



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 22, 2016 – 04:06 AM EDT

PDB ID : 5F2W
Title : Crystal structure of human KDM4A in complex with compound 16
Authors : Le Bihan, Y.-V.; Dempster, S.; Westwood, I.M.; van Montfort, R.L.M.
Deposited on : 2015-12-02
Resolution : 2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

i

X-RAY DIFFRACTION

A.



R_{free}
Clashscore
Ramachandran outliers
Sidechain outliers
RSRZ outliers

		g
1	A	360
1	B	360
1	C	360
1	D	360

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	5UP	A	403	-	-	-	X
4	EDO	A	404	-	-	-	X
4	EDO	B	405	-	-	-	X
4	EDO	B	407	-	-	-	X
4	EDO	B	408	-	-	-	X
4	EDO	B	412	-	-	-	X
4	EDO	B	413	-	-	-	X
4	EDO	C	408	-	-	-	X
4	EDO	D	404	-	-	-	X
4	EDO	D	407	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11386 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 Lysine-specific demethylase 4A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2675	1736	436	490	13			
1	B	350	Total	C	N	O	S	0	1	0
			2769	1789	454	511	15			
1	C	350	Total	C	N	O	S	0	0	0
			2757	1792	447	504	14			
1	D	350	Total	C	N	O	S	0	0	0
			2782	1805	461	501	15			

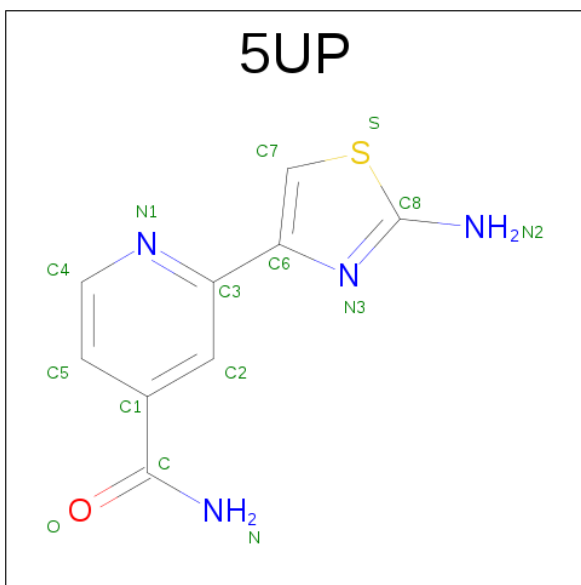
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP O75164
B	0	SER	-	expression tag	UNP O75164
C	0	SER	-	expression tag	UNP O75164
D	0	SER	-	expression tag	UNP O75164

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

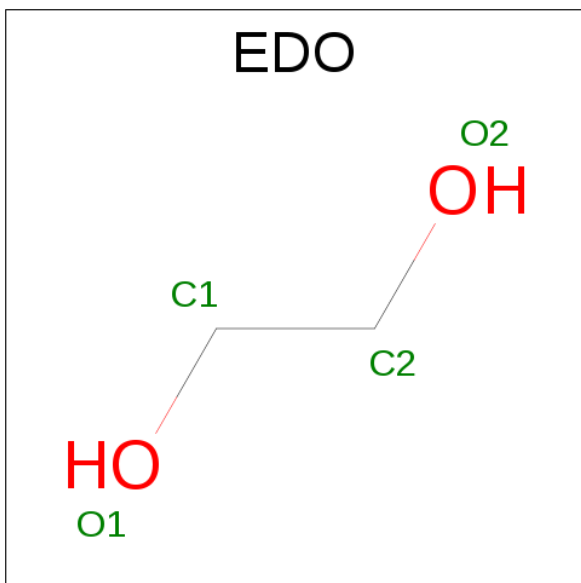
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Zn	0	0
			2	2		
2	A	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		

- Molecule 3 is 2-(2-azanyl-1,3-thiazol-4-yl)pyridine-4-carboxamide (three-letter code: 5UP) (formula: C₉H₈N₄OS).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			15	9	4	1	1		
3	B	1	Total	C	N	O	S	0	0
			15	9	4	1	1		
3	C	1	Total	C	N	O	S	0	0
			15	9	4	1	1		
3	D	1	Total	C	N	O	S	0	0
			15	9	4	1	1		

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



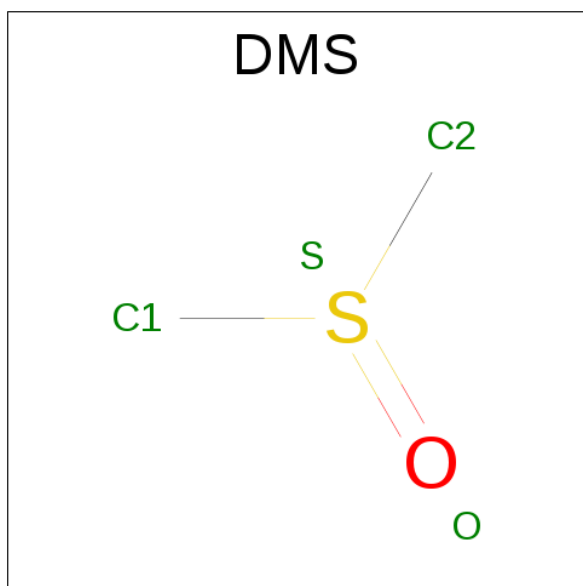
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0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

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

- Molecule 5 is DIMETHYL SULFOXIDE (three-letter code: DMS) (formula: C₂H₆OS).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	45	Total	O	0	0
			45	45		

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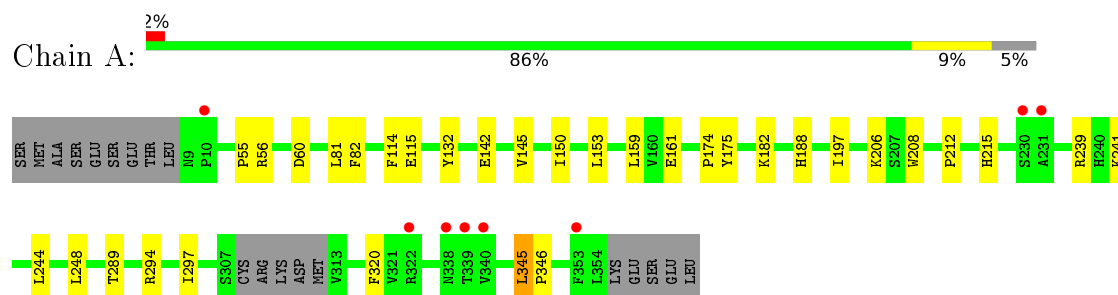
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	57	Total 57	O 57	0	0
6	C	51	Total 51	O 51	0	0
6	D	62	Total 62	O 62	0	0

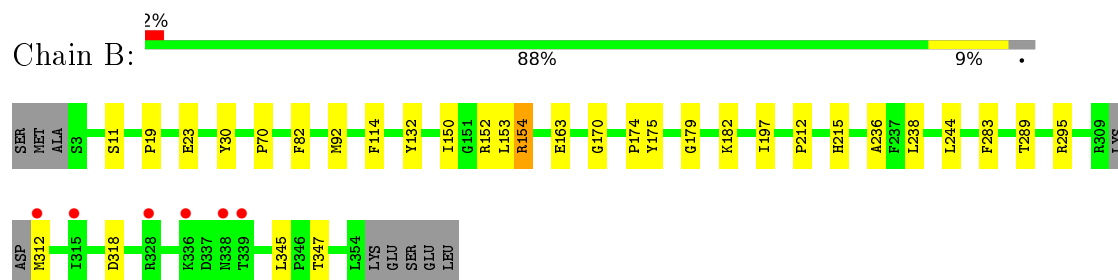
3 Residue-property plots [i](#)

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.

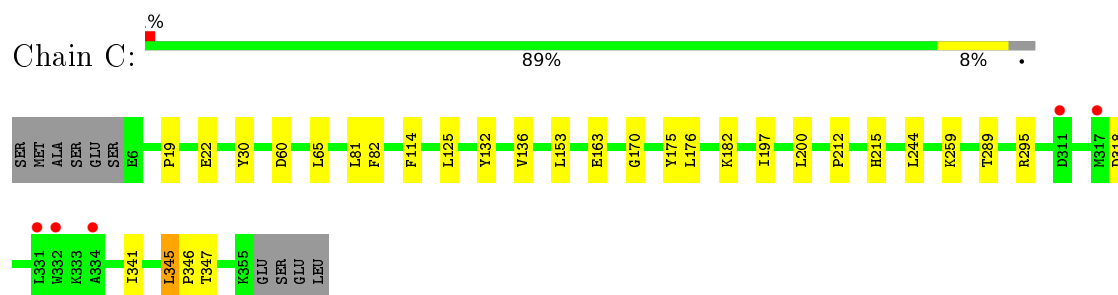
• Molecule 1: Lysine-specific demethylase 4A



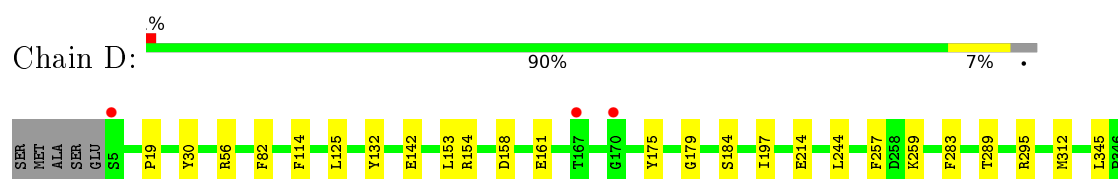
• Molecule 1: Lysine-specific demethylase 4A



• Molecule 1: Lysine-specific demethylase 4A



• Molecule 1: Lysine-specific demethylase 4A



T347	L354
LYS	GLU
GLU	SER
LEU	LEU

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.39Å 102.04Å 142.15Å 90.00° 99.70° 90.00°	Depositor
Resolution (Å)	49.48 – 2.60 49.48 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.7 (49.48-2.60) 95.7 (49.48-2.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.61Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.179 , 0.226 0.193 , 0.240	Depositor DCC
R_{free} test set	2424 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 42.3	EDS
Estimated twinning fraction	0.027 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 47648 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11386	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.57 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.7421e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: ZN, DMS, 5UP, 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	0.51	0/2760	0.67	0/3763
1	B	0.52	0/2854	0.67	0/3887
1	C	0.52	0/2844	0.68	0/3876
1	D	0.53	0/2869	0.69	0/3901
All	All	0.52	0/11327	0.68	0/15427

There are no bond length outliers.

There are no bond angle outliers.

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	2675	0	2448	13	0
1	B	2769	0	2536	15	0
1	C	2757	0	2537	12	0
1	D	2782	0	2596	11	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	15	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	15	0	8	0	0
3	C	15	0	8	0	0
3	D	15	0	8	0	0
4	A	12	0	18	1	0
4	B	40	0	60	6	0
4	C	32	0	48	3	0
4	D	32	0	48	1	0
5	A	4	0	6	0	0
6	A	45	0	0	0	0
6	B	57	0	0	0	0
6	C	51	0	0	0	0
6	D	62	0	0	2	0
All	All	11386	0	10329	51	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 2.

The worst 5 of 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:B:238:LEU:HB2	4:B:413:EDO:H12	1.67	0.76
1:A:239:ARG:O	1:A:241:LYS:NZ	2.26	0.69
1:A:294:ARG:O	1:A:297:ILE:HG12	1.96	0.65
1:B:150:ILE:HG22	1:B:174:PRO:HB3	1.82	0.60
1:C:153:LEU:HD11	1:C:197:ILE:HG21	1.87	0.57

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	337/360 (94%)	330 (98%)	7 (2%)	0	100	100
1	B	347/360 (96%)	339 (98%)	8 (2%)	0	100	100
1	C	348/360 (97%)	342 (98%)	6 (2%)	0	100	100
1	D	348/360 (97%)	340 (98%)	8 (2%)	0	100	100
All	All	1380/1440 (96%)	1351 (98%)	29 (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	263/316 (83%)	252 (96%)	11 (4%)	36	65
1	B	275/316 (87%)	265 (96%)	10 (4%)	42	71
1	C	272/316 (86%)	260 (96%)	12 (4%)	35	63
1	D	277/316 (88%)	269 (97%)	8 (3%)	50	77
All	All	1087/1264 (86%)	1046 (96%)	41 (4%)	40	68

5 of 41 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	312	MET
1	C	60	ASP
1	D	184	SER
1	B	318	ASP
1	B	345	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	215	HIS
1	C	215	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 42 ligands modelled in this entry, 8 are monoatomic - leaving 34 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	5UP	A	403	2	13,16,16	0.42	0	15,22,22	0.52	0
4	EDO	A	404	-	3,3,3	0.40	0	2,2,2	0.43	0
4	EDO	A	405	-	3,3,3	0.46	0	2,2,2	0.65	0
4	EDO	A	406	-	3,3,3	0.57	0	2,2,2	0.27	0
5	DMS	A	407	-	3,3,3	0.23	0	3,3,3	0.50	0
3	5UP	B	403	2	13,16,16	0.52	0	15,22,22	0.56	0
4	EDO	B	404	-	3,3,3	0.25	0	2,2,2	0.58	0
4	EDO	B	405	-	3,3,3	0.64	0	2,2,2	0.15	0
4	EDO	B	406	-	3,3,3	0.54	0	2,2,2	0.30	0
4	EDO	B	407	-	3,3,3	0.60	0	2,2,2	0.32	0
4	EDO	B	408	-	3,3,3	0.44	0	2,2,2	0.46	0
4	EDO	B	409	-	3,3,3	0.71	0	2,2,2	0.28	0
4	EDO	B	410	-	3,3,3	0.50	0	2,2,2	0.52	0
4	EDO	B	411	-	3,3,3	0.41	0	2,2,2	0.46	0
4	EDO	B	412	-	3,3,3	0.56	0	2,2,2	0.43	0
4	EDO	B	413	-	3,3,3	0.88	0	2,2,2	0.41	0
3	5UP	C	403	2	13,16,16	0.51	0	15,22,22	0.53	0
4	EDO	C	404	-	3,3,3	0.69	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	C	405	-	3,3,3	0.63	0	2,2,2	0.28	0
4	EDO	C	406	-	3,3,3	0.61	0	2,2,2	0.29	0
4	EDO	C	407	-	3,3,3	0.53	0	2,2,2	0.40	0
4	EDO	C	408	-	3,3,3	0.66	0	2,2,2	0.11	0
4	EDO	C	409	-	3,3,3	0.47	0	2,2,2	0.76	0
4	EDO	C	410	-	3,3,3	0.68	0	2,2,2	0.33	0
4	EDO	C	411	-	3,3,3	0.49	0	2,2,2	0.57	0
3	5UP	D	403	2	13,16,16	0.53	0	15,22,22	0.59	0
4	EDO	D	404	-	3,3,3	0.41	0	2,2,2	0.59	0
4	EDO	D	405	-	3,3,3	0.64	0	2,2,2	0.18	0
4	EDO	D	406	-	3,3,3	0.81	0	2,2,2	0.21	0
4	EDO	D	407	-	3,3,3	0.74	0	2,2,2	0.27	0
4	EDO	D	408	-	3,3,3	0.61	0	2,2,2	0.21	0
4	EDO	D	409	-	3,3,3	0.59	0	2,2,2	0.36	0
4	EDO	D	410	-	3,3,3	0.42	0	2,2,2	0.65	0
4	EDO	D	411	-	3,3,3	0.54	0	2,2,2	0.38	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	5UP	A	403	2	-	0/6/8/8	0/2/2/2
4	EDO	A	404	-	-	0/1/1/1	0/0/0/0
4	EDO	A	405	-	-	0/1/1/1	0/0/0/0
4	EDO	A	406	-	-	0/1/1/1	0/0/0/0
5	DMS	A	407	-	-	0/0/0/0	0/0/0/0
3	5UP	B	403	2	-	0/6/8/8	0/2/2/2
4	EDO	B	404	-	-	0/1/1/1	0/0/0/0
4	EDO	B	405	-	-	0/1/1/1	0/0/0/0
4	EDO	B	406	-	-	0/1/1/1	0/0/0/0
4	EDO	B	407	-	-	0/1/1/1	0/0/0/0
4	EDO	B	408	-	-	0/1/1/1	0/0/0/0
4	EDO	B	409	-	-	0/1/1/1	0/0/0/0
4	EDO	B	410	-	-	0/1/1/1	0/0/0/0
4	EDO	B	411	-	-	0/1/1/1	0/0/0/0
4	EDO	B	412	-	-	0/1/1/1	0/0/0/0
4	EDO	B	413	-	-	0/1/1/1	0/0/0/0
3	5UP	C	403	2	-	0/6/8/8	0/2/2/2
4	EDO	C	404	-	-	0/1/1/1	0/0/0/0
4	EDO	C	405	-	-	0/1/1/1	0/0/0/0
4	EDO	C	406	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	C	407	-	-	0/1/1/1	0/0/0/0
4	EDO	C	408	-	-	0/1/1/1	0/0/0/0
4	EDO	C	409	-	-	0/1/1/1	0/0/0/0
4	EDO	C	410	-	-	0/1/1/1	0/0/0/0
4	EDO	C	411	-	-	0/1/1/1	0/0/0/0
3	5UP	D	403	2	-	0/6/8/8	0/2/2/2
4	EDO	D	404	-	-	0/1/1/1	0/0/0/0
4	EDO	D	405	-	-	0/1/1/1	0/0/0/0
4	EDO	D	406	-	-	0/1/1/1	0/0/0/0
4	EDO	D	407	-	-	0/1/1/1	0/0/0/0
4	EDO	D	408	-	-	0/1/1/1	0/0/0/0
4	EDO	D	409	-	-	0/1/1/1	0/0/0/0
4	EDO	D	410	-	-	0/1/1/1	0/0/0/0
4	EDO	D	411	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	405	EDO	1	0
4	B	404	EDO	1	0
4	B	410	EDO	1	0
4	B	411	EDO	1	0
4	B	413	EDO	3	0
4	C	405	EDO	1	0
4	C	408	EDO	1	0
4	C	409	EDO	1	0
4	D	409	EDO	1	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 [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	341/360 (94%)	-0.17	8 (2%) 64 57	32, 56, 93, 122	0
1	B	350/360 (97%)	-0.28	6 (1%) 73 68	26, 52, 90, 111	0
1	C	350/360 (97%)	-0.24	5 (1%) 78 74	29, 51, 94, 113	0
1	D	350/360 (97%)	-0.34	3 (0%) 85 83	26, 46, 76, 98	0
All	All	1391/1440 (96%)	-0.26	22 (1%) 74 69	26, 51, 90, 122	0

The worst 5 of 22 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	338	ASN	5.8
1	D	5	SER	3.9
1	C	311	ASP	3.7
1	B	328	ARG	3.5
1	A	339	THR	3.0

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 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(Å ²)	Q<0.9
4	EDO	B	413	4/4	0.90	0.37	10.62	51,55,55,55	0
4	EDO	A	404	4/4	0.91	0.36	7.85	59,60,60,61	0
4	EDO	D	404	4/4	0.85	0.24	6.65	65,67,72,74	0
4	EDO	D	407	4/4	0.94	0.27	6.43	44,46,46,47	0
4	EDO	B	407	4/4	0.89	0.33	4.34	78,79,80,80	0
4	EDO	C	408	4/4	0.88	0.35	3.88	55,59,61,63	0
4	EDO	B	405	4/4	0.89	0.22	3.55	61,62,65,67	0
4	EDO	B	408	4/4	0.95	0.23	2.95	64,64,66,69	0
4	EDO	B	412	4/4	0.76	0.26	2.78	56,57,57,58	0
3	5UP	A	403	15/15	0.96	0.20	2.46	44,49,55,55	15
4	EDO	D	405	4/4	0.95	0.16	1.92	43,43,44,46	0
4	EDO	B	411	4/4	0.97	0.19	1.69	47,48,49,50	0
3	5UP	B	403	15/15	0.96	0.15	1.68	39,43,52,53	15
5	DMS	A	407	4/4	0.91	0.18	1.57	77,78,79,79	0
3	5UP	D	403	15/15	0.95	0.16	1.19	26,32,46,47	15
4	EDO	C	410	4/4	0.93	0.18	0.93	49,49,52,52	0
4	EDO	D	410	4/4	0.95	0.19	0.27	42,43,47,53	0
4	EDO	C	411	4/4	0.92	0.14	0.07	60,61,61,61	0
4	EDO	A	405	4/4	0.93	0.16	-0.00	41,41,45,48	0
4	EDO	B	409	4/4	0.90	0.16	-0.07	58,59,63,65	0
4	EDO	B	404	4/4	0.96	0.15	-0.13	26,30,35,39	0
4	EDO	D	408	4/4	0.93	0.14	-0.23	58,60,63,64	0
4	EDO	C	407	4/4	0.91	0.16	-0.37	66,67,69,72	0
4	EDO	C	409	4/4	0.96	0.12	-0.69	46,47,48,48	0
4	EDO	A	406	4/4	0.94	0.14	-0.87	48,52,57,61	0
2	ZN	C	402	1/1	0.99	0.10	-0.96	44,44,44,44	1
3	5UP	C	403	15/15	0.97	0.11	-1.51	30,38,46,46	15
4	EDO	C	405	4/4	0.87	0.10	-1.76	74,74,75,75	0
2	ZN	D	402	1/1	1.00	0.10	-1.77	37,37,37,37	1
2	ZN	B	402	1/1	0.97	0.05	-2.77	42,42,42,42	1
4	EDO	D	411	4/4	0.92	0.39	-	68,68,68,69	0
4	EDO	C	404	4/4	0.74	0.18	-	62,63,64,64	0
2	ZN	B	401	1/1	1.00	0.10	-	28,28,28,28	1
2	ZN	C	401	1/1	0.99	0.12	-	50,50,50,50	1
4	EDO	D	409	4/4	0.85	0.19	-	73,73,73,74	0
2	ZN	D	401	1/1	0.99	0.12	-	46,46,46,46	1
4	EDO	B	410	4/4	0.93	0.17	-	60,61,61,63	0
4	EDO	D	406	4/4	0.86	0.20	-	51,53,55,55	0
2	ZN	A	402	1/1	0.88	0.10	-	114,114,114,114	1
2	ZN	A	401	1/1	0.99	0.14	-	50,50,50,50	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	EDO	B	406	4/4	0.93	0.28	-	50,50,52,52	0
4	EDO	C	406	4/4	0.88	0.15	-	59,60,63,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.