Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:01 PM GMT

PDB ID	:	5C9D
Title	:	Crystal structure of a retropepsin-like aspartic protease from Rickettsia conorii
Authors	:	Li, M.; Gustchina, A.; Cruz, R.; Simoes, M.; Curto, P.; Martinez, J.; Faro, C.; Simoes, I.; Wlodawer, A.
Deposited on	:	2015-06-26
Resolution	:	2.59 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
http://wwpdb.org/validation/2016/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:

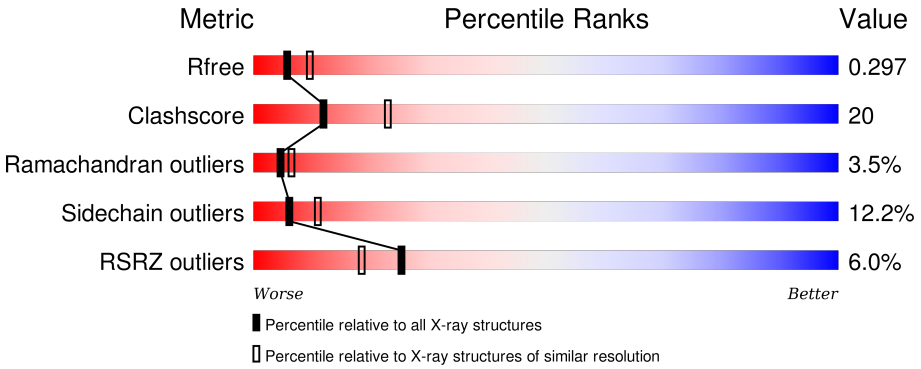
MolProbity	:	4.02b-467
Mogul	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20026982

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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. 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 <=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	139	3%, 53%, 27%, 17%
1	B	139	7%, 46%, 30%, 9%, 14%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1838 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 ApRick protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	115	Total	C	N	O	S	0	0	0
			892	576	151	162	3			
1	B	120	Total	C	N	O	S	0	0	0
			940	605	160	173	2			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	109	MET	-	initiating methionine	UNP Q92FY8
A	232	ALA	-	expression tag	UNP Q92FY8
A	233	ALA	-	expression tag	UNP Q92FY8
A	234	ALA	-	expression tag	UNP Q92FY8
A	235	LEU	-	expression tag	UNP Q92FY8
A	236	GLU	-	expression tag	UNP Q92FY8
A	237	HIS	-	expression tag	UNP Q92FY8
A	238	HIS	-	expression tag	UNP Q92FY8
A	239	HIS	-	expression tag	UNP Q92FY8
A	240	HIS	-	expression tag	UNP Q92FY8
A	241	HIS	-	expression tag	UNP Q92FY8
A	242	HIS	-	expression tag	UNP Q92FY8
B	104	MET	-	initiating methionine	UNP Q92FY8
B	232	ALA	-	expression tag	UNP Q92FY8
B	233	ALA	-	expression tag	UNP Q92FY8
B	234	ALA	-	expression tag	UNP Q92FY8
B	235	LEU	-	expression tag	UNP Q92FY8
B	236	GLU	-	expression tag	UNP Q92FY8
B	237	HIS	-	expression tag	UNP Q92FY8
B	238	HIS	-	expression tag	UNP Q92FY8
B	239	HIS	-	expression tag	UNP Q92FY8
B	240	HIS	-	expression tag	UNP Q92FY8
B	241	HIS	-	expression tag	UNP Q92FY8
B	242	HIS	-	expression tag	UNP Q92FY8

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0

- Molecule 3 is water.

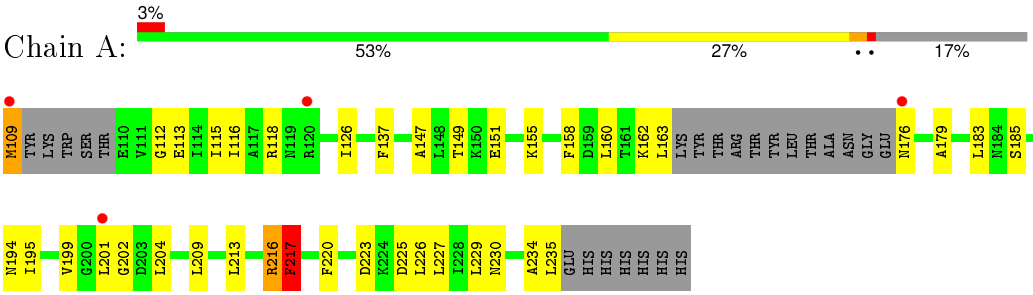
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	B	4	Total O 4 4	0	0

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 errors 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.

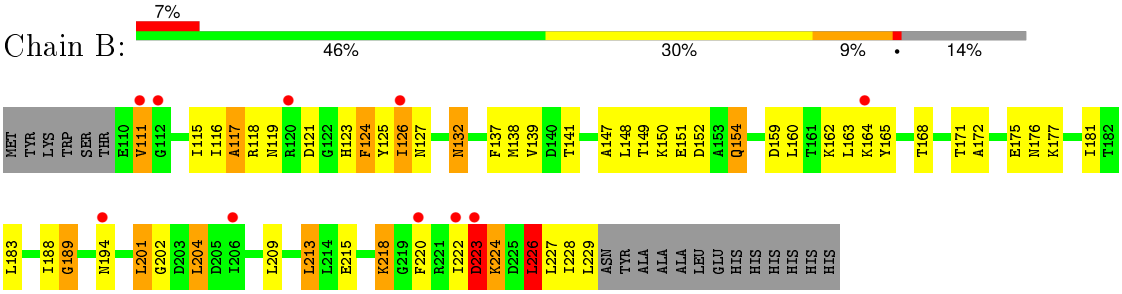
- Molecule 1: ApRick protease

Chain A:



- Molecule 1: ApRick protease

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α, β, γ	105.29Å 105.29Å 91.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.85 – 2.59 41.82 – 2.59	Depositor EDS
% Data completeness (in resolution range)	92.6 (41.85-2.59) 92.7 (41.82-2.59)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.70 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.227 , 0.298 0.234 , 0.297	Depositor DCC
R_{free} test set	353 reflections (4.84%)	DCC
Wilson B-factor (Å^2)	62.0	Xtriage
Anisotropy	0.770	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 94.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 7658 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	1838	wwPDB-VP
Average B, all atoms (Å^2)	90.0	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.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA

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	0.56	0/902	0.80	0/1209
1	B	0.55	0/952	0.84	0/1278
All	All	0.56	0/1854	0.82	0/2487

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	217	PHE	Peptide
1	B	223	ASP	Peptide

5.2 Too-close contacts

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	892	0	932	34	0
1	B	940	0	979	48	1
2	A	1	0	0	0	0
3	A	1	0	0	0	0
3	B	4	0	0	1	0
All	All	1838	0	1911	75	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 20.

All (75) 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:109:MET:HG2	1:B:124:PHE:CZ	2.07	0.90
1:A:109:MET:HG2	1:B:124:PHE:CE2	2.12	0.84
1:A:176:ASN:HD22	1:A:201:LEU:HD22	1.44	0.81
1:B:117:ALA:O	1:B:124:PHE:CD2	2.36	0.79
1:B:171:THR:HG22	1:B:172:ALA:H	1.49	0.76
1:B:150:LYS:HB2	1:B:160:LEU:HD21	1.69	0.75
1:B:119:ASN:HB2	1:B:124:PHE:CD1	2.22	0.73
1:B:117:ALA:O	1:B:124:PHE:CE2	2.44	0.70
1:B:168:THR:HG23	1:B:175:GLU:HG2	1.74	0.69
1:A:179:ALA:HB3	1:A:199:VAL:HG22	1.78	0.65
1:A:209:LEU:HD21	1:A:213:LEU:HD23	1.81	0.63
1:B:147:ALA:HB1	1:B:204:LEU:HD21	1.79	0.63
1:B:118:ARG:HA	1:B:124:PHE:HB2	1.82	0.61
1:A:158:PHE:HD2	1:A:160:LEU:HD12	1.66	0.60
1:A:189:GLY:HA2	1:B:189:GLY:HA2	1.84	0.59
1:B:213:LEU:HD12	1:B:213:LEU:O	2.03	0.59
1:B:226:LEU:CD1	1:B:228:ILE:HG13	2.33	0.59
1:A:189:GLY:CA	1:B:189:GLY:HA2	2.35	0.57
1:B:183:LEU:O	1:B:194:ASN:HA	2.04	0.57
1:A:216:ARG:O	1:A:217:PHE:CD1	2.59	0.56
1:B:171:THR:HG22	1:B:172:ALA:N	2.21	0.56
1:A:192:PHE:CD2	1:A:213:LEU:HD11	2.41	0.55
1:A:149:THR:HG22	1:A:204:LEU:HB3	1.87	0.55
1:A:115:ILE:CG2	1:A:226:LEU:HD22	2.36	0.54
1:A:223:ASP:O	1:A:225:ASP:N	2.42	0.53
1:B:116:ILE:HG22	1:B:124:PHE:HB3	1.91	0.52
1:B:116:ILE:HG22	1:B:124:PHE:CD2	2.45	0.51
1:A:192:PHE:HB3	1:A:195:ILE:HD11	1.94	0.50
1:B:137:PHE:HB3	1:B:209:LEU:HB2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:ASN:HB3	1:B:202:GLY:HA3	1.94	0.49
1:B:117:ALA:C	1:B:124:PHE:CD2	2.86	0.49
1:B:126:ILE:HG13	1:B:139:VAL:HG22	1.94	0.48
1:B:119:ASN:HB2	1:B:124:PHE:HD1	1.72	0.48
1:B:223:ASP:OD1	1:B:224:LYS:N	2.46	0.48
1:B:177:LYS:HG3	1:B:201:LEU:HD22	1.96	0.47
1:B:154:GLN:NE2	1:B:160:LEU:HD22	2.30	0.47
1:B:181:ILE:C	1:B:181:ILE:HD12	2.35	0.47
1:A:126:ILE:HD12	1:A:209:LEU:HD13	1.96	0.47
1:A:209:LEU:CG	1:A:213:LEU:HD23	2.45	0.47
1:B:127:ASN:HB2	3:B:302:HOH:O	2.14	0.47
1:A:220:PHE:HD1	1:A:227:LEU:HD11	1.79	0.46
1:B:151:GLU:HG3	1:B:152:ASP:N	2.30	0.46
1:A:209:LEU:CD2	1:A:213:LEU:HD23	2.46	0.46
1:A:209:LEU:HD11	1:A:213:LEU:HD23	1.98	0.46
1:A:115:ILE:HD12	1:B:111:VAL:HG23	1.98	0.46
1:A:201:LEU:HD23	1:A:202:GLY:N	2.31	0.46
1:B:177:LYS:CG	1:B:201:LEU:HD22	2.46	0.46
1:B:171:THR:CG2	1:B:172:ALA:H	2.24	0.45
1:A:220:PHE:CD1	1:A:227:LEU:HD11	2.52	0.45
1:B:149:THR:HG22	1:B:204:LEU:O	2.16	0.45
1:B:159:ASP:HB3	1:B:162:LYS:HB2	1.98	0.45
1:A:109:MET:HE1	1:B:125:TYR:O	2.18	0.44
1:A:213:LEU:O	1:A:216:ARG:HB2	2.17	0.44
1:B:149:THR:OG1	1:B:151:GLU:HG3	2.18	0.44
1:A:116:ILE:HD13	1:A:126:ILE:HG12	2.00	0.43
1:B:124:PHE:O	1:B:139:VAL:HG23	2.18	0.43
1:B:148:LEU:HD23	1:B:148:LEU:HA	1.91	0.42
1:A:109:MET:CG	1:B:124:PHE:CZ	2.91	0.42
1:B:218:LYS:HD2	1:B:218:LYS:N	2.34	0.42
1:A:162:LYS:CB	1:A:163:LEU:HD23	2.50	0.41
1:B:125:TYR:HA	1:B:137:PHE:O	2.20	0.41
1:B:116:ILE:HG22	1:B:124:PHE:HD2	1.84	0.41
1:B:115:ILE:HG23	1:B:228:ILE:HG12	2.01	0.41
1:A:113:GLU:HA	1:A:229:LEU:O	2.21	0.41
1:A:112:GLY:O	1:A:230:ASN:HA	2.21	0.41
1:A:216:ARG:O	1:A:217:PHE:HD1	2.02	0.41
1:B:163:LEU:HB2	1:B:165:TYR:CE1	2.56	0.41
1:A:137:PHE:HB3	1:A:209:LEU:HB2	2.03	0.41
1:A:109:MET:HB3	1:A:234:ALA:HA	2.03	0.40
1:A:183:LEU:O	1:A:194:ASN:HA	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:ASN:N	1:B:123:HIS:O	2.47	0.40
1:A:147:ALA:CB	1:A:204:LEU:HD21	2.51	0.40
1:B:121:ASP:OD2	1:B:125:TYR:OH	2.38	0.40
1:B:125:TYR:CE1	1:B:138:MET:HG3	2.56	0.40
1:B:209:LEU:C	1:B:209:LEU:HD23	2.41	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:B:220:PHE:O	1:B:220:PHE:O[16_555]	1.89	0.31

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	111/139 (80%)	97 (87%)	13 (12%)	1 (1%)	21	42
1	B	118/139 (85%)	96 (81%)	15 (13%)	7 (6%)	2	2
All	All	229/278 (82%)	193 (84%)	28 (12%)	8 (4%)	4	6

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	117	ALA
1	B	223	ASP
1	B	224	LYS
1	B	111	VAL
1	B	189	GLY
1	B	132	ASN
1	B	226	LEU
1	A	189	GLY

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	95/117 (81%)	87 (92%)	8 (8%)	14 26
1	B	101/117 (86%)	85 (84%)	16 (16%)	3 5
All	All	196/234 (84%)	172 (88%)	24 (12%)	6 11

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	MET
1	A	118	ARG
1	A	151	GLU
1	A	155	LYS
1	A	185	SER
1	A	216	ARG
1	A	217	PHE
1	A	235	LEU
1	B	124	PHE
1	B	126	ILE
1	B	132	ASN
1	B	141	THR
1	B	154	GLN
1	B	164	LYS
1	B	188	ILE
1	B	201	LEU
1	B	204	LEU
1	B	213	LEU
1	B	215	GLU
1	B	218	LYS
1	B	222	ILE
1	B	226	LEU
1	B	227	LEU
1	B	229	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	176	ASN
1	B	119	ASN
1	B	132	ASN
1	B	154	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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, 95^{th} 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	115/139 (82%)	0.19	4 (3%) 48 40	46, 79, 123, 142	0
1	B	120/139 (86%)	0.57	10 (8%) 14 9	59, 90, 147, 185	0
All	All	235/278 (84%)	0.38	14 (5%) 25 18	46, 84, 136, 185	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	223	ASP	3.8
1	A	176	ASN	3.6
1	A	120	ARG	3.1
1	A	201	LEU	2.9
1	B	206	ILE	2.8
1	A	109	MET	2.8
1	B	222	ILE	2.7
1	B	220	PHE	2.4
1	B	120	ARG	2.2
1	B	111	VAL	2.2
1	B	164	LYS	2.2
1	B	126	ILE	2.1
1	B	112	GLY	2.1
1	B	194	ASN	2.1

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 carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	NA	A	301	1/1	0.95	0.11	-1.64	88,88,88,88	0

6.5 Other polymers

There are no such residues in this entry.