



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:18 PM UTC

PDB ID : 9WLP / pdb_00009wlp
Title : Crystal structure of monkeypox virus A30/H2 sub-complex at pH 6.8
Authors : Lin, S.; Jia, X.H.; Lu, G.W.
Deposited on : 2025-09-02
Resolution : 2.50 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

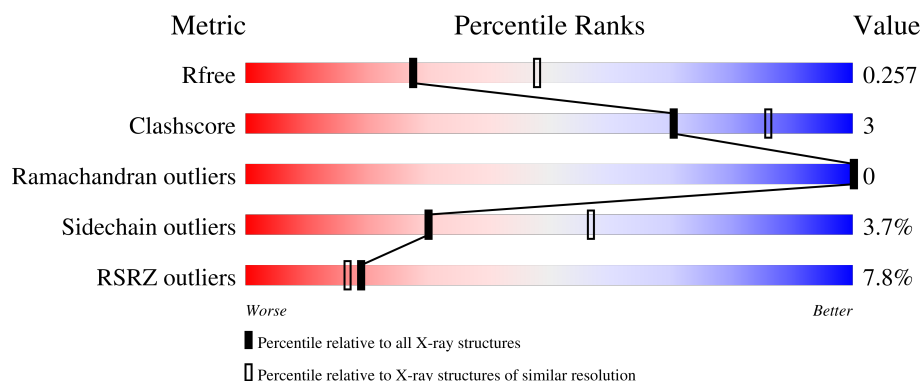
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	168	2% 70% 7% . 22%
1	C	168	4% 65% 7% . 27%
1	E	168	2% 71% . . 24%
2	B	119	11% 69% 12% . 18%
2	D	119	9% 72% 7% 21%

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Mol	Chain	Length	Quality of chain
2	F	119	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5447 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 MPXVgp092.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			1055	668	183	194	10			
1	C	123	Total	C	N	O	S	0	0	0
			981	622	169	182	8			
1	E	128	Total	C	N	O	S	0	0	0
			1030	652	180	189	9			

There are 87 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	GLY	-	expression tag	UNP A0A7H0DN79
A	48	PRO	-	expression tag	UNP A0A7H0DN79
A	49	LEU	-	expression tag	UNP A0A7H0DN79
A	50	GLY	-	expression tag	UNP A0A7H0DN79
A	190	GLY	-	expression tag	UNP A0A7H0DN79
A	191	GLY	-	expression tag	UNP A0A7H0DN79
A	192	GLY	-	expression tag	UNP A0A7H0DN79
A	193	GLY	-	expression tag	UNP A0A7H0DN79
A	194	SER	-	expression tag	UNP A0A7H0DN79
A	195	GLY	-	expression tag	UNP A0A7H0DN79
A	196	GLY	-	expression tag	UNP A0A7H0DN79
A	197	GLY	-	expression tag	UNP A0A7H0DN79
A	198	GLY	-	expression tag	UNP A0A7H0DN79
A	199	SER	-	expression tag	UNP A0A7H0DN79
A	200	GLY	-	expression tag	UNP A0A7H0DN79
A	201	GLY	-	expression tag	UNP A0A7H0DN79
A	202	GLY	-	expression tag	UNP A0A7H0DN79
A	203	GLY	-	expression tag	UNP A0A7H0DN79
A	204	SER	-	expression tag	UNP A0A7H0DN79
A	205	GLY	-	expression tag	UNP A0A7H0DN79
A	206	GLY	-	expression tag	UNP A0A7H0DN79
A	207	GLY	-	expression tag	UNP A0A7H0DN79
A	208	GLY	-	expression tag	UNP A0A7H0DN79

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Chain	Residue	Modelled	Actual	Comment	Reference
A	209	SER	-	expression tag	UNP A0A7H0DN79
A	210	GLY	-	expression tag	UNP A0A7H0DN79
A	211	GLY	-	expression tag	UNP A0A7H0DN79
A	212	GLY	-	expression tag	UNP A0A7H0DN79
A	213	GLY	-	expression tag	UNP A0A7H0DN79
A	214	SER	-	expression tag	UNP A0A7H0DN79
C	47	GLY	-	expression tag	UNP A0A7H0DN79
C	48	PRO	-	expression tag	UNP A0A7H0DN79
C	49	LEU	-	expression tag	UNP A0A7H0DN79
C	50	GLY	-	expression tag	UNP A0A7H0DN79
C	190	GLY	-	expression tag	UNP A0A7H0DN79
C	191	GLY	-	expression tag	UNP A0A7H0DN79
C	192	GLY	-	expression tag	UNP A0A7H0DN79
C	193	GLY	-	expression tag	UNP A0A7H0DN79
C	194	SER	-	expression tag	UNP A0A7H0DN79
C	195	GLY	-	expression tag	UNP A0A7H0DN79
C	196	GLY	-	expression tag	UNP A0A7H0DN79
C	197	GLY	-	expression tag	UNP A0A7H0DN79
C	198	GLY	-	expression tag	UNP A0A7H0DN79
C	199	SER	-	expression tag	UNP A0A7H0DN79
C	200	GLY	-	expression tag	UNP A0A7H0DN79
C	201	GLY	-	expression tag	UNP A0A7H0DN79
C	202	GLY	-	expression tag	UNP A0A7H0DN79
C	203	GLY	-	expression tag	UNP A0A7H0DN79
C	204	SER	-	expression tag	UNP A0A7H0DN79
C	205	GLY	-	expression tag	UNP A0A7H0DN79
C	206	GLY	-	expression tag	UNP A0A7H0DN79
C	207	GLY	-	expression tag	UNP A0A7H0DN79
C	208	GLY	-	expression tag	UNP A0A7H0DN79
C	209	SER	-	expression tag	UNP A0A7H0DN79
C	210	GLY	-	expression tag	UNP A0A7H0DN79
C	211	GLY	-	expression tag	UNP A0A7H0DN79
C	212	GLY	-	expression tag	UNP A0A7H0DN79
C	213	GLY	-	expression tag	UNP A0A7H0DN79
C	214	SER	-	expression tag	UNP A0A7H0DN79
E	47	GLY	-	expression tag	UNP A0A7H0DN79
E	48	PRO	-	expression tag	UNP A0A7H0DN79
E	49	LEU	-	expression tag	UNP A0A7H0DN79
E	50	GLY	-	expression tag	UNP A0A7H0DN79
E	190	GLY	-	expression tag	UNP A0A7H0DN79
E	191	GLY	-	expression tag	UNP A0A7H0DN79
E	192	GLY	-	expression tag	UNP A0A7H0DN79

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Chain	Residue	Modelled	Actual	Comment	Reference
E	193	GLY	-	expression tag	UNP A0A7H0DN79
E	194	SER	-	expression tag	UNP A0A7H0DN79
E	195	GLY	-	expression tag	UNP A0A7H0DN79
E	196	GLY	-	expression tag	UNP A0A7H0DN79
E	197	GLY	-	expression tag	UNP A0A7H0DN79
E	198	GLY	-	expression tag	UNP A0A7H0DN79
E	199	SER	-	expression tag	UNP A0A7H0DN79
E	200	GLY	-	expression tag	UNP A0A7H0DN79
E	201	GLY	-	expression tag	UNP A0A7H0DN79
E	202	GLY	-	expression tag	UNP A0A7H0DN79
E	203	GLY	-	expression tag	UNP A0A7H0DN79
E	204	SER	-	expression tag	UNP A0A7H0DN79
E	205	GLY	-	expression tag	UNP A0A7H0DN79
E	206	GLY	-	expression tag	UNP A0A7H0DN79
E	207	GLY	-	expression tag	UNP A0A7H0DN79
E	208	GLY	-	expression tag	UNP A0A7H0DN79
E	209	SER	-	expression tag	UNP A0A7H0DN79
E	210	GLY	-	expression tag	UNP A0A7H0DN79
E	211	GLY	-	expression tag	UNP A0A7H0DN79
E	212	GLY	-	expression tag	UNP A0A7H0DN79
E	213	GLY	-	expression tag	UNP A0A7H0DN79
E	214	SER	-	expression tag	UNP A0A7H0DN79

- Molecule 2 is a protein called Envelope protein OPG155.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	97	Total	C	N	O	S	0	0	0
			765	480	135	145	5			
2	D	94	Total	C	N	O	S	0	0	0
			750	472	132	141	5			
2	F	94	Total	C	N	O	S	0	0	0
			750	472	132	141	5			

- Molecule 3 is water.

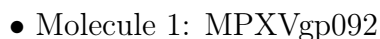
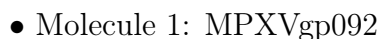
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	17	Total	O	0	0
			17	17		
3	B	21	Total	O	0	0
			21	21		
3	C	23	Total	O	0	0
			23	23		

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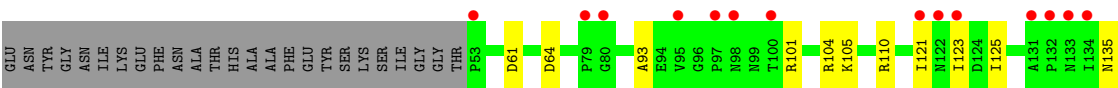
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	12	Total 12	O 12	0	0
3	E	25	Total 25	O 25	0	0
3	F	18	Total 18	O 18	0	0

- Molecule 1: MPXVgp092





● Molecule 2: Envelope protein OPG155



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.41Å 116.21Å 125.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.22 – 2.50 30.22 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.22-2.50) 99.8 (30.22-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.51Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.216 , 0.256 0.220 , 0.257	Depositor DCC
R_{free} test set	1573 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	50.0	Xtrriage
Anisotropy	0.096	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5447	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 45.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2833e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.31	0/1077	0.55	0/1446
1	C	0.32	0/1002	0.57	0/1347
1	E	0.30	0/1052	0.53	0/1413
2	B	0.44	0/780	0.74	0/1057
2	D	0.22	0/765	0.46	0/1036
2	F	0.30	0/765	0.60	0/1036
All	All	0.32	0/5441	0.58	0/7335

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	1055	0	1041	8	0
1	C	981	0	964	8	0
1	E	1030	0	1015	4	0
2	B	765	0	743	10	0
2	D	750	0	731	3	0
2	F	750	0	731	7	0
3	A	17	0	0	1	0
3	B	21	0	0	1	0
3	C	23	0	0	2	0
3	D	12	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	25	0	0	1	0
3	F	18	0	0	2	0
All	All	5447	0	5225	36	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 3.

All (36) 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)
2:B:66:ILE:O	3:B:201:HOH:O	2.00	0.79
1:C:175:TYR:O	3:C:301:HOH:O	2.04	0.74
1:E:145:GLN:NE2	1:E:149:GLN:OE1	2.24	0.69
2:B:119:ASP:OD1	2:B:119:ASP:N	2.25	0.67
2:F:61:ASP:OD2	3:F:201:HOH:O	2.13	0.66
1:A:145:GLN:OE1	1:A:149:GLN:NE2	2.30	0.64
2:F:135:ASN:HB3	2:F:138:GLU:HB2	1.82	0.60
1:A:153:LYS:HE3	2:B:110:ARG:HD2	1.84	0.60
2:B:121:ILE:HG13	2:B:122:ASN:H	1.65	0.60
2:B:81:ASN:HB2	2:B:105:LYS:HD3	1.85	0.59
1:A:112:MET:HE1	1:A:185:LEU:HD11	1.85	0.59
1:E:132:TRP:O	3:E:301:HOH:O	2.17	0.59
1:E:153:LYS:HE3	2:F:110:ARG:HH11	1.69	0.58
2:B:119:ASP:HB2	1:C:116:LYS:HG3	1.87	0.56
2:F:93:ALA:O	2:F:104:ARG:NH1	2.38	0.55
2:D:101:ARG:NH1	2:D:125:ILE:O	2.40	0.55
1:C:80:THR:HG23	1:C:120:LEU:HD11	1.91	0.52
1:C:181:TRP:O	1:C:185:LEU:HD22	2.12	0.50
2:B:101:ARG:NH1	2:B:124:ASP:HB3	2.27	0.49
2:D:135:ASN:HB3	2:D:138:GLU:HB2	1.94	0.49
2:F:110:ARG:CZ	3:F:205:HOH:O	2.61	0.48
2:B:93:ALA:O	2:B:104:ARG:NH1	2.48	0.46
1:A:58:GLU:HG3	1:A:59:LEU:HD13	1.98	0.46
2:D:101:ARG:NH1	2:D:125:ILE:H	2.14	0.45
2:F:101:ARG:NH1	2:F:125:ILE:H	2.15	0.45
1:E:126:ARG:NH2	2:F:64:ASP:OD2	2.47	0.43
1:A:58:GLU:HG3	1:A:59:LEU:N	2.34	0.43
1:C:110:ILE:HD11	1:C:133:MET:HG2	2.01	0.43
1:A:84:SER:OG	1:A:118:THR:O	2.30	0.43
1:C:112:MET:HE3	1:C:112:MET:HB2	1.86	0.42
1:A:98:PHE:O	1:A:113:ARG:HD2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:101:ARG:CZ	2:B:124:ASP:HB3	2.50	0.41
1:C:95:ASP:N	3:C:305:HOH:O	2.54	0.41
1:A:110:ILE:HD11	1:A:133:MET:HG2	2.03	0.41
2:B:135:ASN:HB3	2:B:138:GLU:HB2	2.03	0.41
3:A:302:HOH:O	1:C:153:LYS:NZ	2.54	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	129/168 (77%)	126 (98%)	3 (2%)	0	100	100
1	C	121/168 (72%)	118 (98%)	3 (2%)	0	100	100
1	E	126/168 (75%)	123 (98%)	3 (2%)	0	100	100
2	B	95/119 (80%)	90 (95%)	5 (5%)	0	100	100
2	D	92/119 (77%)	88 (96%)	4 (4%)	0	100	100
2	F	92/119 (77%)	87 (95%)	5 (5%)	0	100	100
All	All	655/861 (76%)	632 (96%)	23 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	114/129 (88%)	111 (97%)	3 (3%)	40	68
1	C	106/129 (82%)	102 (96%)	4 (4%)	29	56
1	E	111/129 (86%)	107 (96%)	4 (4%)	31	58
2	B	87/105 (83%)	84 (97%)	3 (3%)	32	60
2	D	86/105 (82%)	82 (95%)	4 (5%)	23	47
2	F	86/105 (82%)	82 (95%)	4 (5%)	23	47
All	All	590/702 (84%)	568 (96%)	22 (4%)	30	57

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	VAL
1	A	84	SER
1	A	145	GLN
2	B	57	ARG
2	B	119	ASP
2	B	136	ASN
1	C	67	ASN
1	C	116	LYS
1	C	145	GLN
1	C	185	LEU
2	D	100	THR
2	D	121	ILE
2	D	122	ASN
2	D	123	ILE
1	E	69	VAL
1	E	90	LEU
1	E	139	LYS
1	E	145	GLN
2	F	105	LYS
2	F	121	ILE
2	F	123	ILE
2	F	136	ASN

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	70	ASN
1	A	145	GLN
1	A	149	GLN

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Mol	Chain	Res	Type
2	B	92	GLN
2	B	98	ASN
1	C	157	ASN
1	E	187	ASN
2	F	71	GLN
2	F	111	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	131/168 (77%)	0.31	4 (3%)	51	47	36, 50, 84, 103	0
1	C	123/168 (73%)	0.45	6 (4%)	35	31	31, 50, 90, 105	0
1	E	128/168 (76%)	0.23	4 (3%)	51	47	32, 48, 71, 76	0
2	B	97/119 (81%)	0.85	13 (13%)	7	5	30, 49, 103, 118	0
2	D	94/119 (78%)	0.65	11 (11%)	9	7	31, 54, 106, 128	0
2	F	94/119 (78%)	0.78	14 (14%)	5	4	34, 51, 112, 138	0
All	All	667/861 (77%)	0.51	52 (7%)	19	17	30, 50, 92, 138	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	51	GLY	6.1
2	B	50	GLY	5.4
2	F	122	ASN	5.1
2	B	52	THR	5.0
2	B	121	ILE	4.4
2	D	122	ASN	4.2
2	D	97	PRO	4.2
1	C	65	ILE	4.2
2	F	97	PRO	3.9
1	C	92	ALA	3.7
2	F	123	ILE	3.7
2	D	95	VAL	3.6
2	D	123	ILE	3.4
2	D	131	ALA	3.4
2	D	121	ILE	3.4
2	B	97	PRO	3.3
2	B	120	VAL	3.2
2	D	134	ILE	3.2
2	B	119	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	58	GLU	3.0
1	A	59	LEU	2.9
2	D	133	ASN	2.9
2	B	123	ILE	2.9
2	F	133	ASN	2.8
2	F	98	ASN	2.8
2	F	134	ILE	2.8
2	F	131	ALA	2.7
2	F	121	ILE	2.7
2	F	95	VAL	2.7
2	F	100	THR	2.7
1	A	188	MET	2.7
1	C	91	ALA	2.7
1	E	91	ALA	2.6
1	A	89	LEU	2.6
2	F	53	PRO	2.6
2	B	132	PRO	2.5
1	C	69	VAL	2.5
1	C	93	GLY	2.5
1	E	60	THR	2.4
2	F	132	PRO	2.4
1	E	177	MET	2.4
2	D	100	THR	2.3
2	B	122	ASN	2.2
2	D	57	ARG	2.2
2	B	124	ASP	2.2
2	B	98	ASN	2.1
1	C	72	TRP	2.1
1	E	102	CYS	2.1
2	F	80	GLY	2.1
2	F	79	PRO	2.1
2	B	99	ASN	2.0
2	D	132	PRO	2.0

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

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

6.3 Carbohydrates ⓘ

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.