



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 04:21 am GMT

PDB ID : 5GOR  
Title : Crystal structure of alkaline invertase InvA from *Anabaena* sp. PCC 7120  
Authors : Xie, J.; Cai, K.; Hu, H.X.; Jiang, Y.L.; Yang, F.; Hu, P.F.; Chen, Y.; Zhou, C.Z.  
Deposited on : 2016-07-28  
Resolution : 2.67 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

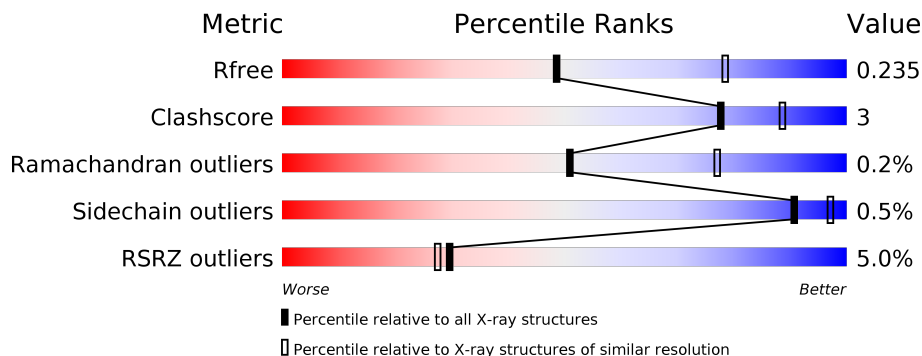
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.67 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))
$R_{free}$	100719	3050 (2.70-2.66)
Clashscore	112137	3418 (2.70-2.66)
Ramachandran outliers	110173	3367 (2.70-2.66)
Sidechain outliers	110143	3367 (2.70-2.66)
RSRZ outliers	101464	3069 (2.70-2.66)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	476	
1	B	476	
1	C	476	
1	D	476	
1	E	476	
1	F	476	

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
2	GOL	A	501	-	-	-	X
2	GOL	C	501	-	-	-	X
2	GOL	D	501	-	-	-	X
2	GOL	E	501	-	-	-	X
2	GOL	F	501	-	-	-	X
3	SO4	C	502	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22286 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 Alkaline Invertase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	457	Total	C	N	O	S	0	0	0
			3666	2380	610	657	19			
1	B	457	Total	C	N	O	S	0	0	0
			3666	2380	610	657	19			
1	C	457	Total	C	N	O	S	0	0	0
			3666	2380	610	657	19			
1	D	455	Total	C	N	O	S	0	0	0
			3652	2371	608	654	19			
1	E	457	Total	C	N	O	S	0	1	0
			3677	2390	611	657	19			
1	F	457	Total	C	N	O	S	0	0	0
			3666	2380	610	657	19			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP Q8YWS9
A	-6	GLY	-	expression tag	UNP Q8YWS9
A	-5	HIS	-	expression tag	UNP Q8YWS9
A	-4	HIS	-	expression tag	UNP Q8YWS9
A	-3	HIS	-	expression tag	UNP Q8YWS9
A	-2	HIS	-	expression tag	UNP Q8YWS9
A	-1	HIS	-	expression tag	UNP Q8YWS9
A	0	HIS	-	expression tag	UNP Q8YWS9
B	-7	MET	-	expression tag	UNP Q8YWS9
B	-6	GLY	-	expression tag	UNP Q8YWS9
B	-5	HIS	-	expression tag	UNP Q8YWS9
B	-4	HIS	-	expression tag	UNP Q8YWS9
B	-3	HIS	-	expression tag	UNP Q8YWS9
B	-2	HIS	-	expression tag	UNP Q8YWS9
B	-1	HIS	-	expression tag	UNP Q8YWS9
B	0	HIS	-	expression tag	UNP Q8YWS9
C	-7	MET	-	expression tag	UNP Q8YWS9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-6	GLY	-	expression tag	UNP Q8YWS9
C	-5	HIS	-	expression tag	UNP Q8YWS9
C	-4	HIS	-	expression tag	UNP Q8YWS9
C	-3	HIS	-	expression tag	UNP Q8YWS9
C	-2	HIS	-	expression tag	UNP Q8YWS9
C	-1	HIS	-	expression tag	UNP Q8YWS9
C	0	HIS	-	expression tag	UNP Q8YWS9
D	-7	MET	-	expression tag	UNP Q8YWS9
D	-6	GLY	-	expression tag	UNP Q8YWS9
D	-5	HIS	-	expression tag	UNP Q8YWS9
D	-4	HIS	-	expression tag	UNP Q8YWS9
D	-3	HIS	-	expression tag	UNP Q8YWS9
D	-2	HIS	-	expression tag	UNP Q8YWS9
D	-1	HIS	-	expression tag	UNP Q8YWS9
D	0	HIS	-	expression tag	UNP Q8YWS9
E	-7	MET	-	expression tag	UNP Q8YWS9
E	-6	GLY	-	expression tag	UNP Q8YWS9
E	-5	HIS	-	expression tag	UNP Q8YWS9
E	-4	HIS	-	expression tag	UNP Q8YWS9
E	-3	HIS	-	expression tag	UNP Q8YWS9
E	-2	HIS	-	expression tag	UNP Q8YWS9
E	-1	HIS	-	expression tag	UNP Q8YWS9
E	0	HIS	-	expression tag	UNP Q8YWS9
F	-7	MET	-	expression tag	UNP Q8YWS9
F	-6	GLY	-	expression tag	UNP Q8YWS9
F	-5	HIS	-	expression tag	UNP Q8YWS9
F	-4	HIS	-	expression tag	UNP Q8YWS9
F	-3	HIS	-	expression tag	UNP Q8YWS9
F	-2	HIS	-	expression tag	UNP Q8YWS9
F	-1	HIS	-	expression tag	UNP Q8YWS9
F	0	HIS	-	expression tag	UNP Q8YWS9

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		

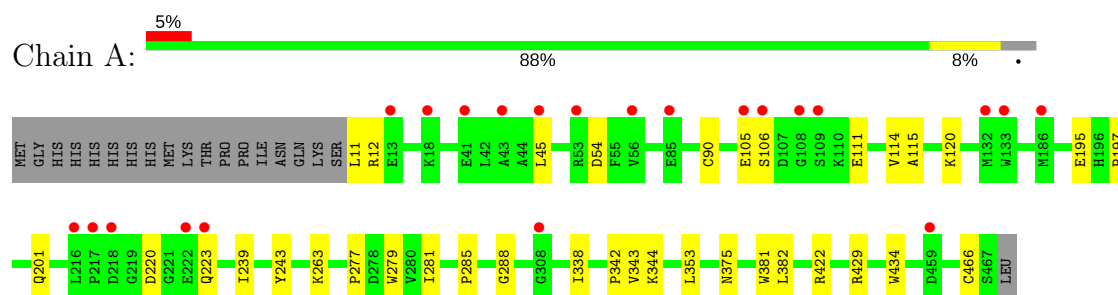
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	48	Total	O	0	0
			48	48		
4	B	46	Total	O	0	0
			46	46		
4	C	39	Total	O	0	0
			39	39		
4	D	36	Total	O	0	0
			36	36		
4	E	44	Total	O	0	0
			44	44		
4	F	39	Total	O	0	0
			39	39		

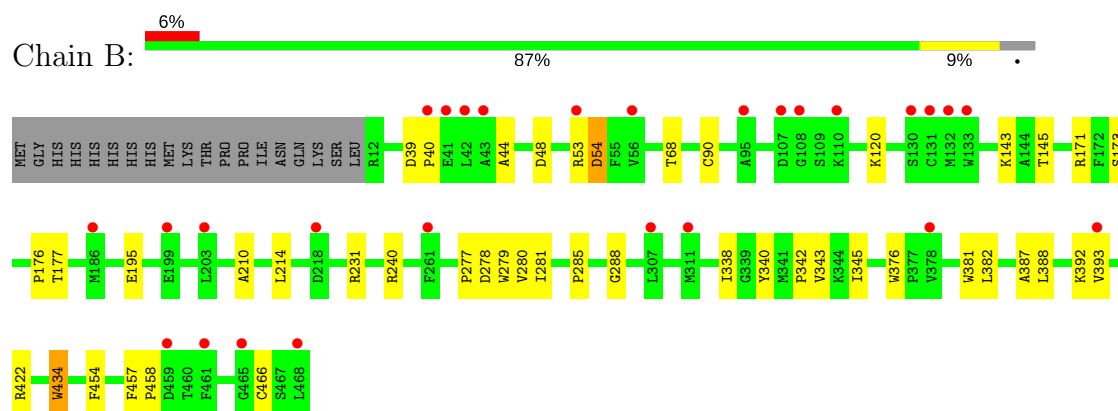
### 3 Residue-property plots [i](#)

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.

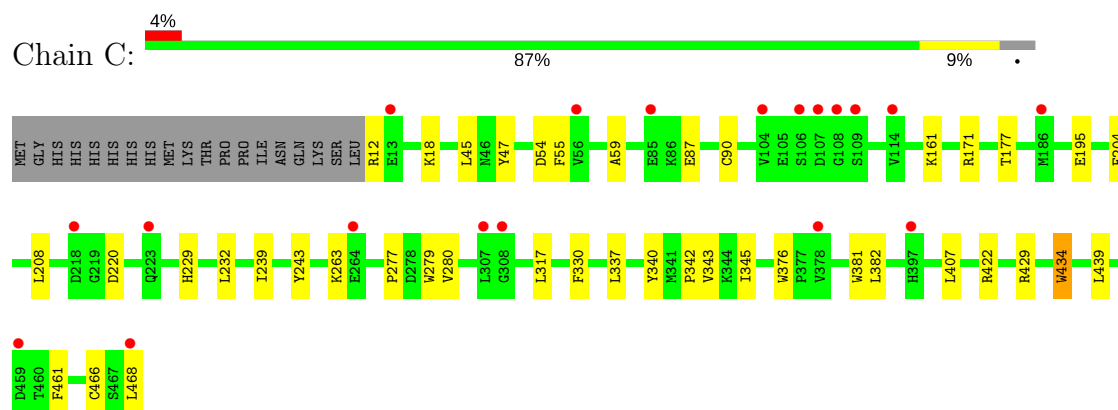
#### • Molecule 1: Alkaline Invertase



#### • Molecule 1: Alkaline Invertase

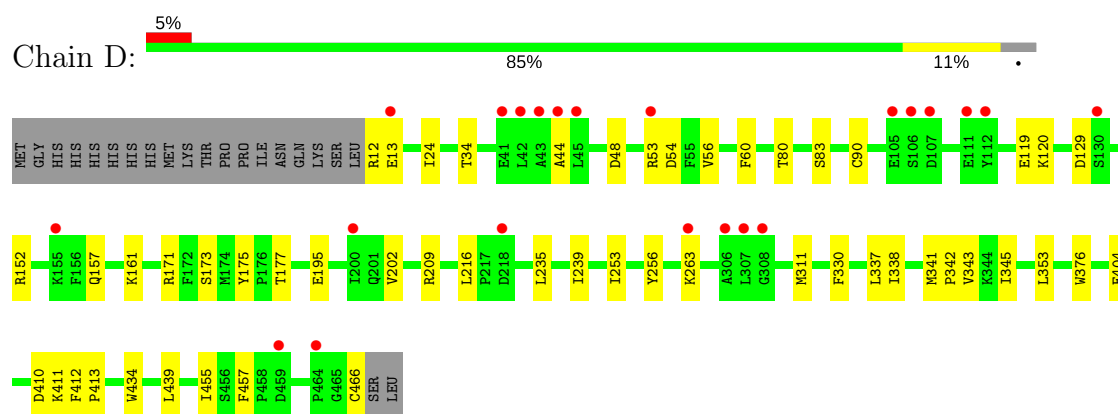


#### • Molecule 1: Alkaline Invertase

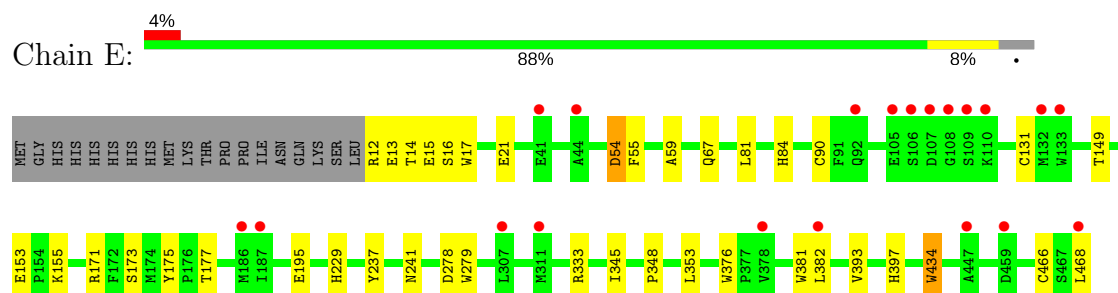


#### • Molecule 1: Alkaline Invertase

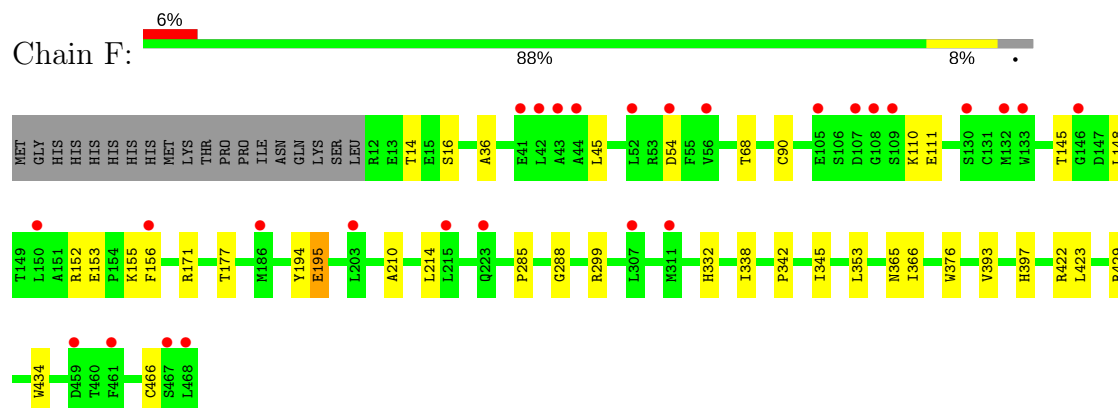




• Molecule 1: Alkaline Invertase



• Molecule 1: Alkaline Invertase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.17Å 179.28Å 96.36Å 90.00° 105.92° 90.00°	Depositor
Resolution (Å)	30.50 – 2.67 30.50 – 2.67	Depositor EDS
% Data completeness (in resolution range)	95.2 (30.50-2.67) 95.2 (30.50-2.67)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.20 (at 2.68Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.192 , 0.235 0.190 , 0.235	Depositor DCC
$R_{free}$ test set	4214 reflections (5.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	56.0	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 42.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	22286	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4

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	0.33	0/3768	0.49	0/5112
1	B	0.34	0/3768	0.49	0/5112
1	C	0.34	0/3768	0.48	0/5112
1	D	0.33	0/3754	0.48	0/5093
1	E	0.34	0/3784	0.49	0/5135
1	F	0.33	0/3768	0.49	0/5112
All	All	0.34	0/22610	0.48	0/30676

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	3666	0	3602	23	0
1	B	3666	0	3602	28	0
1	C	3666	0	3602	26	0
1	D	3652	0	3586	32	0
1	E	3677	0	3612	30	0
1	F	3666	0	3602	24	0
2	A	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	6	0	8	2	0
2	C	6	0	8	0	0
2	D	6	0	8	0	0
2	E	6	0	8	0	0
2	F	6	0	8	0	0
3	C	5	0	0	0	0
4	A	48	0	0	0	0
4	B	46	0	0	1	0
4	C	39	0	0	0	0
4	D	36	0	0	0	0
4	E	44	0	0	3	0
4	F	39	0	0	2	0
All	All	22286	0	21654	148	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 (148) 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:E:14:THR:HB	1:E:17[A]:TRP:HD1	1.51	0.76
1:A:466:CYS:HB2	1:D:90:CYS:SG	2.26	0.75
1:E:90:CYS:SG	1:F:466:CYS:HB2	2.30	0.72
1:B:466:CYS:HB2	1:C:90:CYS:SG	2.30	0.71
1:F:148:LEU:O	1:F:152:ARG:HG2	1.90	0.70
1:A:11:LEU:HG	1:A:12:ARG:H	1.56	0.70
1:A:45:LEU:HD12	1:A:429:ARG:HH12	1.57	0.70
1:B:171:ARG:NH2	1:B:177:THR:O	2.25	0.69
1:D:12:ARG:HG3	1:D:13:GLU:H	1.57	0.69
1:A:281:ILE:HD13	1:F:423:LEU:HD11	1.76	0.67
1:C:171:ARG:NH2	1:C:177:THR:O	2.28	0.66
1:D:455:ILE:O	1:D:455:ILE:HG13	1.97	0.65
1:E:171:ARG:NH2	1:E:177:THR:O	2.30	0.65
1:C:161:LYS:NZ	1:C:220:ASP:OD1	2.28	0.64
1:C:232:LEU:HB3	1:C:461:PHE:HE2	1.63	0.62
1:E:173:SER:HG	1:E:175:TYR:HD1	1.46	0.62
1:E:153:GLU:OE1	1:E:155:LYS:HE2	2.01	0.61
1:F:153:GLU:OE1	1:F:155:LYS:HE2	2.01	0.61
1:D:157:GLN:O	1:D:161:LYS:HG3	2.00	0.60
1:E:466:CYS:HB2	1:F:90:CYS:SG	2.41	0.60
1:C:87:GLU:OE2	1:C:90:CYS:HA	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:119:GLU:OE1	1:D:120:LYS:NZ	2.35	0.58
1:A:115:ALA:O	1:A:120:LYS:NZ	2.37	0.58
1:B:90:CYS:SG	1:C:466:CYS:HB2	2.44	0.57
1:A:12:ARG:O	1:A:12:ARG:HG3	2.04	0.57
1:E:171:ARG:HD2	4:E:643:HOH:O	2.04	0.56
1:B:240:ARG:NH2	4:B:601:HOH:O	2.39	0.56
1:E:278:ASP:OD1	1:E:279:TRP:N	2.39	0.56
1:D:410:ASP:O	1:D:411:LYS:HB2	2.05	0.55
1:F:36:ALA:O	4:F:601:HOH:O	2.18	0.55
1:E:12:ARG:HB3	1:E:17[A]:TRP:CD1	2.43	0.54
1:F:171:ARG:NH2	1:F:177:THR:O	2.40	0.54
1:B:53:ARG:HH11	2:B:501:GOL:H12	1.72	0.54
1:A:381:TRP:CE2	1:A:382:LEU:HD22	2.42	0.54
1:B:381:TRP:CE2	1:B:382:LEU:HD22	2.42	0.54
1:D:44:ALA:HB1	1:D:48:ASP:HB3	1.91	0.53
1:B:210:ALA:O	1:B:214:LEU:HD12	2.08	0.53
1:E:381:TRP:CE2	1:E:382:LEU:HD22	2.43	0.53
1:F:45:LEU:HD12	1:F:429:ARG:HH12	1.72	0.53
1:B:176:PRO:O	1:B:231:ARG:NH1	2.42	0.53
1:B:338:ILE:HG23	1:B:342:PRO:HB3	1.89	0.53
1:D:12:ARG:HG3	1:D:13:GLU:N	2.23	0.53
1:E:54:ASP:HA	1:E:434:TRP:CE3	2.44	0.53
1:C:381:TRP:CE2	1:C:382:LEU:HD22	2.44	0.52
1:B:44:ALA:HB1	1:B:48:ASP:HB3	1.93	0.51
1:D:171:ARG:NH2	1:D:177:THR:O	2.44	0.51
1:D:338:ILE:HG23	1:D:342:PRO:HB3	1.91	0.51
1:B:345:ILE:HB	1:B:376:TRP:CD1	2.46	0.50
1:C:47:TYR:CE1	1:C:429:ARG:NH2	2.79	0.50
1:D:24:ILE:HA	1:D:34:THR:HG22	1.94	0.50
1:D:345:ILE:HB	1:D:376:TRP:CD1	2.47	0.50
1:B:388:LEU:HD23	1:B:393:VAL:HG22	1.94	0.49
1:D:53:ARG:NH2	1:D:129:ASP:OD1	2.39	0.49
1:A:344:LYS:NZ	1:A:375:ASN:OD1	2.44	0.49
1:F:14:THR:HG22	1:F:16:SER:H	1.76	0.49
1:C:239:ILE:O	1:C:243:TYR:HB2	2.14	0.48
1:D:152:ARG:HG2	1:D:216:LEU:HD21	1.95	0.48
1:D:235:LEU:O	1:D:239:ILE:HG12	2.12	0.48
1:A:338:ILE:HG23	1:A:342:PRO:HB3	1.96	0.48
1:B:277:PRO:O	1:B:280:VAL:HG12	2.13	0.48
1:C:330:PHE:HA	1:C:337:LEU:HD12	1.94	0.48
1:F:299:ARG:HD3	4:F:608:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:209:ARG:HB3	1:D:457:PHE:CE2	2.49	0.47
1:E:13:GLU:OE1	4:E:601:HOH:O	2.20	0.47
1:F:285:PRO:HB3	1:F:332:HIS:CG	2.49	0.47
1:B:277:PRO:HB2	1:B:279:TRP:CD1	2.48	0.47
1:B:278:ASP:OD1	1:B:278:ASP:N	2.47	0.47
1:C:204:PHE:O	1:C:208:LEU:HG	2.15	0.47
1:E:173:SER:OG	1:E:175:TYR:HD1	1.98	0.47
1:A:90:CYS:HB2	1:D:466:CYS:CB	2.45	0.47
1:E:17[B]:TRP:NE1	1:E:21:GLU:OE2	2.48	0.47
1:E:81:LEU:O	1:E:84:HIS:HB2	2.15	0.47
1:A:281:ILE:O	1:F:422:ARG:NH1	2.48	0.47
1:C:12:ARG:O	1:C:18:LYS:NZ	2.48	0.46
1:A:106:SER:HB3	1:A:111:GLU:HG2	1.98	0.46
1:D:80:THR:O	1:D:83:SER:OG	2.28	0.46
1:E:14:THR:HG22	1:E:15:GLU:N	2.30	0.46
1:B:342:PRO:HA	1:B:343:VAL:HA	1.77	0.46
1:A:120:LYS:HG2	1:D:263:LYS:HG3	1.98	0.46
1:E:229:HIS:NE2	1:E:468:LEU:HD22	2.29	0.46
1:F:110:LYS:HG2	1:F:111:GLU:N	2.31	0.46
1:E:237:TYR:O	1:E:241:ASN:ND2	2.37	0.46
1:D:404:GLU:HG2	1:D:439:LEU:HD21	1.98	0.46
1:E:55:PHE:CE2	1:E:59:ALA:HB2	2.51	0.46
1:A:342:PRO:HA	1:A:343:VAL:HA	1.76	0.46
1:C:277:PRO:HB2	1:C:279:TRP:CD1	2.51	0.45
1:E:131:CYS:HB3	4:E:611:HOH:O	2.16	0.45
1:F:155:LYS:HE3	1:F:156:PHE:CE1	2.51	0.45
1:E:353:LEU:HB3	1:F:353:LEU:HD23	1.98	0.45
1:F:393:VAL:HG12	1:F:397:HIS:CD2	2.51	0.45
1:C:54:ASP:HA	1:C:434:TRP:CE3	2.52	0.45
1:A:422:ARG:NH1	1:B:281:ILE:O	2.48	0.45
1:A:263:LYS:HE2	1:D:119:GLU:HG3	1.99	0.45
1:B:68:THR:HG21	1:B:145:THR:HG22	1.97	0.45
1:C:345:ILE:HB	1:C:376:TRP:CD1	2.52	0.45
1:A:220:ASP:O	1:A:223:GLN:HG2	2.17	0.45
1:B:340:TYR:CE2	1:B:422:ARG:HG3	2.53	0.44
1:C:340:TYR:CD2	1:C:422:ARG:HG3	2.52	0.44
1:C:342:PRO:HA	1:C:343:VAL:HA	1.78	0.44
1:C:47:TYR:CZ	1:C:429:ARG:NH2	2.85	0.44
1:C:239:ILE:HG21	1:C:317:LEU:HD21	2.00	0.44
1:C:407:LEU:HD12	1:C:439:LEU:HD22	2.00	0.44
1:D:202:VAL:HG11	1:D:311:MET:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:67:GLN:O	1:E:67:GLN:HG3	2.18	0.44
1:B:457:PHE:HA	1:B:458:PRO:HD3	1.86	0.44
1:F:210:ALA:O	1:F:214:LEU:HG	2.18	0.44
1:B:120:LYS:HE3	1:C:263:LYS:HE3	2.00	0.44
1:B:285:PRO:HG2	1:B:288:GLY:HA3	2.00	0.44
1:E:149:THR:O	1:E:153:GLU:HG3	2.18	0.43
1:D:253:ILE:HA	1:D:256:TYR:CD2	2.53	0.43
1:D:341:MET:SD	1:D:342:PRO:HD2	2.58	0.43
1:B:53:ARG:HD2	2:B:501:GOL:O1	2.17	0.43
1:E:345:ILE:HB	1:E:376:TRP:CD1	2.53	0.43
1:F:155:LYS:HE3	1:F:156:PHE:HE1	1.82	0.43
1:E:279:TRP:CZ2	1:E:348:PRO:HD3	2.53	0.43
1:F:338:ILE:HG23	1:F:342:PRO:HB3	2.01	0.43
1:A:197:PRO:O	1:A:201:GLN:HG2	2.19	0.43
1:A:239:ILE:O	1:A:243:TYR:HB2	2.18	0.43
1:B:120:LYS:HG2	1:C:263:LYS:HG3	2.01	0.43
1:C:229:HIS:CE1	1:C:468:LEU:HD23	2.54	0.42
1:A:353:LEU:HD23	1:D:353:LEU:HB3	2.01	0.42
1:E:12:ARG:HG3	1:E:13:GLU:H	1.84	0.42
1:B:143:LYS:HD3	1:B:454:PHE:HA	2.02	0.42
1:C:55:PHE:CE2	1:C:59:ALA:HB2	2.55	0.42
1:F:194:TYR:CD1	1:F:195:GLU:HB2	2.54	0.42
1:A:105:GLU:HG3	1:A:114:VAL:HG21	2.01	0.42
1:D:173:SER:HB3	1:D:175:TYR:HD1	1.84	0.42
1:F:285:PRO:HG2	1:F:288:GLY:HA3	2.01	0.42
1:C:45:LEU:HD12	1:C:429:ARG:HH11	1.85	0.42
1:C:277:PRO:O	1:C:280:VAL:HG12	2.20	0.41
1:D:412:PHE:N	1:D:413:PRO:HD3	2.35	0.41
1:D:56:VAL:HG12	1:D:60:PHE:CE2	2.56	0.41
1:B:54:ASP:HA	1:B:434:TRP:CE3	2.55	0.41
1:D:152:ARG:HD3	1:D:216:LEU:HD21	2.03	0.41
1:B:387:ALA:HB1	1:B:392:LYS:O	2.21	0.41
1:D:342:PRO:HA	1:D:343:VAL:HA	1.73	0.41
1:D:152:ARG:CD	1:D:216:LEU:HD21	2.50	0.41
1:E:333:ARG:HA	1:E:333:ARG:HD3	1.85	0.41
1:F:345:ILE:HB	1:F:376:TRP:CD1	2.55	0.41
1:E:393:VAL:HG12	1:E:397:HIS:CD2	2.56	0.41
1:A:277:PRO:HB2	1:A:279:TRP:CD1	2.56	0.41
1:D:330:PHE:HA	1:D:337:LEU:HD12	2.03	0.41
1:E:14:THR:HB	1:E:17[A]:TRP:CD1	2.43	0.41
1:E:14:THR:HG22	1:E:16:SER:H	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:365:ASN:O	1:F:366:ILE:O	2.39	0.41
1:B:39:ASP:HA	1:B:40:PRO:HD3	1.95	0.40
1:F:68:THR:HG21	1:F:145:THR:HG22	2.02	0.40
1:A:285:PRO:HG2	1:A:288:GLY:HA3	2.04	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	455/476 (96%)	443 (97%)	11 (2%)	1 (0%)	51	77
1	B	455/476 (96%)	445 (98%)	9 (2%)	1 (0%)	51	77
1	C	455/476 (96%)	444 (98%)	10 (2%)	1 (0%)	51	77
1	D	453/476 (95%)	445 (98%)	7 (2%)	1 (0%)	51	77
1	E	456/476 (96%)	445 (98%)	10 (2%)	1 (0%)	51	77
1	F	455/476 (96%)	446 (98%)	8 (2%)	1 (0%)	51	77
All	All	2729/2856 (96%)	2668 (98%)	55 (2%)	6 (0%)	51	77

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	195	GLU
1	E	195	GLU
1	F	195	GLU
1	A	195	GLU
1	B	195	GLU
1	D	195	GLU



### 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	380/398 (96%)	378 (100%)	2 (0%)	91	97
1	B	380/398 (96%)	377 (99%)	3 (1%)	85	95
1	C	380/398 (96%)	379 (100%)	1 (0%)	94	98
1	D	378/398 (95%)	376 (100%)	2 (0%)	91	97
1	E	381/398 (96%)	379 (100%)	2 (0%)	91	97
1	F	380/398 (96%)	378 (100%)	2 (0%)	91	97
All	All	2279/2388 (95%)	2267 (100%)	12 (0%)	91	97

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ASP
1	A	434	TRP
1	B	54	ASP
1	B	173	SER
1	B	434	TRP
1	C	434	TRP
1	D	54	ASP
1	D	434	TRP
1	E	54	ASP
1	E	434	TRP
1	F	54	ASP
1	F	434	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

7 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	GOL	A	501	-	5,5,5	0.15	0	5,5,5	0.30	0
2	GOL	B	501	-	5,5,5	0.19	0	5,5,5	0.28	0
2	GOL	C	501	-	5,5,5	0.16	0	5,5,5	0.28	0
3	SO4	C	502	-	4,4,4	0.34	0	6,6,6	0.29	0
2	GOL	D	501	-	5,5,5	0.14	0	5,5,5	0.30	0
2	GOL	E	501	-	5,5,5	0.17	0	5,5,5	0.30	0
2	GOL	F	501	-	5,5,5	0.16	0	5,5,5	0.35	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	GOL	A	501	-	-	0/4/4/4	0/0/0/0
2	GOL	B	501	-	-	0/4/4/4	0/0/0/0
2	GOL	C	501	-	-	0/4/4/4	0/0/0/0
3	SO4	C	502	-	-	0/0/0/0	0/0/0/0
2	GOL	D	501	-	-	0/4/4/4	0/0/0/0
2	GOL	E	501	-	-	0/4/4/4	0/0/0/0
2	GOL	F	501	-	-	0/4/4/4	0/0/0/0

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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	GOL	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

### 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	457/476 (96%)	0.07	22 (4%)	31	29	42, 60, 88, 119	1 (0%)
1	B	457/476 (96%)	0.16	27 (5%)	23	21	41, 63, 87, 115	1 (0%)
1	C	457/476 (96%)	0.06	19 (4%)	37	35	42, 60, 88, 122	1 (0%)
1	D	455/476 (95%)	0.15	22 (4%)	31	29	42, 65, 93, 121	1 (0%)
1	E	457/476 (96%)	0.08	20 (4%)	35	33	43, 63, 87, 109	0
1	F	457/476 (96%)	0.13	27 (5%)	23	21	42, 65, 95, 116	1 (0%)
All	All	2740/2856 (95%)	0.11	137 (5%)	30	27	41, 63, 90, 122	5 (0%)

All (137) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	107	ASP	6.1
1	C	109	SER	5.4
1	F	43	ALA	5.3
1	D	41	GLU	5.0
1	F	109	SER	4.5
1	F	56	VAL	4.1
1	E	109	SER	3.9
1	A	459	ASP	3.9
1	E	468	LEU	3.9
1	C	218	ASP	3.9
1	A	43	ALA	3.8
1	B	133	TRP	3.8
1	F	108	GLY	3.8
1	B	186	MET	3.8
1	B	261	PHE	3.8
1	E	108	GLY	3.8
1	B	132	MET	3.7
1	C	13	GLU	3.7
1	D	464	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	E	107	ASP	3.6
1	D	106	SER	3.6
1	B	107	ASP	3.5
1	D	43	ALA	3.5
1	A	108	GLY	3.5
1	B	41	GLU	3.4
1	F	468	LEU	3.3
1	B	56	VAL	3.3
1	F	44	ALA	3.3
1	A	41	GLU	3.3
1	D	45	LEU	3.3
1	C	106	SER	3.3
1	E	459	ASP	3.2
1	B	40	PRO	3.2
1	A	218	ASP	3.2
1	F	307	LEU	3.2
1	D	44	ALA	3.1
1	C	108	GLY	3.1
1	A	85	GLU	3.1
1	F	133	TRP	3.1
1	D	307	LEU	3.1
1	D	107	ASP	3.1
1	F	459	ASP	3.1
1	F	41	GLU	3.1
1	A	109	SER	3.0
1	B	307	LEU	3.0
1	B	465	GLY	2.9
1	A	45	LEU	2.9
1	A	132	MET	2.9
1	D	13	GLU	2.9
1	F	223	GLN	2.8
1	C	307	LEU	2.8
1	F	107	ASP	2.8
1	F	186	MET	2.8
1	A	106	SER	2.8
1	A	18	LYS	2.8
1	A	133	TRP	2.8
1	D	112	TYR	2.8
1	E	307	LEU	2.8
1	D	459	ASP	2.7
1	E	106	SER	2.7
1	B	130	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	311	MET	2.7
1	B	461	PHE	2.7
1	A	56	VAL	2.6
1	C	378	VAL	2.6
1	D	42	LEU	2.6
1	E	447	ALA	2.6
1	B	378	VAL	2.5
1	C	468	LEU	2.5
1	D	111	GLU	2.5
1	D	105	GLU	2.5
1	E	105	GLU	2.5
1	B	110	LYS	2.5
1	F	105	GLU	2.5
1	B	468	LEU	2.4
1	C	85	GLU	2.4
1	F	54	ASP	2.4
1	E	41	GLU	2.4
1	C	186	MET	2.4
1	F	203	LEU	2.4
1	E	133	TRP	2.4
1	B	95	ALA	2.4
1	F	52	LEU	2.4
1	F	311	MET	2.4
1	D	218	ASP	2.4
1	B	53	ARG	2.4
1	B	203	LEU	2.3
1	F	461	PHE	2.3
1	E	186	MET	2.3
1	B	199	GLU	2.3
1	B	218	ASP	2.3
1	D	306	ALA	2.3
1	C	56	VAL	2.3
1	A	308	GLY	2.3
1	A	186	MET	2.3
1	B	42	LEU	2.3
1	C	223	GLN	2.3
1	F	42	LEU	2.3
1	A	13	GLU	2.3
1	B	108	GLY	2.2
1	F	156	PHE	2.2
1	C	104	VAL	2.2
1	C	264	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	155	LYS	2.2
1	B	43	ALA	2.2
1	C	459	ASP	2.2
1	B	459	ASP	2.2
1	F	146	GLY	2.2
1	F	130	SER	2.2
1	F	215	LEU	2.2
1	D	263	LYS	2.2
1	F	467	SER	2.2
1	E	110	LYS	2.2
1	D	130	SER	2.2
1	D	308	GLY	2.2
1	F	132	MET	2.2
1	E	378	VAL	2.1
1	A	223	GLN	2.1
1	C	397	HIS	2.1
1	A	216	LEU	2.1
1	D	53	ARG	2.1
1	E	311	MET	2.1
1	B	393	VAL	2.1
1	C	114	VAL	2.1
1	C	308	GLY	2.1
1	A	222	GLU	2.1
1	E	382	LEU	2.1
1	F	150	LEU	2.1
1	A	105	GLU	2.1
1	E	92	GLN	2.1
1	A	217	PRO	2.0
1	B	131	CYS	2.0
1	E	132	MET	2.0
1	A	53	ARG	2.0
1	D	200	ILE	2.0
1	E	187	ILE	2.0
1	E	44	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 [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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	SO4	C	502	5/5	0.88	0.29	3.86	67,68,75,92	0
2	GOL	D	501	6/6	0.94	0.44	3.09	59,60,62,63	0
2	GOL	A	501	6/6	0.93	0.39	2.49	56,58,60,60	0
2	GOL	E	501	6/6	0.92	0.36	2.46	54,59,59,61	0
2	GOL	F	501	6/6	0.94	0.42	2.29	59,61,64,68	0
2	GOL	C	501	6/6	0.93	0.32	2.20	59,59,64,66	0
2	GOL	B	501	6/6	0.94	0.35	1.62	54,60,63,64	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.