



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 11:37 am BST

PDB ID : 6ABX
Title : Crystal structure of citrate synthase (Msed_1522) from Metallosphaera sedula
in complex with citrate
Authors : Lee, S.-H.; Son, H.-F.; Kim, K.-J.
Deposited on : 2018-07-24
Resolution : 1.70 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
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

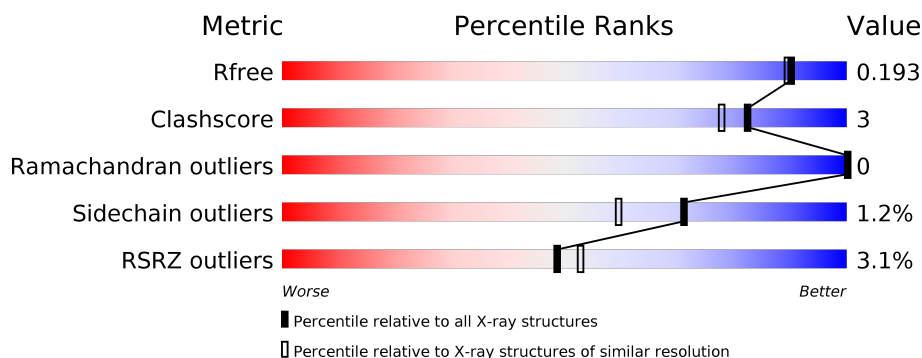
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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	378	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>7%</div> </div> </div>
1	B	378	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6555 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 Citrate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			2959	1885	503	555	16			
1	B	370	Total	C	N	O	S	0	0	0
			2959	1885	503	555	16			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	371	LEU	-	expression tag	UNP A4YGX6
A	372	GLU	-	expression tag	UNP A4YGX6
A	373	HIS	-	expression tag	UNP A4YGX6
A	374	HIS	-	expression tag	UNP A4YGX6
A	375	HIS	-	expression tag	UNP A4YGX6
A	376	HIS	-	expression tag	UNP A4YGX6
A	377	HIS	-	expression tag	UNP A4YGX6
A	378	HIS	-	expression tag	UNP A4YGX6
B	371	LEU	-	expression tag	UNP A4YGX6
B	372	GLU	-	expression tag	UNP A4YGX6
B	373	HIS	-	expression tag	UNP A4YGX6
B	374	HIS	-	expression tag	UNP A4YGX6
B	375	HIS	-	expression tag	UNP A4YGX6
B	376	HIS	-	expression tag	UNP A4YGX6
B	377	HIS	-	expression tag	UNP A4YGX6
B	378	HIS	-	expression tag	UNP A4YGX6

- Molecule 2 is CITRATE ANION (three-letter code: FLC) (formula: C₆H₅O₇).



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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

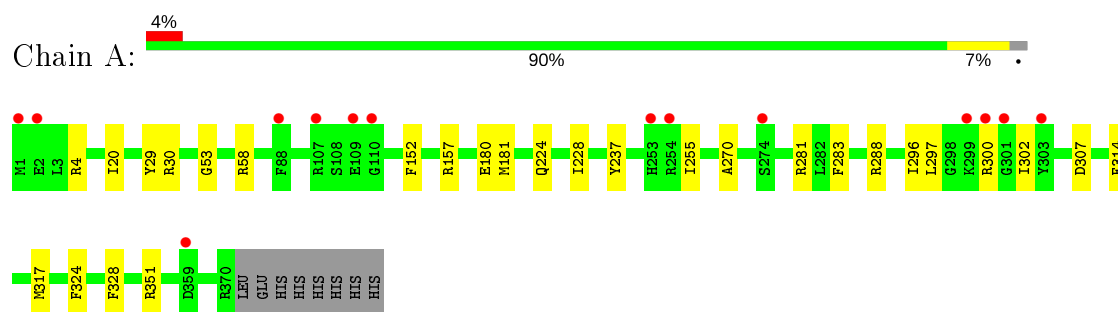
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	293	Total	O	0	0
			293	293		
4	B	300	Total	O	0	0
			300	300		

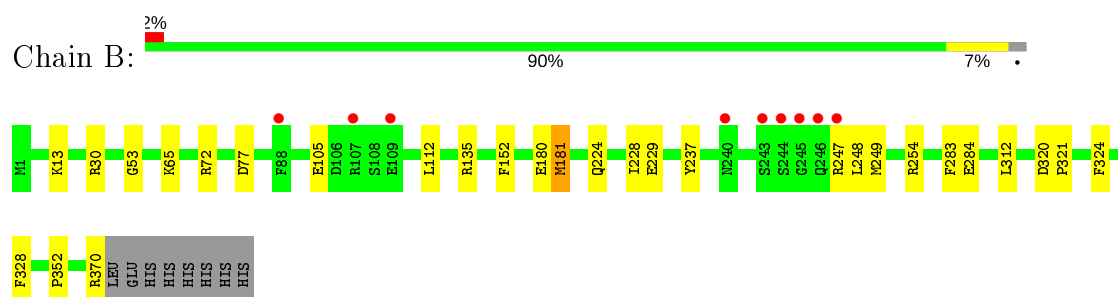
3 Residue-property plots [i](#)

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: Citrate synthase



• Molecule 1: Citrate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	50.18Å 53.52Å 76.28Å 93.56° 105.73° 102.16°	Depositor
Resolution (Å)	29.15 – 1.70 29.13 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.8 (29.15-1.70) 96.8 (29.13-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.144 , 0.180 0.158 , 0.193	Depositor DCC
R_{free} test set	3916 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	16.0	Xtriage
Anisotropy	0.369	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.9	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	0.97	EDS
Total number of atoms	6555	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.45% 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

5.1 Standard geometry

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

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	0.74	0/3017	0.84	3/4064 (0.1%)
1	B	0.69	0/3017	0.80	2/4064 (0.0%)
All	All	0.71	0/6034	0.82	5/8128 (0.1%)

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	5
1	B	0	2
All	All	0	7

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	GLU	CB-CA-C	5.63	121.65	110.40
1	B	72	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	B	370	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	317	MET	CA-CB-CG	-5.37	104.17	113.30
1	A	29	TYR	CB-CG-CD1	5.05	124.03	121.00

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	157	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	A	281	ARG	Sidechain
1	A	288	ARG	Sidechain
1	A	4	ARG	Sidechain
1	A	58	ARG	Sidechain
1	B	247	ARG	Sidechain
1	B	254	ARG	Sidechain

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	2959	0	2982	17	0
1	B	2959	0	2982	15	0
2	A	13	0	5	0	0
2	B	13	0	5	0	0
3	A	12	0	16	1	0
3	B	6	0	8	0	0
4	A	293	0	0	0	0
4	B	300	0	0	2	0
All	All	6555	0	5998	30	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 (30) 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:224:GLN:HE22	1:B:249:MET:H	1.28	0.78
1:A:314:PHE:CE2	1:A:324:PHE:HE1	2.08	0.72
1:A:181:MET:O	1:A:181:MET:HG3	2.05	0.57
1:A:314:PHE:CD2	1:A:324:PHE:CE1	2.94	0.56
1:B:135:ARG:NH2	1:B:320:ASP:OD1	2.39	0.56
1:A:181:MET:HG3	1:B:352:PRO:HG2	1.88	0.56
1:A:314:PHE:CE2	1:A:324:PHE:CE1	2.91	0.55
1:B:224:GLN:NE2	1:B:249:MET:H	2.05	0.49
1:A:181:MET:O	1:A:181:MET:CG	2.60	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LYS:HD2	4:B:609:HOH:O	2.12	0.48
1:B:321:PRO:HA	1:B:324:PHE:CZ	2.48	0.48
1:A:314:PHE:HE2	1:A:324:PHE:HE1	1.56	0.48
1:A:270:ALA:HB1	1:A:283:PHE:CE1	2.49	0.47
1:B:284:GLU:HG3	4:B:514:HOH:O	2.15	0.47
1:B:30:ARG:CD	1:B:53:GLY:HA2	2.46	0.46
1:A:307:ASP:OD2	3:A:403:GOL:H2	2.16	0.46
1:B:228:ILE:HG12	1:B:237:TYR:CG	2.50	0.45
1:A:224:GLN:HG3	1:A:237:TYR:CE1	2.51	0.45
1:A:314:PHE:CD2	1:A:324:PHE:CD1	3.05	0.45
1:B:112:LEU:C	1:B:112:LEU:HD23	2.38	0.43
1:A:296:ILE:HG22	1:A:300:ARG:NH2	2.34	0.43
1:A:30:ARG:HD2	1:A:53:GLY:HA2	2.00	0.43
1:A:181:MET:CG	1:B:352:PRO:HG2	2.49	0.43
1:B:77:ASP:OD2	1:B:105:GLU:OE2	2.37	0.42
1:A:297:LEU:HB3	1:A:302:ILE:HB	2.01	0.42
1:B:13:LYS:HD2	1:B:181:MET:HE1	2.01	0.42
1:B:283:PHE:CE2	1:B:312:LEU:HG	2.55	0.41
1:A:228:ILE:HG12	1:A:237:TYR:CG	2.56	0.41
1:B:248:LEU:HD23	1:B:248:LEU:HA	1.96	0.40
1:A:20:ILE:HG21	1:A:255:ILE:HG21	2.04	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	368/378 (97%)	359 (98%)	9 (2%)	0	100	100
1	B	368/378 (97%)	356 (97%)	12 (3%)	0	100	100
All	All	736/756 (97%)	715 (97%)	21 (3%)	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	323/331 (98%)	320 (99%)	3 (1%)	78	70
1	B	323/331 (98%)	318 (98%)	5 (2%)	65	51
All	All	646/662 (98%)	638 (99%)	8 (1%)	71	59

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	PHE
1	A	328	PHE
1	A	351	ARG
1	B	152	PHE
1	B	180	GLU
1	B	181	MET
1	B	229	GLU
1	B	328	PHE

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

Mol	Chain	Res	Type
1	B	224	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

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
2	FLC	B	401	-	3,12,12	0.98	0	3,17,17	1.65	1 (33%)
2	FLC	A	401	-	3,12,12	7.27	3 (100%)	3,17,17	5.69	3 (100%)
3	GOL	B	402	-	5,5,5	0.47	0	5,5,5	0.86	0
3	GOL	A	403	-	5,5,5	0.52	0	5,5,5	0.38	0
3	GOL	A	402	-	5,5,5	0.55	0	5,5,5	0.43	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
2	FLC	B	401	-	-	0/6/16/16	-
2	FLC	A	401	-	-	4/6/16/16	-
3	GOL	B	402	-	-	4/4/4/4	-
3	GOL	A	403	-	-	2/4/4/4	-
3	GOL	A	402	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FLC	CG-CB	-11.86	1.38	1.54
2	A	401	FLC	OHB-CB	3.26	1.48	1.43
2	A	401	FLC	CA-CB	-2.72	1.51	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FLC	CB-CA-CAC	7.68	127.28	114.98
2	A	401	FLC	CG-CB-CA	5.56	124.20	109.33
2	A	401	FLC	CB-CG-CGC	-2.71	110.65	114.98
2	B	401	FLC	CB-CG-CGC	-2.48	111.01	114.98

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	FLC	CAC-CA-CB-CBC
2	A	401	FLC	CAC-CA-CB-OHB
2	A	401	FLC	CAC-CA-CB-CG
3	B	402	GOL	O2-C2-C3-O3
3	B	402	GOL	O1-C1-C2-C3
3	B	402	GOL	C1-C2-C3-O3
3	A	403	GOL	O1-C1-C2-C3
3	A	402	GOL	C1-C2-C3-O3
3	A	403	GOL	O1-C1-C2-O2
2	A	401	FLC	CA-CB-CG-CGC
3	B	402	GOL	O1-C1-C2-O2
3	A	402	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	GOL	1	0

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	370/378 (97%)	-0.09	14 (3%)	40 45	9, 17, 38, 70	0
1	B	370/378 (97%)	-0.12	9 (2%)	59 63	11, 19, 38, 60	0
All	All	740/756 (97%)	-0.10	23 (3%)	49 53	9, 18, 39, 70	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	5.7
1	B	88	PHE	4.2
1	A	254	ARG	4.2
1	A	109	GLU	4.0
1	B	243	SER	3.5
1	A	299	LYS	3.5
1	B	109	GLU	3.4
1	A	359	ASP	3.2
1	B	245	GLY	3.1
1	A	107	ARG	3.1
1	A	301	GLY	2.9
1	A	274	SER	2.8
1	B	247	ARG	2.4
1	B	240	ASN	2.4
1	A	303	TYR	2.4
1	A	88	PHE	2.3
1	A	110	GLY	2.3
1	A	2	GLU	2.2
1	B	246	GLN	2.2
1	B	244	SER	2.1
1	A	253	HIS	2.1
1	A	300	ARG	2.1
1	B	107	ARG	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](#)

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, 95th 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(\AA^2)	Q<0.9
3	GOL	A	403	6/6	0.70	0.18	44,50,50,52	0
3	GOL	B	402	6/6	0.88	0.10	33,45,46,50	0
2	FLC	A	401	13/13	0.89	0.12	19,26,38,46	0
3	GOL	A	402	6/6	0.93	0.11	19,28,30,35	0
2	FLC	B	401	13/13	0.97	0.07	13,14,22,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.