



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 03:39 PM UTC

PDB ID : 3ZLI / pdb_00003zli
Title : Crystal structure of JmjC domain of human histone demethylase UTY
Authors : Vollmar, M.; Gileadi, C.; Shrestha, L.; Goubin, S.; Johansson, C.; Krojer, T.; Raynor, J.W.; Bradley, A.; von Delft, F.; Arrowsmith, C.H.; Bountra, C.; Edwards, A.; Oppermann, U.
Deposited on : 2013-01-31
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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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

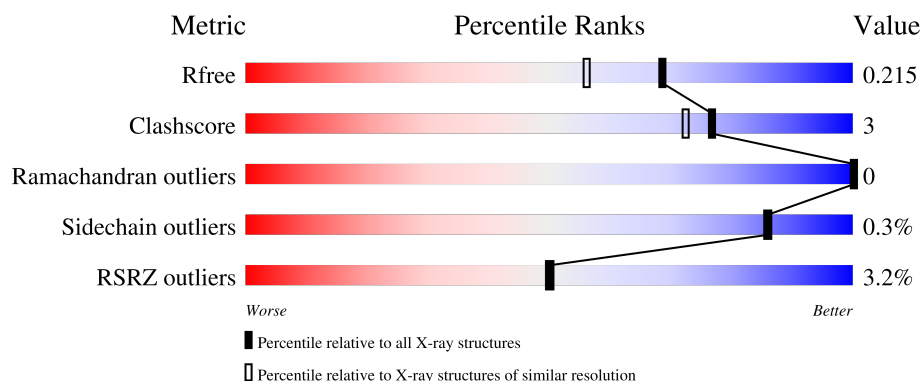
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

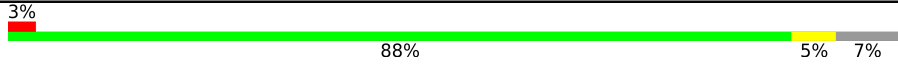
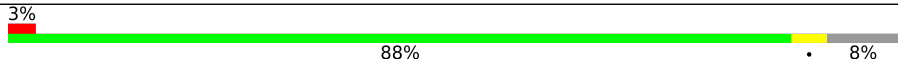
The reported resolution of this entry is 1.80 Å.

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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	478	
1	B	478	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7747 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 HISTONE DEMETHYLASE UTY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	11	1
			3573	2295	607	650	21			
1	B	439	Total	C	N	O	S	0	9	1
			3545	2278	601	646	20			

There are 16 discrepancies between the modelled and reference sequences:

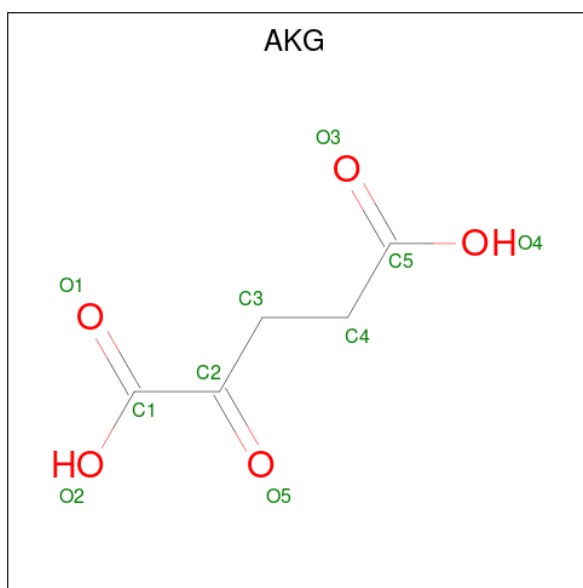
Chain	Residue	Modelled	Actual	Comment	Reference
A	877	MET	-	expression tag	UNP O14607
A	1348	ALA	-	expression tag	UNP O14607
A	1349	GLU	-	expression tag	UNP O14607
A	1350	ASN	-	expression tag	UNP O14607
A	1351	LEU	-	expression tag	UNP O14607
A	1352	TYR	-	expression tag	UNP O14607
A	1353	PHE	-	expression tag	UNP O14607
A	1354	GLN	-	expression tag	UNP O14607
B	877	MET	-	expression tag	UNP O14607
B	1348	ALA	-	expression tag	UNP O14607
B	1349	GLU	-	expression tag	UNP O14607
B	1350	ASN	-	expression tag	UNP O14607
B	1351	LEU	-	expression tag	UNP O14607
B	1352	TYR	-	expression tag	UNP O14607
B	1353	PHE	-	expression tag	UNP O14607
B	1354	GLN	-	expression tag	UNP O14607

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



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

- Molecule 3 is 2-OXOGLUTARIC ACID (CCD ID: AKG) (formula: $C_5H_6O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	5	5		
3	B	1	Total	C	O	0	0
			10	5	5		

- Molecule 4 is FE (II) ION (CCD ID: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Fe	0	0
			1	1		
4	B	1	Total	Fe	0	0
			1	1		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		
5	B	1	Total	Zn	0	0
			1	1		

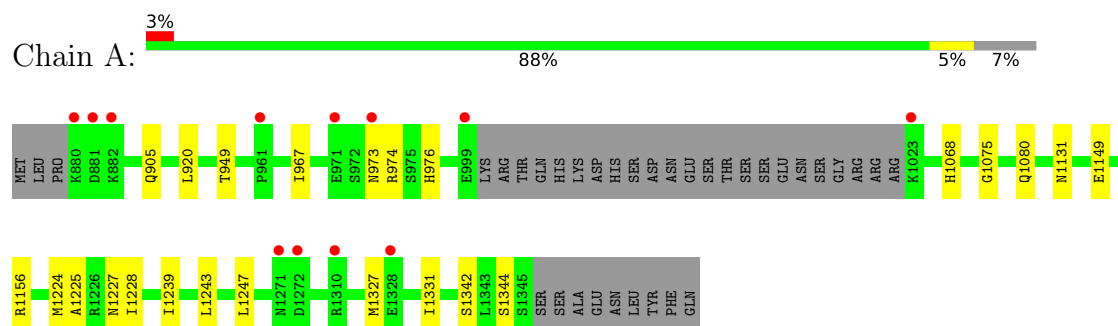
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	286	Total 286	O 286	0	0
6	B	263	Total 263	O 263	0	0

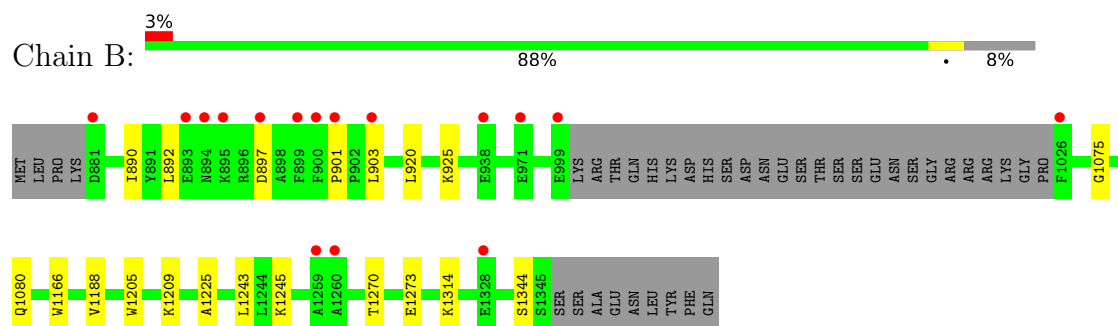
3 Residue-property plots [i](#)

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: HISTONE DEMETHYLASE UTY



• Molecule 1: HISTONE DEMETHYLASE UTY



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.58Å 110.25Å 118.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.01 – 1.80 50.01 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.01-1.80) 99.3 (50.01-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.176 , 0.204 (Not available) , 0.215	Depositor DCC
R_{free} test set	5494 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.517	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 26.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7747	wwPDB-VP
Average B, all atoms (Å ²)	33.0	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, 0.333 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: AKG, FE2, ZN, EDO

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	0.92	2/3701 (0.1%)	0.91	0/5047
1	B	0.89	1/3666 (0.0%)	0.88	1/4999 (0.0%)
All	All	0.90	3/7367 (0.0%)	0.90	1/10046 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1344	SER	C-N	-7.93	1.22	1.33
1	B	1344	SER	C-N	-7.41	1.23	1.33
1	A	1156	ARG	N-CA	-5.08	1.39	1.46

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	901	PRO	N-CA-C	5.48	117.38	110.70

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	A	3573	0	3473	24	0
1	B	3545	0	3461	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	32	0	48	4	0
2	B	24	0	36	1	0
3	A	10	0	4	0	0
3	B	10	0	4	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	286	0	0	4	0
6	B	263	0	0	1	0
All	All	7747	0	7026	41	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (41) 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:890:ILE:HD12	1:B:903:LEU:CD1	2.02	0.89
1:A:905[B]:GLN:OE1	6:A:2030:HOH:O	1.89	0.88
1:A:1224[B]:MET:SD	1:A:1228[B]:ILE:HG12	2.24	0.77
1:A:1247:LEU:HD21	1:A:1331:ILE:HD13	1.70	0.72
1:A:1068:HIS:ND1	1:A:1228[B]:ILE:HD11	2.10	0.66
1:B:890:ILE:CD1	1:B:903:LEU:CD1	2.75	0.65
1:A:1068:HIS:CE1	1:A:1228[B]:ILE:HD11	2.32	0.64
1:B:1075:GLY:HA2	1:B:1080:GLN:HG3	1.82	0.62
1:A:1224[B]:MET:HE1	1:A:1228[B]:ILE:HG21	1.82	0.61
1:B:890:ILE:HD12	1:B:903:LEU:HD12	1.81	0.58
2:B:2350:EDO:H12	6:B:2114:HOH:O	2.04	0.58
1:A:1131:ASN:HB3	2:A:2345:EDO:H12	1.87	0.57
1:A:1075:GLY:HA2	1:A:1080:GLN:HG3	1.87	0.57
2:A:2351:EDO:H12	6:A:2133:HOH:O	2.06	0.54
1:B:1166:TRP:CH2	1:B:1188[A]:VAL:HG21	2.45	0.52
1:A:1224[B]:MET:CE	1:A:1228[B]:ILE:HG12	2.41	0.51
1:A:1342:SER:O	1:B:1245:LYS:HE3	2.11	0.49
1:A:1225:ALA:HB2	1:A:1243:LEU:HD11	1.94	0.49
1:B:1225:ALA:HB2	1:B:1243:LEU:HD11	1.94	0.49
1:B:890:ILE:CD1	1:B:903:LEU:HD12	2.41	0.49
1:A:1327:MET:O	1:A:1331:ILE:HG12	2.14	0.48
1:A:949:THR:HB	1:A:973[B]:ASN:OD1	2.14	0.47
1:A:920:LEU:C	1:A:920:LEU:HD13	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1227:ASN:C	1:A:1228[B]:ILE:CD1	2.89	0.46
1:A:1075:GLY:O	2:A:2351:EDO:H11	2.16	0.46
1:B:1270:THR:HG22	1:B:1273:GLU:HB2	1.98	0.45
1:A:1224[A]:MET:HE2	1:A:1239:ILE:CD1	2.46	0.45
1:A:976:HIS:HE1	6:A:2053:HOH:O	2.00	0.44
1:B:920:LEU:C	1:B:920:LEU:HD13	2.42	0.44
1:B:892:LEU:HD22	1:B:897:ASP:HB3	2.00	0.44
1:A:1342:SER:O	1:B:1245:LYS:CE	2.66	0.43
1:B:1270:THR:CG2	1:B:1273:GLU:HB2	2.49	0.43
1:A:967:ILE:HB	2:A:2345:EDO:H21	2.01	0.43
1:A:1227:ASN:C	1:A:1228[B]:ILE:HD12	2.44	0.43
1:B:890:ILE:HD12	1:B:903:LEU:HD11	1.96	0.42
1:B:1166:TRP:CZ3	1:B:1188[A]:VAL:HG21	2.55	0.42
1:A:974:ARG:NH2	6:A:2053:HOH:O	2.53	0.41
1:B:1314:LYS:HA	1:B:1314:LYS:HD3	1.79	0.41
1:B:1205:TRP:CZ2	1:B:1209:LYS:HE3	2.57	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	450/478 (94%)	441 (98%)	9 (2%)	0	100	100
1	B	444/478 (93%)	436 (98%)	8 (2%)	0	100	100
All	All	894/956 (94%)	877 (98%)	17 (2%)	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	391/432 (90%)	390 (100%)	1 (0%)	86	86
1	B	391/432 (90%)	390 (100%)	1 (0%)	86	86
All	All	782/864 (90%)	780 (100%)	2 (0%)	86	86

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

Mol	Chain	Res	Type
1	A	1149	GLU
1	B	925	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	958	ASN
1	A	1040	ASN
1	A	1048	HIS
1	A	1080	GLN
1	A	1096	ASN
1	A	1183	ASN
1	A	1296	ASN
1	A	1318	ASN
1	B	986	GLN
1	B	1040	ASN
1	B	1080	GLN
1	B	1182	ASN
1	B	1183	ASN
1	B	1296	ASN
1	B	1306	HIS
1	B	1318	ASN

5.3.3 RNA

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 20 ligands modelled in this entry, 4 are monoatomic - leaving 16 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	$\# Z > 2$
2	EDO	A	2348	-	3,3,3	0.36	0	2,2,2	0.31	0
2	EDO	B	2348	-	3,3,3	0.30	0	2,2,2	0.12	0
3	AKG	A	4000	4	9,9,9	1.43	1 (11%)	11,11,11	2.18	6 (54%)
2	EDO	A	2345	-	3,3,3	0.40	0	2,2,2	0.07	0
2	EDO	B	2350	-	3,3,3	0.83	0	2,2,2	0.24	0
2	EDO	B	2345	-	3,3,3	0.44	0	2,2,2	0.08	0
2	EDO	B	2349	-	3,3,3	0.38	0	2,2,2	0.28	0
2	EDO	B	2346	-	3,3,3	0.35	0	2,2,2	0.43	0
2	EDO	A	2352	-	3,3,3	0.32	0	2,2,2	0.63	0
2	EDO	A	2351	-	3,3,3	0.76	0	2,2,2	0.77	0
3	AKG	B	4000	4	9,9,9	1.65	1 (11%)	11,11,11	2.54	5 (45%)
2	EDO	A	2347	-	3,3,3	0.53	0	2,2,2	0.69	0
2	EDO	A	2346	-	3,3,3	0.32	0	2,2,2	0.60	0
2	EDO	A	2350	-	3,3,3	0.29	0	2,2,2	0.60	0
2	EDO	B	2347	-	3,3,3	0.72	0	2,2,2	0.19	0
2	EDO	A	2349	-	3,3,3	0.42	0	2,2,2	0.63	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
2	EDO	A	2348	-	-	0/1/1/1	-
2	EDO	B	2348	-	-	1/1/1/1	-
3	AKG	A	4000	4	-	4/9/9/9	-
2	EDO	A	2345	-	-	0/1/1/1	-
2	EDO	B	2350	-	-	1/1/1/1	-
2	EDO	B	2345	-	-	0/1/1/1	-
2	EDO	B	2349	-	-	0/1/1/1	-
2	EDO	B	2346	-	-	1/1/1/1	-
2	EDO	A	2352	-	-	1/1/1/1	-
2	EDO	A	2351	-	-	0/1/1/1	-
3	AKG	B	4000	4	-	4/9/9/9	-
2	EDO	A	2347	-	-	1/1/1/1	-
2	EDO	A	2346	-	-	0/1/1/1	-
2	EDO	A	2350	-	-	0/1/1/1	-
2	EDO	B	2347	-	-	0/1/1/1	-
2	EDO	A	2349	-	-	1/1/1/1	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	4000	AKG	C2-C1	-3.57	1.48	1.53
3	A	4000	AKG	C2-C1	-3.56	1.48	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	4000	AKG	C3-C2-C1	4.65	123.76	115.86
3	B	4000	AKG	O2-C1-C2	3.69	123.81	113.59
3	B	4000	AKG	O5-C2-C3	-3.38	113.92	121.17
3	A	4000	AKG	C3-C2-C1	3.23	121.35	115.86
3	A	4000	AKG	O2-C1-C2	3.01	121.94	113.59
3	B	4000	AKG	O2-C1-O1	-2.81	117.22	123.90
3	A	4000	AKG	O1-C1-C2	-2.73	118.42	121.81
3	A	4000	AKG	O3-C5-C4	-2.46	115.30	123.09
3	B	4000	AKG	O1-C1-C2	-2.33	118.92	121.81
3	A	4000	AKG	O4-C5-C4	2.26	121.13	114.00
3	A	4000	AKG	O5-C2-C3	-2.22	116.40	121.17

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	4000	AKG	C1-C2-C3-C4
3	B	4000	AKG	C1-C2-C3-C4
2	A	2347	EDO	O1-C1-C2-O2
3	A	4000	AKG	C3-C4-C5-O3
3	A	4000	AKG	C3-C4-C5-O4
3	B	4000	AKG	C3-C4-C5-O4
3	B	4000	AKG	C3-C4-C5-O3
3	B	4000	AKG	O5-C2-C3-C4
2	A	2352	EDO	O1-C1-C2-O2
3	A	4000	AKG	O5-C2-C3-C4
2	B	2346	EDO	O1-C1-C2-O2
2	B	2348	EDO	O1-C1-C2-O2
2	B	2350	EDO	O1-C1-C2-O2
2	A	2349	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2345	EDO	2	0
2	B	2350	EDO	1	0
2	A	2351	EDO	2	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	443/478 (92%)	-0.05	12 (2%)	56 56	13, 30, 48, 79	11 (2%)
1	B	439/478 (91%)	0.05	16 (3%)	46 46	15, 32, 54, 80	9 (2%)
All	All	882/956 (92%)	-0.00	28 (3%)	50 50	13, 31, 51, 80	20 (2%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	900	PHE	5.3
1	B	999	GLU	4.5
1	A	880	LYS	4.4
1	A	999	GLU	3.9
1	B	901	PRO	3.9
1	B	971	GLU	3.7
1	B	899	PHE	3.5
1	B	1328	GLU	3.3
1	B	881	ASP	3.1
1	B	894	ASN	3.1
1	A	1272	ASP	3.0
1	A	1310[A]	ARG	2.9
1	B	897	ASP	2.8
1	A	1271	ASN	2.8
1	A	1023	LYS	2.7
1	A	961	PRO	2.6
1	A	1328	GLU	2.6
1	B	1259	ALA	2.5
1	B	895	LYS	2.5
1	B	1026	PHE	2.5
1	B	1260	ALA	2.4
1	A	881	ASP	2.4
1	A	882	LYS	2.3
1	A	971	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	973[A]	ASN	2.2
1	B	938	GLU	2.2
1	B	903	LEU	2.2
1	B	893	GLU	2.1

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
2	EDO	A	2352	4/4	0.77	0.15	45,51,51,58	0
2	EDO	B	2350	4/4	0.80	0.20	45,46,47,59	0
2	EDO	A	2345	4/4	0.81	0.15	43,47,48,53	0
2	EDO	A	2349	4/4	0.82	0.16	47,49,49,52	0
2	EDO	B	2348	4/4	0.84	0.19	53,57,58,60	0
2	EDO	A	2351	4/4	0.85	0.19	34,45,48,50	0
2	EDO	B	2347	4/4	0.85	0.16	33,41,42,46	0
2	EDO	A	2347	4/4	0.86	0.15	32,43,45,53	0
2	EDO	B	2349	4/4	0.91	0.11	45,46,50,52	0
2	EDO	A	2348	4/4	0.93	0.11	43,46,50,51	0
3	AKG	A	4000	10/10	0.93	0.09	24,29,33,33	0
2	EDO	B	2346	4/4	0.94	0.09	47,48,50,52	0
2	EDO	A	2346	4/4	0.94	0.10	31,31,34,34	0
2	EDO	A	2350	4/4	0.94	0.09	42,43,45,47	0
2	EDO	B	2345	4/4	0.95	0.08	30,32,35,35	0
3	AKG	B	4000	10/10	0.96	0.07	25,32,35,36	0
5	ZN	B	4002	1/1	0.98	0.03	29,29,29,29	0
5	ZN	A	4002	1/1	0.99	0.02	27,27,27,27	0
4	FE2	A	4001	1/1	1.00	0.04	24,24,24,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FE2	B	4001	1/1	1.00	0.04	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.