



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2023 – 12:07 PM JST

PDB ID : 8GWS
Title : SARS-CoV-2 Mpro 1-302 c145a in complex with peptide 4
Authors : Liu, M.; Fu, Z.; Huang, H.
Deposited on : 2022-09-17
Resolution : 2.90 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

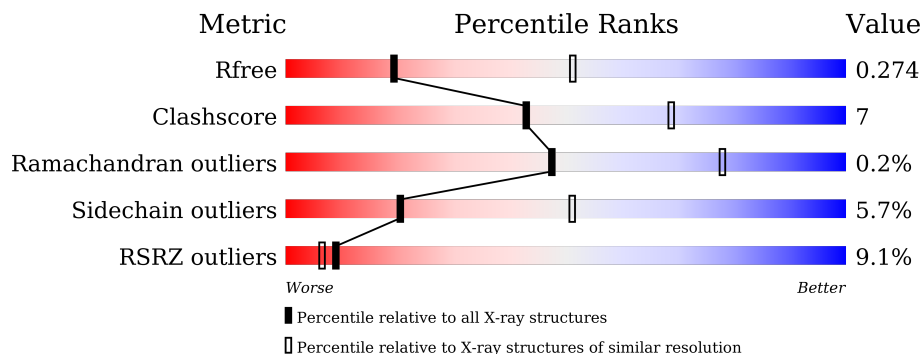
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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	302	
1	B	302	
2	C	8	
2	D	8	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4768 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 Replicase polyprotein 1ab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	300	2307	1461	392	433	21	0	0	0
1	B	302	2325	1470	397	437	21	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	145	ALA	CYS	conflict	UNP P0DTD1
B	145	ALA	CYS	conflict	UNP P0DTD1

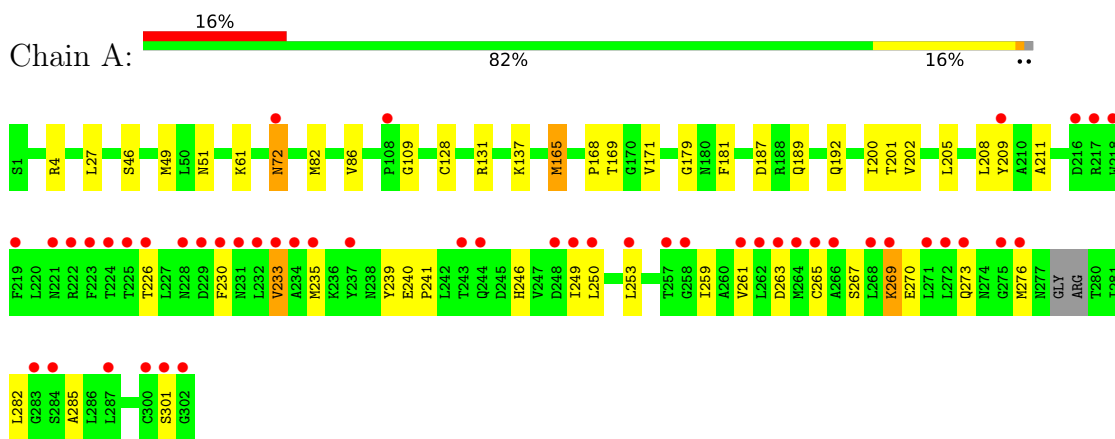
- Molecule 2 is a protein called VAL-LYS-LEU-GLN-ALA-ILE-PHE-ARG.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	8	68	46	13	9	0	0	0
2	D	8	68	46	13	9	0	0	0

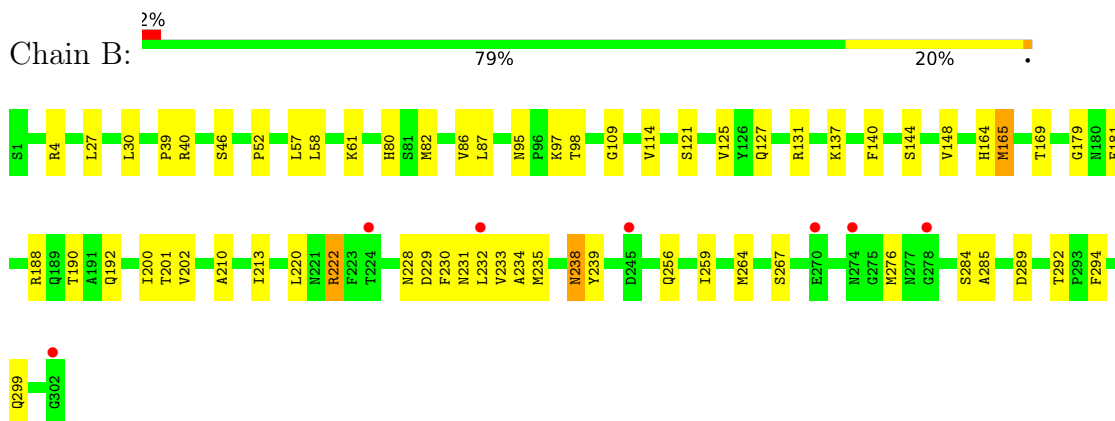
3 Residue-property plots [i](#)

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: Replicase polyprotein 1ab



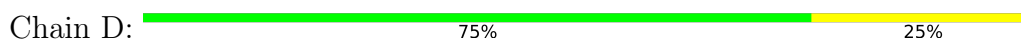
- Molecule 1: Replicase polyprotein 1ab



- Molecule 2: VAL-LYS-LEU-GLN-ALA-ILE-PHE-ARG



- Molecule 2: VAL-LYS-LEU-GLN-ALA-ILE-PHE-ARG





4 Data and refinement statistics i

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	67.03Å 67.03Å 241.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	22.42 – 2.90 22.42 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (22.42-2.90) 99.3 (22.42-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	388.82 (at 2.89Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.222 , 0.253 0.241 , 0.274	Depositor DCC
R_{free} test set	647 reflections (4.44%)	wwPDB-VP
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 0.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.108 for -h,-k,l	Xtriage
Reported twinning fraction	0.502 for H, K, L 0.498 for -h,-k,l	Depositor
Outliers	0 of 14571 reflections	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	4768	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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	0.25	0/2357	0.49	0/3203
1	B	0.25	0/2376	0.48	0/3229
2	C	0.28	0/68	0.59	0/89
2	D	0.26	0/68	0.54	0/89
All	All	0.25	0/4869	0.49	0/6610

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	2307	0	2255	33	0
1	B	2325	0	2274	32	0
2	C	68	0	78	7	0
2	D	68	0	78	2	0
All	All	4768	0	4685	64	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 7.

All (64) 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:49:MET:HB3	1:A:189:GLN:HG3	1.56	0.86
1:A:109:GLY:HA2	1:A:200:ILE:HD13	1.60	0.83
1:B:222:ARG:HD3	1:B:222:ARG:H	1.46	0.80
1:A:72:ASN:N	1:A:72:ASN:OD1	2.20	0.72
1:B:109:GLY:HA2	1:B:200:ILE:HD13	1.75	0.69
1:B:235:MET:HA	1:B:238:ASN:HA	1.75	0.68
1:B:292:THR:HG22	1:B:294:PHE:H	1.58	0.68
1:B:30:LEU:HD22	1:B:148:VAL:HG11	1.75	0.66
1:A:86:VAL:HG13	1:A:179:GLY:HA2	1.82	0.62
1:A:189:GLN:NE2	2:C:4:LEU:H	1.98	0.62
1:A:205:LEU:HA	1:A:208:LEU:HD12	1.82	0.60
1:B:86:VAL:HG13	1:B:179:GLY:HA2	1.86	0.58
1:A:201:THR:HG22	1:A:239:TYR:HD2	1.68	0.57
1:A:202:VAL:HG23	1:A:250:LEU:HD11	1.85	0.57
1:A:269:LYS:NZ	1:A:270:GLU:HG2	2.19	0.57
1:B:192:GLN:HB2	2:D:2:VAL:HG21	1.86	0.56
1:A:137:LYS:HG3	1:A:171:VAL:HG12	1.90	0.54
1:A:187:ASP:OD2	1:A:187:ASP:N	2.39	0.54
1:A:4:ARG:NH2	1:B:127:GLN:O	2.31	0.54
1:B:58:LEU:HD22	1:B:82:MET:HB2	1.91	0.52
1:A:301:SER:O	1:A:301:SER:OG	2.26	0.52
1:B:165:MET:HB3	2:D:4:LEU:HD13	1.91	0.51
1:A:240:GLU:HG3	1:A:241:PRO:HD2	1.91	0.51
1:A:201:THR:HG22	1:A:239:TYR:CD2	2.46	0.51
1:B:231:ASN:HA	1:B:234:ALA:HB3	1.93	0.49
1:A:211:ALA:HA	1:A:282:LEU:HD11	1.94	0.49
1:A:189:GLN:HE22	2:C:4:LEU:H	1.60	0.49
1:A:269:LYS:HB3	1:A:269:LYS:HE3	1.63	0.49
1:A:259:ILE:HG23	1:A:263:ASP:HB3	1.95	0.48
1:B:58:LEU:HD11	1:B:80:HIS:HD2	1.78	0.48
1:B:131:ARG:HD3	1:B:137:LYS:HD2	1.95	0.47
1:A:246:HIS:O	1:A:249:ILE:N	2.42	0.47
1:B:57:LEU:O	1:B:61:LYS:HG2	2.14	0.47
1:B:220:LEU:HD11	1:B:259:ILE:HD13	1.96	0.47
1:A:233:VAL:HG11	1:A:269:LYS:HG2	1.96	0.47
1:A:285:ALA:HB3	1:B:285:ALA:HB3	1.97	0.47
1:B:228:ASN:O	1:B:232:LEU:N	2.34	0.47
1:A:189:GLN:HE21	2:C:4:LEU:HD23	1.79	0.46
1:A:165:MET:HB3	2:C:4:LEU:HD13	1.97	0.46
1:A:168:PRO:HD3	2:C:2:VAL:HG22	1.97	0.46
1:B:95:ASN:O	1:B:98:THR:OG1	2.29	0.46
1:B:58:LEU:HD13	1:B:87:LEU:HD21	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:230:PHE:O	1:B:234:ALA:N	2.49	0.46
1:B:52:PRO:HD2	1:B:188:ARG:CZ	2.47	0.45
1:A:269:LYS:HZ2	1:A:270:GLU:HG2	1.83	0.44
1:B:229:ASP:O	1:B:233:VAL:HG23	2.16	0.44
1:B:40:ARG:HA	1:B:87:LEU:HD12	1.98	0.44
1:A:269:LYS:HD2	1:A:273:GLN:HE22	1.83	0.44
1:B:201:THR:HG22	1:B:239:TYR:HD2	1.82	0.44
1:B:210:ALA:HA	1:B:213:ILE:HD12	2.01	0.43
1:A:61:LYS:HA	1:A:61:LYS:HD3	1.66	0.42
1:B:4:ARG:O	1:B:299:GLN:NE2	2.46	0.42
1:B:188:ARG:HG3	1:B:190:THR:HG23	2.01	0.42
1:A:192:GLN:H	2:C:2:VAL:HG21	1.84	0.42
2:C:7:ILE:HD13	2:C:7:ILE:HA	1.90	0.41
1:B:292:THR:HG22	1:B:294:PHE:N	2.30	0.41
1:A:253:LEU:HD23	1:A:253:LEU:HA	1.90	0.41
1:A:230:PHE:HD1	1:A:265:CYS:HB3	1.85	0.41
1:B:140:PHE:HB3	1:B:144:SER:OG	2.20	0.41
1:A:246:HIS:O	1:A:250:LEU:HD12	2.20	0.41
1:B:39:PRO:HB3	1:B:164:HIS:CD2	2.56	0.41
1:A:209:TYR:OH	1:A:261:VAL:HA	2.21	0.41
1:B:200:ILE:HD12	1:B:289:ASP:O	2.21	0.40
1:B:114:VAL:O	1:B:125:VAL:HA	2.21	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	296/302 (98%)	278 (94%)	18 (6%)	0	100	100
1	B	300/302 (99%)	287 (96%)	12 (4%)	1 (0%)	41	71
2	C	6/8 (75%)	5 (83%)	1 (17%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	6/8 (75%)	6 (100%)	0	0	100	100
All	All	608/620 (98%)	576 (95%)	31 (5%)	1 (0%)	47	78

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	238	ASN

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	255/258 (99%)	239 (94%)	16 (6%)	18	46
1	B	257/258 (100%)	243 (95%)	14 (5%)	22	54
2	C	7/7 (100%)	7 (100%)	0	100	100
2	D	7/7 (100%)	7 (100%)	0	100	100
All	All	526/530 (99%)	496 (94%)	30 (6%)	20	51

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	46	SER
1	A	51	ASN
1	A	72	ASN
1	A	82	MET
1	A	128	CYS
1	A	131	ARG
1	A	165	MET
1	A	169	THR
1	A	181	PHE
1	A	226	THR
1	A	233	VAL
1	A	235	MET

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Mol	Chain	Res	Type
1	A	267	SER
1	A	269	LYS
1	A	276	MET
1	B	27	LEU
1	B	46	SER
1	B	97	LYS
1	B	121	SER
1	B	165	MET
1	B	169	THR
1	B	181	PHE
1	B	202	VAL
1	B	222	ARG
1	B	256	GLN
1	B	264	MET
1	B	267	SER
1	B	276	MET
1	B	284	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	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 monosaccharides 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	300/302 (99%)	0.77	49 (16%) 1 1	11, 29, 102, 123	0
1	B	302/302 (100%)	0.13	7 (2%) 60 58	6, 24, 55, 70	0
2	C	8/8 (100%)	0.39	0 100 100	15, 20, 25, 34	0
2	D	8/8 (100%)	0.28	0 100 100	18, 24, 27, 55	0
All	All	618/620 (99%)	0.45	56 (9%) 9 6	6, 26, 92, 123	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	LEU	9.2
1	A	272	LEU	6.5
1	A	226	THR	6.3
1	A	233	VAL	6.2
1	A	266	ALA	5.2
1	A	221	ASN	5.0
1	A	248	ASP	4.6
1	A	225	THR	4.6
1	A	230	PHE	4.5
1	A	273	GLN	4.5
1	A	224	THR	4.4
1	A	271	LEU	4.4
1	A	72	ASN	4.3
1	A	237	TYR	4.3
1	A	228	ASN	4.3
1	B	302	GLY	4.1
1	A	234	ALA	4.0
1	A	287	LEU	3.6
1	A	263	ASP	3.6
1	A	265	CYS	3.6
1	A	229	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	262	LEU	3.3
1	A	235	MET	3.3
1	A	219	PHE	3.2
1	A	217	ARG	3.2
1	A	268	LEU	3.1
1	A	222	ARG	2.9
1	A	302	GLY	2.9
1	A	284	SER	2.9
1	A	253	LEU	2.9
1	A	301	SER	2.9
1	B	224	THR	2.8
1	B	232	LEU	2.8
1	A	264	MET	2.7
1	A	243	THR	2.7
1	A	249	ILE	2.6
1	A	261	VAL	2.6
1	A	216	ASP	2.6
1	A	209	TYR	2.5
1	B	274	ASN	2.5
1	A	218	TRP	2.5
1	A	283	GLY	2.4
1	A	244	GLN	2.4
1	B	278	GLY	2.4
1	A	258	GLY	2.3
1	A	300	CYS	2.3
1	A	223	PHE	2.2
1	A	275	GLY	2.2
1	A	269	LYS	2.2
1	B	245	ASP	2.2
1	A	108	PRO	2.2
1	A	231	ASN	2.1
1	A	250	LEU	2.1
1	A	257	THR	2.1
1	B	270	GLU	2.0
1	A	276	MET	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 [i](#)

There are no monosaccharides 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.