



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:40 PM GMT

PDB ID : 1Y7B
Title : BETA-D-XYLOSIDASE, A FAMILY 43 GLYCOSIDE HYDROLASE
Authors : Teplyakov, A.; Fedorov, E.; Gilliland, G.L.; Almo, S.C.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2004-12-08
Resolution : 1.60 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
<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 : 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

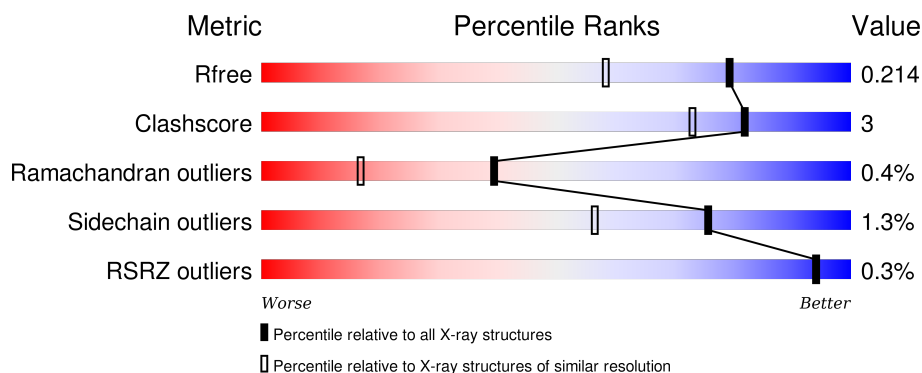
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 1.60 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}	91344	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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 $\leq 5\%$. 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	542	 90% 8% ..
1	B	542	 90% 8% .
1	C	542	 87% 10% ..
1	D	542	 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
2	CA	A	3001	-	-	-	X
2	CA	C	3003	-	-	-	X
2	CA	D	3004	-	-	-	X
3	SO4	A	3027	-	-	-	X
6	GOL	A	3017	-	-	-	X
6	GOL	B	3018	-	-	-	X
6	GOL	C	3014	-	-	-	X
6	GOL	C	3020	-	-	-	X
6	GOL	D	3019	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 20552 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 Beta-xylosidase, family 43 glycosyl hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	534	Total	C	N	O	S	0	0	0
			4341	2783	723	821	14			
1	B	534	Total	C	N	O	S	0	0	0
			4341	2783	723	821	14			
1	C	534	Total	C	N	O	S	0	0	0
			4341	2783	723	821	14			
1	D	534	Total	C	N	O	S	0	0	0
			4341	2783	723	821	14			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	-	CLONING ARTIFACT	UNP Q97DM1
A	3	LEU	-	CLONING ARTIFACT	UNP Q97DM1
A	535	GLU	LEU	CLONING ARTIFACT	UNP Q97DM1
A	536	GLY	-	CLONING ARTIFACT	UNP Q97DM1
A	537	HIS	-	EXPRESSION TAG	UNP Q97DM1
A	538	HIS	-	EXPRESSION TAG	UNP Q97DM1
A	539	HIS	-	EXPRESSION TAG	UNP Q97DM1
A	540	HIS	-	EXPRESSION TAG	UNP Q97DM1
A	541	HIS	-	EXPRESSION TAG	UNP Q97DM1
A	542	HIS	-	EXPRESSION TAG	UNP Q97DM1
B	2	SER	-	CLONING ARTIFACT	UNP Q97DM1
B	3	LEU	-	CLONING ARTIFACT	UNP Q97DM1
B	535	GLU	LEU	CLONING ARTIFACT	UNP Q97DM1
B	536	GLY	-	CLONING ARTIFACT	UNP Q97DM1
B	537	HIS	-	EXPRESSION TAG	UNP Q97DM1
B	538	HIS	-	EXPRESSION TAG	UNP Q97DM1
B	539	HIS	-	EXPRESSION TAG	UNP Q97DM1
B	540	HIS	-	EXPRESSION TAG	UNP Q97DM1
B	541	HIS	-	EXPRESSION TAG	UNP Q97DM1
B	542	HIS	-	EXPRESSION TAG	UNP Q97DM1
C	2	SER	-	CLONING ARTIFACT	UNP Q97DM1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	3	LEU	-	CLONING ARTIFACT	UNP Q97DM1
C	535	GLU	LEU	CLONING ARTIFACT	UNP Q97DM1
C	536	GLY	-	CLONING ARTIFACT	UNP Q97DM1
C	537	HIS	-	EXPRESSION TAG	UNP Q97DM1
C	538	HIS	-	EXPRESSION TAG	UNP Q97DM1
C	539	HIS	-	EXPRESSION TAG	UNP Q97DM1
C	540	HIS	-	EXPRESSION TAG	UNP Q97DM1
C	541	HIS	-	EXPRESSION TAG	UNP Q97DM1
C	542	HIS	-	EXPRESSION TAG	UNP Q97DM1
D	2	SER	-	CLONING ARTIFACT	UNP Q97DM1
D	3	LEU	-	CLONING ARTIFACT	UNP Q97DM1
D	535	GLU	LEU	CLONING ARTIFACT	UNP Q97DM1
D	536	GLY	-	CLONING ARTIFACT	UNP Q97DM1
D	537	HIS	-	EXPRESSION TAG	UNP Q97DM1
D	538	HIS	-	EXPRESSION TAG	UNP Q97DM1
D	539	HIS	-	EXPRESSION TAG	UNP Q97DM1
D	540	HIS	-	EXPRESSION TAG	UNP Q97DM1
D	541	HIS	-	EXPRESSION TAG	UNP Q97DM1
D	542	HIS	-	EXPRESSION TAG	UNP Q97DM1

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

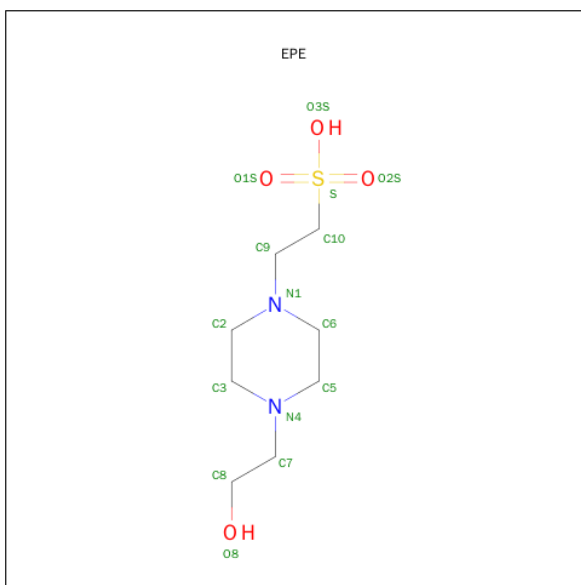
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0
2	A	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0
2	C	1	Total Ca 1 1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



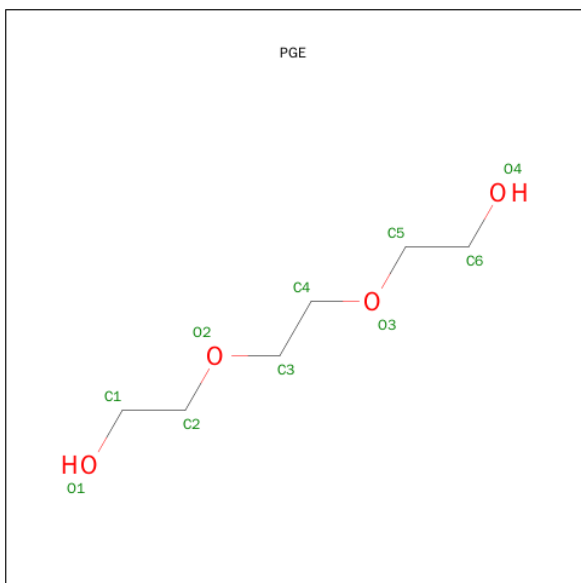
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



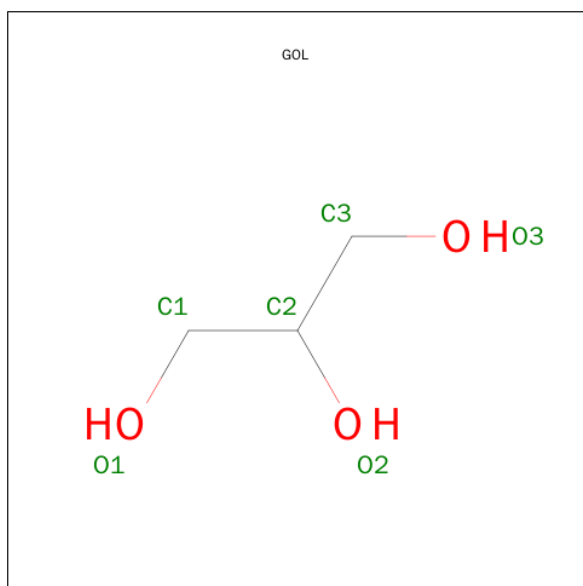
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

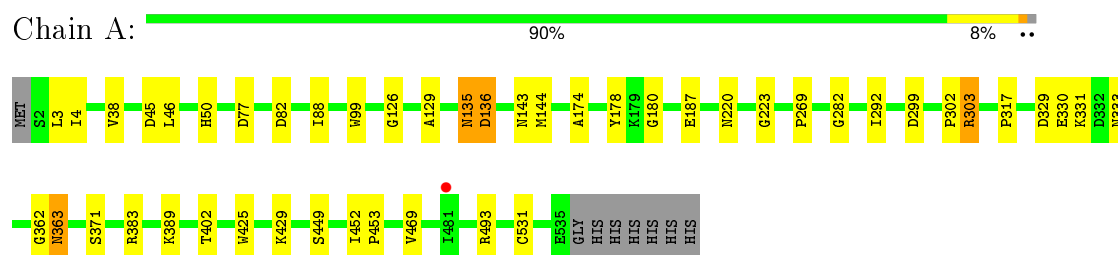
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	800	Total	O	0	0
			800	800		
7	B	704	Total	O	0	0
			704	704		
7	C	745	Total	O	0	0
			745	745		
7	D	751	Total	O	0	0
			751	751		

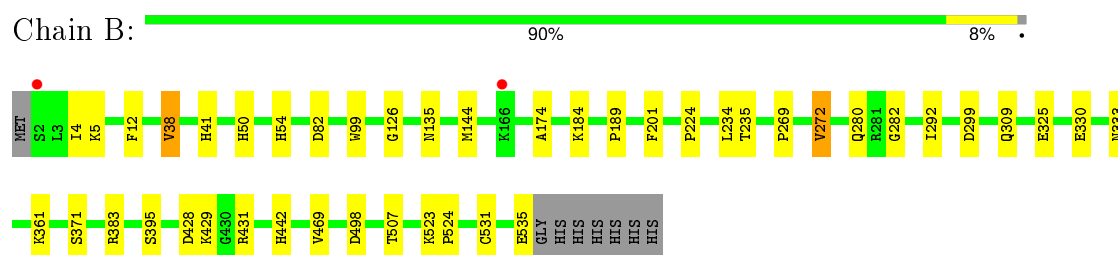
3 Residue-property plots

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.

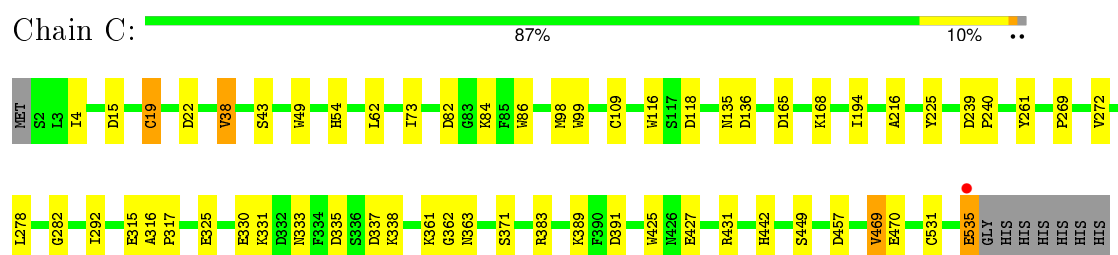
- Molecule 1: Beta-xylosidase, family 43 glycosyl hydrolase



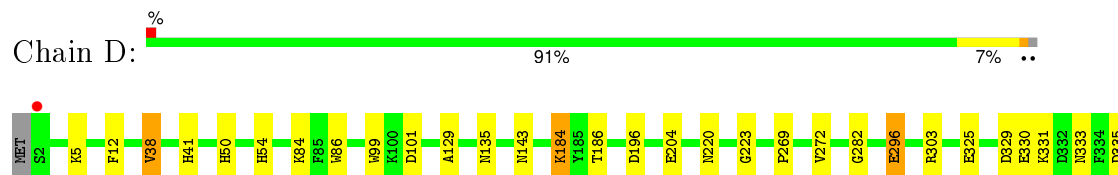
- Molecule 1: Beta-xylosidase, family 43 glycosyl hydrolase

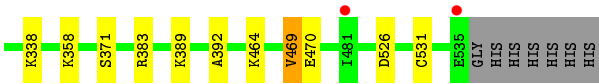


- Molecule 1: Beta-xylosidase, family 43 glycosyl hydrolase



- Molecule 1: Beta-xylosidase, family 43 glycosyl hydrolase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	94.48Å 136.15Å 183.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.60 36.47 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-1.60) 93.6 (36.47-1.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.157 , 0.206 0.164 , 0.214	Depositor DCC
R_{free} test set	5898 reflections (2.07%)	DCC
Wilson B-factor (Å ²)	14.6	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 290830 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	20552	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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, 0.375 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: EPE, GOL, CA, PGE, 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	0.71	0/4474	0.83	6/6089 (0.1%)
1	B	0.69	0/4474	0.81	4/6089 (0.1%)
1	C	0.68	1/4474 (0.0%)	0.81	9/6089 (0.1%)
1	D	0.68	0/4474	0.81	4/6089 (0.1%)
All	All	0.69	1/17896 (0.0%)	0.82	23/24356 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	19	CYS	CB-SG	-5.27	1.73	1.81

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	15	ASP	CB-CG-OD2	6.60	124.24	118.30
1	C	136	ASP	CB-CG-OD2	6.13	123.82	118.30
1	C	22	ASP	CB-CG-OD2	5.86	123.57	118.30
1	A	136	ASP	CB-CG-OD2	5.64	123.37	118.30
1	A	82	ASP	CB-CG-OD2	5.62	123.36	118.30
1	A	329	ASP	CB-CG-OD2	5.58	123.32	118.30
1	A	299	ASP	CB-CG-OD2	5.56	123.31	118.30
1	B	428	ASP	CB-CG-OD2	5.50	123.25	118.30
1	C	82	ASP	CB-CG-OD2	5.49	123.24	118.30
1	D	526	ASP	CB-CG-OD2	5.43	123.19	118.30
1	D	196	ASP	CB-CG-OD2	5.34	123.11	118.30
1	A	45	ASP	CB-CG-OD2	5.34	123.11	118.30
1	C	239	ASP	CB-CG-OD2	5.20	122.98	118.30
1	C	118	ASP	CB-CG-OD2	5.11	122.90	118.30
1	A	493	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	D	329	ASP	CB-CG-OD2	5.07	122.86	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	101	ASP	CB-CG-OD2	5.07	122.86	118.30
1	B	82	ASP	CB-CG-OD2	5.06	122.85	118.30
1	B	299	ASP	CB-CG-OD2	5.05	122.85	118.30
1	C	391	ASP	CB-CG-OD1	5.03	122.82	118.30
1	C	457	ASP	CB-CG-OD2	5.01	122.81	118.30
1	C	335	ASP	CB-CG-OD2	5.01	122.81	118.30
1	B	498	ASP	CB-CG-OD1	5.00	122.80	118.30

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	4341	0	4127	31	0
1	B	4341	0	4127	26	0
1	C	4341	0	4127	33	0
1	D	4341	0	4127	25	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	0	0
3	D	10	0	0	0	0
4	A	15	0	17	1	0
4	B	15	0	17	1	0
4	C	15	0	17	0	0
4	D	15	0	17	0	0
5	A	10	0	14	0	0
6	A	24	0	32	2	0
6	B	24	0	32	0	0
6	C	18	0	24	1	0
6	D	18	0	24	0	0
7	A	800	0	0	4	0
7	B	704	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	C	745	0	0	9	0
7	D	751	0	0	10	0
All	All	20552	0	16702	109	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 (109) 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:C:261:TYR:HB3	1:C:292:ILE:HD11	1.52	0.90
1:C:315:GLU:HG3	7:C:3300:HOH:O	1.73	0.88
1:C:442:HIS:HB2	7:C:3656:HOH:O	1.79	0.83
1:B:5:LYS:HG3	7:B:3299:HOH:O	1.81	0.81
1:A:303:ARG:HH11	1:A:303:ARG:HG3	1.51	0.75
1:D:335:ASP:HB2	7:D:3747:HOH:O	1.85	0.75
1:D:5:LYS:HG2	7:D:3707:HOH:O	1.89	0.71
1:B:442:HIS:HB2	7:B:3686:HOH:O	1.90	0.70
1:D:272:VAL:HG12	7:D:3376:HOH:O	1.90	0.70
1:C:389:LYS:HG2	1:C:535:GLU:HG3	1.74	0.69
1:C:338:LYS:HE2	7:C:3604:HOH:O	1.93	0.67
1:D:41:HIS:ND1	7:D:3341:HOH:O	2.11	0.67
1:D:269:PRO:HB3	1:D:282:GLY:HA3	1.75	0.66
1:C:449:SER:HA	7:C:3729:HOH:O	1.96	0.65
1:C:331:LYS:HE2	1:C:333:ASN:HD21	1.61	0.65
1:A:449:SER:HA	7:A:3625:HOH:O	1.97	0.65
1:B:272:VAL:HG23	7:B:3420:HOH:O	1.98	0.63
1:C:337:ASP:HB2	7:C:3752:HOH:O	2.01	0.59
1:C:84:LYS:HE3	1:C:86:TRP:CZ2	2.38	0.58
1:A:269:PRO:HB3	1:A:282:GLY:HA3	1.86	0.58
1:C:194:ILE:HG22	6:C:3020:GOL:H31	1.86	0.57
1:A:4:ILE:HG21	1:A:292:ILE:CG2	2.35	0.57
1:C:261:TYR:CB	1:C:292:ILE:HD11	2.33	0.57
1:C:269:PRO:HB3	1:C:282:GLY:HA3	1.87	0.57
1:C:272:VAL:HG23	7:C:3754:HOH:O	2.05	0.57
1:A:303:ARG:HH11	1:A:303:ARG:CG	2.17	0.57
1:B:361:LYS:HE3	7:B:3609:HOH:O	2.06	0.56
1:A:362:GLY:C	1:A:363:ASN:HD22	2.09	0.56
1:C:361:LYS:HD3	7:C:3761:HOH:O	2.07	0.55
1:A:99:TRP:CD2	1:B:371:SER:HB2	2.42	0.55
1:D:220:ASN:ND2	1:D:223:GLY:O	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:PRO:HB3	1:B:282:GLY:HA3	1.90	0.52
1:D:296:GLU:HG3	7:D:3664:HOH:O	2.09	0.52
1:A:187:GLU:OE2	6:A:3017:GOL:H12	2.10	0.52
1:D:358:LYS:HG3	7:D:3404:HOH:O	2.10	0.52
1:A:331:LYS:HE2	1:A:333:ASN:HD21	1.76	0.51
1:D:470:GLU:H	1:D:470:GLU:CD	2.12	0.51
1:C:4:ILE:HG21	1:C:292:ILE:HG23	1.94	0.50
1:A:129:ALA:HA	1:A:143:ASN:HB3	1.94	0.50
1:C:4:ILE:HG21	1:C:292:ILE:CG2	2.42	0.49
1:C:165:ASP:O	1:C:168:LYS:HD3	2.12	0.49
1:C:330:GLU:O	1:C:531:CYS:HA	2.13	0.49
1:B:126:GLY:HA3	1:B:144:MET:O	2.13	0.49
1:B:280:GLN:HB3	7:B:3543:HOH:O	2.13	0.48
1:A:302:PRO:O	1:A:303:ARG:HD3	2.14	0.48
1:B:523:LYS:HG3	1:B:524:PRO:HD2	1.96	0.47
1:A:371:SER:HB2	1:B:99:TRP:CD2	2.50	0.47
1:A:126:GLY:HA3	1:A:144:MET:O	2.14	0.47
1:B:12:PHE:CE1	1:B:507:THR:HA	2.49	0.47
1:C:371:SER:HB2	1:D:99:TRP:CD2	2.50	0.47
1:A:174:ALA:HB3	4:A:3006:EPE:H101	1.97	0.47
1:B:234:LEU:O	1:B:235:THR:HB	2.15	0.46
1:C:38:VAL:O	1:C:54:HIS:HA	2.16	0.46
1:C:389:LYS:HE3	1:C:469:VAL:HG23	1.98	0.46
1:A:99:TRP:CE3	1:B:371:SER:HB2	2.50	0.46
1:D:389:LYS:HB2	1:D:389:LYS:HE3	1.72	0.46
1:B:174:ALA:HB3	4:B:3005:EPE:H101	1.98	0.46
1:B:41:HIS:ND1	7:B:3343:HOH:O	2.27	0.46
1:A:220:ASN:ND2	1:A:223:GLY:O	2.40	0.46
1:D:184:LYS:NZ	7:D:3625:HOH:O	2.47	0.46
1:D:129:ALA:HA	1:D:143:ASN:HB3	1.98	0.45
1:C:316:ALA:HB1	1:C:317:PRO:HD2	1.99	0.45
1:A:371:SER:HB2	1:B:99:TRP:CE3	2.52	0.45
1:C:98:MET:HG3	1:C:99:TRP:CE3	2.52	0.45
1:C:325:GLU:HG2	7:C:3671:HOH:O	2.17	0.45
1:A:3:LEU:O	1:A:3:LEU:HD12	2.17	0.45
1:A:46:LEU:HD12	1:A:317:PRO:HG3	1.97	0.45
1:A:135:ASN:HD22	1:A:135:ASN:C	2.20	0.44
1:D:41:HIS:CE1	7:D:3341:HOH:O	2.66	0.44
1:A:330:GLU:HG3	7:A:3476:HOH:O	2.17	0.44
1:B:330:GLU:O	1:B:531:CYS:HA	2.18	0.44
1:C:43:SER:HB2	1:C:49:TRP:CD2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:THR:HA	1:A:425:TRP:CE3	2.52	0.44
1:D:84:LYS:HE3	1:D:86:TRP:CZ2	2.53	0.44
1:D:389:LYS:HE2	1:D:469:VAL:HG23	1.99	0.43
1:C:470:GLU:HG2	7:C:3502:HOH:O	2.17	0.43
1:C:362:GLY:C	1:C:363:ASN:HD22	2.20	0.43
1:A:429:LYS:NZ	1:A:449:SER:O	2.51	0.43
1:B:224:PRO:HD2	7:B:3520:HOH:O	2.19	0.43
1:A:77:ASP:HB3	1:A:88:ILE:HB	1.99	0.43
1:C:62:LEU:HD13	1:C:73:ILE:HD11	1.99	0.43
1:D:38:VAL:O	1:D:54:HIS:HA	2.19	0.43
1:A:389:LYS:HB3	1:A:389:LYS:HE2	1.68	0.42
1:C:99:TRP:CD2	1:D:371:SER:HB2	2.54	0.42
1:D:330:GLU:O	1:D:531:CYS:HA	2.20	0.42
1:D:338:LYS:HE2	7:D:3537:HOH:O	2.18	0.42
1:C:240:PRO:HG3	1:C:278:LEU:HD23	2.01	0.42
1:A:303:ARG:NH2	7:A:3605:HOH:O	2.52	0.42
1:A:99:TRP:CD2	1:B:371:SER:CB	3.02	0.42
1:A:178:TYR:CE2	1:A:180:GLY:HA2	2.55	0.42
1:C:109:CYS:HB2	1:C:116:TRP:CD2	2.55	0.41
1:A:135:ASN:HD22	1:A:136:ASP:N	2.18	0.41
1:D:392:ALA:HA	1:D:531:CYS:O	2.20	0.41
1:C:425:TRP:CH2	1:C:427:GLU:HA	2.56	0.41
1:B:429:LYS:HE3	7:B:3630:HOH:O	2.21	0.41
1:A:452:ILE:HA	1:A:453:PRO:HD3	1.94	0.41
1:D:186:THR:HA	1:D:204:GLU:HG2	2.02	0.41
1:B:38:VAL:O	1:B:54:HIS:HA	2.20	0.41
1:A:330:GLU:O	1:A:531:CYS:HA	2.21	0.40
1:B:333:ASN:ND2	7:B:3165:HOH:O	2.53	0.40
1:B:4:ILE:HG21	1:B:292:ILE:CG2	2.50	0.40
1:D:464:LYS:NZ	7:D:3598:HOH:O	2.51	0.40
1:B:184:LYS:HE3	7:B:3089:HOH:O	2.21	0.40
1:D:303:ARG:HD3	1:D:303:ARG:HA	1.89	0.40
1:B:189:PRO:HA	1:B:201:PHE:O	2.22	0.40
1:B:235:THR:O	1:B:309:GLN:HA	2.21	0.40
6:A:3022:GOL:H11	7:A:3316:HOH:O	2.21	0.40
1:C:216:ALA:HB1	1:C:225:TYR:HB3	2.04	0.40
1:D:331:LYS:HE2	1:D:333:ASN:HD21	1.87	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	532/542 (98%)	502 (94%)	28 (5%)	2 (0%)	39	17
1	B	532/542 (98%)	509 (96%)	21 (4%)	2 (0%)	39	17
1	C	532/542 (98%)	505 (95%)	25 (5%)	2 (0%)	39	17
1	D	532/542 (98%)	504 (95%)	25 (5%)	3 (1%)	30	9
All	All	2128/2168 (98%)	2020 (95%)	99 (5%)	9 (0%)	39	17

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	469	VAL
1	A	469	VAL
1	C	469	VAL
1	D	469	VAL
1	D	12	PHE
1	A	38	VAL
1	B	38	VAL
1	C	38	VAL
1	D	38	VAL

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	475/482 (98%)	470 (99%)	5 (1%)	80	63
1	B	475/482 (98%)	467 (98%)	8 (2%)	68	44

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	475/482 (98%)	470 (99%)	5 (1%)	80	63
1	D	475/482 (98%)	469 (99%)	6 (1%)	76	56
All	All	1900/1928 (98%)	1876 (99%)	24 (1%)	76	56

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

Mol	Chain	Res	Type
1	A	50	HIS
1	A	135	ASN
1	A	303	ARG
1	A	363	ASN
1	A	383	ARG
1	B	50	HIS
1	B	135	ASN
1	B	272	VAL
1	B	325	GLU
1	B	383	ARG
1	B	395	SER
1	B	431	ARG
1	B	535	GLU
1	C	19	CYS
1	C	135	ASN
1	C	383	ARG
1	C	431	ARG
1	C	535	GLU
1	D	50	HIS
1	D	135	ASN
1	D	184	LYS
1	D	296	GLU
1	D	325	GLU
1	D	383	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	333	ASN
1	B	135	ASN
1	B	333	ASN
1	C	135	ASN

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Mol	Chain	Res	Type
1	C	333	ASN
1	D	81	HIS
1	D	135	ASN
1	D	333	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](#)

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](#)

Of 29 ligands modelled in this entry, 4 are monoatomic - leaving 25 for Mogul analysis.

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	$\# Z > 2$
4	EPE	A	3006	-	14,15,15	0.52	0	18,20,20	1.74	4 (22%)
6	GOL	A	3011	-	5,5,5	0.42	0	5,5,5	1.35	0
6	GOL	A	3012	-	5,5,5	0.40	0	5,5,5	0.47	0
6	GOL	A	3017	-	5,5,5	0.46	0	5,5,5	0.93	0
6	GOL	A	3022	-	5,5,5	0.36	0	5,5,5	0.78	0
5	PGE	A	3023	-	9,9,9	0.47	0	8,8,8	0.37	0
3	SO4	A	3027	-	4,4,4	0.22	0	6,6,6	0.33	0
3	SO4	A	3029	-	4,4,4	0.17	0	6,6,6	0.23	0
4	EPE	B	3005	-	14,15,15	0.43	0	18,20,20	1.53	4 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GOL	B	3009	-	5,5,5	0.35	0	5,5,5	0.27	0
6	GOL	B	3010	-	5,5,5	0.35	0	5,5,5	1.37	1 (20%)
6	GOL	B	3018	-	5,5,5	0.33	0	5,5,5	0.48	0
6	GOL	B	3021	-	5,5,5	0.19	0	5,5,5	0.56	0
3	SO4	B	3025	-	4,4,4	0.12	0	6,6,6	0.68	0
3	SO4	B	3026	-	4,4,4	0.15	0	6,6,6	0.38	0
4	EPE	C	3008	-	14,15,15	0.52	0	18,20,20	1.55	3 (16%)
6	GOL	C	3013	-	5,5,5	0.35	0	5,5,5	0.48	0
6	GOL	C	3014	-	5,5,5	0.49	0	5,5,5	1.14	0
6	GOL	C	3020	-	5,5,5	0.25	0	5,5,5	0.43	0
4	EPE	D	3007	-	14,15,15	0.45	0	18,20,20	1.60	4 (22%)
6	GOL	D	3015	-	5,5,5	0.53	0	5,5,5	1.09	0
6	GOL	D	3016	-	5,5,5	0.30	0	5,5,5	0.41	0
6	GOL	D	3019	-	5,5,5	0.28	0	5,5,5	0.24	0
3	SO4	D	3024	-	4,4,4	0.40	0	6,6,6	0.56	0
3	SO4	D	3028	-	4,4,4	0.18	0	6,6,6	0.28	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
4	EPE	A	3006	-	-	0/9/19/19	0/1/1/1
6	GOL	A	3011	-	-	0/4/4/4	0/0/0/0
6	GOL	A	3012	-	-	0/4/4/4	0/0/0/0
6	GOL	A	3017	-	-	0/4/4/4	0/0/0/0
6	GOL	A	3022	-	-	0/4/4/4	0/0/0/0
5	PGE	A	3023	-	-	0/7/7/7	0/0/0/0
3	SO4	A	3027	-	-	0/0/0/0	0/0/0/0
3	SO4	A	3029	-	-	0/0/0/0	0/0/0/0
4	EPE	B	3005	-	-	0/9/19/19	0/1/1/1
6	GOL	B	3009	-	-	0/4/4/4	0/0/0/0
6	GOL	B	3010	-	-	0/4/4/4	0/0/0/0
6	GOL	B	3018	-	-	0/4/4/4	0/0/0/0
6	GOL	B	3021	-	-	0/4/4/4	0/0/0/0
3	SO4	B	3025	-	-	0/0/0/0	0/0/0/0
3	SO4	B	3026	-	-	0/0/0/0	0/0/0/0
4	EPE	C	3008	-	-	0/9/19/19	0/1/1/1
6	GOL	C	3013	-	-	0/4/4/4	0/0/0/0
6	GOL	C	3014	-	-	0/4/4/4	0/0/0/0
6	GOL	C	3020	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EPE	D	3007	-	-	0/9/19/19	0/1/1/1
6	GOL	D	3015	-	-	0/4/4/4	0/0/0/0
6	GOL	D	3016	-	-	0/4/4/4	0/0/0/0
6	GOL	D	3019	-	-	0/4/4/4	0/0/0/0
3	SO4	D	3024	-	-	0/0/0/0	0/0/0/0
3	SO4	D	3028	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	3010	GOL	C3-C2-C1	-2.72	100.46	111.12
4	D	3007	EPE	O2S-S-C10	-2.70	104.60	106.91
4	C	3008	EPE	C9-N1-C2	-2.60	104.61	111.27
4	B	3005	EPE	C9-N1-C6	-2.11	105.84	111.27
4	A	3006	EPE	C7-N4-C5	2.07	116.56	111.27
4	B	3005	EPE	O1S-S-C10	2.10	108.70	106.91
4	A	3006	EPE	C7-N4-C3	2.13	116.74	111.27
4	D	3007	EPE	C6-N1-C2	2.15	113.56	108.90
4	B	3005	EPE	C2-C3-N4	2.54	115.17	110.63
4	B	3005	EPE	C5-N4-C3	2.97	115.34	108.90
4	C	3008	EPE	C2-C3-N4	3.04	116.06	110.63
4	A	3006	EPE	C2-C3-N4	3.12	116.21	110.63
4	D	3007	EPE	C6-C5-N4	3.18	116.32	110.63
4	D	3007	EPE	C5-N4-C3	3.31	116.07	108.90
4	C	3008	EPE	C5-N4-C3	3.59	116.67	108.90
4	A	3006	EPE	C5-N4-C3	4.52	118.70	108.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	3006	EPE	1	0
6	A	3017	GOL	1	0
6	A	3022	GOL	1	0
4	B	3005	EPE	1	0
6	C	3020	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, 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	534/542 (98%)	-0.39	1 (0%) 95 94	13, 18, 35, 70	0
1	B	534/542 (98%)	-0.31	2 (0%) 93 93	14, 20, 39, 75	0
1	C	534/542 (98%)	-0.33	1 (0%) 95 94	13, 19, 37, 72	0
1	D	534/542 (98%)	-0.36	3 (0%) 90 90	14, 19, 38, 77	0
All	All	2136/2168 (98%)	-0.35	7 (0%) 94 94	13, 19, 37, 77	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	535	GLU	2.7
1	A	481	ILE	2.6
1	C	535	GLU	2.4
1	D	481	ILE	2.3
1	D	2	SER	2.2
1	B	166	LYS	2.1
1	B	2	SER	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 [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
6	GOL	B	3018	6/6	0.83	0.18	7.71	32,38,53,69	0
2	CA	A	3001	1/1	0.99	0.15	4.63	27,27,27,27	0
6	GOL	A	3017	6/6	0.87	0.16	4.56	29,47,48,107	0
2	CA	C	3003	1/1	0.99	0.20	4.50	27,27,27,27	0
2	CA	D	3004	1/1	0.99	0.15	3.09	27,27,27,27	0
6	GOL	C	3014	6/6	0.98	0.14	3.02	14,20,23,26	0
3	SO4	A	3027	5/5	0.96	0.18	2.91	32,36,43,54	0
6	GOL	D	3019	6/6	0.94	0.12	2.66	31,33,48,86	0
6	GOL	C	3020	6/6	0.78	0.13	2.60	37,48,57,63	0
6	GOL	D	3015	6/6	0.97	0.13	1.94	15,20,23,24	0
6	GOL	A	3011	6/6	0.97	0.14	1.92	14,21,25,35	0
6	GOL	A	3022	6/6	0.97	0.08	1.75	21,32,36,52	0
6	GOL	B	3010	6/6	0.97	0.14	1.45	18,18,23,29	0
4	EPE	D	3007	15/15	0.92	0.10	1.43	19,27,38,40	0
3	SO4	D	3028	5/5	0.94	0.16	1.12	29,38,61,62	0
6	GOL	B	3021	6/6	0.96	0.07	1.06	25,32,36,39	0
3	SO4	B	3026	5/5	0.97	0.16	1.04	28,34,51,56	0
4	EPE	B	3005	15/15	0.93	0.09	0.54	18,25,51,68	0
4	EPE	C	3008	15/15	0.91	0.10	0.39	20,23,49,52	0
4	EPE	A	3006	15/15	0.93	0.09	0.28	18,24,46,73	0
5	PGE	A	3023	10/10	0.93	0.09	-0.26	25,32,40,46	0
2	CA	B	3002	1/1	0.99	0.08	-0.42	26,26,26,26	0
6	GOL	A	3012	6/6	0.96	0.06	-0.83	14,17,18,19	0
6	GOL	D	3016	6/6	0.98	0.05	-1.12	15,16,17,19	0
6	GOL	B	3009	6/6	0.97	0.05	-1.29	13,17,21,22	0
6	GOL	C	3013	6/6	0.98	0.05	-1.45	14,17,19,22	0
3	SO4	A	3029	5/5	0.96	0.18	-	29,47,50,75	0
3	SO4	D	3024	5/5	0.93	0.17	-	32,32,65,85	0
3	SO4	B	3025	5/5	0.93	0.20	-	29,30,50,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.