Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 06:42 PM GMT

PDB ID	:	4MHS
Title	:	Crystal structure of Gpb in complex with SUGAR (N-[(2E)-3-(BIPHENYL-4 -YL)PROP-2-ENOYL]-BETA-D-GLUCOPYRANOSYLAMINE
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Deposited on	:	2013-08-30
Resolution	:	2.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

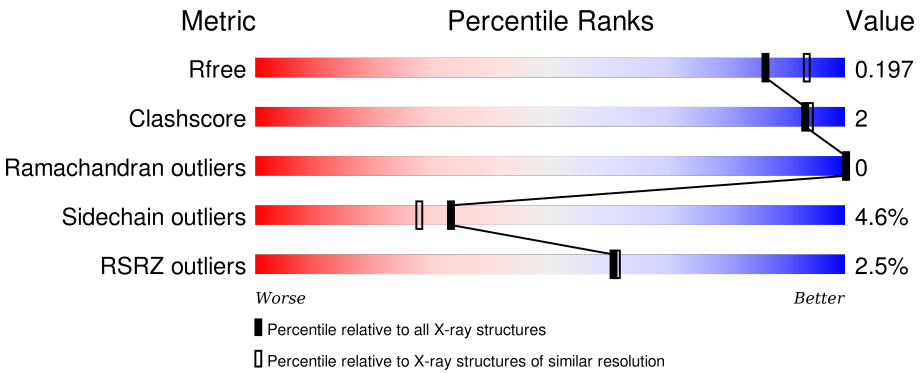
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

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 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	825	2% 91% 7% • •

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	DMS	A	902	-	-	-	X

2 Entry composition (i)

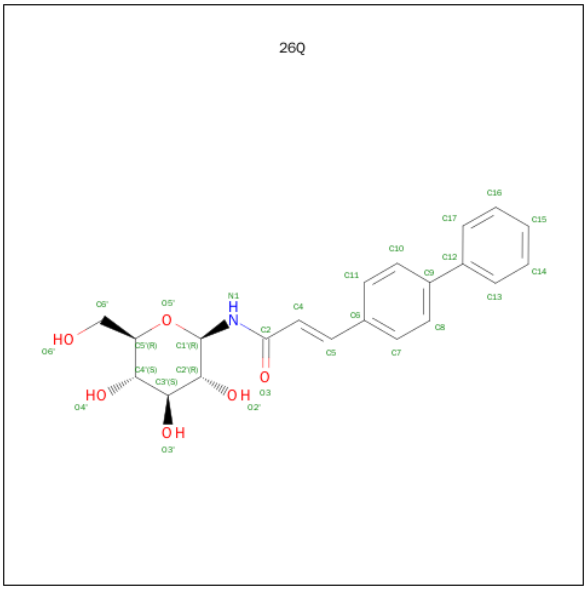
There are 4 unique types of molecules in this entry. The entry contains 6870 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycogen phosphorylase, muscle form.

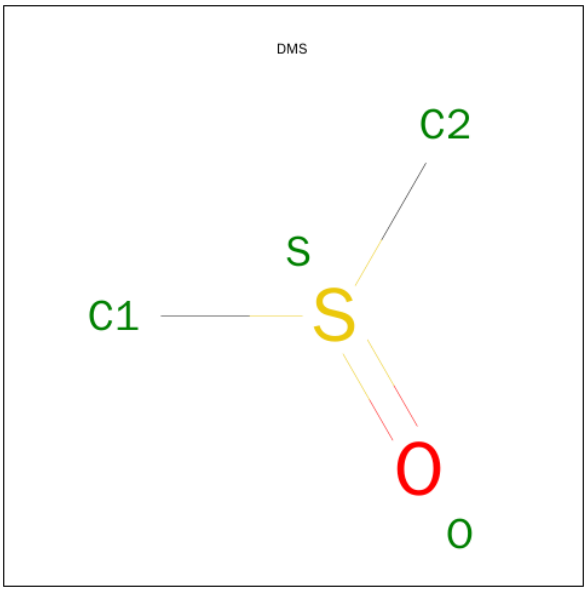
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	810	6615	4216	1163	1205	1	30	0	3	0

- Molecule 2 is SUGAR (N-[(2E)-3-(BIPHENYL-4-YL)PROP-2-ENOYL]-BETA-D-GLUCO PYRANOSYLAMINE) (three-letter code: 26Q) (formula: $C_{21}H_{23}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	28	21	1	6	0	0

- Molecule 3 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
3	A	1	4	2	1	1	0	0

- Molecule 4 is water.

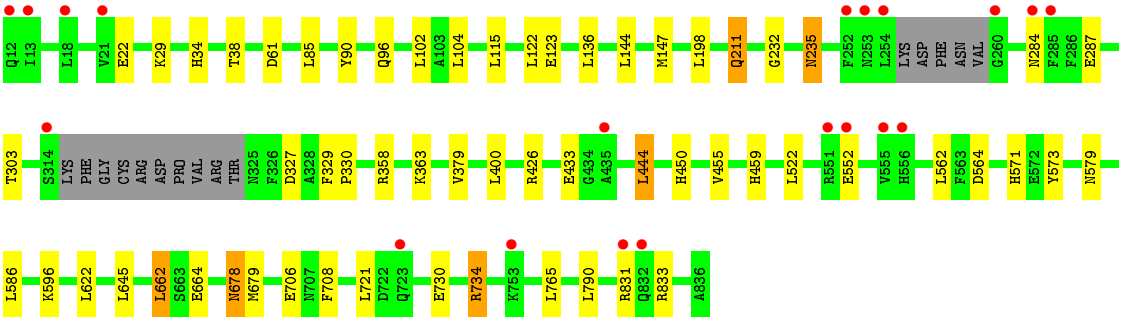
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	223	223	223	0	0

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: Glycogen phosphorylase, muscle form

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	128.87Å 128.87Å 116.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.15 – 2.00 27.55 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.4 (28.15-2.00) 95.5 (27.55-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.16 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.165 , 0.189 0.174 , 0.197	Depositor DCC
R_{free} test set	3236 reflections (5.37%)	DCC
Wilson B-factor (Å^2)	32.0	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 41.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 63546 reflections	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	6870	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 2.62% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 26Q, DMS, LLP

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	0.43	0/6748	0.64	1/9131 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	444	LEU	CA-CB-CG	5.19	127.24	115.30

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	6615	0	6549	25	0
2	A	28	0	23	0	0
3	A	4	0	6	2	0
4	A	223	0	0	4	0
All	All	6870	0	6578	25	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 2.

All (25) 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:85:LEU:HD11	1:A:303:THR:HG21	1.82	0.61
1:A:678:ASN:HD22	1:A:679:MET:H	1.49	0.60
1:A:358:ARG:NH1	4:A:1191:HOH:O	2.34	0.59
1:A:455:VAL:H	1:A:459:HIS:HD2	1.52	0.56
1:A:730:GLU:O	1:A:734:ARG:HG3	2.08	0.54
1:A:327:ASP:OD1	1:A:363:LYS:NZ	2.29	0.53
1:A:235:ASN:H	1:A:235:ASN:HD22	1.55	0.53
1:A:571[A]:HIS:HD2	1:A:573:TYR:H	1.56	0.51
1:A:34:HIS:HE1	1:A:61:ASP:OD2	1.94	0.51
1:A:379:VAL:HG22	4:A:1066:HOH:O	2.11	0.50
1:A:450:HIS:HD2	4:A:1163:HOH:O	1.96	0.48
1:A:34:HIS:HD2	1:A:38:THR:OG1	1.98	0.47
1:A:450:HIS:HE1	4:A:1146:HOH:O	1.95	0.47
1:A:329:PHE:HB3	1:A:330:PRO:HD3	1.99	0.44
1:A:284:ASN:HB2	1:A:571[B]:HIS:CE1	2.51	0.44
1:A:232:GLY:HA3	1:A:235:ASN:HD21	1.83	0.43
1:A:96:GLN:HB2	3:A:902:DMS:S	2.59	0.43
1:A:211:GLN:O	1:A:358:ARG:NH1	2.50	0.42
1:A:123:GLU:O	3:A:902:DMS:H11	2.19	0.42
1:A:136:LEU:HD23	1:A:136:LEU:C	2.40	0.42
1:A:678:ASN:HD22	1:A:679:MET:N	2.16	0.42
1:A:144:LEU:HD23	1:A:147:MET:CE	2.51	0.41
1:A:564:ASP:OD1	1:A:664:GLU:OE1	2.39	0.41
1:A:235:ASN:HD22	1:A:235:ASN:N	2.17	0.40
1:A:562:LEU:HD21	1:A:662:LEU:HB2	2.02	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	806/825 (98%)	778 (96%)	28 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	702/713 (98%)	670 (95%)	32 (5%)	33	28

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	29	LYS
1	A	90	TYR
1	A	102	LEU
1	A	104	LEU
1	A	115	LEU
1	A	122	LEU
1	A	198	LEU
1	A	211	GLN
1	A	235	ASN
1	A	287	GLU
1	A	400	LEU
1	A	426	ARG
1	A	433	GLU
1	A	444	LEU
1	A	522	LEU
1	A	552	GLU
1	A	579	ASN
1	A	586	LEU
1	A	596	LYS
1	A	622	LEU
1	A	645	LEU
1	A	662	LEU

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Mol	Chain	Res	Type
1	A	678	ASN
1	A	706	GLU
1	A	708	PHE
1	A	721	LEU
1	A	734	ARG
1	A	765	LEU
1	A	790	LEU
1	A	831	ARG
1	A	833	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	235	ASN
1	A	270	ASN
1	A	284	ASN
1	A	412	ASN
1	A	450	HIS
1	A	459	HIS
1	A	481	ASN
1	A	484	ASN
1	A	560	ASN
1	A	566	GLN
1	A	579	ASN
1	A	678	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	LLP	A	680	1	23,24,25	2.78	5 (21%)	28,32,34	1.44	5 (17%)

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
1	LLP	A	680	1	-	0/15/17/19	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	680	LLP	C4-C4'	2.37	1.50	1.46
1	A	680	LLP	C4-C5	4.88	1.48	1.42
1	A	680	LLP	C4'-NZ	5.25	1.43	1.27
1	A	680	LLP	C4-C3	5.54	1.47	1.40
1	A	680	LLP	C3-C2	9.04	1.47	1.40

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	680	LLP	OP4-P-OP1	-3.52	98.19	107.14
1	A	680	LLP	C5-C4-C4'	-2.04	118.58	121.52
1	A	680	LLP	OP3-P-OP2	2.19	115.71	107.38
1	A	680	LLP	C6-N1-C2	2.28	123.93	119.28
1	A	680	LLP	C3-C4-C4'	2.75	123.72	120.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

2 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	26Q	A	901	-	30,30,30	2.01	8 (26%)	40,41,41	0.86	0
3	DMS	A	902	-	3,3,3	0.36	0	3,3,3	0.82	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	26Q	A	901	-	-	0/15/35/35	0/3/3/3
3	DMS	A	902	-	-	0/0/0/0	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	26Q	C12-C9	-2.65	1.42	1.49
2	A	901	26Q	C6-C5	-2.19	1.40	1.47
2	A	901	26Q	C2'-C3'	2.04	1.57	1.52
2	A	901	26Q	O5'-C1'	2.15	1.46	1.43
2	A	901	26Q	C4'-C5'	2.54	1.58	1.53
2	A	901	26Q	C4-C2	3.27	1.54	1.48
2	A	901	26Q	C1'-N1	3.74	1.47	1.43
2	A	901	26Q	C4-C5	7.46	1.53	1.32

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	DMS	2	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 (i)

6.1 Protein, DNA and RNA chains (i)

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	809/825 (98%)	-0.16	20 (2%) 61 61	20, 33, 58, 108	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	252	PHE	6.3
1	A	260	GLY	5.0
1	A	253	ASN	5.0
1	A	556	HIS	5.0
1	A	285	PHE	3.8
1	A	314	SER	3.6
1	A	435	ALA	3.4
1	A	551	ARG	3.2
1	A	254	LEU	2.9
1	A	555	VAL	2.8
1	A	753	LYS	2.6
1	A	832	GLN	2.6
1	A	284	ASN	2.5
1	A	552	GLU	2.4
1	A	12	GLN	2.4
1	A	21	VAL	2.3
1	A	13	ILE	2.2
1	A	831	ARG	2.2
1	A	18	LEU	2.1
1	A	723	GLN	2.1

6.2 Non-standard residues in protein, DNA, RNA chains (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. 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, 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	LLDF	B-factors(Å^2)	Q<0.9
1	LLP	A	680	24/25	0.98	0.11	-	21,23,24,25	0

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. 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, 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	LLDF	B-factors(Å^2)	Q<0.9
3	DMS	A	902	4/4	0.67	0.37	18.61	65,69,71,73	0
2	26Q	A	901	28/28	0.90	0.18	1.01	30,37,44,45	0

6.5 Other polymers (i)

There are no such residues in this entry.