



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:00 PM UTC

PDB ID : 5FY1 / pdb_00005fy1
Title : Crystal structure of JmjC domain of human histone demethylase UTY in complex with N08619b
Authors : Nowak, R.; Krojer, T.; Johansson, C.; Gileadi, C.; Kupinska, K.; Pearce, N.M.; von Delft, F.; Arrowsmith, C.H.; Bountra, C.; Edwards, A.; Oppermann, U.
Deposited on : 2016-03-03
Resolution : 1.78 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

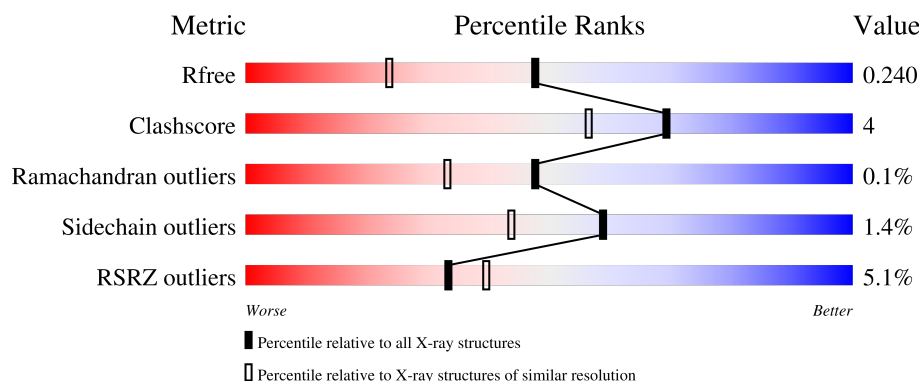
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	5% 85% 7% 8%
1	B	478	4% 87% 6% 7%

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
5	DMS	A	2350	-	-	X	-
5	DMS	B	2354	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7902 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	442	Total	C	N	O	S	0	7	0
			3576	2292	611	651	22			
1	B	446	Total	C	N	O	S	0	7	0
			3576	2291	610	652	23			

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

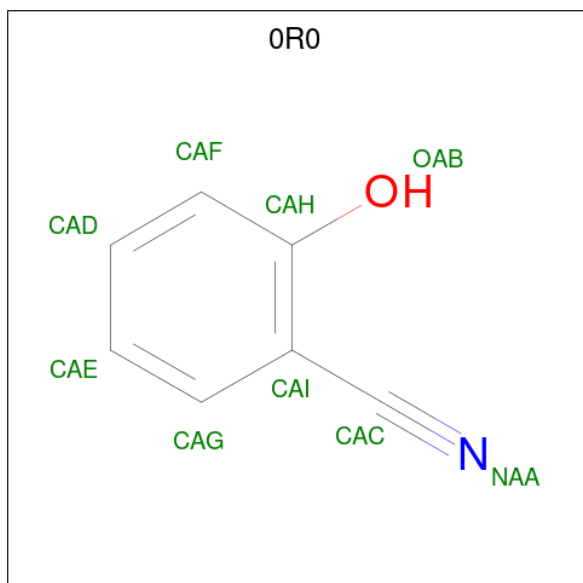
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	B	1	Total	Mn	0	0
			1	1		

- Molecule 4 is 2-hydroxybenzonitrile (CCD ID: 0R0) (formula: C₇H₅NO).



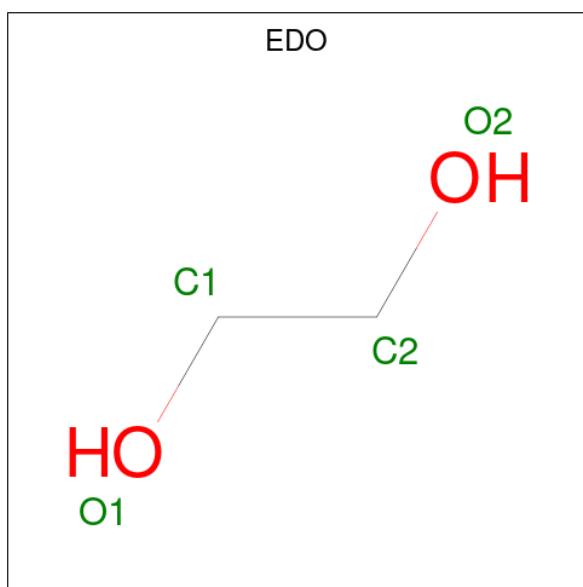
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			9	7	1	1		
4	A	1	Total	C	N	O	0	0
			9	7	1	1		

- Molecule 5 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



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

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

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

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

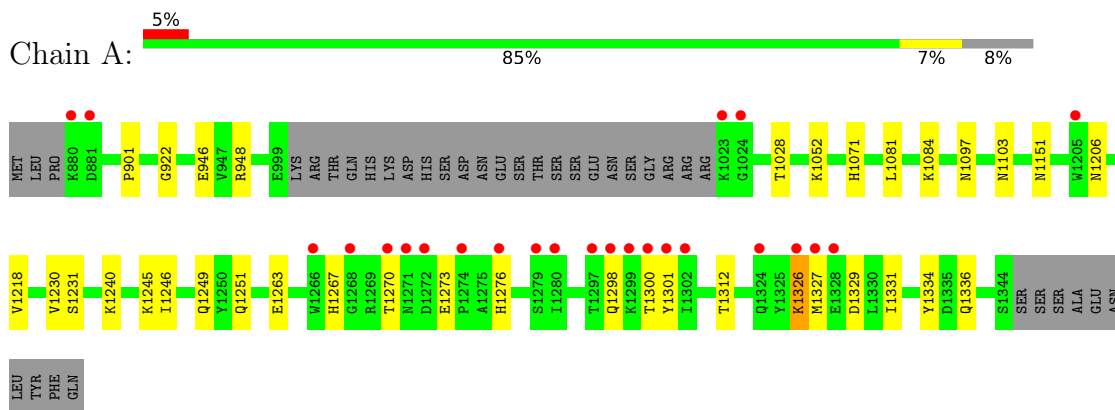
- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	306	Total O 306 306	0	0
7	B	286	Total O 286 286	0	0

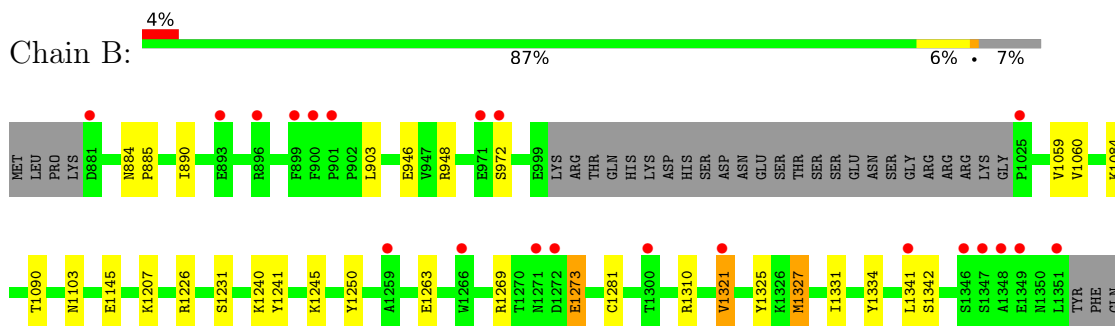
3 Residue-property plots [i](#)

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	91.71Å 110.08Å 118.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.19 – 1.78 47.19 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.9 (47.19-1.78) 99.9 (47.19-1.78)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.78Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.201 , 0.239 0.205 , 0.240	Depositor DCC
R_{free} test set	5722 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.5	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	0.96	EDS
Total number of atoms	7902	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% 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: DMS, OR0, EDO, MN, ZN

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	0.56	0/3671	0.78	2/5004 (0.0%)
1	B	0.57	0/3670	0.79	2/5005 (0.0%)
All	All	0.56	0/7341	0.79	4/10009 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	901	PRO	CA-C-N	-5.84	114.10	119.82
1	A	901	PRO	C-N-CA	-5.84	114.10	119.82
1	B	885	PRO	CA-C-N	5.34	125.28	119.78
1	B	885	PRO	C-N-CA	5.34	125.28	119.78

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	3576	0	3454	27	1
1	B	3576	0	3446	26	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
4	A	18	0	10	0	0
5	A	8	0	12	7	0
5	B	4	0	6	6	0
6	A	56	0	84	5	0
6	B	68	0	102	4	0
7	A	306	0	0	4	0
7	B	286	0	0	5	1
All	All	7902	0	7114	51	1

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 (51) 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:1206:ASN:HD21	5:A:2350:DMS:H22	1.22	1.03
1:A:1097:ASN:H	5:A:2350:DMS:H21	1.25	1.02
1:B:1250:TYR:OH	7:B:2228:HOH:O	2.01	0.79
1:B:1103:ASN:HD21	5:B:2354:DMS:H12	1.47	0.79
1:B:1325:TYR:O	7:B:2228:HOH:O	2.01	0.79
1:A:946:GLU:OE2	1:A:948:ARG:NH1	2.16	0.74
1:A:1251:GLN:HE21	1:A:1327[A]:MET:HE3	1.51	0.74
1:A:1084:LYS:HZ3	5:A:2349:DMS:H21	1.57	0.70
1:A:1326:LYS:HE2	1:A:1329:ASP:CB	2.22	0.69
1:A:1097:ASN:H	5:A:2350:DMS:C2	2.04	0.68
1:A:1071:HIS:ND1	7:A:2132:HOH:O	2.25	0.68
1:B:1145:GLU:OE1	7:B:2174:HOH:O	2.13	0.67
1:A:1052[B]:LYS:NZ	7:A:2111:HOH:O	2.29	0.66
1:B:884[A]:ASN:ND2	7:B:2006:HOH:O	2.23	0.63
1:A:1251:GLN:HG3	1:A:1327[B]:MET:HE3	1.81	0.62
1:B:946:GLU:OE2	1:B:948[A]:ARG:NH1	2.34	0.61
1:B:948[B]:ARG:NH1	1:B:972:SER:O	2.34	0.61
1:A:1052[A]:LYS:NZ	7:A:2034:HOH:O	2.36	0.59
1:B:1226:ARG:NH1	7:B:2215:HOH:O	2.36	0.58
1:B:1084:LYS:HZ3	5:B:2354:DMS:H21	1.70	0.57
1:B:1231:SER:HB2	1:B:1342:SER:HB3	1.88	0.54
1:B:1084:LYS:HZ3	5:B:2354:DMS:C2	2.22	0.53
1:A:1270:THR:HG23	1:A:1273:GLU:HB2	1.91	0.53
1:B:1103:ASN:HD21	5:B:2354:DMS:C1	2.20	0.53
1:A:1267:HIS:NE2	6:A:2354:EDO:O1	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1245:LYS:O	1:A:1249[B]:GLN:OE1	2.28	0.51
1:B:1241:TYR:CZ	1:B:1245:LYS:HD2	2.46	0.51
1:A:1240:LYS:HE3	1:A:1334:TYR:CZ	2.47	0.50
1:A:1103:ASN:HD21	5:A:2349:DMS:C1	2.24	0.50
1:B:890:ILE:HD13	1:B:903:LEU:HD13	1.94	0.49
1:A:1097:ASN:N	5:A:2350:DMS:H21	2.09	0.49
1:A:1231:SER:H	6:A:2361:EDO:H21	1.78	0.48
1:A:1336:GLN:HA	1:B:1310:ARG:NH1	2.28	0.48
1:A:1336:GLN:HA	1:B:1310:ARG:HH11	1.79	0.48
1:B:1240:LYS:HE3	1:B:1334:TYR:CZ	2.49	0.48
1:A:1103:ASN:HD21	5:A:2349:DMS:H12	1.80	0.47
1:B:1090:THR:OG1	5:B:2354:DMS:H23	2.16	0.46
1:B:1207:LYS:HA	6:B:2361:EDO:H12	1.97	0.46
1:A:1218:VAL:HG13	1:A:1246:ILE:HD13	1.98	0.45
1:B:1263:GLU:O	1:B:1321:VAL:HG22	2.16	0.45
1:B:1327:MET:HE2	1:B:1331:ILE:HD11	1.98	0.45
1:A:948:ARG:NE	7:A:2053:HOH:O	2.50	0.43
1:A:922:GLY:H	6:A:2355:EDO:C2	2.30	0.43
1:A:1230:VAL:HA	6:A:2361:EDO:H21	2.00	0.43
1:B:1059:VAL:HB	6:B:2359:EDO:H22	2.00	0.43
1:A:1276:HIS:CE1	6:A:2354:EDO:H21	2.54	0.43
1:A:1298:GLN:HB2	1:A:1300:THR:HG23	2.01	0.43
1:B:1269:ARG:NH1	1:B:1273:GLU:O	2.37	0.42
1:B:1103:ASN:ND2	5:B:2354:DMS:H12	2.25	0.41
1:B:1060:VAL:H	6:B:2359:EDO:H22	1.85	0.41
1:B:1281:CYS:O	6:B:2361:EDO:H21	2.20	0.40

All (1) 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)
1:A:1249[B]:GLN:NE2	7:B:2219:HOH:O[4_555]	1.91	0.29

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	445/478 (93%)	436 (98%)	8 (2%)	1 (0%)	43	29
1	B	449/478 (94%)	439 (98%)	10 (2%)	0	100	100
All	All	894/956 (94%)	875 (98%)	18 (2%)	1 (0%)	48	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1301	TYR

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	388/432 (90%)	381 (98%)	7 (2%)	51	33
1	B	386/432 (89%)	382 (99%)	4 (1%)	68	55
All	All	774/864 (90%)	763 (99%)	11 (1%)	59	44

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1028	THR
1	A	1081	LEU
1	A	1151	ASN
1	A	1263	GLU
1	A	1312	THR
1	A	1326	LYS
1	A	1331	ILE
1	B	1273	GLU
1	B	1321	VAL
1	B	1327	MET
1	B	1341	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1096	ASN
1	A	1183	ASN
1	A	1206	ASN
1	A	1251	GLN
1	A	1304	HIS
1	A	1318	ASN
1	B	991	GLN
1	B	1080	GLN
1	B	1318	ASN

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 40 ligands modelled in this entry, 4 are monoatomic - leaving 36 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	$\# Z > 2$
6	EDO	B	2363	-	3,3,3	0.41	0	2,2,2	0.73	0
6	EDO	A	2353	-	3,3,3	0.46	0	2,2,2	0.34	0
6	EDO	B	2364	-	3,3,3	0.39	0	2,2,2	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	B	2359	-	3,3,3	0.40	0	2,2,2	0.34	0
5	DMS	A	2350	-	3,3,3	0.52	0	3,3,3	0.55	0
5	DMS	A	2349	-	3,3,3	0.56	0	3,3,3	0.76	0
6	EDO	B	2367	-	3,3,3	0.43	0	2,2,2	0.54	0
4	OR0	A	2348	-	9,9,9	1.90	2 (22%)	10,11,11	0.64	0
6	EDO	A	2351	-	3,3,3	0.31	0	2,2,2	0.47	0
4	OR0	A	2347	-	9,9,9	1.88	2 (22%)	10,11,11	0.50	0
6	EDO	A	2352	-	3,3,3	0.44	0	2,2,2	0.36	0
6	EDO	B	2356	-	3,3,3	0.46	0	2,2,2	0.38	0
6	EDO	A	2358	-	3,3,3	0.38	0	2,2,2	0.56	0
6	EDO	A	2355	-	3,3,3	0.50	0	2,2,2	0.25	0
6	EDO	B	2360	-	3,3,3	0.47	0	2,2,2	0.45	0
6	EDO	B	2368	-	3,3,3	0.49	0	2,2,2	0.25	0
6	EDO	B	2365	-	3,3,3	0.42	0	2,2,2	0.43	0
6	EDO	B	2371	-	3,3,3	0.43	0	2,2,2	0.61	0
6	EDO	B	2362	-	3,3,3	0.44	0	2,2,2	0.51	0
6	EDO	B	2370	-	3,3,3	0.43	0	2,2,2	0.39	0
6	EDO	A	2356	-	3,3,3	0.34	0	2,2,2	0.98	0
6	EDO	B	2361	-	3,3,3	0.44	0	2,2,2	0.48	0
6	EDO	B	2358	-	3,3,3	0.34	0	2,2,2	0.84	0
6	EDO	B	2366	-	3,3,3	0.54	0	2,2,2	0.38	0
6	EDO	A	2361	-	3,3,3	0.36	0	2,2,2	0.61	0
6	EDO	B	2369	-	3,3,3	0.38	0	2,2,2	0.41	0
6	EDO	B	2355	-	3,3,3	0.48	0	2,2,2	0.30	0
6	EDO	B	2357	-	3,3,3	0.41	0	2,2,2	0.38	0
6	EDO	A	2364	-	3,3,3	0.44	0	2,2,2	0.46	0
6	EDO	A	2362	-	3,3,3	0.51	0	2,2,2	0.26	0
6	EDO	A	2354	-	3,3,3	0.42	0	2,2,2	0.39	0
6	EDO	A	2363	-	3,3,3	0.42	0	2,2,2	0.44	0
6	EDO	A	2359	-	3,3,3	0.47	0	2,2,2	0.73	0
6	EDO	A	2360	-	3,3,3	0.43	0	2,2,2	0.52	0
5	DMS	B	2354	-	3,3,3	0.45	0	3,3,3	0.69	0
6	EDO	A	2357	-	3,3,3	0.43	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
6	EDO	B	2363	-	-	1/1/1/1	-
6	EDO	A	2353	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	EDO	B	2364	-	-	0/1/1/1	-
6	EDO	B	2359	-	-	1/1/1/1	-
6	EDO	B	2367	-	-	0/1/1/1	-
4	OR0	A	2348	-	-	0/2/2/2	0/1/1/1
6	EDO	A	2351	-	-	1/1/1/1	-
4	OR0	A	2347	-	-	1/2/2/2	0/1/1/1
6	EDO	A	2352	-	-	0/1/1/1	-
6	EDO	B	2356	-	-	0/1/1/1	-
6	EDO	A	2358	-	-	0/1/1/1	-
6	EDO	A	2355	-	-	1/1/1/1	-
6	EDO	B	2360	-	-	1/1/1/1	-
6	EDO	B	2368	-	-	0/1/1/1	-
6	EDO	B	2365	-	-	1/1/1/1	-
6	EDO	B	2371	-	-	0/1/1/1	-
6	EDO	B	2362	-	-	0/1/1/1	-
6	EDO	B	2370	-	-	0/1/1/1	-
6	EDO	A	2356	-	-	1/1/1/1	-
6	EDO	B	2361	-	-	0/1/1/1	-
6	EDO	B	2358	-	-	1/1/1/1	-
6	EDO	B	2366	-	-	0/1/1/1	-
6	EDO	A	2361	-	-	1/1/1/1	-
6	EDO	B	2369	-	-	0/1/1/1	-
6	EDO	B	2355	-	-	0/1/1/1	-
6	EDO	B	2357	-	-	1/1/1/1	-
6	EDO	A	2364	-	-	0/1/1/1	-
6	EDO	A	2362	-	-	1/1/1/1	-
6	EDO	A	2354	-	-	0/1/1/1	-
6	EDO	A	2363	-	-	1/1/1/1	-
6	EDO	A	2359	-	-	0/1/1/1	-
6	EDO	A	2360	-	-	0/1/1/1	-
6	EDO	A	2357	-	-	0/1/1/1	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2348	OR0	CAI-CAC	5.11	1.51	1.44
4	A	2347	OR0	CAI-CAC	4.91	1.51	1.44
4	A	2347	OR0	OAB-CAH	2.65	1.41	1.36
4	A	2348	OR0	OAB-CAH	2.28	1.40	1.36

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	2359	EDO	O1-C1-C2-O2
6	B	2360	EDO	O1-C1-C2-O2
6	A	2355	EDO	O1-C1-C2-O2
6	A	2361	EDO	O1-C1-C2-O2
6	A	2362	EDO	O1-C1-C2-O2
6	A	2353	EDO	O1-C1-C2-O2
6	A	2356	EDO	O1-C1-C2-O2
6	A	2363	EDO	O1-C1-C2-O2
6	B	2365	EDO	O1-C1-C2-O2
4	A	2347	OR0	NAA-CAC-CAI-CAH
6	B	2357	EDO	O1-C1-C2-O2
6	B	2358	EDO	O1-C1-C2-O2
6	B	2363	EDO	O1-C1-C2-O2
6	A	2351	EDO	O1-C1-C2-O2

There are no ring outliers.

8 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	2359	EDO	2	0
5	A	2350	DMS	4	0
5	A	2349	DMS	3	0
6	A	2355	EDO	1	0
6	B	2361	EDO	2	0
6	A	2361	EDO	2	0
6	A	2354	EDO	2	0
5	B	2354	DMS	6	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	442/478 (92%)	0.25	24 (5%)	31 37	12, 32, 56, 88	7 (1%)
1	B	446/478 (93%)	0.28	21 (4%)	36 42	14, 33, 56, 82	7 (1%)
All	All	888/956 (92%)	0.27	45 (5%)	33 39	12, 32, 56, 88	14 (1%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1341	LEU	3.9
1	A	1270	THR	3.8
1	B	1025	PRO	3.5
1	A	1301	TYR	3.4
1	A	1302	ILE	3.4
1	B	1347	SER	3.3
1	B	900	PHE	3.2
1	B	1349	GLU	3.1
1	B	1351	LEU	3.1
1	A	1279	SER	3.1
1	A	1327[A]	MET	3.0
1	B	1271	ASN	3.0
1	B	971	GLU	2.8
1	B	1266	TRP	2.7
1	A	1023	LYS	2.7
1	A	1299	LYS	2.7
1	A	1205	TRP	2.6
1	B	1259	ALA	2.6
1	A	880	LYS	2.6
1	A	1300	THR	2.6
1	B	1346	SER	2.5
1	A	881	ASP	2.5
1	B	1272	ASP	2.5
1	A	1266	TRP	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	1321	VAL	2.5
1	A	1280	ILE	2.4
1	A	1268	GLY	2.4
1	B	1348	ALA	2.4
1	A	1326	LYS	2.4
1	A	1272	ASP	2.3
1	B	896	ARG	2.3
1	A	1324	GLN	2.3
1	B	899	PHE	2.3
1	B	893	GLU	2.3
1	A	1271	ASN	2.3
1	B	972	SER	2.3
1	A	1276	HIS	2.2
1	A	1297	THR	2.2
1	B	881	ASP	2.1
1	B	1300	THR	2.1
1	A	1298	GLN	2.1
1	A	1274	PRO	2.1
1	B	901	PRO	2.1
1	A	1024	GLY	2.1
1	A	1328	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($\text{\AA}^2$)	Q<0.9
6	EDO	B	2360	4/4	0.79	0.16	33,59,60,60	0
6	EDO	B	2371	4/4	0.80	0.20	42,52,53,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	OR0	A	2348	9/9	0.81	0.20	36,42,46,46	9
6	EDO	A	2363	4/4	0.81	0.15	52,56,57,58	0
6	EDO	A	2364	4/4	0.83	0.13	56,57,60,68	0
6	EDO	A	2361	4/4	0.83	0.17	27,37,47,47	0
4	OR0	A	2347	9/9	0.83	0.16	31,36,40,47	9
6	EDO	B	2369	4/4	0.84	0.15	36,36,53,59	0
6	EDO	B	2363	4/4	0.86	0.14	42,46,47,59	0
5	DMS	A	2350	4/4	0.87	0.14	32,41,55,72	0
5	DMS	A	2349	4/4	0.88	0.15	33,42,49,61	0
6	EDO	B	2370	4/4	0.88	0.12	30,36,41,43	0
6	EDO	B	2365	4/4	0.88	0.13	42,43,51,52	0
6	EDO	B	2366	4/4	0.89	0.17	32,33,38,39	0
6	EDO	B	2362	4/4	0.89	0.10	28,30,44,46	0
6	EDO	A	2360	4/4	0.89	0.12	33,38,43,47	0
6	EDO	B	2361	4/4	0.89	0.13	30,41,48,58	0
6	EDO	B	2358	4/4	0.90	0.16	30,37,40,49	0
6	EDO	A	2358	4/4	0.90	0.14	27,28,37,68	0
6	EDO	A	2352	4/4	0.91	0.10	38,42,46,52	0
6	EDO	B	2357	4/4	0.91	0.10	36,40,47,51	0
6	EDO	A	2362	4/4	0.91	0.11	38,44,44,47	0
6	EDO	A	2356	4/4	0.91	0.12	39,39,43,44	0
6	EDO	B	2364	4/4	0.92	0.12	39,47,48,50	0
5	DMS	B	2354	4/4	0.92	0.11	30,34,46,60	0
6	EDO	A	2353	4/4	0.92	0.09	44,44,49,57	0
6	EDO	A	2354	4/4	0.92	0.11	44,50,53,57	0
6	EDO	B	2356	4/4	0.92	0.13	34,36,46,52	0
6	EDO	A	2355	4/4	0.92	0.20	20,24,40,48	0
6	EDO	A	2359	4/4	0.93	0.09	29,33,35,40	0
6	EDO	B	2359	4/4	0.93	0.25	33,40,41,43	0
6	EDO	A	2351	4/4	0.93	0.09	25,30,37,43	0
6	EDO	B	2355	4/4	0.95	0.12	34,35,39,44	0
6	EDO	B	2367	4/4	0.95	0.10	26,28,29,33	0
6	EDO	B	2368	4/4	0.95	0.09	24,30,32,35	0
6	EDO	A	2357	4/4	0.97	0.06	24,25,27,29	0
2	ZN	A	2345	1/1	0.98	0.03	38,38,38,38	0
3	MN	A	2346	1/1	0.99	0.04	30,30,30,30	0
2	ZN	B	2352	1/1	1.00	0.02	27,27,27,27	0
3	MN	B	2353	1/1	1.00	0.04	29,29,29,29	0

6.5 Other polymers ⓘ

There are no such residues in this entry.