



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 04:34 am BST

PDB ID : 3EU8
Title : Crystal structure of putative glucoamylase (YP_210071.1) from *Bacteroides fragilis* NCTC 9343 at 2.12 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2008-10-09
Resolution : 2.12 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

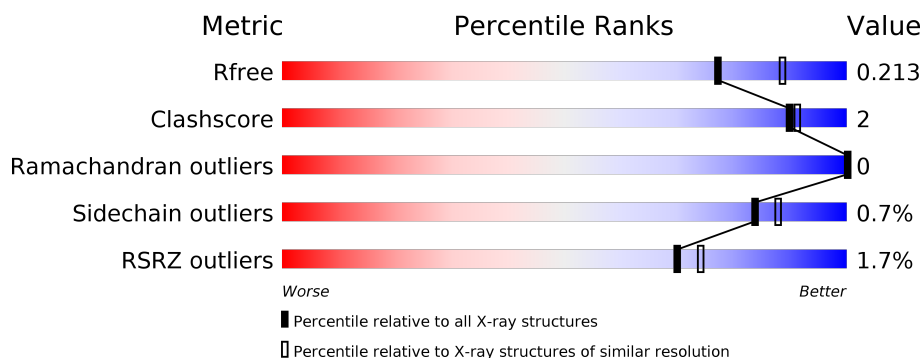
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.12 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	6241 (2.14-2.10)
Clashscore	141614	6778 (2.14-2.10)
Ramachandran outliers	138981	6705 (2.14-2.10)
Sidechain outliers	138945	6706 (2.14-2.10)
RSRZ outliers	127900	6112 (2.14-2.10)

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	430	<div> <div>%</div> <div> <div></div> <div>95%</div> <div>7%</div> <div>• •</div> </div> </div>
1	B	430	<div> <div>%</div> <div> <div></div> <div>92%</div> <div>7%</div> <div>•</div> </div> </div>
1	C	430	<div> <div>3%</div> <div> <div></div> <div>91%</div> <div>7%</div> <div>•</div> </div> </div>
1	D	430	<div> <div>%</div> <div> <div></div> <div>94%</div> <div>• •</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15148 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 putative glucoamylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	422	Total	C	N	O	S	Se	0	0	0
			3362	2160	568	618	4	12			
1	B	422	Total	C	N	O	S	Se	0	4	0
			3393	2177	578	622	4	12			
1	C	422	Total	C	N	O	S	Se	0	5	0
			3401	2184	581	620	4	12			
1	D	422	Total	C	N	O	S	Se	0	7	0
			3402	2186	576	624	4	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	GLY	-	leader sequence	UNP Q5LIB7
B	22	GLY	-	leader sequence	UNP Q5LIB7
C	22	GLY	-	leader sequence	UNP Q5LIB7
D	22	GLY	-	leader sequence	UNP Q5LIB7

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

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

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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

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

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

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

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0

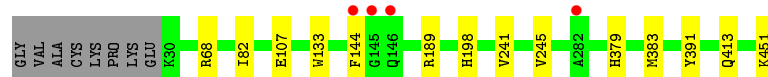
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	599	Total O 600 600	0	1
5	B	249	Total O 249 249	0	0
5	C	180	Total O 181 181	0	1
5	D	218	Total O 218 218	0	1

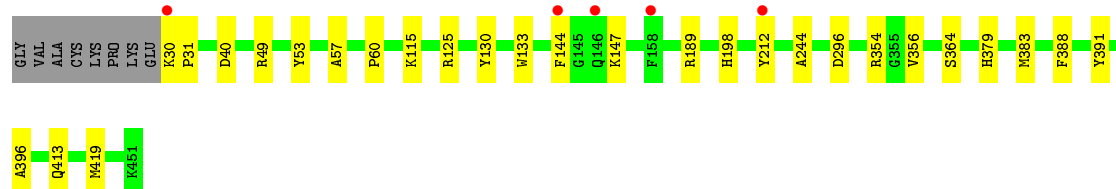
3 Residue-property plots [i](#)

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.

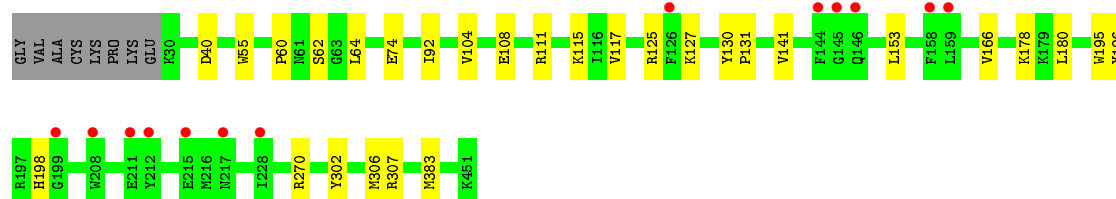
- Molecule 1: putative glucoamylase



- Molecule 1: putative glucoamylase



- Molecule 1: putative glucoamylase



- Molecule 1: putative glucoamylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.18 Å 71.76 Å 221.79 Å 90.00° 92.30° 90.00°	Depositor
Resolution (Å)	29.59 – 2.12 29.60 – 2.12	Depositor EDS
% Data completeness (in resolution range)	98.1 (29.59-2.12) 98.2 (29.60-2.12)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.12 Å)	Xtriage
Refinement program	REFMAC 5.2.0019, PHENIX	Depositor
R, R_{free}	0.158 , 0.213 0.162 , 0.213	Depositor DCC
R_{free} test set	5086 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.349	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15148	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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: K, EDO, CL

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	0.77	1/3461 (0.0%)	0.74	3/4684 (0.1%)
1	B	0.75	0/3504	0.73	2/4741 (0.0%)
1	C	0.69	1/3515 (0.0%)	0.68	0/4753
1	D	0.74	0/3522	0.71	3/4766 (0.1%)
All	All	0.74	2/14002 (0.0%)	0.72	8/18944 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	383	MSE	SE-CE	-6.75	1.55	1.95
1	A	107	GLU	CG-CD	5.90	1.60	1.51

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	189	ARG	NE-CZ-NH1	7.01	123.80	120.30
1	B	189	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	D	68	ARG	NE-CZ-NH1	6.61	123.60	120.30
1	D	68	ARG	NE-CZ-NH2	-6.57	117.02	120.30
1	A	68	ARG	NE-CZ-NH1	6.07	123.34	120.30
1	A	189	ARG	NE-CZ-NH2	-5.73	117.44	120.30
1	B	296	ASP	CB-CG-OD1	5.33	123.10	118.30
1	D	385	ASP	CB-CA-C	-5.32	99.75	110.40

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	3362	0	3146	7	0
1	B	3393	0	3185	15	0
1	C	3401	0	3212	16	0
1	D	3402	0	3201	9	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
3	A	124	0	184	1	0
3	B	84	0	125	2	0
3	C	72	0	107	3	0
3	D	56	0	84	0	0
4	B	1	0	0	0	0
5	A	600	0	0	2	0
5	B	249	0	0	1	0
5	C	181	0	0	2	0
5	D	218	0	0	0	0
All	All	15148	0	13244	48	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 (48) 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:133:TRP:HZ3	1:A:144:PHE:HB2	1.30	0.95
1:A:133:TRP:CZ3	1:A:144:PHE:HB2	2.09	0.88
1:D:133:TRP:HZ3	1:D:144:PHE:HB2	1.45	0.82
1:B:133:TRP:HZ3	1:B:144:PHE:HB2	1.43	0.81
1:D:133:TRP:CZ3	1:D:144:PHE:HB2	2.24	0.73
1:B:133:TRP:CZ3	1:B:144:PHE:HB2	2.25	0.71
1:C:60:PRO:O	1:C:115:LYS:NZ	2.19	0.59
1:B:125:ARG:HD3	1:B:130:TYR:CE2	2.40	0.56
1:A:379:HIS:CD2	1:A:383:MSE:HE3	2.42	0.55
1:B:356:VAL:CG1	1:B:396:ALA:HB1	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:53:TYR:O	1:B:57:ALA:HB3	2.10	0.51
1:A:391:TYR:O	1:A:413:GLN:NE2	2.44	0.50
1:B:40:ASP:HB3	5:B:707:HOH:O	2.12	0.50
1:B:60:PRO:O	1:B:115:LYS:NZ	2.36	0.49
1:D:379:HIS:CD2	1:D:383:MSE:HE3	2.46	0.49
1:D:220:VAL:HG21	1:D:229:MSE:HE1	1.94	0.49
1:B:49:ARG:HD2	1:B:388:PHE:CZ	2.47	0.49
1:B:49:ARG:HH21	3:B:9:EDO:H12	1.78	0.48
1:B:147:LYS:HA	1:B:212:TYR:CD2	2.50	0.47
1:B:391:TYR:O	1:B:413:GLN:NE2	2.47	0.47
1:A:82:ILE:HG21	3:A:472:EDO:C1	2.45	0.47
1:B:379:HIS:CD2	1:B:383:MSE:HE3	2.49	0.46
1:C:92:ILE:CG2	1:C:166:VAL:CG2	2.94	0.46
1:D:387[A]:VAL:HG12	1:D:393:PHE:HD1	1.81	0.46
1:C:125:ARG:HD3	1:C:130:TYR:CZ	2.52	0.45
1:C:117:VAL:HG21	1:C:180:LEU:HD21	1.98	0.45
1:D:92:ILE:HG21	1:D:92:ILE:HD12	1.74	0.45
1:D:395:ASP:OD1	1:D:412:ASP:OD2	2.34	0.45
1:C:131:PRO:CB	1:C:141:VAL:HG23	2.46	0.44
1:C:108:GLU:HG3	5:C:471:HOH:O	2.18	0.44
1:C:125:ARG:HD3	1:C:130:TYR:CE2	2.53	0.44
1:B:364:SER:O	1:B:419:MSE:HG3	2.17	0.43
1:C:62:SER:HB2	1:C:64:LEU:HD13	2.00	0.43
1:D:246:TYR:CE1	1:D:250:TRP:CD1	3.07	0.43
1:C:153:LEU:HD22	1:C:196:TYR:CB	2.49	0.43
1:D:125:ARG:HD3	1:D:130:TYR:CE2	2.55	0.42
1:B:244:ALA:HB3	3:B:454:EDO:H11	2.02	0.42
1:C:178:LYS:O	3:C:457:EDO:H22	2.20	0.42
1:C:127:LYS:HG3	1:C:195:TRP:CD1	2.55	0.42
1:C:111:ARG:HD3	5:C:612:HOH:O	2.18	0.42
3:C:16:EDO:O2	3:C:459:EDO:H22	2.20	0.41
1:A:241:VAL:HB	1:A:245:VAL:HG21	2.02	0.41
1:B:30:LYS:HA	1:B:31:PRO:HD3	1.97	0.41
1:C:307[A]:ARG:NH1	3:C:452:EDO:O2	2.50	0.41
1:C:55:TRP:CE2	1:C:104:VAL:HG11	2.55	0.41
1:C:302:TYR:O	1:C:306:MSE:HG2	2.21	0.41
5:A:637:HOH:O	1:C:307[A]:ARG:HD3	2.21	0.40
1:A:451:LYS:NZ	5:A:954:HOH:O	2.49	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	420/430 (98%)	405 (96%)	15 (4%)	0	100	100
1	B	424/430 (99%)	406 (96%)	18 (4%)	0	100	100
1	C	425/430 (99%)	413 (97%)	12 (3%)	0	100	100
1	D	427/430 (99%)	412 (96%)	15 (4%)	0	100	100
All	All	1696/1720 (99%)	1636 (96%)	60 (4%)	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	337/336 (100%)	336 (100%)	1 (0%)	92	95
1	B	342/336 (102%)	339 (99%)	3 (1%)	78	83
1	C	344/336 (102%)	340 (99%)	4 (1%)	71	77
1	D	344/336 (102%)	340 (99%)	4 (1%)	71	77
All	All	1367/1344 (102%)	1355 (99%)	12 (1%)	84	83

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

Mol	Chain	Res	Type
1	A	198	HIS
1	B	198	HIS

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Mol	Chain	Res	Type
1	B	354[A]	ARG
1	B	354[B]	ARG
1	C	40	ASP
1	C	74	GLU
1	C	198	HIS
1	C	270	ARG
1	D	92	ILE
1	D	198	HIS
1	D	354[A]	ARG
1	D	354[B]	ARG

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

Mol	Chain	Res	Type
1	A	146	GLN

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 90 ligands modelled in this entry, 6 are monoatomic - leaving 84 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
3	EDO	C	17	-	3,3,3	0.39	0	2,2,2	0.54	0
3	EDO	C	460	-	3,3,3	0.27	0	2,2,2	0.63	0
3	EDO	B	8	2	3,3,3	0.30	0	2,2,2	0.67	0
3	EDO	B	456	-	3,3,3	0.56	0	2,2,2	0.19	0
3	EDO	C	456	-	3,3,3	0.37	0	2,2,2	0.48	0
3	EDO	D	452	-	3,3,3	0.34	0	2,2,2	0.41	0
3	EDO	C	457	-	3,3,3	0.46	0	2,2,2	0.23	0
3	EDO	C	452	-	3,3,3	0.39	0	2,2,2	0.68	0
3	EDO	B	9	-	3,3,3	0.42	0	2,2,2	0.37	0
3	EDO	B	455	-	3,3,3	0.35	0	2,2,2	0.50	0
3	EDO	C	453	-	3,3,3	0.43	0	2,2,2	0.34	0
3	EDO	A	469	-	3,3,3	0.49	0	2,2,2	0.10	0
3	EDO	A	477	-	3,3,3	0.51	0	2,2,2	0.09	0
3	EDO	A	470	-	3,3,3	0.52	0	2,2,2	0.13	0
3	EDO	A	453	-	3,3,3	0.40	0	2,2,2	0.44	0
3	EDO	A	459	2	3,3,3	0.26	0	2,2,2	0.96	0
3	EDO	D	458	-	3,3,3	0.38	0	2,2,2	0.46	0
3	EDO	C	19	-	3,3,3	0.24	0	2,2,2	0.97	0
3	EDO	D	453	-	3,3,3	0.44	0	2,2,2	0.38	0
3	EDO	A	21	-	3,3,3	0.46	0	2,2,2	0.25	0
3	EDO	B	457	-	3,3,3	0.60	0	2,2,2	0.33	0
3	EDO	C	7	-	3,3,3	0.49	0	2,2,2	0.74	0
3	EDO	B	466	-	3,3,3	0.50	0	2,2,2	0.17	0
3	EDO	C	464	-	3,3,3	0.37	0	2,2,2	0.37	0
3	EDO	A	463	-	3,3,3	0.47	0	2,2,2	0.06	0
3	EDO	B	467	-	3,3,3	0.52	0	2,2,2	0.02	0
3	EDO	A	452	-	3,3,3	0.66	0	2,2,2	0.16	0
3	EDO	C	15	-	3,3,3	0.60	0	2,2,2	0.10	0
3	EDO	C	458	-	3,3,3	0.37	0	2,2,2	0.23	0
3	EDO	A	462	-	3,3,3	0.29	0	2,2,2	0.67	0
3	EDO	A	476	-	3,3,3	0.45	0	2,2,2	0.44	0
3	EDO	D	462	-	3,3,3	0.49	0	2,2,2	0.25	0
3	EDO	A	11	-	3,3,3	0.25	0	2,2,2	0.41	0
3	EDO	C	461	-	3,3,3	0.40	0	2,2,2	0.77	0
3	EDO	D	461	-	3,3,3	0.45	0	2,2,2	0.04	0
3	EDO	A	461	-	3,3,3	0.35	0	2,2,2	0.61	0
3	EDO	B	461	-	3,3,3	0.32	0	2,2,2	0.16	0
3	EDO	A	455	-	3,3,3	0.32	0	2,2,2	0.56	0
3	EDO	B	465	-	3,3,3	0.25	0	2,2,2	0.75	0
3	EDO	D	455	-	3,3,3	0.36	0	2,2,2	0.59	0
3	EDO	A	464	-	3,3,3	0.50	0	2,2,2	0.10	0
3	EDO	C	455	-	3,3,3	0.51	0	2,2,2	0.24	0
3	EDO	D	457	-	3,3,3	0.42	0	2,2,2	0.53	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	B	463	-	3,3,3	0.57	0	2,2,2	0.18	0
3	EDO	A	472	-	3,3,3	0.59	0	2,2,2	0.15	0
3	EDO	A	473	-	3,3,3	0.45	0	2,2,2	0.32	0
3	EDO	C	463	-	3,3,3	0.51	0	2,2,2	0.30	0
3	EDO	C	462	-	3,3,3	0.44	0	2,2,2	0.32	0
3	EDO	D	456	-	3,3,3	0.39	0	2,2,2	0.03	0
3	EDO	B	460	-	3,3,3	0.57	0	2,2,2	0.26	0
3	EDO	D	463	-	3,3,3	0.37	0	2,2,2	0.46	0
3	EDO	D	465	-	3,3,3	0.61	0	2,2,2	0.04	0
3	EDO	A	458	-	3,3,3	0.39	0	2,2,2	0.35	0
3	EDO	A	466	-	3,3,3	0.42	0	2,2,2	0.39	0
3	EDO	B	14	-	3,3,3	0.35	0	2,2,2	0.69	0
3	EDO	D	459	-	3,3,3	0.41	0	2,2,2	0.48	0
3	EDO	A	475	-	3,3,3	0.36	0	2,2,2	0.48	0
3	EDO	B	10	-	3,3,3	0.41	0	2,2,2	0.44	0
3	EDO	A	12	-	3,3,3	0.46	0	2,2,2	0.40	0
3	EDO	A	471	-	3,3,3	0.65	0	2,2,2	0.37	0
3	EDO	A	13	-	3,3,3	0.35	0	2,2,2	0.33	0
3	EDO	C	16	-	3,3,3	0.73	0	2,2,2	0.04	0
3	EDO	B	453	-	3,3,3	0.40	0	2,2,2	0.65	0
3	EDO	B	458	-	3,3,3	0.34	0	2,2,2	0.87	0
3	EDO	A	456	-	3,3,3	0.30	0	2,2,2	0.52	0
3	EDO	B	452	-	3,3,3	0.46	0	2,2,2	0.09	0
3	EDO	B	462	-	3,3,3	0.61	0	2,2,2	0.17	0
3	EDO	A	467	-	3,3,3	0.41	0	2,2,2	0.85	0
3	EDO	B	18	-	3,3,3	0.41	0	2,2,2	0.41	0
3	EDO	A	474	-	3,3,3	0.46	0	2,2,2	0.52	0
3	EDO	D	454	-	3,3,3	0.51	0	2,2,2	0.47	0
3	EDO	D	464	-	3,3,3	0.63	0	2,2,2	0.02	0
3	EDO	A	457	-	3,3,3	0.38	0	2,2,2	0.40	0
3	EDO	B	464	-	3,3,3	0.51	0	2,2,2	0.17	0
3	EDO	A	468	-	3,3,3	0.48	0	2,2,2	0.30	0
3	EDO	D	460	-	3,3,3	0.45	0	2,2,2	0.50	0
3	EDO	C	454	-	3,3,3	0.43	0	2,2,2	0.64	0
3	EDO	A	20	-	3,3,3	0.38	0	2,2,2	0.69	0
3	EDO	A	454	-	3,3,3	0.37	0	2,2,2	0.58	0
3	EDO	B	454	-	3,3,3	0.44	0	2,2,2	0.34	0
3	EDO	B	459	-	3,3,3	0.59	0	2,2,2	0.23	0
3	EDO	A	460	-	3,3,3	0.45	0	2,2,2	0.30	0
3	EDO	A	465	-	3,3,3	0.48	0	2,2,2	0.21	0
3	EDO	C	459	2	3,3,3	0.26	0	2,2,2	0.76	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	C	17	-	-	0/1/1/1	-
3	EDO	C	460	-	-	1/1/1/1	-
3	EDO	B	8	2	-	1/1/1/1	-
3	EDO	B	456	-	-	1/1/1/1	-
3	EDO	C	456	-	-	0/1/1/1	-
3	EDO	D	452	-	-	0/1/1/1	-
3	EDO	C	457	-	-	1/1/1/1	-
3	EDO	C	452	-	-	0/1/1/1	-
3	EDO	B	9	-	-	1/1/1/1	-
3	EDO	B	455	-	-	1/1/1/1	-
3	EDO	C	453	-	-	0/1/1/1	-
3	EDO	A	469	-	-	0/1/1/1	-
3	EDO	A	477	-	-	1/1/1/1	-
3	EDO	A	470	-	-	0/1/1/1	-
3	EDO	A	453	-	-	0/1/1/1	-
3	EDO	A	459	2	-	0/1/1/1	-
3	EDO	D	458	-	-	0/1/1/1	-
3	EDO	C	19	-	-	0/1/1/1	-
3	EDO	D	453	-	-	0/1/1/1	-
3	EDO	A	21	-	-	1/1/1/1	-
3	EDO	B	457	-	-	1/1/1/1	-
3	EDO	C	7	-	-	0/1/1/1	-
3	EDO	B	466	-	-	0/1/1/1	-
3	EDO	C	464	-	-	1/1/1/1	-
3	EDO	A	463	-	-	0/1/1/1	-
3	EDO	B	467	-	-	0/1/1/1	-
3	EDO	A	452	-	-	0/1/1/1	-
3	EDO	C	15	-	-	0/1/1/1	-
3	EDO	C	458	-	-	1/1/1/1	-
3	EDO	A	462	-	-	1/1/1/1	-
3	EDO	A	476	-	-	1/1/1/1	-
3	EDO	D	462	-	-	0/1/1/1	-
3	EDO	A	11	-	-	0/1/1/1	-
3	EDO	C	461	-	-	0/1/1/1	-
3	EDO	D	461	-	-	1/1/1/1	-
3	EDO	A	461	-	-	0/1/1/1	-
3	EDO	B	461	-	-	1/1/1/1	-
3	EDO	A	455	-	-	1/1/1/1	-
3	EDO	B	465	-	-	0/1/1/1	-
3	EDO	D	455	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	464	-	-	0/1/1/1	-
3	EDO	C	455	-	-	1/1/1/1	-
3	EDO	D	457	-	-	0/1/1/1	-
3	EDO	B	463	-	-	0/1/1/1	-
3	EDO	A	472	-	-	0/1/1/1	-
3	EDO	A	473	-	-	1/1/1/1	-
3	EDO	C	463	-	-	1/1/1/1	-
3	EDO	C	462	-	-	0/1/1/1	-
3	EDO	D	456	-	-	0/1/1/1	-
3	EDO	B	460	-	-	1/1/1/1	-
3	EDO	D	463	-	-	0/1/1/1	-
3	EDO	D	465	-	-	0/1/1/1	-
3	EDO	A	458	-	-	0/1/1/1	-
3	EDO	A	466	-	-	0/1/1/1	-
3	EDO	B	14	-	-	0/1/1/1	-
3	EDO	D	459	-	-	1/1/1/1	-
3	EDO	A	475	-	-	1/1/1/1	-
3	EDO	B	10	-	-	1/1/1/1	-
3	EDO	A	12	-	-	0/1/1/1	-
3	EDO	A	471	-	-	1/1/1/1	-
3	EDO	A	13	-	-	0/1/1/1	-
3	EDO	C	16	-	-	1/1/1/1	-
3	EDO	B	453	-	-	1/1/1/1	-
3	EDO	B	458	-	-	0/1/1/1	-
3	EDO	A	456	-	-	1/1/1/1	-
3	EDO	B	452	-	-	0/1/1/1	-
3	EDO	B	462	-	-	1/1/1/1	-
3	EDO	A	467	-	-	1/1/1/1	-
3	EDO	B	18	-	-	0/1/1/1	-
3	EDO	A	474	-	-	0/1/1/1	-
3	EDO	D	454	-	-	0/1/1/1	-
3	EDO	D	464	-	-	1/1/1/1	-
3	EDO	A	457	-	-	1/1/1/1	-
3	EDO	B	464	-	-	0/1/1/1	-
3	EDO	A	468	-	-	1/1/1/1	-
3	EDO	D	460	-	-	1/1/1/1	-
3	EDO	C	454	-	-	0/1/1/1	-
3	EDO	A	20	-	-	0/1/1/1	-
3	EDO	A	454	-	-	1/1/1/1	-
3	EDO	B	454	-	-	1/1/1/1	-
3	EDO	B	459	-	-	0/1/1/1	-
3	EDO	A	460	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	465	-	-	1/1/1/1	-
3	EDO	C	459	2	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	456	EDO	O1-C1-C2-O2
3	A	454	EDO	O1-C1-C2-O2
3	B	461	EDO	O1-C1-C2-O2
3	A	473	EDO	O1-C1-C2-O2
3	C	460	EDO	O1-C1-C2-O2
3	B	9	EDO	O1-C1-C2-O2
3	A	471	EDO	O1-C1-C2-O2
3	A	462	EDO	O1-C1-C2-O2
3	C	16	EDO	O1-C1-C2-O2
3	D	461	EDO	O1-C1-C2-O2
3	B	453	EDO	O1-C1-C2-O2
3	D	464	EDO	O1-C1-C2-O2
3	B	460	EDO	O1-C1-C2-O2
3	B	462	EDO	O1-C1-C2-O2
3	A	457	EDO	O1-C1-C2-O2
3	A	468	EDO	O1-C1-C2-O2
3	C	457	EDO	O1-C1-C2-O2
3	B	457	EDO	O1-C1-C2-O2
3	D	459	EDO	O1-C1-C2-O2
3	C	455	EDO	O1-C1-C2-O2
3	B	456	EDO	O1-C1-C2-O2
3	A	475	EDO	O1-C1-C2-O2
3	C	463	EDO	O1-C1-C2-O2
3	B	10	EDO	O1-C1-C2-O2
3	B	454	EDO	O1-C1-C2-O2
3	A	477	EDO	O1-C1-C2-O2
3	B	8	EDO	O1-C1-C2-O2
3	C	464	EDO	O1-C1-C2-O2
3	C	458	EDO	O1-C1-C2-O2
3	A	21	EDO	O1-C1-C2-O2
3	A	476	EDO	O1-C1-C2-O2
3	A	455	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
3	B	455	EDO	O1-C1-C2-O2
3	A	460	EDO	O1-C1-C2-O2
3	A	467	EDO	O1-C1-C2-O2
3	A	465	EDO	O1-C1-C2-O2
3	C	459	EDO	O1-C1-C2-O2
3	D	460	EDO	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	457	EDO	1	0
3	C	452	EDO	1	0
3	B	9	EDO	1	0
3	A	472	EDO	1	0
3	C	16	EDO	1	0
3	B	454	EDO	1	0
3	C	459	EDO	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	410/430 (95%)	-0.36	4 (0%) 82 85	19, 24, 35, 58	0
1	B	410/430 (95%)	-0.26	5 (1%) 79 82	18, 24, 36, 58	0
1	C	410/430 (95%)	0.02	13 (3%) 47 54	19, 24, 36, 57	0
1	D	410/430 (95%)	-0.26	6 (1%) 73 77	18, 24, 36, 58	0
All	All	1640/1720 (95%)	-0.21	28 (1%) 70 74	18, 24, 36, 58	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	144	PHE	5.9
1	C	145	GLY	5.9
1	B	144	PHE	4.8
1	C	146	GLN	3.9
1	A	145	GLY	3.8
1	D	145	GLY	3.8
1	D	212	TYR	3.4
1	A	144	PHE	3.3
1	C	212	TYR	3.3
1	A	146	GLN	3.2
1	D	146	GLN	3.1
1	C	215	GLU	3.1
1	B	212	TYR	2.9
1	C	211	GLU	2.9
1	D	144	PHE	2.6
1	C	158	PHE	2.5
1	B	158	PHE	2.5
1	C	217	ASN	2.4
1	C	208	TRP	2.3
1	C	159	LEU	2.3
1	B	30	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	146	GLN	2.3
1	D	158	PHE	2.2
1	C	126	PHE	2.1
1	C	199	GLY	2.1
1	A	282	ALA	2.1
1	C	228	ILE	2.1
1	D	159	LEU	2.0

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(Å ²)	Q<0.9
3	EDO	A	454	4/4	0.69	0.34	56,56,60,61	0
3	EDO	A	466	4/4	0.70	0.30	70,73,74,78	0
3	EDO	B	459	4/4	0.76	0.27	42,48,48,49	0
3	EDO	D	454	4/4	0.78	0.31	38,43,44,55	0
3	EDO	C	454	4/4	0.79	0.28	61,64,65,66	0
3	EDO	B	464	4/4	0.79	0.22	50,53,57,59	0
3	EDO	B	460	4/4	0.81	0.22	47,51,51,53	0
3	EDO	B	9	4/4	0.83	0.22	43,43,43,48	0
3	EDO	A	463	4/4	0.83	0.29	48,51,52,54	0
3	EDO	D	462	4/4	0.83	0.23	50,53,54,55	0
3	EDO	D	464	4/4	0.83	0.20	39,41,48,51	0
3	EDO	A	473	4/4	0.84	0.13	47,49,51,56	0
3	EDO	A	462	4/4	0.84	0.26	44,50,52,55	0
3	EDO	A	467	4/4	0.84	0.18	40,40,46,52	0
3	EDO	A	464	4/4	0.85	0.20	56,57,61,62	0
3	EDO	A	470	4/4	0.85	0.17	36,43,48,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	EDO	B	457	4/4	0.85	0.15	44,51,55,57	0
3	EDO	D	455	4/4	0.85	0.24	44,45,46,49	0
3	EDO	A	471	4/4	0.85	0.21	45,46,49,51	0
3	EDO	C	462	4/4	0.86	0.27	45,49,50,54	0
3	EDO	D	453	4/4	0.86	0.15	42,47,48,49	0
3	EDO	C	464	4/4	0.86	0.26	57,57,59,63	0
3	EDO	C	461	4/4	0.87	0.16	46,50,52,52	0
3	EDO	A	461	4/4	0.88	0.26	46,52,57,58	0
3	EDO	D	457	4/4	0.88	0.38	54,55,57,61	0
3	EDO	C	463	4/4	0.88	0.14	67,70,70,71	0
3	EDO	D	458	4/4	0.88	0.21	45,48,50,53	0
3	EDO	A	452	4/4	0.88	0.13	29,31,38,50	0
3	EDO	B	454	4/4	0.88	0.36	51,52,53,55	0
3	EDO	A	465	4/4	0.88	0.32	61,62,63,64	0
3	EDO	B	10	4/4	0.89	0.16	52,53,57,58	0
3	EDO	B	456	4/4	0.89	0.33	47,57,62,63	0
3	EDO	D	463	4/4	0.89	0.18	50,51,60,65	0
4	CL	B	6	1/1	0.89	0.10	64,64,64,64	0
3	EDO	A	477	4/4	0.89	0.17	43,46,47,49	0
3	EDO	C	457	4/4	0.90	0.25	39,39,46,48	0
3	EDO	A	460	4/4	0.90	0.15	42,45,49,49	0
3	EDO	A	472	4/4	0.90	0.22	29,38,42,47	0
3	EDO	D	459	4/4	0.91	0.16	52,54,54,56	0
3	EDO	A	456	4/4	0.91	0.25	42,43,44,47	0
3	EDO	B	18	4/4	0.91	0.17	23,26,35,47	0
3	EDO	B	461	4/4	0.91	0.14	26,34,43,43	0
3	EDO	A	476	4/4	0.91	0.15	36,41,46,55	0
3	EDO	C	455	4/4	0.92	0.16	39,58,62,70	0
3	EDO	C	16	4/4	0.92	0.17	37,40,43,46	0
3	EDO	A	457	4/4	0.92	0.16	32,38,39,51	0
3	EDO	C	452	4/4	0.92	0.13	33,35,36,40	0
3	EDO	A	469	4/4	0.92	0.25	36,40,44,48	0
3	EDO	B	8	4/4	0.92	0.12	24,27,31,34	0
3	EDO	D	452	4/4	0.93	0.10	32,34,35,39	0
3	EDO	B	467	4/4	0.93	0.13	38,39,42,42	0
3	EDO	A	458	4/4	0.93	0.18	32,32,39,40	0
3	EDO	A	475	4/4	0.93	0.19	47,47,48,57	0
3	EDO	D	461	4/4	0.93	0.19	39,42,43,45	0
3	EDO	A	11	4/4	0.93	0.13	21,27,27,28	0
3	EDO	B	453	4/4	0.93	0.10	28,32,36,36	0
3	EDO	C	15	4/4	0.93	0.11	30,33,34,40	0
3	EDO	A	453	4/4	0.93	0.15	54,55,59,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	EDO	B	463	4/4	0.94	0.14	37,41,41,43	0
3	EDO	A	459	4/4	0.94	0.21	33,37,38,38	0
3	EDO	C	453	4/4	0.94	0.18	29,30,34,36	0
3	EDO	B	465	4/4	0.94	0.13	41,48,49,51	0
3	EDO	D	460	4/4	0.94	0.13	46,49,50,52	0
3	EDO	D	456	4/4	0.94	0.12	29,30,36,42	0
3	EDO	D	465	4/4	0.94	0.13	36,45,45,47	0
3	EDO	B	452	4/4	0.95	0.16	31,34,35,40	0
3	EDO	C	458	4/4	0.95	0.22	30,39,42,45	0
3	EDO	C	456	4/4	0.95	0.14	45,46,46,47	0
3	EDO	B	462	4/4	0.95	0.15	26,29,33,34	0
3	EDO	B	466	4/4	0.95	0.29	29,32,34,37	0
3	EDO	A	21	4/4	0.95	0.09	30,32,38,39	0
3	EDO	B	455	4/4	0.95	0.15	32,37,37,42	0
3	EDO	A	468	4/4	0.95	0.21	57,58,58,58	0
3	EDO	C	19	4/4	0.95	0.15	30,38,39,43	0
3	EDO	C	459	4/4	0.95	0.17	31,37,46,47	0
3	EDO	C	7	4/4	0.96	0.11	20,21,21,21	0
3	EDO	A	455	4/4	0.96	0.18	43,44,48,49	0
3	EDO	B	14	4/4	0.96	0.07	25,25,27,33	0
3	EDO	C	460	4/4	0.96	0.09	26,30,34,45	0
2	K	C	5	1/1	0.96	0.08	60,60,60,60	0
3	EDO	A	474	4/4	0.96	0.14	19,28,33,42	0
3	EDO	A	12	4/4	0.97	0.10	29,29,31,36	0
3	EDO	B	458	4/4	0.97	0.08	25,27,29,31	0
3	EDO	C	17	4/4	0.97	0.07	17,27,29,34	0
3	EDO	A	20	4/4	0.97	0.07	16,17,21,28	0
2	K	C	3	1/1	0.97	0.13	41,41,41,41	0
3	EDO	A	13	4/4	0.97	0.12	29,33,34,35	0
2	K	D	2	1/1	0.98	0.14	40,40,40,40	0
2	K	A	1	1/1	0.98	0.10	36,36,36,36	0
2	K	B	4	1/1	0.99	0.10	36,36,36,36	0

6.5 Other polymers ⓘ

There are no such residues in this entry.