Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 02:35 pm BST

PDB ID	:	6FXI
Title	:	Human PARP10 (ARTD10), catalytic fragment in complex with 3-aminobenzamide and citrate
Authors	:	Karlberg, T.; Thorsell, A.G.; Schuler, H.
Deposited on	:	2018-03-09
Resolution	:	2.60 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 ①) were used in the production of this report:

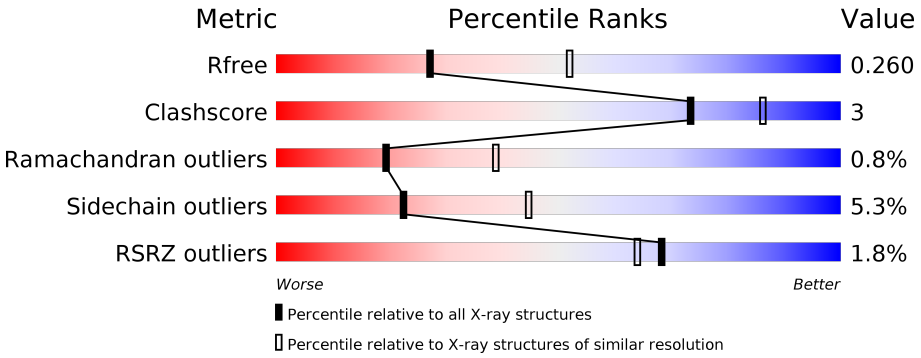
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (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 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 <=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	191	3% 88% 10%
1	B	191	% 88% 10%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3119 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.

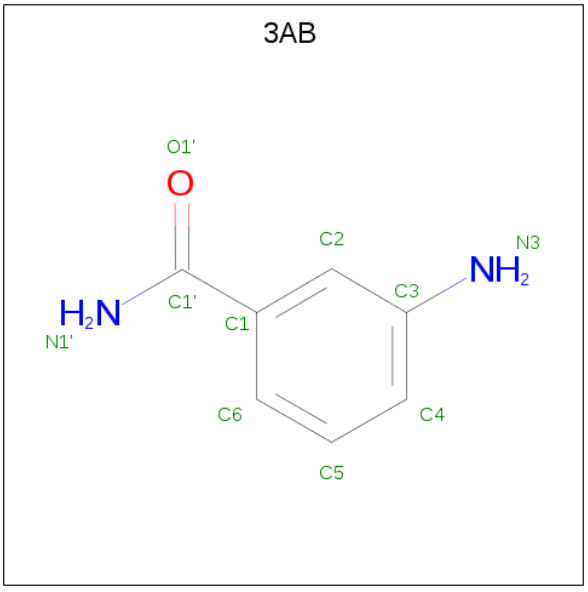
- Molecule 1 is a protein called Poly [ADP-ribose] polymerase 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	190	Total	C	N	O	S	0	0	0
			1515	949	289	270	7			
1	B	189	Total	C	N	O	S	0	0	0
			1507	944	288	269	6			

There are 4 discrepancies between the modelled and reference sequences:

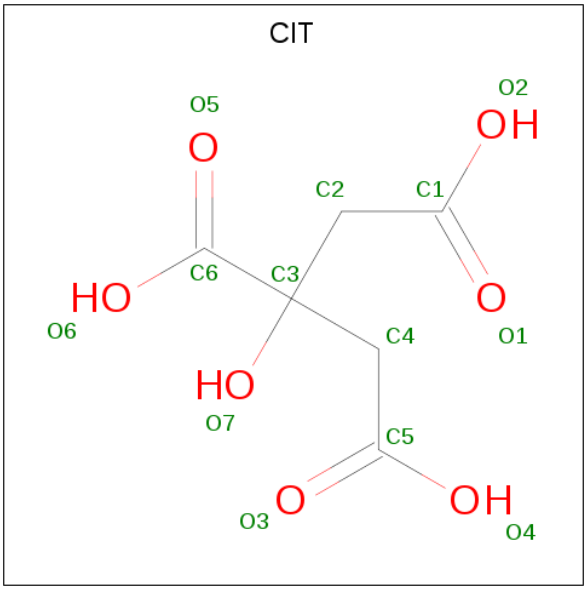
Chain	Residue	Modelled	Actual	Comment	Reference
A	817	SER	-	expression tag	UNP Q53GL7
A	818	MET	-	expression tag	UNP Q53GL7
B	817	SER	-	expression tag	UNP Q53GL7
B	818	MET	-	expression tag	UNP Q53GL7

- Molecule 2 is 3-aminobenzamide (three-letter code: 3AB) (formula: $C_7H_8N_2O$).



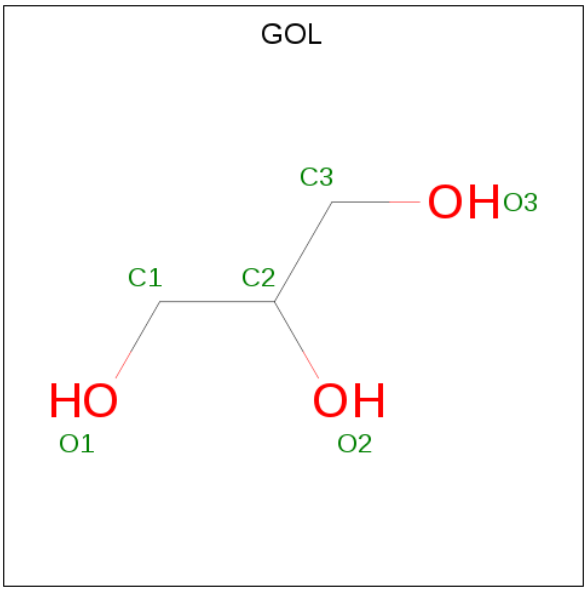
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 10 7 2 1	0	0
2	B	1	Total C N O 10 7 2 1	0	0

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 13 6 7	0	0
3	B	1	Total C O 13 6 7	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0

- Molecule 5 is water.

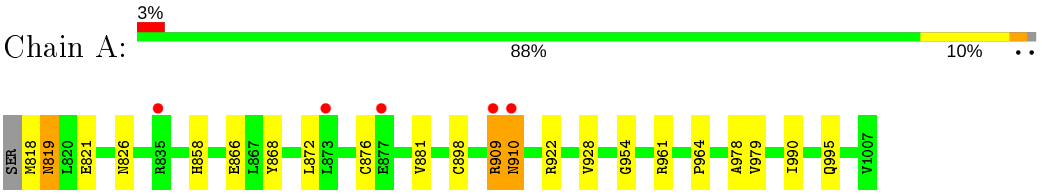
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	24	Total O 24 24	0	0
5	B	21	Total O 21 21	0	0

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 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.

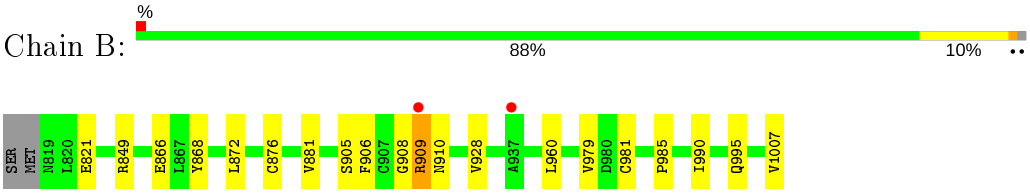
- Molecule 1: Poly [ADP-ribose] polymerase 10

Chain A:



- Molecule 1: Poly [ADP-ribose] polymerase 10

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	51.60Å 86.60Å 57.46Å 90.00° 104.98° 90.00°	Depositor
Resolution (Å)	49.85 – 2.60 49.85 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.85-2.60) 99.7 (49.85-2.60)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.26 (at 2.61Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.210 , 0.245 0.216 , 0.260	Depositor DCC
R_{free} test set	756 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	29.2	Xtriage
Anisotropy	0.864	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 37.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	3119	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: 3AB, GOL, CIT

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	0.52	0/1550	0.71	0/2103
1	B	0.50	0/1542	0.74	0/2093
All	All	0.51	0/3092	0.72	0/4196

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

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	1515	0	1493	10	0
1	B	1507	0	1484	11	0
2	A	10	0	8	2	0
2	B	10	0	8	0	0
3	A	13	0	5	0	0
3	B	13	0	5	2	0
4	A	6	0	8	1	0
5	A	24	0	0	0	0
5	B	21	0	0	0	0
All	All	3119	0	3011	20	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 (20) 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:B:979:VAL:HG21	1:B:985:PRO:HB3	1.47	0.97
1:B:979:VAL:CG2	1:B:985:PRO:HB3	2.01	0.90
1:A:990:ILE:HD13	1:A:995:GLN:HB3	1.78	0.66
1:B:990:ILE:HD13	1:B:995:GLN:HB3	1.78	0.65
1:B:909:ARG:HD2	1:B:910:ASN:HB2	1.82	0.62
1:B:909:ARG:HG3	3:B:1102:CIT:O2	2.04	0.57
1:A:909:ARG:NH1	2:A:1101:3AB:HN1'	2.06	0.54
1:B:979:VAL:HG22	1:B:985:PRO:HB3	1.90	0.54
1:A:909:ARG:HG3	1:A:910:ASN:H	1.75	0.50
1:A:964:PRO:HD3	1:A:978:ALA:HB2	1.95	0.48
1:B:872:LEU:HG	1:B:876:CYS:SG	2.54	0.47
1:B:906:PHE:HB2	3:B:1102:CIT:H42	1.97	0.47
1:A:819:ASN:HB2	1:A:898:CYS:HB3	1.97	0.46
1:A:872:LEU:HG	1:A:876:CYS:SG	2.56	0.46
1:A:954:GLY:O	1:A:979:VAL:HG22	2.18	0.43
1:B:849:ARG:HB2	1:B:1007:VAL:CG2	2.49	0.43
1:B:979:VAL:HG21	1:B:985:PRO:CB	2.32	0.42
1:A:909:ARG:HH12	2:A:1101:3AB:HN1'	1.67	0.41
1:A:858:HIS:CE1	4:A:1103:GOL:H32	2.56	0.41
1:A:961:ARG:HD2	1:B:905:SER:O	2.21	0.41

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	188/191 (98%)	179 (95%)	8 (4%)	1 (0%)	29 52
1	B	187/191 (98%)	179 (96%)	6 (3%)	2 (1%)	14 30

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	375/382 (98%)	358 (96%)	14 (4%)	3 (1%)	19 39

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	909	ARG
1	B	909	ARG
1	B	908	GLY

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	161/162 (99%)	151 (94%)	10 (6%)	18 37
1	B	160/162 (99%)	153 (96%)	7 (4%)	28 53
All	All	321/324 (99%)	304 (95%)	17 (5%)	22 45

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	818	MET
1	A	819	ASN
1	A	821	GLU
1	A	826	ASN
1	A	866	GLU
1	A	868	TYR
1	A	881	VAL
1	A	910	ASN
1	A	922	ARG
1	A	928	VAL
1	B	821	GLU
1	B	866	GLU
1	B	868	TYR
1	B	881	VAL
1	B	928	VAL

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Mol	Chain	Res	Type
1	B	960	LEU
1	B	981	CYS

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

Mol	Chain	Res	Type
1	A	910	ASN

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

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	CIT	B	1102	-	3,12,12	0.59	0	3,17,17	1.92	2 (66%)
2	3AB	A	1101	-	10,10,10	1.71	2 (20%)	13,13,13	1.46	2 (15%)
3	CIT	A	1102	-	3,12,12	0.39	0	3,17,17	1.94	1 (33%)
2	3AB	B	1101	-	10,10,10	1.81	3 (30%)	13,13,13	2.00	4 (30%)
4	GOL	A	1103	-	5,5,5	0.18	0	5,5,5	0.56	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CIT	B	1102	-	-	3/6/16/16	-
2	3AB	A	1101	-	-	0/4/4/4	0/1/1/1
3	CIT	A	1102	-	-	3/6/16/16	-
2	3AB	B	1101	-	-	1/4/4/4	0/1/1/1
4	GOL	A	1103	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1101	3AB	O1'-C1'	-3.98	1.16	1.24
2	B	1101	3AB	O1'-C1'	-3.63	1.17	1.24
2	B	1101	3AB	C1-C1'	-2.29	1.47	1.50
2	A	1101	3AB	C6-C1	-2.27	1.35	1.39
2	B	1101	3AB	C6-C1	-2.25	1.35	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1101	3AB	C1-C1'-N1'	3.89	122.42	117.75
2	A	1101	3AB	C1-C1'-N1'	3.54	122.00	117.75
2	B	1101	3AB	O1'-C1'-C1	-3.20	115.81	119.63
3	A	1102	CIT	C3-C2-C1	2.76	119.40	114.98
2	A	1101	3AB	O1'-C1'-C1	-2.42	116.74	119.63
2	B	1101	3AB	C4-C3-N3	-2.40	116.45	120.91
3	B	1102	CIT	C4-C3-C2	-2.10	103.72	109.33
2	B	1101	3AB	C5-C6-C1	2.07	122.79	120.34
3	B	1102	CIT	C3-C2-C1	2.00	118.19	114.98

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1102	CIT	C2-C3-C4-C5
3	B	1102	CIT	C6-C3-C4-C5
3	A	1102	CIT	C1-C2-C3-O7
3	A	1102	CIT	C1-C2-C3-C4
3	A	1102	CIT	C1-C2-C3-C6

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Mol	Chain	Res	Type	Atoms
4	A	1103	GOL	O1-C1-C2-C3
3	B	1102	CIT	O7-C3-C4-C5
4	A	1103	GOL	O1-C1-C2-O2
2	B	1101	3AB	C6-C1-C1'-N1'

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1102	CIT	2	0
2	A	1101	3AB	2	0
4	A	1103	GOL	1	0

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	190/191 (99%)	0.12	5 (2%)	56	50	18, 32, 63, 81	0
1	B	189/191 (98%)	0.12	2 (1%)	80	78	14, 34, 62, 76	0
All	All	379/382 (99%)	0.12	7 (1%)	68	64	14, 33, 63, 81	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	910	ASN	3.2
1	A	877	GLU	2.8
1	A	909	ARG	2.7
1	B	909	ARG	2.4
1	A	835	ARG	2.3
1	A	873	LEU	2.3
1	B	937	ALA	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 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. 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	B-factors(Å^2)	Q<0.9
3	CIT	A	1102	13/13	0.86	0.23	42,50,54,57	0
4	GOL	A	1103	6/6	0.89	0.16	25,27,28,29	0
3	CIT	B	1102	13/13	0.93	0.28	22,37,43,50	0
2	3AB	B	1101	10/10	0.97	0.18	18,26,28,29	0
2	3AB	A	1101	10/10	0.97	0.20	18,21,25,27	0

6.5 Other polymers (i)

There are no such residues in this entry.