



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:14 PM UTC

PDB ID : 7PW1 / pdb_00007pw1
Title : Crystal structure of ancestral haloalkane dehalogenase AncLinB-DmbA
Authors : Mazur, A.; Grinkevich, P.; Prudnikova, T.
Deposited on : 2021-10-05
Resolution : 1.50 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 types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

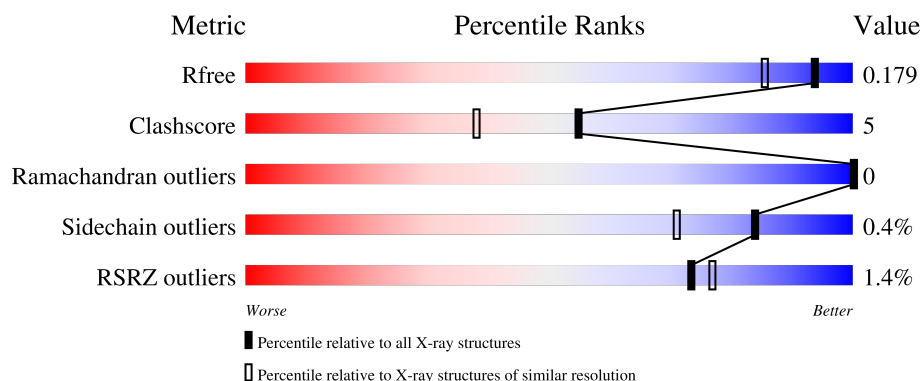
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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}	180053	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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 ≥ 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	296	% 69% 26% ..

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
4	EDO	A	406[B]	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2747 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 Haloalkane dehalogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	11	0
			2417	1531	421	455	10			

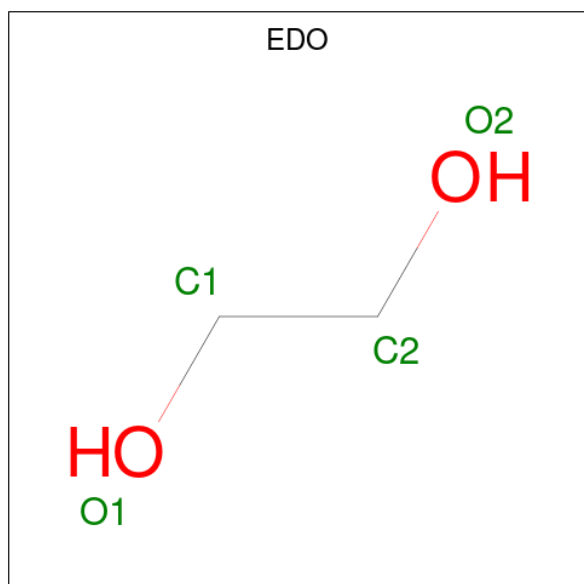
- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	Na	0	0
			4	4		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	1
			8	4	4		

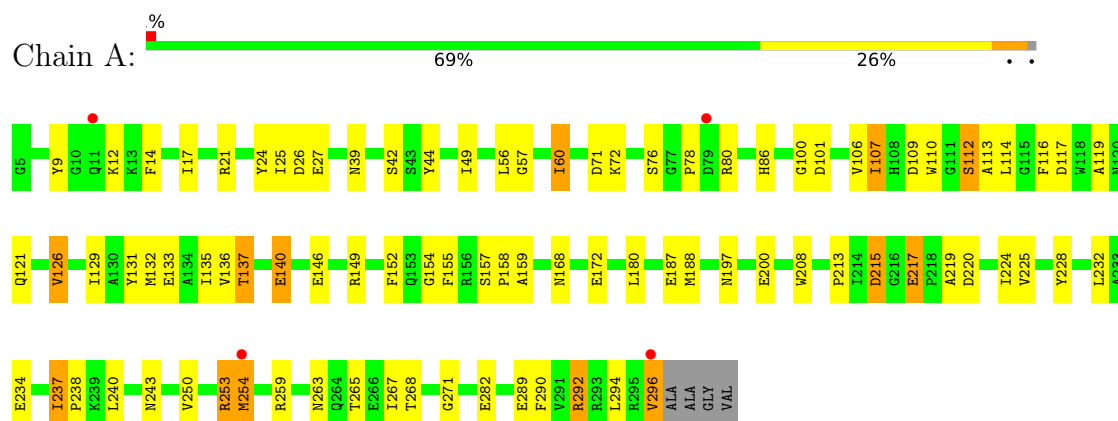
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	317	Total	O	0	0
			317	317		

3 Residue-property plots

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: Haloalkane dehalogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	68.90Å 68.90Å 156.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.04 – 1.50 63.04 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (63.04-1.50) 100.0 (63.04-1.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, R_{free}	0.138 , 0.174 0.146 , 0.179	Depositor DCC
R_{free} test set	3065 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2747	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: EDO, CL, NA

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	2.07	83/2489 (3.3%)	1.53	21/3380 (0.6%)

All (83) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	137	THR	N-CA	9.62	1.55	1.46
1	A	157	SER	CA-C	-8.28	1.44	1.52
1	A	119	ALA	N-CA	-8.10	1.36	1.46
1	A	187	GLU	CA-C	-7.90	1.42	1.52
1	A	265	THR	CA-CB	-7.70	1.41	1.53
1	A	135	ILE	CA-C	-7.46	1.45	1.53
1	A	234	GLU	CA-C	-7.33	1.42	1.52
1	A	107	ILE	C-O	7.29	1.31	1.24
1	A	187	GLU	N-CA	7.21	1.55	1.46
1	A	290	PHE	N-CA	7.08	1.54	1.46
1	A	49	ILE	CB-CG1	6.85	1.67	1.53
1	A	237	ILE	N-CA	-6.72	1.39	1.47
1	A	113	ALA	CA-CB	-6.69	1.43	1.53
1	A	113	ALA	N-CA	6.69	1.54	1.46
1	A	140	GLU	CA-C	-6.60	1.44	1.52
1	A	116	PHE	C-O	-6.51	1.16	1.24
1	A	159	ALA	CA-C	6.48	1.61	1.52
1	A	219	ALA	C-O	-6.40	1.16	1.24
1	A	168	ASN	N-CA	6.39	1.56	1.46
1	A	57	GLY	N-CA	6.33	1.51	1.45
1	A	228	TYR	C-O	-6.33	1.16	1.24
1	A	200	GLU	C-O	6.32	1.32	1.24
1	A	78	PRO	N-CA	-6.26	1.39	1.47
1	A	253[A]	ARG	CA-C	6.13	1.61	1.52
1	A	253[B]	ARG	CA-C	6.13	1.61	1.52
1	A	187	GLU	C-O	6.07	1.31	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	157	SER	N-CA	6.06	1.51	1.46
1	A	232	LEU	CA-C	6.05	1.60	1.52
1	A	263	ASN	CA-C	6.04	1.60	1.53
1	A	44	TYR	C-O	-6.03	1.16	1.24
1	A	259	ARG	C-O	-5.97	1.16	1.24
1	A	213	PRO	N-CA	5.90	1.54	1.47
1	A	271	GLY	CA-C	-5.88	1.44	1.51
1	A	76	SER	CA-C	-5.87	1.45	1.52
1	A	289	GLU	C-O	-5.81	1.17	1.24
1	A	24	TYR	C-O	-5.80	1.16	1.23
1	A	60	ILE	CA-CB	-5.77	1.47	1.54
1	A	243	ASN	C-O	5.77	1.31	1.23
1	A	126	VAL	CA-CB	-5.71	1.47	1.54
1	A	131	TYR	C-O	-5.68	1.16	1.23
1	A	220	ASP	N-CA	-5.66	1.39	1.46
1	A	158	PRO	CA-CB	5.63	1.62	1.53
1	A	197	ASN	CA-C	-5.62	1.45	1.52
1	A	86	HIS	CA-C	-5.61	1.45	1.52
1	A	225	VAL	C-O	-5.56	1.17	1.24
1	A	26	ASP	C-O	-5.52	1.17	1.23
1	A	106	VAL	C-O	5.52	1.30	1.24
1	A	254	MET	C-O	-5.52	1.17	1.24
1	A	136	VAL	C-N	-5.49	1.27	1.33
1	A	180	LEU	N-CA	5.43	1.53	1.46
1	A	240	LEU	CA-CB	-5.43	1.45	1.53
1	A	220	ASP	CA-C	5.42	1.59	1.52
1	A	72	LYS	CA-C	5.39	1.59	1.52
1	A	117	ASP	CA-CB	-5.37	1.45	1.53
1	A	224	ILE	CA-C	5.37	1.59	1.52
1	A	159	ALA	C-O	-5.35	1.17	1.24
1	A	154	GLY	CA-C	5.34	1.57	1.52
1	A	9	TYR	C-O	-5.33	1.17	1.24
1	A	224	ILE	CA-CB	-5.32	1.48	1.54
1	A	250	VAL	CA-CB	-5.31	1.47	1.54
1	A	107	ILE	CA-C	-5.30	1.46	1.52
1	A	100	GLY	CA-C	-5.30	1.45	1.52
1	A	80	ARG	C-O	-5.28	1.17	1.23
1	A	114	LEU	N-CA	-5.27	1.40	1.46
1	A	238	PRO	CA-CB	-5.25	1.46	1.53
1	A	12	LYS	CA-CB	-5.23	1.44	1.53
1	A	129	ILE	CA-CB	-5.21	1.47	1.54
1	A	117	ASP	C-O	5.21	1.30	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	155	PHE	C-O	-5.20	1.18	1.24
1	A	121	GLN	CA-C	-5.16	1.45	1.52
1	A	292	ARG	C-O	-5.15	1.18	1.24
1	A	149	ARG	N-CA	-5.15	1.40	1.46
1	A	268	THR	C-O	5.15	1.30	1.24
1	A	215	ASP	C-O	-5.12	1.17	1.23
1	A	146	GLU	CA-C	5.11	1.59	1.52
1	A	56	LEU	N-CA	-5.11	1.39	1.46
1	A	217	GLU	C-O	-5.08	1.17	1.24
1	A	149	ARG	CD-NE	5.05	1.53	1.46
1	A	159	ALA	N-CA	-5.04	1.39	1.46
1	A	27	GLU	C-O	-5.03	1.17	1.23
1	A	42	SER	C-O	5.03	1.28	1.24
1	A	180	LEU	C-O	5.02	1.30	1.24
1	A	294	LEU	CA-C	-5.02	1.46	1.52

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	56	LEU	O-C-N	7.29	131.46	122.34
1	A	136	VAL	O-C-N	6.73	129.05	122.25
1	A	152	PHE	N-CA-C	6.57	119.44	111.82
1	A	101	ASP	N-CA-C	6.43	120.43	112.59
1	A	172	GLU	CB-CG-CD	6.37	123.43	112.60
1	A	296	VAL	CA-C-O	-6.21	110.25	120.80
1	A	137	THR	N-CA-CB	-5.95	105.67	111.27
1	A	119	ALA	CA-C-O	-5.84	114.35	120.55
1	A	112	SER	CA-C-O	-5.83	113.77	120.24
1	A	17	ILE	CA-C-O	5.76	126.32	120.33
1	A	219	ALA	N-CA-C	5.61	117.39	111.28
1	A	121	GLN	O-C-N	-5.50	114.87	122.46
1	A	267	ILE	O-C-N	-5.40	117.05	123.04
1	A	208	TRP	CA-C-N	-5.34	114.17	119.56
1	A	208	TRP	C-N-CA	-5.34	114.17	119.56
1	A	253[A]	ARG	N-CA-C	5.33	119.04	112.54
1	A	253[B]	ARG	N-CA-C	5.33	119.04	112.54
1	A	137	THR	O-C-N	-5.24	116.67	122.17
1	A	39	ASN	CA-C-O	-5.20	115.82	120.19
1	A	107	ILE	N-CA-C	5.17	117.28	108.86
1	A	282	GLU	O-C-N	5.11	127.98	122.15

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	2417	0	2320	23	0
2	A	1	0	0	0	0
3	A	4	0	0	0	0
4	A	8	0	12	0	0
5	A	317	0	0	2	0
All	All	2747	0	2332	23	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 5.

All (23) 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:253[A]:ARG:CZ	1:A:254:MET:HE3	2.09	0.82
1:A:292:ARG:O	1:A:296:VAL:HG13	1.86	0.76
1:A:254:MET:HE2	1:A:254:MET:HA	1.76	0.67
1:A:253[A]:ARG:CZ	1:A:254:MET:CE	2.75	0.64
1:A:137:THR:HG21	1:A:254:MET:HE1	1.83	0.61
1:A:137:THR:CG2	1:A:254:MET:HE1	2.36	0.56
1:A:109:ASP:CG	5:A:507:HOH:O	2.47	0.55
1:A:253[A]:ARG:NH1	1:A:254:MET:CE	2.70	0.54
1:A:253[A]:ARG:NH1	1:A:254:MET:HE1	2.23	0.54
1:A:21[B]:ARG:HB3	1:A:71:ASP:CG	2.40	0.47
1:A:126:VAL:HG12	1:A:237:ILE:HD11	1.96	0.47
1:A:132:MET:O	1:A:133:GLU:C	2.58	0.46
1:A:14:PHE:CD1	1:A:21[B]:ARG:HD3	2.52	0.45
1:A:254:MET:HE2	1:A:254:MET:CA	2.46	0.44
1:A:25:ILE:O	1:A:60:ILE:HA	2.16	0.44
1:A:107:ILE:HD12	1:A:112:SER:HA	1.99	0.44
1:A:296:VAL:O	1:A:296:VAL:HG23	2.18	0.44
1:A:140:GLU:OE1	1:A:215:ASP:OD1	2.38	0.41
1:A:107:ILE:HB	1:A:112:SER:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:VAL:CG1	1:A:237:ILE:HD11	2.51	0.41
1:A:253[A]:ARG:NE	1:A:254:MET:HE3	2.35	0.41
1:A:109:ASP:CG	1:A:110:TRP:H	2.29	0.40
1:A:217:GLU:OE1	5:A:501:HOH:O	2.22	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	301/296 (102%)	287 (95%)	14 (5%)	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	A	253/243 (104%)	252 (100%)	1 (0%)	84	71

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	188	MET

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	86	HIS
1	A	122	HIS
1	A	182	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 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	A	406[A]	-	3,3,3	0.43	0	2,2,2	0.39	0
4	EDO	A	406[B]	-	3,3,3	2.70	2 (66%)	2,2,2	0.31	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	EDO	A	406[A]	-	-	0/1/1/1	-
4	EDO	A	406[B]	-	-	1/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	406[B]	EDO	O1-C1	3.93	1.62	1.42
4	A	406[B]	EDO	O2-C2	2.52	1.54	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	406[B]	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	292/296 (98%)	-0.17	4 (1%) 73 77	8, 19, 34, 51	11 (3%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	VAL	4.4
1	A	79	ASP	3.6
1	A	254	MET	2.6
1	A	11	GLN	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 oligosaccharides 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
4	EDO	A	406[A]	4/4	0.65	0.18	15,18,19,25	4
4	EDO	A	406[B]	4/4	0.65	0.18	14,17,20,28	4
3	NA	A	405	1/1	0.99	0.10	22,22,22,22	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	A	402	1/1	0.99	0.11	21,21,21,21	0
3	NA	A	404	1/1	0.99	0.05	21,21,21,21	0
2	CL	A	401	1/1	1.00	0.03	18,18,18,18	0
3	NA	A	403	1/1	1.00	0.05	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.