



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:08 PM GMT

PDB ID : 3WA2
Title : High resolution crystal structure of copper amine oxidase from arthrobacter globiformis
Authors : Murakawa, T.; Hayashi, H.; Sunami, T.; Kurihara, K.; Tamada, T.; Kuroki, R.; Suzuki, M.; Tanizawa, K.; Okajima, T.
Deposited on : 2013-04-22
Resolution : 1.08 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

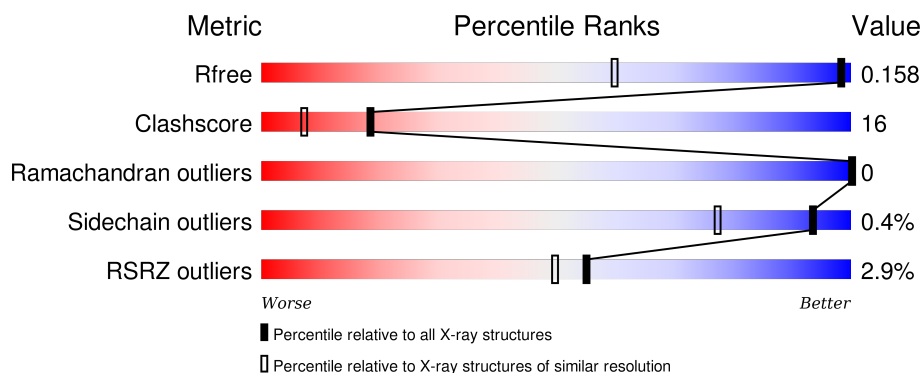
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.08 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	1120 (1.14-1.02)
Clashscore	102246	1181 (1.14-1.02)
Ramachandran outliers	100387	1134 (1.14-1.02)
Sidechain outliers	100360	1132 (1.14-1.02)
RSRZ outliers	91569	1124 (1.14-1.02)

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	X	621	<div> <div>3%</div> <div>87%</div> <div>13%</div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NA	X	702	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	EDO	X	703	-	-	X	X
4	EDO	X	705	-	-	-	X
4	EDO	X	707	-	-	-	X
4	EDO	X	708	-	-	X	X
4	EDO	X	709	-	-	-	X
4	EDO	X	711	-	-	X	X
4	EDO	X	712	-	-	-	X
4	EDO	X	714	-	-	-	X
5	PEG	X	715	-	-	-	X
5	PEG	X	716	-	-	-	X
5	PEG	X	717	-	-	X	X
5	PEG	X	718	-	-	-	X
5	PEG	X	719	-	-	X	X
5	PEG	X	721	-	-	X	X
5	PEG	X	722	-	-	X	X
5	PEG	X	723	-	-	X	X
6	PGE	X	724	-	-	X	X
6	PGE	X	725	-	-	-	X
6	PGE	X	727	-	-	-	X
6	PGE	X	728	-	-	X	X
7	PG4	X	729	-	-	-	X
8	1PE	X	730	-	-	X	X
8	1PE	X	731	-	-	-	X
9	OXY	X	732	-	-	-	X
9	OXY	X	733	-	-	-	X
9	OXY	X	734	-	-	-	X
9	OXY	X	735	-	-	-	X

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 12203 atoms, of which 5706 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 Phenylethylamine oxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	X	621	Total	C	H	N	O	S	0	91	0
			10768	3422	5412	933	988	13			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	X	1	Total	Cu	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

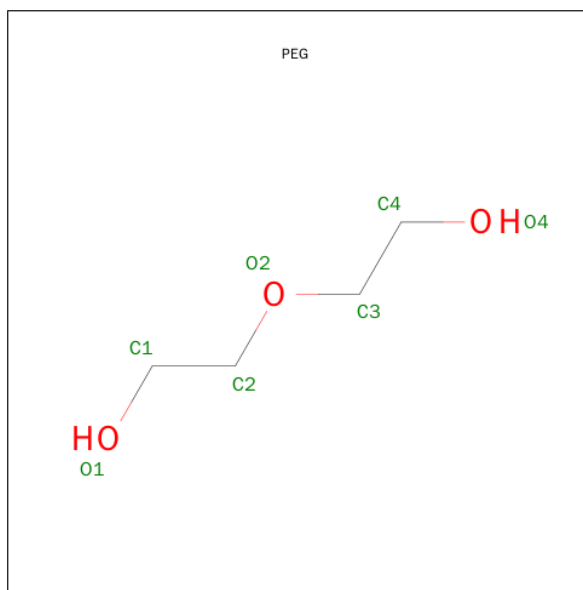
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	1	Total	Na	0	0
			1	1		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



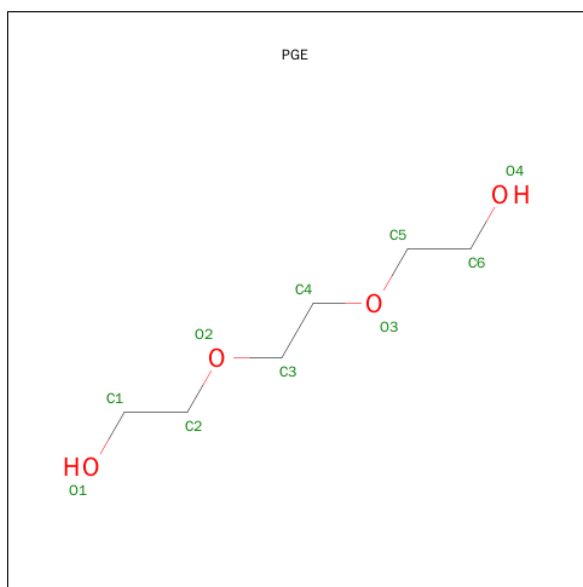
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0
4	X	1	Total C H O 10 2 6 2	0	0

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



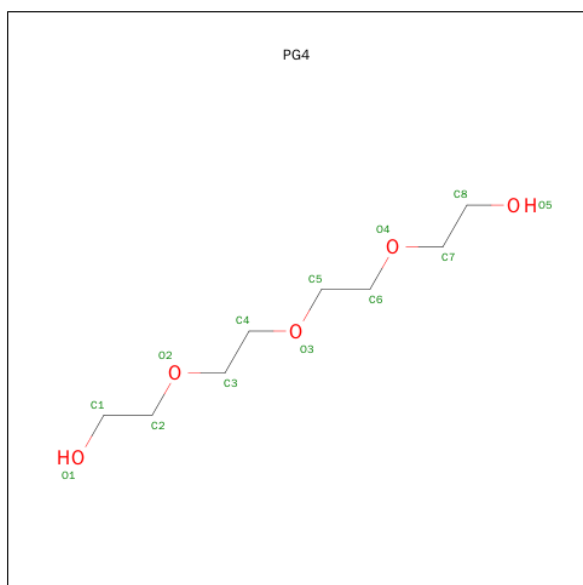
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	X	1	Total	C	H	O	0	0
			24	6	14	4		
6	X	1	Total	C	H	O	0	0
			24	6	14	4		
6	X	1	Total	C	H	O	0	0
			24	6	14	4		

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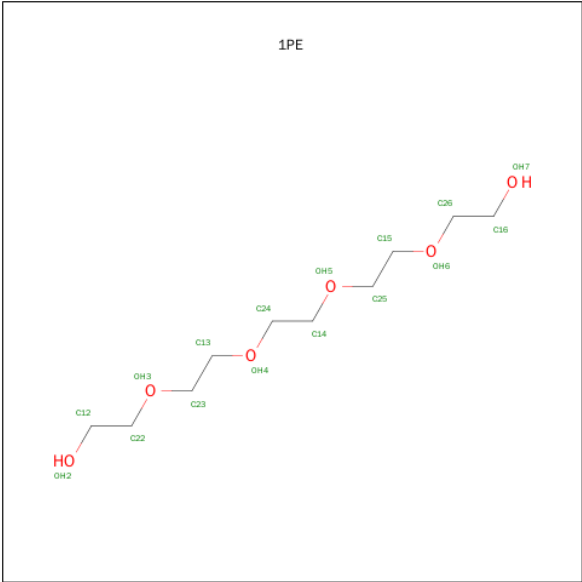
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	X	1	Total	C	H	O	0	0
			24	6	14	4		
6	X	1	Total	C	H	O	0	0
			24	6	14	4		

- Molecule 7 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



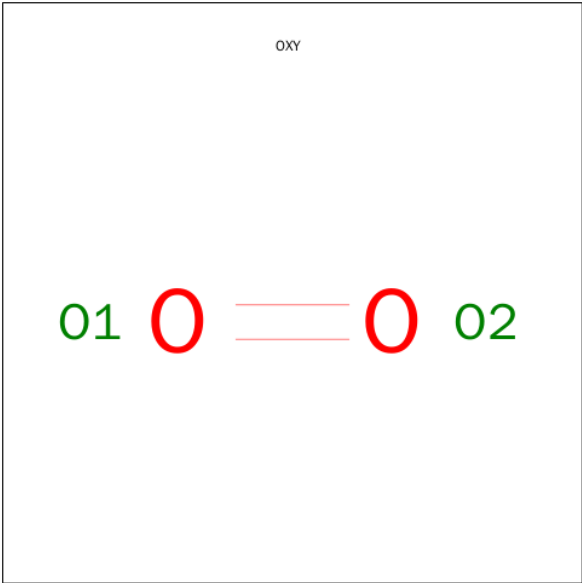
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	X	1	Total	C	H	O	0	0
			31	8	18	5		

- Molecule 8 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	X	1	Total	C	H	O	0	0
			38	10	22	6		
8	X	1	Total	C	H	O	0	0
			38	10	22	6		

- Molecule 9 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	X	1	Total	O	0	0
			2	2		
9	X	1	Total	O	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	X	1	Total	O	0	0
			2	2		
9	X	1	Total	O	0	0
			2	2		

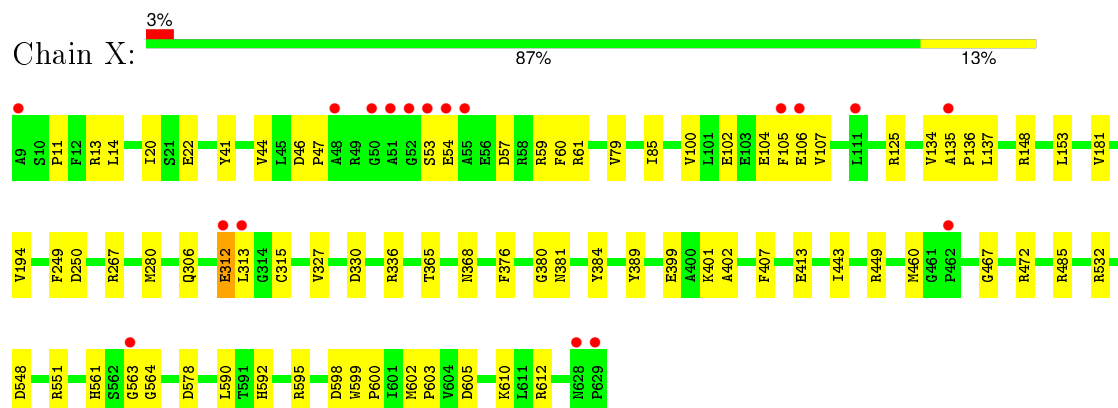
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	X	905	Total	O	0	20
			925	925		

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Phenylethylamine oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.75Å 62.38Å 92.08Å 90.00° 112.10° 90.00°	Depositor
Resolution (Å)	31.97 – 1.08 31.97 – 1.08	Depositor EDS
% Data completeness (in resolution range)	86.5 (31.97-1.08) 86.6 (31.97-1.08)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.03 (at 1.08Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.2_1309)	Depositor
R, R_{free}	0.130 , 0.150 0.137 , 0.158	Depositor DCC
R_{free} test set	15335 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	9.9	Xtriage
Anisotropy	0.079	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 56.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 305305 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	12203	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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: PGE, OXY, NA, EDO, 1PE, PG4, TPQ, PEG, CU

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	X	0.49	0/5741	0.72	1/7810 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	X	612	ARG	NE-CZ-NH1	5.25	122.92	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	5356	5412	5255	139	0
2	X	1	0	0	0	0
3	X	1	0	0	0	0
4	X	48	72	72	20	0
5	X	63	90	85	42	0
6	X	50	70	70	19	0
7	X	13	18	18	4	0
8	X	32	44	40	40	0
9	X	8	0	0	0	0
10	X	925	0	0	44	3

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6497	5706	5540	174	3

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 16.

All (174) 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:X:602[B]:MET:HG3	8:X:730:1PE:H221	1.40	1.04
1:X:595[B]:ARG:NE	8:X:730:1PE:H261	1.72	1.02
1:X:368:ASN:HD22	4:X:705:EDO:H12	1.28	0.96
1:X:602[B]:MET:CG	8:X:730:1PE:H221	1.95	0.94
5:X:719:PEG:O1	10:X:1591:HOH:O	1.88	0.91
1:X:134[A]:VAL:O	10:X:1466:HOH:O	1.88	0.90
1:X:44:VAL:HG12	4:X:711:EDO:H21	1.54	0.88
1:X:602[B]:MET:SD	8:X:730:1PE:C23	2.63	0.86
1:X:380:GLY:HA2	5:X:718:PEG:H41	1.58	0.84
1:X:368:ASN:ND2	4:X:705:EDO:H12	1.96	0.81
1:X:148[B]:ARG:NH1	10:X:1491:HOH:O	2.14	0.80
1:X:61[A]:ARG:NH1	4:X:703:EDO:O2	2.12	0.79
1:X:485[B]:ARG:NH1	10:X:1646:HOH:O	2.14	0.79
1:X:563[B]:GLY:O	10:X:947:HOH:O	2.01	0.79
1:X:602[B]:MET:SD	8:X:730:1PE:H232	2.23	0.78
1:X:602[B]:MET:SD	8:X:730:1PE:OH3	2.42	0.78
4:X:705:EDO:O1	10:X:929:HOH:O	2.02	0.76
5:X:722:PEG:O1	10:X:1238[A]:HOH:O	2.03	0.76
1:X:603[B]:PRO:HG2	8:X:730:1PE:OH2	1.86	0.75
1:X:135[B]:ALA:O	10:X:1478:HOH:O	2.02	0.75
1:X:399:GLU:OE1	5:X:721:PEG:C1	2.35	0.74
1:X:605[B]:ASP:OD2	8:X:730:1PE:OH5	2.07	0.72
1:X:402:ALA:HB3	8:X:730:1PE:H131	1.72	0.71
1:X:595[B]:ARG:HE	8:X:730:1PE:H261	1.56	0.71
1:X:592[A]:HIS:CD2	8:X:730:1PE:H161	2.26	0.70
7:X:729:PG4:H42	10:X:1424:HOH:O	1.91	0.70
1:X:14:LEU:O	5:X:719:PEG:H41	1.92	0.69
1:X:41:TYR:OH	4:X:703:EDO:H12	1.93	0.69
1:X:590[B]:LEU:HD13	8:X:730:1PE:H241	1.72	0.69
1:X:399:GLU:OE1	5:X:721:PEG:H12	1.94	0.68
1:X:315:CYS:HB2	10:X:1583:HOH:O	1.93	0.68
1:X:602[B]:MET:CG	8:X:730:1PE:C22	2.73	0.67
5:X:721:PEG:H22	10:X:1695:HOH:O	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:104:GLU:HG2	1:X:181[A]:VAL:HG11	1.78	0.67
1:X:57[A]:ASP:OD2	10:X:1559:HOH:O	2.13	0.66
1:X:106:GLU:OE1	10:X:1185:HOH:O	2.12	0.66
1:X:595[B]:ARG:HG3	8:X:730:1PE:C16	2.26	0.66
1:X:250:ASP:OD1	4:X:711:EDO:H11	1.96	0.66
1:X:22[B]:GLU:OE2	10:X:1024:HOH:O	2.14	0.65
1:X:100:VAL:HG12	6:X:724:PGE:H1	1.78	0.64
1:X:380:GLY:H	6:X:724:PGE:H62	1.64	0.63
5:X:722:PEG:H11	10:X:1312:HOH:O	1.99	0.62
1:X:551:ARG:HG2	5:X:722:PEG:H22	1.81	0.62
1:X:605[B]:ASP:HB2	8:X:730:1PE:H231	1.81	0.61
1:X:595[B]:ARG:HG3	8:X:730:1PE:H162	1.81	0.61
1:X:135[B]:ALA:HB1	6:X:728:PGE:H6	1.81	0.61
1:X:102:GLU:HA	6:X:724:PGE:H22	1.83	0.60
1:X:605[B]:ASP:OD2	8:X:730:1PE:C25	2.51	0.58
1:X:599:TRP:CD2	1:X:600:PRO:HA	2.39	0.58
1:X:249:PHE:O	4:X:711:EDO:H22	2.02	0.58
1:X:134[B]:VAL:O	10:X:1452[B]:HOH:O	2.17	0.58
1:X:105:PHE:CA	6:X:724:PGE:H2	2.34	0.57
1:X:44:VAL:HG12	4:X:711:EDO:C2	2.32	0.57
7:X:729:PG4:H21	10:X:1424:HOH:O	2.03	0.57
1:X:401:LYS:NZ	5:X:721:PEG:H42	2.19	0.57
1:X:460[B]:MET:HG3	1:X:467:GLY:N	2.20	0.56
1:X:595[B]:ARG:CD	8:X:730:1PE:H261	2.35	0.56
1:X:313[B]:LEU:H	1:X:313[B]:LEU:HD12	1.70	0.56
1:X:59:ARG:NH2	10:X:1559:HOH:O	2.38	0.56
1:X:137:LEU:HG	10:X:1478:HOH:O	2.04	0.56
1:X:595[A]:ARG:HG2	8:X:730:1PE:H162	1.88	0.56
1:X:551:ARG:HD3	5:X:722:PEG:O1	2.06	0.55
1:X:47:PRO:O	5:X:723:PEG:H12	2.06	0.55
1:X:551:ARG:CG	5:X:722:PEG:H22	2.37	0.55
5:X:722:PEG:C1	10:X:1238[A]:HOH:O	2.54	0.55
1:X:13:ARG:CD	5:X:719:PEG:H42	2.36	0.55
1:X:280[B]:MET:HE1	1:X:384:TYR:CE1	2.42	0.54
1:X:610[B]:LYS:NZ	10:X:1230:HOH:O	2.41	0.54
1:X:595[B]:ARG:HG3	8:X:730:1PE:OH7	2.07	0.54
6:X:728:PGE:H42	6:X:728:PGE:C1	2.37	0.54
8:X:730:1PE:H222	8:X:730:1PE:OH6	2.09	0.53
1:X:136:PRO:HG3	6:X:724:PGE:H12	1.91	0.53
5:X:722:PEG:H42	5:X:723:PEG:H22	1.91	0.52
4:X:703:EDO:H21	10:X:1238[B]:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:148[A]:ARG:NH2	10:X:1457:HOH:O	2.43	0.52
5:X:720:PEG:H32	10:X:1701:HOH:O	2.10	0.52
1:X:136:PRO:CG	6:X:724:PGE:H12	2.40	0.52
1:X:148[B]:ARG:CZ	10:X:1491:HOH:O	2.54	0.52
1:X:389:TYR:CE2	5:X:721:PEG:H12	2.45	0.52
1:X:401:LYS:HZ3	5:X:721:PEG:H42	1.75	0.51
5:X:722:PEG:C4	5:X:723:PEG:H22	2.40	0.51
1:X:460[B]:MET:HG3	1:X:467:GLY:CA	2.40	0.51
1:X:592[A]:HIS:HD2	8:X:730:1PE:H161	1.75	0.51
1:X:595[B]:ARG:HE	8:X:730:1PE:C26	2.22	0.51
1:X:605[B]:ASP:CB	8:X:730:1PE:H231	2.41	0.50
1:X:449[A]:ARG:HD2	1:X:578:ASP:OD1	2.12	0.50
5:X:717:PEG:H11	10:X:942:HOH:O	2.10	0.50
1:X:485[A]:ARG:CZ	10:X:1485[A]:HOH:O	2.59	0.50
1:X:46:ASP:HB2	5:X:723:PEG:H21	1.94	0.50
1:X:602[B]:MET:HG2	1:X:603[B]:PRO:O	2.12	0.49
1:X:603[A]:PRO:O	8:X:730:1PE:H122	2.11	0.49
1:X:181[A]:VAL:O	1:X:181[A]:VAL:HG12	2.12	0.49
1:X:60:PHE:CD1	1:X:79[B]:VAL:HG21	2.48	0.49
1:X:595[B]:ARG:NE	8:X:730:1PE:C26	2.60	0.49
1:X:532[A]:ARG:HB3	1:X:563[A]:GLY:HA3	1.94	0.49
1:X:61[A]:ARG:HE	4:X:703:EDO:H11	1.77	0.49
5:X:717:PEG:H12	10:X:887:HOH:O	2.13	0.49
1:X:61[A]:ARG:HE	4:X:703:EDO:C2	2.26	0.48
1:X:306:GLN:NE2	10:X:1392:HOH:O	2.30	0.48
1:X:148[B]:ARG:NE	10:X:1491:HOH:O	2.46	0.48
1:X:153:LEU:HD12	1:X:181[A]:VAL:HG21	1.96	0.48
1:X:125:ARG:HG2	1:X:194[A]:VAL:HG23	1.95	0.48
1:X:381:ASN:OD1	6:X:728:PGE:H22	2.14	0.47
1:X:551:ARG:NH2	10:X:1284:HOH:O	2.47	0.47
1:X:107[A]:VAL:HG11	1:X:181[A]:VAL:HG13	1.96	0.47
1:X:413:GLU:H	5:X:717:PEG:C2	2.28	0.47
1:X:136:PRO:HG3	6:X:724:PGE:C1	2.44	0.47
1:X:595[B]:ARG:NH2	8:X:730:1PE:OH2	2.48	0.46
1:X:610[B]:LYS:NZ	10:X:1444:HOH:O	2.43	0.46
1:X:381:ASN:OD1	6:X:728:PGE:C2	2.63	0.46
1:X:399:GLU:OE1	5:X:721:PEG:H11	2.13	0.46
7:X:729:PG4:H82	10:X:1465:HOH:O	2.15	0.46
1:X:59:ARG:CZ	1:X:85[B]:ILE:HD13	2.45	0.46
8:X:730:1PE:OH4	8:X:730:1PE:C25	2.63	0.46
1:X:380:GLY:HA3	6:X:728:PGE:H42	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:532[A]:ARG:HA	4:X:708:EDO:H12	1.98	0.46
1:X:13:ARG:NE	5:X:719:PEG:H42	2.31	0.46
1:X:598:ASP:OD2	8:X:730:1PE:H162	2.16	0.45
1:X:105:PHE:HA	6:X:724:PGE:H2	1.97	0.45
1:X:532[B]:ARG:HA	4:X:708:EDO:H12	1.99	0.45
1:X:602[B]:MET:SD	8:X:730:1PE:C22	3.04	0.45
1:X:561:HIS:HD1	1:X:563[B]:GLY:H	1.64	0.45
1:X:11:PRO:HB2	1:X:47:PRO:HG3	1.98	0.45
1:X:564[B]:GLY:HA2	4:X:708:EDO:H11	1.98	0.45
1:X:44:VAL:N	4:X:711:EDO:H12	2.31	0.45
1:X:61[B]:ARG:HB3	4:X:703:EDO:H11	1.99	0.45
1:X:134[A]:VAL:O	1:X:135[A]:ALA:CB	2.65	0.44
1:X:60:PHE:CD1	1:X:79[B]:VAL:CG2	3.01	0.44
7:X:729:PG4:H61	10:X:1424:HOH:O	2.16	0.44
1:X:564[A]:GLY:HA2	4:X:708:EDO:H11	2.00	0.44
1:X:20[B]:ILE:HD12	1:X:327:VAL:HG12	2.00	0.44
1:X:267[B]:ARG:NH2	10:X:1469:HOH:O	2.48	0.44
1:X:564[B]:GLY:HA2	4:X:708:EDO:C1	2.48	0.43
1:X:330:ASP:OD2	1:X:336[B]:ARG:NH2	2.50	0.43
1:X:312:GLU:O	10:X:1371:HOH:O	2.21	0.43
1:X:14:LEU:O	5:X:719:PEG:C4	2.65	0.43
5:X:717:PEG:HO4	6:X:726:PGE:H3	1.83	0.43
1:X:402:ALA:HB3	8:X:730:1PE:C13	2.46	0.43
8:X:730:1PE:OH3	8:X:730:1PE:H252	2.19	0.43
1:X:365[A]:THR:HG22	10:X:1399:HOH:O	2.18	0.43
1:X:472:ARG:NH1	5:X:715:PEG:H21	2.33	0.43
8:X:730:1PE:C22	8:X:730:1PE:C25	2.97	0.43
1:X:137:LEU:CD2	6:X:728:PGE:H1	2.48	0.43
8:X:731:1PE:H241	10:X:1408:HOH:O	2.18	0.42
1:X:592[B]:HIS:HD2	10:X:1192:HOH:O	2.02	0.42
1:X:563[B]:GLY:HA2	10:X:902:HOH:O	2.18	0.42
1:X:590[A]:LEU:HD21	10:X:1179:HOH:O	2.19	0.42
1:X:595[A]:ARG:HD3	8:X:730:1PE:H261	2.00	0.42
1:X:605[A]:ASP:OD2	8:X:730:1PE:H151	2.20	0.42
1:X:13:ARG:HG3	5:X:719:PEG:H42	2.02	0.42
1:X:53:SER:HA	1:X:54:GLU:HB2	2.02	0.42
1:X:605[B]:ASP:HB2	8:X:730:1PE:H252	2.01	0.41
8:X:730:1PE:OH6	8:X:730:1PE:H121	2.19	0.41
1:X:548[B]:ASP:HB2	5:X:723:PEG:O2	2.20	0.41
1:X:407:PHE:HE2	6:X:728:PGE:H12	1.85	0.41
5:X:721:PEG:C1	10:X:1695:HOH:O	2.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:105:PHE:N	6:X:724:PGE:H2	2.35	0.41
1:X:598:ASP:OD1	8:X:730:1PE:C22	2.69	0.41
1:X:380:GLY:H	6:X:724:PGE:C6	2.32	0.41
5:X:722:PEG:O4	5:X:723:PEG:H22	2.21	0.41
1:X:315:CYS:HB2	10:X:1585:HOH:O	2.19	0.41
1:X:148[B]:ARG:HB3	1:X:148[B]:ARG:HH21	1.85	0.40
1:X:602[B]:MET:SD	8:X:730:1PE:C13	3.08	0.40
5:X:718:PEG:O4	6:X:728:PGE:C6	2.69	0.40
1:X:564[B]:GLY:HA2	4:X:708:EDO:O1	2.21	0.40
5:X:721:PEG:O4	10:X:1534:HOH:O	2.22	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:X:1279:HOH:O	10:X:1295:HOH:O[2_556]	1.73	0.47
10:X:1177:HOH:O	10:X:1177:HOH:O[2_555]	1.91	0.29
10:X:1534:HOH:O	10:X:1547:HOH:O[2_556]	2.08	0.12

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	X	711/621 (114%)	689 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	X	601/514 (117%)	599 (100%)	2 (0%)	94 78

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	312	GLU
1	X	376	PHE

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	TPQ	X	382	1	13,14,15	2.49	4 (30%)	15,19,21	1.73	2 (13%)

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	TPQ	X	382	1	-	0/4/22/24	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	382	TPQ	C1-C2	-2.42	1.46	1.49
1	X	382	TPQ	CB-CA	2.94	1.59	1.53
1	X	382	TPQ	C6-C1	4.65	1.46	1.34
1	X	382	TPQ	C3-C4	6.35	1.46	1.35

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	382	TPQ	C1-C6-C5	-4.18	120.42	122.97
1	X	382	TPQ	C6-C1-C2	3.80	121.12	118.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 35 ligands modelled in this entry, 2 are monoatomic - leaving 33 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	EDO	X	703	-	3,3,3	0.42	0	2,2,2	1.09	0
4	EDO	X	704	-	3,3,3	0.47	0	2,2,2	0.45	0
4	EDO	X	705	-	3,3,3	0.85	0	2,2,2	0.72	0
4	EDO	X	706	-	3,3,3	0.38	0	2,2,2	0.64	0
4	EDO	X	707	-	3,3,3	0.48	0	2,2,2	0.36	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	X	708	-	3,3,3	0.39	0	2,2,2	0.19	0
4	EDO	X	709	-	3,3,3	0.48	0	2,2,2	0.42	0
4	EDO	X	710	-	3,3,3	0.48	0	2,2,2	0.23	0
4	EDO	X	711	-	3,3,3	0.56	0	2,2,2	0.81	0
4	EDO	X	712	-	3,3,3	0.51	0	2,2,2	0.19	0
4	EDO	X	713	-	3,3,3	0.46	0	2,2,2	0.31	0
4	EDO	X	714	-	3,3,3	0.46	0	2,2,2	0.39	0
5	PEG	X	715	-	6,6,6	0.71	0	5,5,5	1.21	1 (20%)
5	PEG	X	716	-	6,6,6	0.80	0	5,5,5	1.43	1 (20%)
5	PEG	X	717	-	6,6,6	0.64	0	5,5,5	0.76	0
5	PEG	X	718	-	6,6,6	0.82	0	5,5,5	1.27	0
5	PEG	X	719	-	6,6,6	0.87	0	5,5,5	0.71	0
5	PEG	X	720	-	6,6,6	0.63	0	5,5,5	0.82	0
5	PEG	X	721	-	6,6,6	1.06	0	5,5,5	1.47	1 (20%)
5	PEG	X	722	-	6,6,6	1.03	0	5,5,5	1.77	1 (20%)
5	PEG	X	723	-	6,6,6	0.65	0	5,5,5	0.93	0
6	PGE	X	724	-	9,9,9	0.86	0	8,8,8	0.90	0
6	PGE	X	725	-	9,9,9	0.71	0	8,8,8	0.80	0
6	PGE	X	726	-	9,9,9	0.60	0	8,8,8	0.75	0
6	PGE	X	727	-	9,9,9	0.71	0	8,8,8	0.76	0
6	PGE	X	728	-	9,9,9	0.64	0	8,8,8	1.37	1 (12%)
7	PG4	X	729	-	12,12,12	0.70	0	11,11,11	1.12	1 (9%)
8	1PE	X	730	-	15,15,15	1.17	1 (6%)	14,14,14	1.83	4 (28%)
8	1PE	X	731	-	15,15,15	0.63	0	14,14,14	1.27	3 (21%)
9	OXY	X	732	-	1,1,1	0.36	0	0,0,0	0.00	-
9	OXY	X	733	-	1,1,1	0.29	0	0,0,0	0.00	-
9	OXY	X	734	-	1,1,1	0.24	0	0,0,0	0.00	-
9	OXY	X	735	-	1,1,1	0.28	0	0,0,0	0.00	-

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	EDO	X	703	-	-	0/1/1/1	0/0/0/0
4	EDO	X	704	-	-	0/1/1/1	0/0/0/0
4	EDO	X	705	-	-	0/1/1/1	0/0/0/0
4	EDO	X	706	-	-	0/1/1/1	0/0/0/0
4	EDO	X	707	-	-	0/1/1/1	0/0/0/0
4	EDO	X	708	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	X	709	-	-	0/1/1/1	0/0/0/0
4	EDO	X	710	-	-	0/1/1/1	0/0/0/0
4	EDO	X	711	-	-	0/1/1/1	0/0/0/0
4	EDO	X	712	-	-	0/1/1/1	0/0/0/0
4	EDO	X	713	-	-	0/1/1/1	0/0/0/0
4	EDO	X	714	-	-	0/1/1/1	0/0/0/0
5	PEG	X	715	-	-	0/4/4/4	0/0/0/0
5	PEG	X	716	-	-	0/4/4/4	0/0/0/0
5	PEG	X	717	-	-	0/4/4/4	0/0/0/0
5	PEG	X	718	-	-	0/4/4/4	0/0/0/0
5	PEG	X	719	-	-	0/4/4/4	0/0/0/0
5	PEG	X	720	-	-	0/4/4/4	0/0/0/0
5	PEG	X	721	-	-	0/4/4/4	0/0/0/0
5	PEG	X	722	-	-	0/4/4/4	0/0/0/0
5	PEG	X	723	-	-	0/4/4/4	0/0/0/0
6	PGE	X	724	-	-	0/7/7/7	0/0/0/0
6	PGE	X	725	-	-	0/7/7/7	0/0/0/0
6	PGE	X	726	-	-	0/7/7/7	0/0/0/0
6	PGE	X	727	-	-	0/7/7/7	0/0/0/0
6	PGE	X	728	-	-	0/7/7/7	0/0/0/0
7	PG4	X	729	-	-	0/10/10/10	0/0/0/0
8	1PE	X	730	-	-	0/13/13/13	0/0/0/0
8	1PE	X	731	-	-	0/13/13/13	0/0/0/0
9	OXY	X	732	-	-	0/0/0/0	0/0/0/0
9	OXY	X	733	-	-	0/0/0/0	0/0/0/0
9	OXY	X	734	-	-	0/0/0/0	0/0/0/0
9	OXY	X	735	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	X	730	1PE	OH6-C26	-2.21	1.32	1.42

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	X	730	1PE	OH7-C16-C26	-3.33	91.50	112.03
5	X	722	PEG	O2-C2-C1	-3.06	96.36	110.43
8	X	730	1PE	C26-OH6-C15	-2.68	101.78	113.31
8	X	730	1PE	OH3-C23-C13	-2.51	99.20	110.36
8	X	731	1PE	C24-OH4-C13	2.06	122.16	113.31
8	X	731	1PE	C26-OH6-C15	2.52	124.16	113.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	X	731	1PE	C25-OH5-C14	2.56	124.31	113.31
5	X	715	PEG	C3-O2-C2	2.68	124.84	113.31
5	X	716	PEG	C3-O2-C2	2.70	124.92	113.31
6	X	728	PGE	O2-C2-C1	2.75	123.10	110.43
7	X	729	PG4	C7-O4-C6	2.96	126.03	113.31
5	X	721	PEG	O2-C2-C1	3.12	124.79	110.43
8	X	730	1PE	C23-OH3-C22	3.29	127.45	113.31

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

18 monomers are involved in 123 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	X	703	EDO	6	0
4	X	705	EDO	3	0
4	X	708	EDO	6	0
4	X	711	EDO	5	0
5	X	715	PEG	1	0
5	X	717	PEG	4	0
5	X	718	PEG	2	0
5	X	719	PEG	6	0
5	X	720	PEG	1	0
5	X	721	PEG	16	0
5	X	722	PEG	9	0
5	X	723	PEG	6	0
6	X	724	PGE	10	0
6	X	726	PGE	1	0
6	X	728	PGE	8	0
7	X	729	PG4	4	0
8	X	730	1PE	39	0
8	X	731	1PE	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	X	620/621 (99%)	-0.22	18 (2%) 55 50	6, 13, 33, 144	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	53	SER	12.1
1	X	54	GLU	8.0
1	X	52	GLY	8.0
1	X	105	PHE	6.1
1	X	55	ALA	4.3
1	X	51	ALA	4.2
1	X	628	ASN	4.2
1	X	313[A]	LEU	3.7
1	X	629	PRO	3.7
1	X	9	ALA	3.1
1	X	462	PRO	2.9
1	X	50	GLY	2.7
1	X	312	GLU	2.5
1	X	135[A]	ALA	2.4
1	X	48	ALA	2.4
1	X	563[A]	GLY	2.4
1	X	111[A]	LEU	2.2
1	X	106	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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(\AA^2)	Q<0.9
1	TPQ	X	382	14/15	0.95	0.08	-	8,12,14,14	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	PEG	X	721	7/7	0.77	0.40	28.33	8,23,31,31	17
6	PGE	X	725	10/10	0.84	0.25	27.76	20,29,36,71	24
5	PEG	X	722	7/7	0.70	0.55	26.13	10,18,39,39	17
5	PEG	X	719	7/7	0.65	0.59	21.17	16,22,33,39	17
8	1PE	X	730	16/16	0.83	0.28	20.00	11,26,34,37	38
4	EDO	X	703	4/4	0.80	0.27	15.00	19,23,28,28	10
4	EDO	X	705	4/4	0.88	0.24	14.14	10,15,22,26	10
5	PEG	X	716	7/7	0.69	0.35	13.24	20,26,35,42	17
5	PEG	X	723	7/7	0.67	0.44	12.27	10,19,32,32	17
4	EDO	X	711	4/4	0.93	0.13	8.30	14,21,43,43	10
9	OXY	X	734	2/2	0.99	0.28	7.92	21,21,21,21	0
5	PEG	X	717	7/7	0.81	0.28	7.90	15,23,49,59	17
7	PG4	X	729	13/13	0.90	0.13	6.59	9,19,35,62	31
6	PGE	X	724	10/10	0.65	0.43	6.43	31,80,109,123	24
4	EDO	X	707	4/4	0.80	0.20	5.72	57,69,88,105	0
8	1PE	X	731	16/16	0.87	0.17	5.70	37,77,100,111	0
4	EDO	X	712	4/4	0.74	0.14	5.38	21,26,29,30	10
3	NA	X	702	1/1	1.00	0.10	4.52	7,7,7,7	0
5	PEG	X	718	7/7	0.73	0.29	4.14	9,16,24,25	17
4	EDO	X	708	4/4	0.82	0.28	4.02	22,27,34,40	10
5	PEG	X	715	7/7	0.87	0.14	3.77	22,32,46,56	17
4	EDO	X	709	4/4	0.90	0.25	3.64	20,29,40,40	10
9	OXY	X	735	2/2	0.99	0.14	3.27	11,11,11,27	0
9	OXY	X	732	2/2	0.98	0.12	3.00	16,16,16,22	0
6	PGE	X	727	10/10	0.71	0.25	2.96	21,26,36,42	24

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	EDO	X	714	4/4	0.86	0.13	2.61	20,24,28,28	10
9	OXY	X	733	2/2	0.98	0.10	2.61	14,14,14,31	0
6	PGE	X	728	10/10	0.89	0.19	2.20	7,13,27,32	24
6	PGE	X	726	10/10	0.95	0.10	1.94	19,30,62,75	0
4	EDO	X	713	4/4	0.76	0.13	1.93	23,30,36,36	10
4	EDO	X	710	4/4	0.85	0.11	1.92	18,22,30,36	10
5	PEG	X	720	7/7	0.74	0.22	1.62	17,26,31,33	17
4	EDO	X	706	4/4	0.85	0.12	0.68	22,27,28,32	10
2	CU	X	701	1/1	1.00	0.03	-4.74	9,9,9,9	0
4	EDO	X	704	4/4	0.91	0.16	-	44,53,66,79	0

6.5 Other polymers

There are no such residues in this entry.