



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:33 PM UTC

PDB ID : 5YXA / pdb_00005yxa
Title : Crystal structure of the C-terminal fragment of NS1 protein from yellow fever virus
Authors : Wang, H.; Song, H.; Qi, J.; Shi, Y.; Gao, G.F.
Deposited on : 2017-12-04
Resolution : 2.10 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

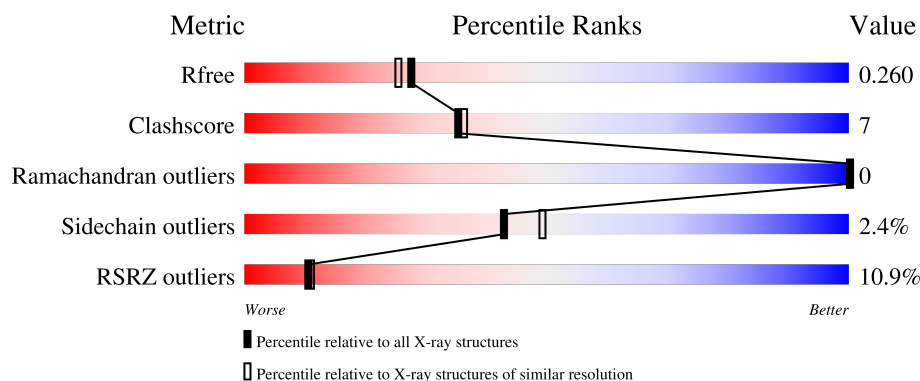
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 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	181	
1	B	181	
1	C	181	
1	D	181	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5540 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 Non-structural protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	177	Total	C	N	O	S	0	0	0
			1355	842	237	260	16			
1	B	177	Total	C	N	O	S	0	0	0
			1355	842	237	260	16			
1	C	177	Total	C	N	O	S	0	0	0
			1355	842	237	260	16			
1	D	177	Total	C	N	O	S	0	0	0
			1355	842	237	260	16			

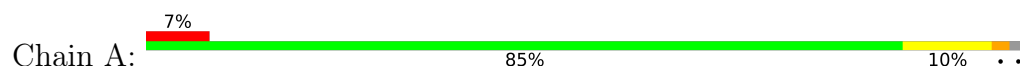
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	49	Total	O	0	0
			49	49		
2	B	37	Total	O	0	0
			37	37		
2	C	22	Total	O	0	0
			22	22		
2	D	12	Total	O	0	0
			12	12		

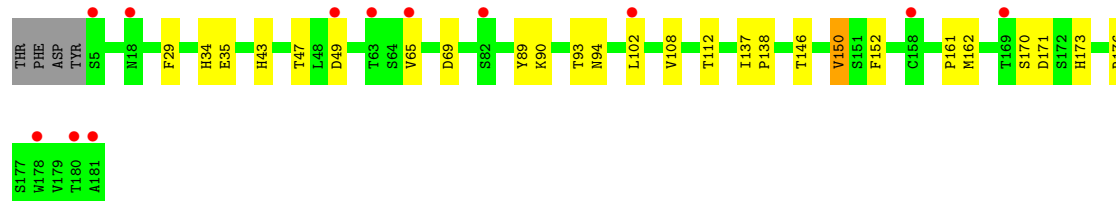
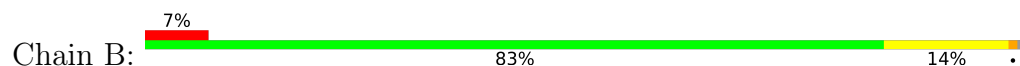
3 Residue-property plots [i](#)

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.

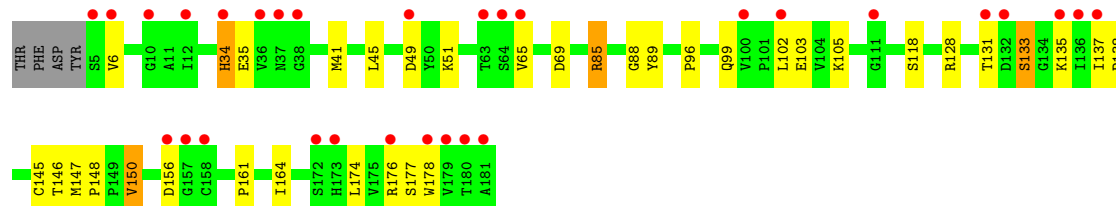
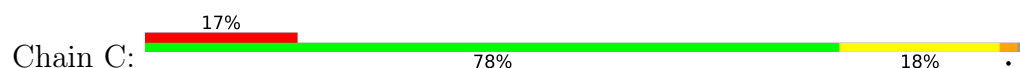
- Molecule 1: Non-structural protein 1



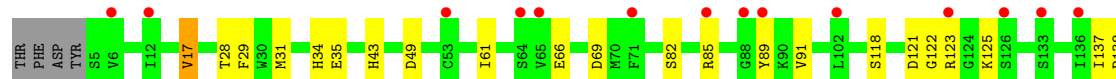
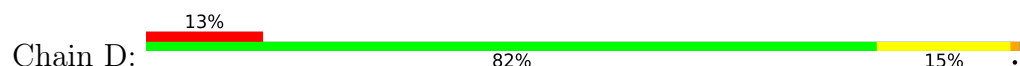
- Molecule 1: Non-structural protein 1

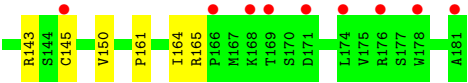


- Molecule 1: Non-structural protein 1



- Molecule 1: Non-structural protein 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.29Å 114.10Å 139.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.61 – 2.10 41.61 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (41.61-2.10) 100.0 (41.61-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.223 , 0.258 0.227 , 0.260	Depositor DCC
R_{free} test set	2808 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	1.190	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.6	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	5540	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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	0.39	0/1394	0.78	1/1897 (0.1%)
1	B	0.38	0/1394	0.85	1/1897 (0.1%)
1	C	0.38	0/1394	0.74	0/1897
1	D	0.35	0/1394	0.80	0/1897
All	All	0.37	0/5576	0.79	2/7588 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	65	VAL	N-CA-C	8.05	120.51	109.80
1	B	65	VAL	N-CA-C	5.33	118.09	110.09

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	1355	0	1287	24	0
1	B	1355	0	1287	15	0
1	C	1355	0	1287	23	1
1	D	1355	0	1287	17	0
2	A	49	0	0	2	0
2	B	37	0	0	2	0
2	C	22	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	12	0	0	2	0
All	All	5540	0	5148	78	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 7.

All (78) 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:34:HIS:HD2	1:A:43:HIS:HD2	1.15	0.94
1:B:49:ASP:OD2	2:B:201:HOH:O	1.87	0.91
1:B:171:ASP:OD1	1:B:176:ARG:NH2	2.04	0.90
1:A:34:HIS:HD2	1:A:43:HIS:CD2	1.93	0.87
1:C:65:VAL:HG11	1:C:85:ARG:HD3	1.66	0.78
1:C:51:LYS:NZ	2:C:202:HOH:O	2.14	0.76
1:C:177:SER:O	2:C:201:HOH:O	2.08	0.72
1:C:176:ARG:NH1	2:C:204:HOH:O	2.24	0.70
1:A:34:HIS:CD2	1:A:43:HIS:HD2	2.06	0.70
1:A:34:HIS:HE2	1:A:43:HIS:HB3	1.57	0.69
1:A:85:ARG:HH12	1:A:87:PRO:HG3	1.58	0.69
1:A:62:GLY:O	1:A:64:SER:OG	2.12	0.68
1:A:34:HIS:CD2	1:A:43:HIS:CD2	2.81	0.66
1:C:131:THR:HG22	1:C:135:LYS:H	1.60	0.66
1:C:35:GLU:OE2	2:C:203:HOH:O	2.15	0.65
1:C:128:ARG:NH2	1:C:131:THR:O	2.30	0.65
1:D:35:GLU:OE2	2:D:201:HOH:O	2.15	0.64
1:A:133:SER:OG	2:A:201:HOH:O	2.14	0.64
1:A:170:SER:HB3	1:A:173:HIS:CE1	2.32	0.63
1:C:131:THR:HG23	1:C:133:SER:H	1.65	0.62
1:A:17:VAL:HG22	1:A:61:ILE:HD13	1.83	0.61
1:A:34:HIS:NE2	1:A:43:HIS:HB3	2.16	0.61
1:D:34:HIS:CE1	1:D:43:HIS:HB3	2.37	0.60
1:B:93:THR:HB	1:B:94:ASN:HD22	1.71	0.55
1:D:66:GLU:O	1:D:69:ASP:HB2	2.06	0.55
1:D:91:VAL:O	1:D:123:ARG:NH1	2.41	0.54
1:A:34:HIS:ND1	1:A:36:VAL:HG23	2.23	0.54
1:C:69:ASP:O	1:C:85:ARG:HG2	2.08	0.53
1:D:29:PHE:HZ	1:D:31:MET:HE3	1.73	0.53
1:A:85:ARG:HG2	1:A:85:ARG:HH11	1.74	0.52
1:D:125:LYS:O	1:D:165:ARG:NH1	2.26	0.51
1:A:108:VAL:HG11	1:A:112:THR:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:89:TYR:CZ	1:D:145:CYS:HB2	2.45	0.50
1:C:99:GLN:NE2	1:C:156:ASP:OD1	2.39	0.50
1:D:17:VAL:CG2	1:D:61:ILE:HG21	2.42	0.50
1:A:170:SER:HB3	1:A:173:HIS:HE1	1.77	0.50
1:C:150:VAL:HG22	1:C:161:PRO:HG3	1.95	0.49
1:B:89:TYR:CZ	1:B:150:VAL:HG13	2.48	0.48
1:B:35:GLU:OE2	2:B:202:HOH:O	2.20	0.48
1:B:34:HIS:CE1	1:B:43:HIS:HB3	2.48	0.48
1:B:29:PHE:HD2	1:B:47:THR:HG22	1.80	0.47
1:B:94:ASN:HB2	1:B:162:MET:SD	2.54	0.47
1:A:131:THR:HB	1:A:173:HIS:HD2	1.79	0.47
1:C:103:GLU:OE2	1:C:105:LYS:NZ	2.45	0.46
1:C:174:LEU:HD12	1:C:176:ARG:HH21	1.80	0.46
1:A:131:THR:CG2	1:A:173:HIS:HD2	2.29	0.46
1:D:17:VAL:HG21	1:D:61:ILE:HG21	1.98	0.45
1:D:161:PRO:HD2	1:D:164:ILE:HD12	1.98	0.45
1:A:9:ASP:O	1:A:13:LEU:HG	2.16	0.45
1:C:89:TYR:CZ	1:C:145:CYS:HB2	2.51	0.45
1:B:69:ASP:OD1	1:B:90:LYS:NZ	2.48	0.45
1:A:34:HIS:CD2	1:A:43:HIS:HB3	2.51	0.44
1:D:17:VAL:CG2	1:D:61:ILE:HD13	2.47	0.44
1:D:137:ILE:HA	1:D:138:PRO:HD3	1.81	0.44
1:A:118:SER:HB3	1:B:146:THR:HB	1.98	0.44
1:A:85:ARG:HG2	1:A:85:ARG:NH1	2.33	0.44
1:B:170:SER:OG	1:B:173:HIS:CE1	2.71	0.44
1:C:178:TRP:HA	2:C:201:HOH:O	2.18	0.44
1:C:131:THR:HG22	1:C:135:LYS:N	2.28	0.44
1:D:17:VAL:HG22	1:D:61:ILE:HD13	1.99	0.44
1:D:122:GLY:N	1:D:143:ARG:HH11	2.16	0.44
1:D:28:THR:HB	1:D:49:ASP:HB2	1.99	0.43
1:B:108:VAL:HG11	1:B:112:THR:O	2.17	0.43
1:C:131:THR:CG2	1:C:135:LYS:H	2.29	0.43
1:A:108:VAL:HG12	1:A:109:CYS:O	2.19	0.42
1:C:161:PRO:HD2	1:C:164:ILE:HD12	2.00	0.42
1:A:129:SER:OG	2:A:202:HOH:O	2.21	0.42
1:B:102:LEU:HD22	1:B:152:PHE:HB3	2.00	0.42
1:C:88:GLY:HA3	1:C:146:THR:HG23	2.02	0.41
1:C:147:MET:HA	1:C:148:PRO:C	2.45	0.41
1:A:171:ASP:OD1	1:A:176:ARG:NH2	2.54	0.41
1:C:34:HIS:CE1	1:C:41:MET:HG3	2.56	0.41
1:B:137:ILE:HA	1:B:138:PRO:HD3	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:VAL:HG22	1:B:161:PRO:HG3	2.03	0.41
1:D:121:ASP:C	1:D:143:ARG:HE	2.29	0.41
1:C:96:PRO:O	1:C:102:LEU:HD11	2.20	0.40
1:C:137:ILE:HA	1:C:138:PRO:HD3	1.87	0.40
1:D:82:SER:OG	2:D:202:HOH:O	2.22	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)
1:C:49:ASP:OD1	1:C:118:SER:OG[4_455]	2.12	0.08

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	175/181 (97%)	169 (97%)	6 (3%)	0	100	100
1	B	175/181 (97%)	168 (96%)	7 (4%)	0	100	100
1	C	175/181 (97%)	168 (96%)	7 (4%)	0	100	100
1	D	175/181 (97%)	165 (94%)	10 (6%)	0	100	100
All	All	700/724 (97%)	670 (96%)	30 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	155/159 (98%)	151 (97%)	4 (3%)	40	46
1	B	155/159 (98%)	154 (99%)	1 (1%)	78	86
1	C	155/159 (98%)	149 (96%)	6 (4%)	28	31
1	D	155/159 (98%)	151 (97%)	4 (3%)	40	46
All	All	620/636 (98%)	605 (98%)	15 (2%)	43	49

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

Mol	Chain	Res	Type
1	A	17	VAL
1	A	34	HIS
1	A	64	SER
1	A	176	ARG
1	B	150	VAL
1	C	6	VAL
1	C	34	HIS
1	C	45	LEU
1	C	85	ARG
1	C	133	SER
1	C	150	VAL
1	D	17	VAL
1	D	85	ARG
1	D	118	SER
1	D	150	VAL

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	34	HIS
1	A	173	HIS
1	B	94	ASN
1	B	173	HIS
1	C	173	HIS
1	D	37	ASN

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	177/181 (97%)	0.53	12 (6%)	23 25	32, 44, 71, 118	0
1	B	177/181 (97%)	0.64	12 (6%)	23 25	35, 45, 75, 112	0
1	C	177/181 (97%)	1.07	30 (16%)	4 4	33, 53, 92, 116	0
1	D	177/181 (97%)	1.06	23 (12%)	7 7	36, 63, 97, 122	0
All	All	708/724 (97%)	0.82	77 (10%)	10 11	32, 50, 91, 122	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	63	THR	5.7
1	C	63	THR	5.7
1	B	169	THR	5.0
1	B	181	ALA	4.5
1	D	178	TRP	4.5
1	C	6	VAL	4.4
1	D	65	VAL	4.4
1	C	178	TRP	4.2
1	C	64	SER	4.1
1	C	65	VAL	4.1
1	A	63	THR	3.9
1	C	131	THR	3.9
1	C	49	ASP	3.8
1	A	65	VAL	3.8
1	C	5	SER	3.5
1	B	65	VAL	3.5
1	C	181	ALA	3.5
1	B	102	LEU	3.4
1	C	173	HIS	3.4
1	C	37	ASN	3.4
1	D	102	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	172	SER	3.2
1	A	102	LEU	3.1
1	C	136	ILE	3.0
1	D	71	PHE	3.0
1	D	166	PRO	3.0
1	A	62	GLY	2.9
1	B	5	SER	2.9
1	D	176	ARG	2.8
1	C	36	VAL	2.8
1	B	18	ASN	2.8
1	D	169	THR	2.8
1	C	12	ILE	2.8
1	A	34	HIS	2.8
1	C	137	ILE	2.8
1	A	5	SER	2.7
1	D	171	ASP	2.7
1	B	82	SER	2.7
1	C	180	THR	2.6
1	A	171	ASP	2.6
1	C	156	ASP	2.6
1	A	61	ILE	2.6
1	C	100	VAL	2.5
1	B	178	TRP	2.5
1	D	64	SER	2.4
1	A	64	SER	2.4
1	C	10	GLY	2.4
1	D	85	ARG	2.4
1	D	145	CYS	2.4
1	B	180	THR	2.3
1	D	6	VAL	2.3
1	B	49	ASP	2.3
1	A	18	ASN	2.3
1	C	111	GLY	2.3
1	D	126	SER	2.2
1	C	158	CYS	2.2
1	A	173	HIS	2.2
1	C	38	GLY	2.2
1	B	158	CYS	2.2
1	C	34	HIS	2.2
1	C	176	ARG	2.2
1	C	179	VAL	2.1
1	C	157	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	181	ALA	2.1
1	D	53	CYS	2.1
1	C	135	LYS	2.1
1	D	136	ILE	2.1
1	D	88	GLY	2.1
1	A	176	ARG	2.1
1	C	132	ASP	2.1
1	D	12	ILE	2.1
1	D	174	LEU	2.1
1	D	89	TYR	2.1
1	C	102	LEU	2.0
1	D	123	ARG	2.0
1	D	133	SER	2.0
1	D	168	LYS	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.