



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 03:55 pm BST

PDB ID : 4M9M
Title : NS2B-NS3 protease from dengue virus at pH 8.5
Authors : Yildiz, M.; Hardy, J.A.
Deposited on : 2013-08-14
Resolution : 1.53 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

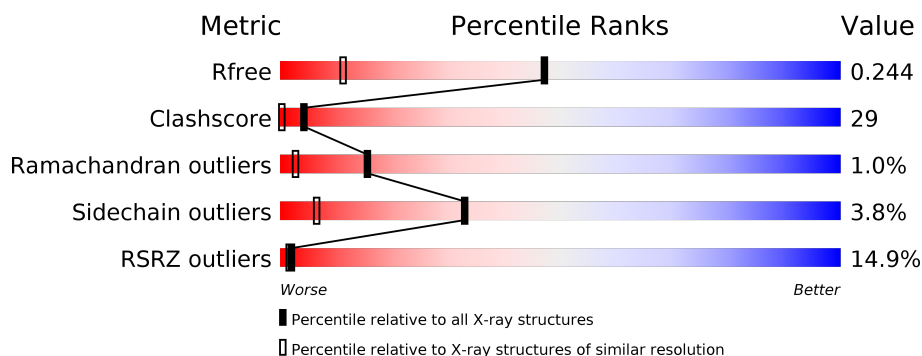
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.53 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}	130704	2556 (1.56-1.52)
Clashscore	141614	2634 (1.56-1.52)
Ramachandran outliers	138981	2580 (1.56-1.52)
Sidechain outliers	138945	2577 (1.56-1.52)
RSRZ outliers	127900	2524 (1.56-1.52)

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 on 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	247	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 1760 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 NS2B-NS3 protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	12	0
			1591	1014	270	303	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	GLY	-	EXPRESSION TAG	UNP Q91H74
A	44	SER	-	EXPRESSION TAG	UNP Q91H74
A	45	HIS	-	EXPRESSION TAG	UNP Q91H74
A	46	MET	-	EXPRESSION TAG	UNP Q91H74
A	47	LEU	-	EXPRESSION TAG	UNP Q91H74
A	48	GLU	-	EXPRESSION TAG	UNP Q91H74
A	96	GLY	-	LINKER	UNP Q91H74
A	993	GLY	-	LINKER	UNP Q91H74
A	994	GLY	-	LINKER	UNP Q91H74
A	995	GLY	-	LINKER	UNP Q91H74
A	996	SER	-	LINKER	UNP Q91H74
A	997	GLY	-	LINKER	UNP Q91H74
A	998	GLY	-	LINKER	UNP Q91H74
A	999	GLY	-	LINKER	UNP Q91H74
A	1000	GLY	-	LINKER	UNP Q91H74
A	1167	ASN	GLN	CONFLICT	UNP Q91H74

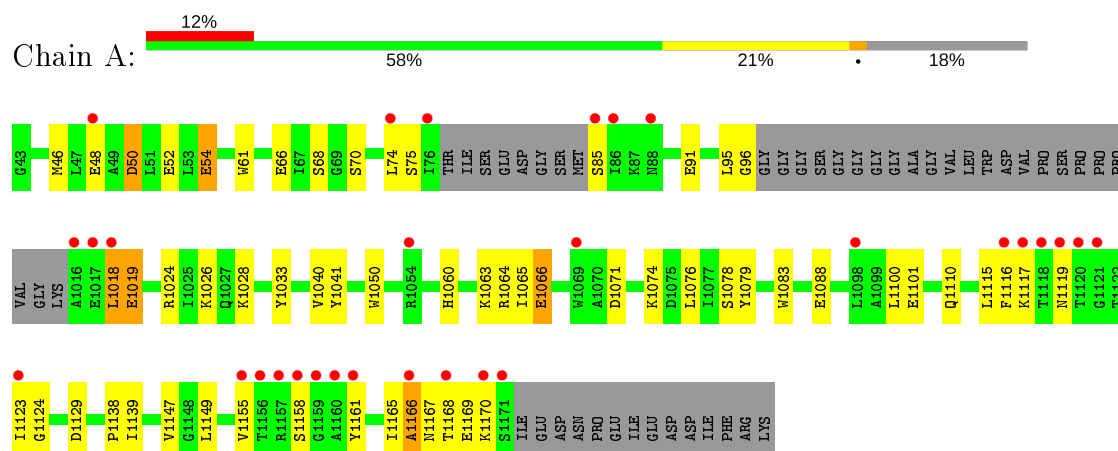
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	169	Total	O	0	0
			169	169		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: NS2B-NS3 protease



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	59.91Å 62.46Å 114.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.95 – 1.53 29.95 – 1.53	Depositor EDS
% Data completeness (in resolution range)	99.1 (29.95-1.53) 99.1 (29.95-1.53)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 1.53Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.203 , 0.250 0.199 , 0.244	Depositor DCC
R_{free} test set	1616 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 49.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.036 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	1760	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.99% 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	1.23	7/1657 (0.4%)	1.20	7/2234 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	61	TRP	CD2-CE2	8.30	1.51	1.41
1	A	91	GLU	CD-OE2	6.58	1.32	1.25
1	A	1101	GLU	CD-OE1	6.02	1.32	1.25
1	A	61	TRP	CG-CD1	5.97	1.45	1.36
1	A	1050	TRP	CD2-CE2	5.83	1.48	1.41
1	A	1079	TYR	CG-CD1	5.71	1.46	1.39
1	A	1050	TRP	CG-CD1	5.19	1.44	1.36

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1129	ASP	CB-CG-OD2	-8.79	110.39	118.30
1	A	1033	TYR	CB-CG-CD1	-8.63	115.82	121.00
1	A	1147	VAL	CG1-CB-CG2	-7.26	99.28	110.90
1	A	1024	ARG	NE-CZ-NH2	-6.17	117.22	120.30
1	A	54	GLU	OE1-CD-OE2	6.13	130.66	123.30
1	A	95	LEU	CB-CG-CD2	-5.85	101.05	111.00
1	A	50	ASP	CB-CG-OD1	5.40	123.16	118.30

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	1591	0	1607	94	0
2	A	169	0	0	54	0
All	All	1760	0	1607	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 29.

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:A:1155[B]:VAL:HG23	1:A:1161[B]:TYR:CE2	1.68	1.29
1:A:96:GLY:C	2:A:1286:HOH:O	1.70	1.27
1:A:1100:LEU:HG	2:A:1322:HOH:O	1.33	1.26
1:A:1155[B]:VAL:CG2	1:A:1161[B]:TYR:CZ	2.22	1.22
1:A:1123:ILE:HG22	2:A:1282:HOH:O	1.37	1.20
1:A:1139:ILE:HG13	2:A:1315:HOH:O	1.38	1.19
1:A:1018:LEU:HD13	1:A:1060:HIS:ND1	1.59	1.17
1:A:1155[B]:VAL:CG2	1:A:1161[B]:TYR:CE2	2.28	1.17
1:A:1123:ILE:HG13	2:A:1270:HOH:O	1.47	1.13
1:A:1123:ILE:HA	2:A:1270:HOH:O	1.48	1.09
1:A:1066:GLU:HB2	2:A:1289:HOH:O	1.52	1.08
1:A:1168:THR:HG23	2:A:1270:HOH:O	1.59	1.03
1:A:54:GLU:HB3	2:A:1345:HOH:O	1.60	1.02
1:A:1168:THR:HB	2:A:1292:HOH:O	0.82	0.98
1:A:1071:ASP:HB3	2:A:1301:HOH:O	1.61	0.98
1:A:1165:ILE:O	1:A:1168:THR:HG22	1.65	0.96
1:A:1139:ILE:CG1	2:A:1315:HOH:O	2.05	0.94
1:A:1155[B]:VAL:HG21	1:A:1161[B]:TYR:CZ	2.01	0.93
1:A:1063:LYS:HB2	2:A:1366:HOH:O	1.69	0.92
1:A:52:GLU:HB2	2:A:1306:HOH:O	1.72	0.87
1:A:1018:LEU:CD1	1:A:1060:HIS:ND1	2.39	0.86
1:A:1026:LYS:HB2	2:A:1306:HOH:O	1.75	0.86
1:A:1078:SER:CB	2:A:1303:HOH:O	2.25	0.84
1:A:1078:SER:HB3	2:A:1303:HOH:O	1.76	0.83
1:A:1018:LEU:HG	1:A:1041[B]:TYR:CZ	2.14	0.82
1:A:1100:LEU:CD2	2:A:1322:HOH:O	2.20	0.82
1:A:1155[B]:VAL:CG2	1:A:1161[B]:TYR:CE1	2.63	0.82
1:A:1155[B]:VAL:HG23	1:A:1161[B]:TYR:CZ	2.06	0.79
1:A:1124:GLY:H	1:A:1168:THR:HG21	1.47	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1124:GLY:H	1:A:1168:THR:CG2	1.98	0.77
1:A:1018:LEU:HD12	1:A:1041[B]:TYR:CE2	2.19	0.77
1:A:1074:LYS:HB2	2:A:1301:HOH:O	1.84	0.76
1:A:1166:ALA:O	1:A:1168:THR:N	2.20	0.74
1:A:50:ASP:OD2	1:A:1028:LYS:HE3	1.86	0.73
1:A:1040:VAL:HB	2:A:1323:HOH:O	1.86	0.72
1:A:1155[B]:VAL:HG22	1:A:1161[B]:TYR:CE1	2.23	0.72
1:A:1155[B]:VAL:HG22	1:A:1161[B]:TYR:CG	2.25	0.72
1:A:1155[B]:VAL:HG22	1:A:1161[B]:TYR:CD1	2.26	0.71
1:A:1155[B]:VAL:HG22	1:A:1161[B]:TYR:CZ	2.24	0.71
1:A:1018:LEU:CD1	1:A:1041[B]:TYR:CE2	2.73	0.71
1:A:1155[B]:VAL:HG22	1:A:1161[B]:TYR:CD2	2.25	0.71
1:A:1155[B]:VAL:CG2	1:A:1161[B]:TYR:CD2	2.74	0.71
1:A:54:GLU:CG	2:A:1345:HOH:O	2.39	0.70
1:A:1115:LEU:HG	2:A:1360:HOH:O	1.92	0.70
1:A:54:GLU:CB	2:A:1345:HOH:O	2.25	0.69
1:A:1155[B]:VAL:HG22	1:A:1161[B]:TYR:CE2	2.25	0.69
1:A:1018:LEU:HG	1:A:1041[B]:TYR:OH	1.93	0.68
1:A:1083:TRP:HA	2:A:1303:HOH:O	1.93	0.67
1:A:1169:GLU:HA	2:A:1270:HOH:O	1.94	0.67
1:A:1088:GLU:HG3	2:A:1241:HOH:O	1.96	0.66
1:A:1041[A]:TYR:HE1	2:A:1235:HOH:O	1.78	0.66
1:A:85:SER:HB2	2:A:1331:HOH:O	1.95	0.65
1:A:1165:ILE:O	1:A:1166:ALA:O	2.16	0.64
1:A:1116:PHE:HB2	2:A:1282:HOH:O	1.99	0.61
1:A:1018:LEU:HD13	1:A:1060:HIS:CE1	2.36	0.60
1:A:1018:LEU:HD23	1:A:1063:LYS:HD3	1.84	0.60
1:A:1155[B]:VAL:HG21	1:A:1161[B]:TYR:CE1	2.35	0.60
1:A:1064:ARG:HD3	2:A:1289:HOH:O	2.02	0.59
1:A:1076:LEU:HB2	2:A:1301:HOH:O	2.01	0.59
1:A:1119:ASN:HB2	2:A:1265:HOH:O	2.03	0.58
1:A:1019:GLU:HG2	2:A:1294:HOH:O	2.03	0.58
1:A:1149:LEU:HD21	2:A:1323:HOH:O	2.03	0.57
1:A:1065:ILE:HG23	2:A:1314:HOH:O	2.05	0.56
1:A:1018:LEU:CD2	1:A:1063:LYS:HD3	2.36	0.56
1:A:1170:LYS:HD3	2:A:1308:HOH:O	2.05	0.55
1:A:1066:GLU:CB	2:A:1289:HOH:O	2.30	0.54
1:A:1161[A]:TYR:CE1	2:A:1304:HOH:O	2.54	0.53
1:A:46[A]:MET:HE3	2:A:1357:HOH:O	2.08	0.53
1:A:1018:LEU:HG	1:A:1041[B]:TYR:CE2	2.45	0.51
1:A:1026:LYS:CB	2:A:1306:HOH:O	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1040:VAL:CB	2:A:1323:HOH:O	2.53	0.50
1:A:1115:LEU:CG	2:A:1360:HOH:O	2.55	0.49
1:A:1169:GLU:N	2:A:1270:HOH:O	2.45	0.48
1:A:1018:LEU:CD1	1:A:1041[B]:TYR:HE2	2.26	0.48
1:A:1168:THR:C	2:A:1270:HOH:O	2.52	0.48
1:A:1166:ALA:C	1:A:1168:THR:H	2.17	0.47
1:A:1165:ILE:C	1:A:1166:ALA:O	2.49	0.47
1:A:1100:LEU:HD23	2:A:1322:HOH:O	2.03	0.46
1:A:1124:GLY:N	1:A:1168:THR:HG21	2.25	0.46
1:A:48:GLU:HA	2:A:1325:HOH:O	2.16	0.45
1:A:1083:TRP:HE3	2:A:1303:HOH:O	2.00	0.45
1:A:1018:LEU:HB3	1:A:1041[A]:TYR:CZ	2.52	0.44
1:A:1115:LEU:CD2	2:A:1360:HOH:O	2.64	0.44
1:A:68[B]:SER:OG	1:A:70:SER:O	2.36	0.44
1:A:1117:LYS:CE	2:A:1330:HOH:O	2.66	0.43
1:A:1169:GLU:CA	2:A:1270:HOH:O	2.60	0.43
1:A:1064:ARG:HD2	2:A:1320:HOH:O	2.19	0.42
1:A:1124:GLY:H	1:A:1168:THR:HG23	1.78	0.42
1:A:66:GLU:HA	1:A:1110:GLN:OE1	2.19	0.42
1:A:85:SER:N	2:A:1245:HOH:O	2.52	0.42
1:A:1040:VAL:HG21	1:A:1138:PRO:HB3	2.01	0.41
1:A:1083:TRP:CA	2:A:1303:HOH:O	2.62	0.41
1:A:1040:VAL:CG2	2:A:1323:HOH:O	2.68	0.41
1:A:1018:LEU:CG	1:A:1041[B]:TYR:CE2	3.03	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	209/247 (85%)	204 (98%)	3 (1%)	2 (1%)	15 3

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1167	ASN
1	A	1166	ALA

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	168/194 (87%)	162 (96%)	6 (4%)	35 8

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	LEU
1	A	75	SER
1	A	1018	LEU
1	A	1019	GLU
1	A	1066	GLU
1	A	1158	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates 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	202/247 (81%)	0.80	30 (14%) 2 1	13, 23, 50, 106	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1016	ALA	17.7
1	A	1017	GLU	8.4
1	A	1018	LEU	6.1
1	A	1121	GLY	5.2
1	A	1166	ALA	4.8
1	A	86	ILE	4.7
1	A	1159	GLY	4.7
1	A	1158	SER	4.7
1	A	1117	LYS	4.5
1	A	1119	ASN	4.2
1	A	74	LEU	4.1
1	A	76	ILE	4.1
1	A	85	SER	4.0
1	A	1161[A]	TYR	4.0
1	A	1170	LYS	3.7
1	A	1168	THR	3.7
1	A	1120	THR	3.1
1	A	1118	THR	3.1
1	A	1157	ARG	3.0
1	A	1098[A]	LEU	2.9
1	A	1171	SER	2.9
1	A	1123	ILE	2.9
1	A	1054[A]	ARG	2.6
1	A	48	GLU	2.6
1	A	1069	TRP	2.3
1	A	1156	THR	2.3
1	A	88	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1160	ALA	2.2
1	A	1155[A]	VAL	2.2
1	A	1116	PHE	2.1

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 carbohydrates 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.