2
2	F	211	ASN	2.1
1	A	100	ALA	2.1
1	B	200	LEU	2.0
1	A	74	THR	2.0
2	F	158	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
2	F	210	PHE	2.0
1	A	141	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 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.