



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 20, 2023 – 12:04 pm GMT

PDB ID : 8B31  
Title : Crystal structure of UDP-glucose pyrophosphorylase from *Thermocristipum agreste* DSM 44070  
Authors : Laustsen, J.U.; Kumpf, A.; Bento, I.  
Deposited on : 2022-09-15  
Resolution : 1.75 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

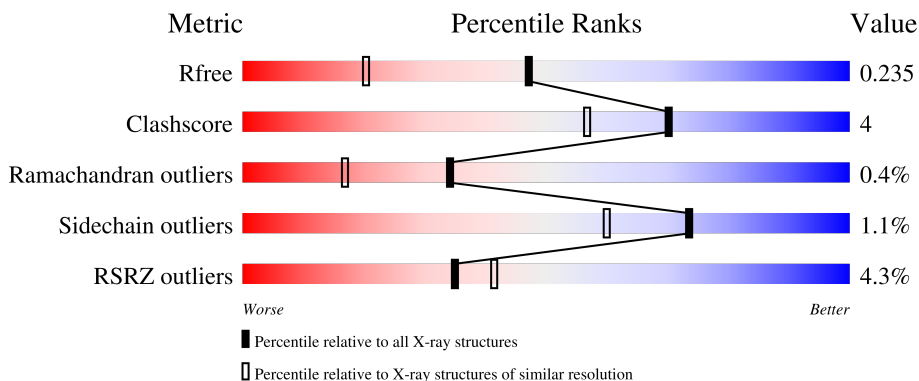
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	299	
1	B	299	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9578 atoms, of which 4625 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.

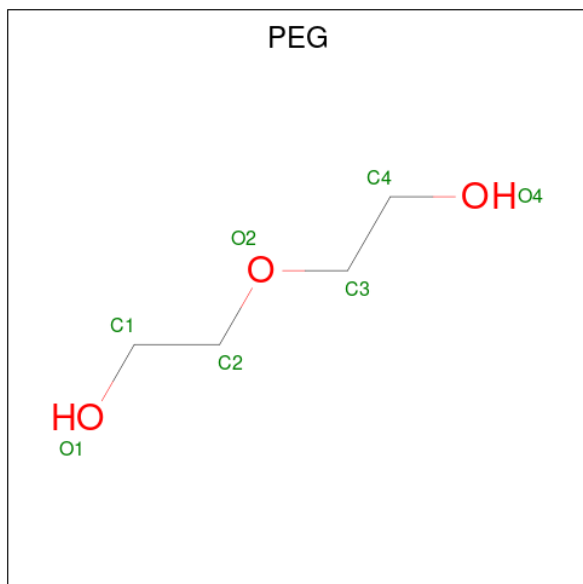
- Molecule 1 is a protein called UTP--glucose-1-phosphate uridylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	288	4463	1403	2246	389	419	6	106	5	0
1	B	289	4448	1397	2237	392	418	4	106	2	0

There are 4 discrepancies between the modelled and reference sequences:

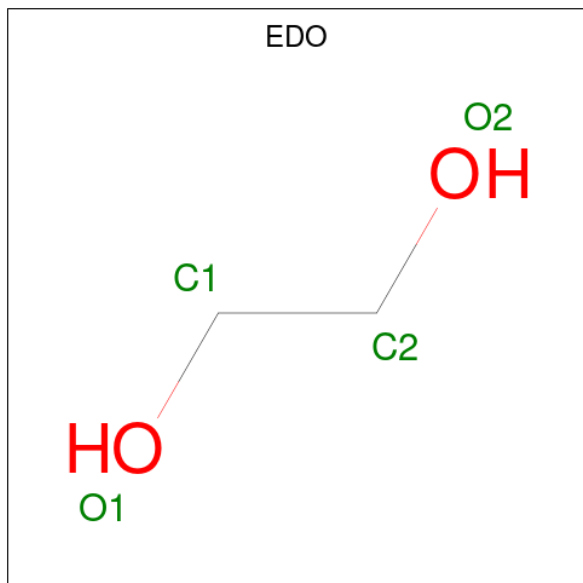
Chain	Residue	Modelled	Actual	Comment	Reference
A	260	ASP	VAL	conflict	UNP A0A2W4LV58
A	269	GLY	ASP	conflict	UNP A0A2W4LV58
B	260	ASP	VAL	conflict	UNP A0A2W4LV58
B	269	GLY	ASP	conflict	UNP A0A2W4LV58

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	A	1	17	4	10	3	1	0

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



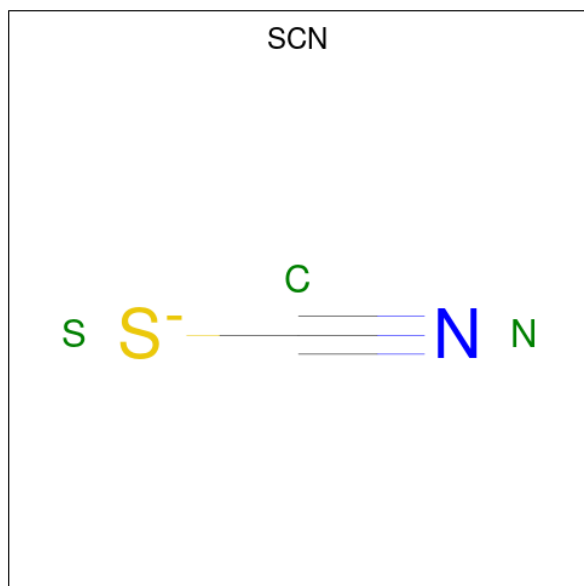
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
3	A	1	10	2	6	2	1	0
3	A	1	10	2	6	2	1	0
3	A	1	10	2	6	2	1	1
3	A	1	10	2	6	2	1	0
3	A	1	10	2	6	2	1	0
3	A	1	10	2	6	2	1	0
3	A	1	20	4	12	4	2	1
3	A	1	10	2	6	2	1	0
3	A	1	10	2	6	2	1	0
3	A	1	10	2	6	2	1	1
3	B	1	10	2	6	2	1	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	H	O	2	1
			20	4	12	4		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		
3	B	1	Total	C	H	O	1	0
			10	2	6	2		

- Molecule 4 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



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

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total K 2 2	0	0

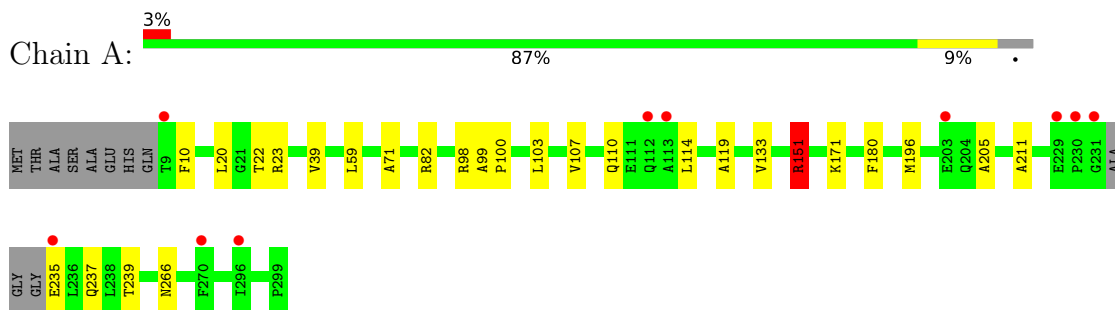
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	224	Total O 229 229	0	8
6	B	193	Total O 196 196	0	3

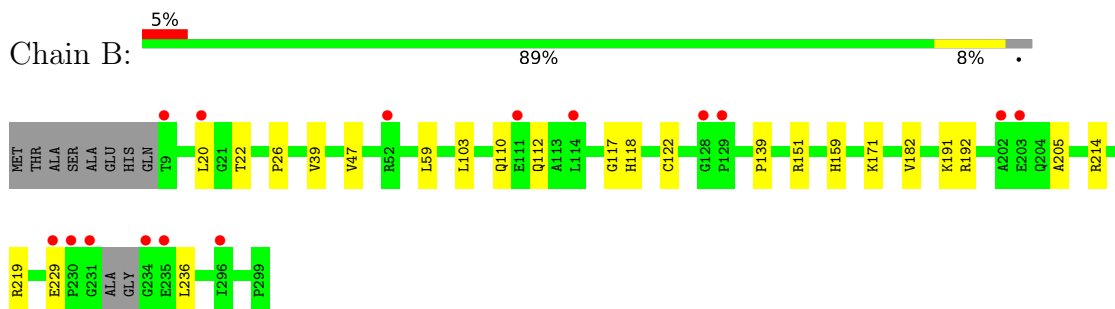
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: UTP--glucose-1-phosphate uridylyltransferase



- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	48.09Å 78.59Å 89.79Å 90.00° 93.89° 90.00°	Depositor
Resolution (Å)	48.02 – 1.75 47.98 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (48.02-1.75) 99.8 (47.98-1.75)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.73 (at 1.75Å)	Xtrriage
Refinement program	REFMAC 5.8.0352	Depositor
R, $R_{free}$	0.196 , 0.226 0.205 , 0.235	Depositor DCC
$R_{free}$ test set	3270 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.5	Xtrriage
Anisotropy	0.542	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 49.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9578	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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: CSX, EDO, SCN, PEG, K

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	0.52	0/2261	0.86	2/3076 (0.1%)
1	B	0.49	0/2250	0.80	2/3060 (0.1%)
All	All	0.51	0/4511	0.83	4/6136 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	151	ARG	NE-CZ-NH1	-9.99	115.31	120.30
1	A	151	ARG	NE-CZ-NH2	7.08	123.84	120.30
1	B	214	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	B	192	ARG	NE-CZ-NH2	-5.35	117.63	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	151	ARG	Sidechain
1	B	219	ARG	Sidechain

## 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	2217	2246	2241	19	0
1	B	2211	2237	2227	13	0
2	A	7	10	10	0	0
3	A	44	66	65	4	0
3	B	44	66	65	3	0
4	A	3	0	0	0	0
5	A	2	0	0	0	0
6	A	229	0	0	2	0
6	B	196	0	0	0	0
All	All	4953	4625	4608	33	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 4.

All (33) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:LEU:HB2	1:B:103:LEU:HD11	1.63	0.78
1:A:59:LEU:HB2	1:A:103:LEU:HD11	1.77	0.67
1:B:20:LEU:HD23	1:B:22:THR:HG22	1.81	0.62
1:A:151:ARG:HB3	3:A:304[A]:EDO:H22	1.82	0.61
1:A:71:ALA:HB2	1:A:107:VAL:HG21	1.84	0.58
1:A:23:ARG:HD2	3:A:303:EDO:H22	1.86	0.58
1:A:10:PHE:HB3	1:A:133[A]:VAL:HG23	1.84	0.57
1:A:180:PHE:CE1	1:A:196[B]:MET:HE3	2.41	0.56
1:B:110:GLN:HG3	1:B:122:CSX:HG	1.75	0.51
1:A:266:ASN:ND2	6:A:408:HOH:O	2.43	0.51
1:B:59:LEU:CB	1:B:103:LEU:HD11	2.38	0.50
1:B:151:ARG:HH11	3:B:303:EDO:H22	1.76	0.50
1:A:59:LEU:CB	1:A:103:LEU:HD11	2.41	0.50
1:A:20:LEU:HD23	1:A:22:THR:HG22	1.94	0.50
1:A:114:LEU:O	1:A:235:GLU:HB2	2.12	0.49
1:B:110:GLN:HE21	1:B:118:HIS:HD2	1.61	0.49
1:A:59:LEU:HB2	1:A:103:LEU:CD1	2.43	0.48
1:A:196[B]:MET:CE	1:A:211:ALA:HB2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ARG:HD3	3:A:310:EDO:H22	1.96	0.47
1:A:180:PHE:CE1	1:A:196[B]:MET:CE	2.98	0.46
1:B:171:LYS:HD3	1:B:205:ALA:HB3	1.99	0.44
6:A:495:HOH:O	3:B:306:EDO:H22	2.17	0.44
1:B:47:VAL:HG11	1:B:139:PRO:HA	2.00	0.43
1:A:239:THR:HG22	3:A:302:EDO:H21	2.01	0.42
1:A:99:ALA:HB3	1:A:100:PRO:HD3	2.00	0.42
1:B:117:GLY:HA3	1:B:236:LEU:O	2.19	0.42
1:A:196[B]:MET:HE1	1:A:211:ALA:HB2	2.00	0.42
1:A:171:LYS:HD3	1:A:205:ALA:HB3	2.02	0.42
1:B:110:GLN:HE21	1:B:118:HIS:CD2	2.39	0.41
1:B:159:HIS:HD2	3:B:307:EDO:C2	2.34	0.41
1:A:110:GLN:HE22	1:A:119:ALA:HA	1.85	0.41
1:B:182:VAL:HB	1:B:191:LYS:HB3	2.02	0.41
1:B:103:LEU:HD12	1:B:103:LEU:C	2.42	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	287/299 (96%)	283 (99%)	3 (1%)	1 (0%)	41	22
1	B	286/299 (96%)	281 (98%)	4 (1%)	1 (0%)	41	22
All	All	573/598 (96%)	564 (98%)	7 (1%)	2 (0%)	34	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	VAL
1	B	39	VAL

### 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	231/233 (99%)	229 (99%)	2 (1%)	78	67
1	B	229/233 (98%)	226 (99%)	3 (1%)	69	54
All	All	460/466 (99%)	455 (99%)	5 (1%)	73	60

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ARG
1	A	237	GLN
1	B	26	PRO
1	B	112	GLN
1	B	229	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	GLN
1	B	118	HIS
1	B	159	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues 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
1	CSX	A	122[B]	-	3,6,7	0.55	0	1,6,8	0.64	0
1	CSX	A	122[A]	-	3,6,7	0.54	0	1,6,8	1.31	0
1	CSX	B	122	1	3,6,7	0.82	0	1,6,8	1.79	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
1	CSX	A	122[B]	-	-	0/1/5/7	-
1	CSX	A	122[A]	-	-	0/1/5/7	-
1	CSX	B	122	1	-	0/1/5/7	-

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 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	122	CSX	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 26 ligands modelled in this entry, 2 are monoatomic - leaving 24 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	B	309	-	3,3,3	0.29	0	2,2,2	0.19	0
3	EDO	B	304	-	3,3,3	0.55	0	2,2,2	0.29	0
3	EDO	B	307	-	3,3,3	0.13	0	2,2,2	0.28	0
3	EDO	B	310	-	3,3,3	0.53	0	2,2,2	0.14	0
3	EDO	A	310	-	3,3,3	0.35	0	2,2,2	0.03	0
3	EDO	B	306	-	3,3,3	0.27	0	2,2,2	0.76	0
3	EDO	B	302[B]	-	3,3,3	0.13	0	2,2,2	0.69	0
3	EDO	A	308[B]	-	3,3,3	0.22	0	2,2,2	1.24	0
3	EDO	B	308	-	3,3,3	0.28	0	2,2,2	0.27	0
3	EDO	A	302	-	3,3,3	0.62	0	2,2,2	0.25	0
3	EDO	B	303	-	3,3,3	0.19	0	2,2,2	0.31	0
3	EDO	B	305	-	3,3,3	0.33	0	2,2,2	0.75	0
3	EDO	A	308[A]	5	3,3,3	0.53	0	2,2,2	0.52	0
3	EDO	B	302[A]	-	3,3,3	0.25	0	2,2,2	0.18	0
3	EDO	A	306	-	3,3,3	0.30	0	2,2,2	0.50	0
3	EDO	A	311[A]	-	3,3,3	0.22	0	2,2,2	0.18	0
3	EDO	A	304[A]	-	3,3,3	0.51	0	2,2,2	0.30	0
3	EDO	A	307	-	3,3,3	0.14	0	2,2,2	0.44	0
2	PEG	A	301	-	6,6,6	0.66	0	5,5,5	0.40	0
4	SCN	A	312	-	1,2,2	0.35	0	0,1,1	-	-
3	EDO	A	305	-	3,3,3	0.14	0	2,2,2	0.20	0
3	EDO	A	303	-	3,3,3	0.69	0	2,2,2	0.33	0
3	EDO	B	301	-	3,3,3	0.89	0	2,2,2	0.98	0
3	EDO	A	309	-	3,3,3	0.26	0	2,2,2	0.36	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	B	309	-	-	0/1/1/1	-
3	EDO	B	304	-	-	0/1/1/1	-
3	EDO	B	307	-	-	1/1/1/1	-
3	EDO	B	310	-	-	0/1/1/1	-
3	EDO	A	310	-	-	0/1/1/1	-
3	EDO	B	306	-	-	0/1/1/1	-
3	EDO	B	302[B]	-	-	1/1/1/1	-
3	EDO	A	308[B]	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	B	308	-	-	0/1/1/1	-
3	EDO	A	302	-	-	1/1/1/1	-
3	EDO	B	303	-	-	1/1/1/1	-
3	EDO	B	305	-	-	1/1/1/1	-
3	EDO	A	308[A]	5	-	1/1/1/1	-
3	EDO	B	302[A]	-	-	1/1/1/1	-
3	EDO	A	306	-	-	0/1/1/1	-
3	EDO	A	311[A]	-	-	1/1/1/1	-
3	EDO	A	304[A]	-	-	0/1/1/1	-
3	EDO	A	307	-	-	1/1/1/1	-
2	PEG	A	301	-	-	1/4/4/4	-
3	EDO	A	305	-	-	0/1/1/1	-
3	EDO	A	303	-	-	1/1/1/1	-
3	EDO	B	301	-	-	0/1/1/1	-
3	EDO	A	309	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	EDO	O1-C1-C2-O2
3	A	308[A]	EDO	O1-C1-C2-O2
3	B	307	EDO	O1-C1-C2-O2
3	B	303	EDO	O1-C1-C2-O2
3	A	302	EDO	O1-C1-C2-O2
2	A	301	PEG	C1-C2-O2-C3
3	A	307	EDO	O1-C1-C2-O2
3	A	308[B]	EDO	O1-C1-C2-O2
3	A	311[A]	EDO	O1-C1-C2-O2
3	B	305	EDO	O1-C1-C2-O2
3	B	302[A]	EDO	O1-C1-C2-O2
3	B	302[B]	EDO	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	307	EDO	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	310	EDO	1	0
3	B	306	EDO	1	0
3	A	302	EDO	1	0
3	B	303	EDO	1	0
3	A	304[A]	EDO	1	0
3	A	303	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, 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	287/299 (95%)	0.35	10 (3%) 44 50	19, 31, 56, 83	0
1	B	288/299 (96%)	0.53	15 (5%) 27 33	20, 35, 60, 88	0
All	All	575/598 (96%)	0.44	25 (4%) 35 41	19, 33, 59, 88	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	231	GLY	6.2
1	A	231	GLY	5.7
1	A	230	PRO	5.5
1	A	235	GLU	4.1
1	B	9	THR	3.8
1	A	112	GLN	3.8
1	B	229	GLU	3.2
1	B	230	PRO	3.0
1	B	52	ARG	3.0
1	A	9	THR	2.7
1	B	114	LEU	2.6
1	A	296	ILE	2.5
1	B	203	GLU	2.4
1	B	129	PRO	2.3
1	A	203	GLU	2.3
1	B	235	GLU	2.3
1	A	113	ALA	2.3
1	B	128	GLY	2.2
1	B	296	ILE	2.2
1	A	229	GLU	2.1
1	B	234	GLY	2.1
1	B	111	GLU	2.1
1	B	20	LEU	2.1
1	A	270	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	202	ALA	2.0

## 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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
1	CSX	A	122[A]	7/8	0.93	0.08	35,39,44,49	8
1	CSX	A	122[B]	7/8	0.93	0.08	35,39,44,49	8
1	CSX	B	122	7/8	0.94	0.09	42,47,57,57	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides 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, 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	B-factors(Å <sup>2</sup> )	Q<0.9
5	K	A	314	1/1	0.52	0.14	87,87,87,87	0
3	EDO	B	305	4/4	0.62	0.23	30,41,42,47	10
3	EDO	A	306	4/4	0.70	0.17	30,57,59,59	1
3	EDO	B	308	4/4	0.71	0.17	30,49,58,58	1
3	EDO	B	306	4/4	0.74	0.37	30,40,42,43	10
3	EDO	B	303	4/4	0.74	0.22	30,49,53,55	10
2	PEG	A	301	7/7	0.74	0.17	30,32,45,46	17
3	EDO	A	309	4/4	0.77	0.23	30,61,63,64	1
3	EDO	A	302	4/4	0.79	0.18	30,49,53,53	1
3	EDO	A	307	4/4	0.79	0.20	30,36,39,41	10
3	EDO	B	310	4/4	0.80	0.16	30,50,52,52	1
3	EDO	A	308[B]	4/4	0.81	0.24	30,35,36,36	10
3	EDO	A	308[A]	4/4	0.81	0.24	21,22,22,30	10
3	EDO	B	307	4/4	0.82	0.18	30,62,64,72	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	EDO	B	309	4/4	0.82	0.16	30,53,58,60	1
3	EDO	A	311[A]	4/4	0.83	0.21	27,29,32,33	10
3	EDO	B	301	4/4	0.85	0.21	30,36,39,39	10
3	EDO	A	310	4/4	0.86	0.27	30,62,66,69	1
3	EDO	B	304	4/4	0.89	0.22	30,40,43,50	10
4	SCN	A	312	3/3	0.90	0.12	46,46,47,55	0
3	EDO	A	305	4/4	0.90	0.17	30,56,57,58	1
3	EDO	B	302[A]	4/4	0.92	0.15	30,34,35,35	10
3	EDO	B	302[B]	4/4	0.92	0.15	30,44,47,47	10
3	EDO	A	304[A]	4/4	0.92	0.21	30,38,40,41	10
3	EDO	A	303	4/4	0.92	0.12	30,37,43,44	1
5	K	A	313	1/1	0.95	0.27	47,47,47,47	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.