



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 05:48 AM UTC

PDB ID : 6H10 / pdb_00006h10
Title : Crystal Structure of KDM4D with tetrazolyhydrazide ligand NR073
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Deposited on : 2018-07-10
Resolution : 1.10 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

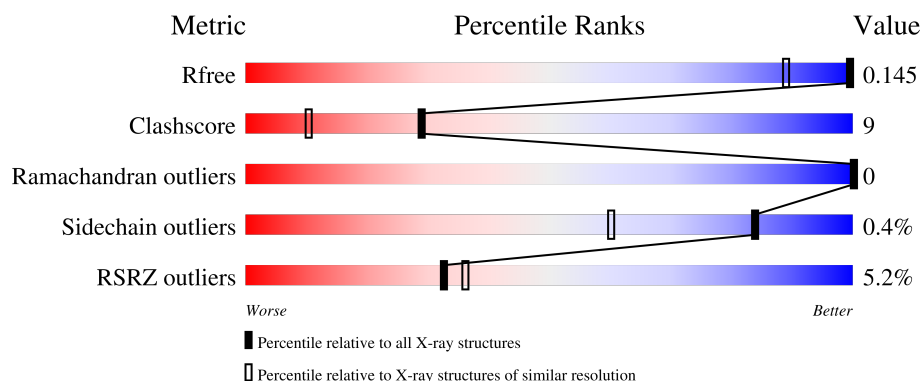
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.10 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	1625 (1.12-1.08)
Clashscore	190562	1656 (1.12-1.08)
Ramachandran outliers	187476	1616 (1.12-1.08)
Sidechain outliers	187428	1612 (1.12-1.08)
RSRZ outliers	180081	1625 (1.12-1.08)

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	342	5% 87% 10%

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 3662 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 4D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	59	0
			3060	1981	523	538	18			

- 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		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	1
			8	4	4		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	1
			4	2	2		
5	A	1	Total	C	O	0	1
			8	4	4		
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

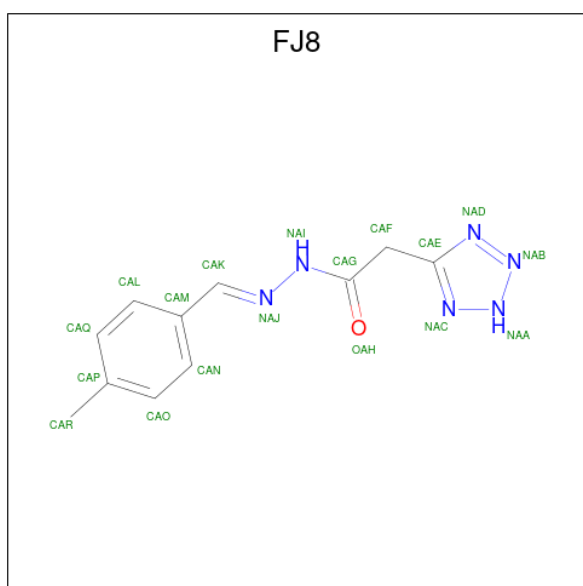


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	1
			5	4	1		

- Molecule 7 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

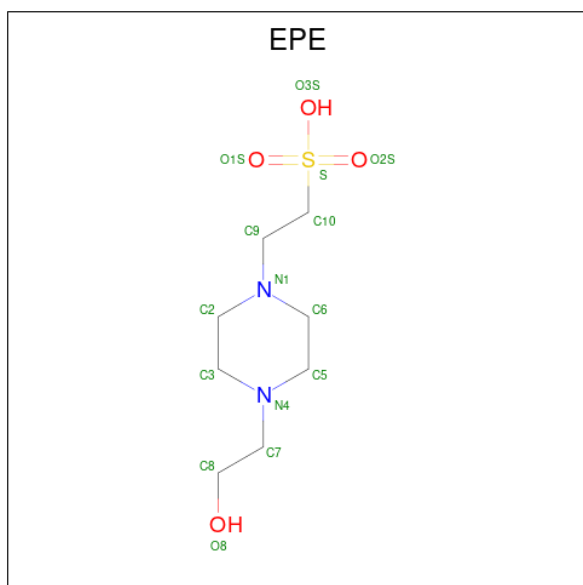
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Ni	0	0
			1	1		

- Molecule 8 is {N}-[({E})-(4-methylphenyl)methylideneamino]-2-(2 {H}-1,2,3,4-tetrazol-5-yl)ethanamide (CCD ID: FJ8) (formula: C₁₁H₁₂N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	1
			10	3	6	1		
8	A	1	Total	C	N	O	0	1
			18	11	6	1		

- Molecule 9 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	N	O	S	0	1
			15	8	2	4	1		

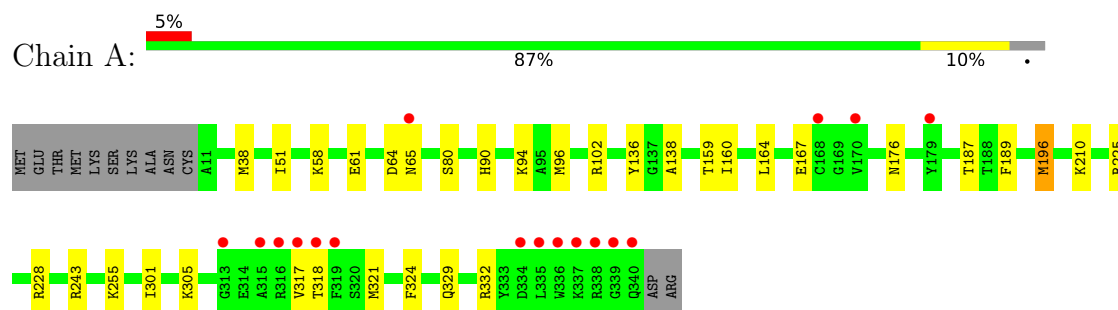
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	483	Total	O	0	62
			506	506		

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: Lysine-specific demethylase 4D



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	71.45Å 71.45Å 150.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.73 – 1.10 35.73 – 1.10	Depositor EDS
% Data completeness (in resolution range)	97.5 (35.73-1.10) 97.5 (35.73-1.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 1.10Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.121 , 0.133 0.133 , 0.145	Depositor DCC
R_{free} test set	1603 reflections (1.05%)	wwPDB-VP
Wilson B-factor (Å ²)	10.9	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.4	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	0.98	EDS
Total number of atoms	3662	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.87% 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: NA, FJ8, NI, EDO, EPE, SO4, ZN, CL

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	0.52	0/3262	0.71	1/4407 (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	A	196	MET	CG-SD-CE	7.85	118.17	100.90

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	3060	0	3057	52	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	44	0	66	3	0
6	A	5	0	0	0	0
7	A	1	0	0	0	0
8	A	28	0	0	0	0
9	A	15	0	17	0	0
10	A	506	0	0	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3662	0	3140	55	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 9.

All (55) 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:136[B]:TYR:HB3	1:A:187[B]:THR:HG21	1.68	0.74
1:A:225[A]:ARG:HG2	1:A:228[A]:ARG:HH22	1.56	0.71
1:A:136[A]:TYR:CD1	1:A:187[A]:THR:HG21	2.25	0.71
1:A:160[B]:ILE:HG12	10:A:520[B]:HOH:O	1.94	0.67
1:A:164[B]:LEU:CD1	1:A:324[B]:PHE:CE2	2.78	0.66
1:A:136[A]:TYR:CE1	1:A:187[A]:THR:HG21	2.31	0.66
1:A:196[MET:HG2	1:A:321[B]:MET:SD	2.35	0.66
1:A:136[B]:TYR:CE1	1:A:138[B]:ALA:HB2	2.32	0.65
1:A:136[B]:TYR:HB3	1:A:187[B]:THR:CG2	2.25	0.65
1:A:317:VAL:HG11	1:A:321[B]:MET:HG3	1.78	0.64
1:A:225[A]:ARG:HG2	1:A:228[A]:ARG:NH2	2.13	0.64
1:A:255[A]:LYS:HD3	10:A:522[A]:HOH:O	1.98	0.62
1:A:176[A]:ASN:HB3	10:A:582:HOH:O	2.00	0.61
1:A:90[B]:HIS:HE1	10:A:523:HOH:O	1.84	0.60
1:A:164[B]:LEU:HD12	1:A:324[B]:PHE:CE2	2.37	0.60
1:A:164[B]:LEU:HD11	1:A:324[B]:PHE:CE2	2.35	0.59
1:A:305[A]:LYS:HD2	10:A:619:HOH:O	2.02	0.59
1:A:160[B]:ILE:HD11	1:A:176[B]:ASN:OD1	2.03	0.58
1:A:321[B]:MET:HE2	1:A:321[B]:MET:HA	1.84	0.58
1:A:136[A]:TYR:CD2	1:A:187[A]:THR:CG2	2.86	0.57
1:A:164[B]:LEU:HD11	1:A:324[B]:PHE:CD2	2.39	0.57
1:A:136[A]:TYR:CD2	1:A:187[A]:THR:HG23	2.40	0.57
1:A:225[B]:ARG:NH2	10:A:507:HOH:O	2.32	0.57
1:A:136[A]:TYR:CG	1:A:187[A]:THR:CG2	2.89	0.55
1:A:317:VAL:HG22	1:A:318:THR:H	1.70	0.55
1:A:176[B]:ASN:HB3	10:A:559:HOH:O	2.08	0.54
1:A:94[B]:LYS:HE2	10:A:578:HOH:O	2.08	0.54
1:A:189[B]:PHE:CE2	1:A:210[B]:LYS:HE3	2.43	0.54
1:A:136[A]:TYR:CG	1:A:187[A]:THR:HG23	2.44	0.53
1:A:64[B]:ASP:OD1	1:A:65:ASN:ND2	2.43	0.52
1:A:317:VAL:HG11	1:A:321[A]:MET:SD	2.49	0.52
1:A:176[B]:ASN:O	10:A:501[B]:HOH:O	2.18	0.49
5:A:408:EDO:H22	10:A:694:HOH:O	2.10	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102[B]:ARG:NH1	10:A:508:HOH:O	2.36	0.49
1:A:94[B]:LYS:HG2	10:A:578:HOH:O	2.13	0.48
1:A:80[B]:SER:HB2	10:A:589:HOH:O	2.13	0.47
1:A:317:VAL:CG1	1:A:321[B]:MET:HG3	2.44	0.47
1:A:136[A]:TYR:CG	1:A:187[A]:THR:HG21	2.50	0.47
1:A:329[B]:GLN:OE1	1:A:332:ARG:HD2	2.15	0.47
1:A:196:MET:HG2	1:A:321[B]:MET:HE1	1.96	0.46
5:A:406:EDO:O1	5:A:408:EDO:H11	2.16	0.46
1:A:90[B]:HIS:CE1	10:A:523:HOH:O	2.66	0.44
1:A:225[A]:ARG:NH2	10:A:503:HOH:O	2.26	0.44
1:A:136[A]:TYR:CE2	1:A:187[A]:THR:CG2	3.01	0.44
5:A:410[A]:EDO:H22	10:A:882:HOH:O	2.18	0.43
1:A:58[A]:LYS:NZ	1:A:61[A]:GLU:OE1	2.43	0.43
1:A:159:THR:HG22	10:A:501[B]:HOH:O	2.19	0.43
1:A:317:VAL:HG22	1:A:318:THR:HG22	2.01	0.43
1:A:243[B]:ARG:HG3	10:A:872[B]:HOH:O	2.19	0.42
1:A:243[B]:ARG:CG	10:A:872[B]:HOH:O	2.67	0.42
1:A:317:VAL:HG22	1:A:318:THR:N	2.35	0.42
1:A:38[A]:MET:HE1	1:A:51:ILE:HD11	2.02	0.41
1:A:136[A]:TYR:HB2	10:A:524[A]:HOH:O	2.19	0.41
1:A:167:GLU:HG3	1:A:301[B]:ILE:HD11	2.03	0.40
1:A:136[B]:TYR:CZ	1:A:138[B]:ALA:HB2	2.56	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	387/342 (113%)	385 (100%)	2 (0%)	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	337/292 (115%)	335 (99%)	2 (1%)	78	49

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

Mol	Chain	Res	Type
1	A	96[A]	MET
1	A	96[B]	MET

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

Mol	Chain	Res	Type
1	A	65	ASN
1	A	262	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 19 ligands modelled in this entry, 4 are monoatomic - leaving 15 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
5	EDO	A	405	-	3,3,3	0.34	0	2,2,2	0.13	0
5	EDO	A	404[A]	-	3,3,3	0.33	0	2,2,2	0.67	0
5	EDO	A	409	-	3,3,3	0.48	0	2,2,2	0.76	0
5	EDO	A	408	-	3,3,3	0.47	0	2,2,2	0.11	0
5	EDO	A	411[A]	-	3,3,3	0.33	0	2,2,2	0.48	0
5	EDO	A	410[A]	-	3,3,3	0.47	0	2,2,2	0.24	0
5	EDO	A	411[B]	-	3,3,3	0.37	0	2,2,2	0.40	0
6	SO4	A	413[A]	-	4,4,4	0.21	0	6,6,6	0.24	0
9	EPE	A	417[A]	-	15,15,15	0.56	0	19,20,20	2.56	7 (36%)
5	EDO	A	407	-	3,3,3	0.40	0	2,2,2	0.65	0
8	FJ8	A	415[A]	7	10,10,19	1.92	5 (50%)	10,12,24	1.97	4 (40%)
8	FJ8	A	416[A]	-	19,19,19	2.02	5 (26%)	23,24,24	1.14	2 (8%)
5	EDO	A	406	-	3,3,3	0.31	0	2,2,2	0.32	0
5	EDO	A	404[B]	-	3,3,3	0.42	0	2,2,2	0.41	0
5	EDO	A	412	-	3,3,3	0.32	0	2,2,2	0.89	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
5	EDO	A	405	-	-	0/1/1/1	-
5	EDO	A	404[A]	-	-	1/1/1/1	-
5	EDO	A	409	-	-	1/1/1/1	-
5	EDO	A	408	-	-	0/1/1/1	-
5	EDO	A	411[A]	-	-	0/1/1/1	-
5	EDO	A	410[A]	-	-	0/1/1/1	-
5	EDO	A	411[B]	-	-	1/1/1/1	-
9	EPE	A	417[A]	-	-	3/9/19/19	0/1/1/1
5	EDO	A	407	-	-	0/1/1/1	-
8	FJ8	A	415[A]	7	-	2/6/6/10	0/1/1/2
8	FJ8	A	416[A]	-	-	2/10/10/10	0/2/2/2
5	EDO	A	406	-	-	0/1/1/1	-
5	EDO	A	404[B]	-	-	1/1/1/1	-
5	EDO	A	412	-	-	0/1/1/1	-

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	416[A]	FJ8	NAA-NAB	5.48	1.39	1.31
8	A	416[A]	FJ8	CAR-CAP	-3.55	1.40	1.51
8	A	415[A]	FJ8	CAE-NAD	-3.47	1.29	1.34
8	A	416[A]	FJ8	CAM-CAK	-3.04	1.40	1.47
8	A	416[A]	FJ8	CAF-CAE	2.71	1.52	1.50
8	A	416[A]	FJ8	CAK-NAJ	2.66	1.31	1.28
8	A	415[A]	FJ8	NAJ-NAI	-2.51	1.38	1.41
8	A	415[A]	FJ8	CAF-CAE	2.48	1.52	1.50
8	A	415[A]	FJ8	NAA-NAB	-2.02	1.28	1.31
8	A	415[A]	FJ8	CAF-CAG	2.01	1.54	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	417[A]	EPE	O2S-S-C10	-7.30	95.70	106.73
9	A	417[A]	EPE	O3S-S-C10	4.62	115.04	106.00
8	A	415[A]	FJ8	CAF-CAE-NAD	3.56	127.53	123.03
9	A	417[A]	EPE	O1S-S-C10	3.35	111.80	106.73
9	A	417[A]	EPE	C5-N4-C3	3.34	116.04	108.84
8	A	415[A]	FJ8	CAE-NAC-NAA	2.72	106.48	102.45
8	A	415[A]	FJ8	NAC-CAE-NAD	-2.57	108.86	111.96
9	A	417[A]	EPE	C9-N1-C6	-2.53	104.50	111.24
8	A	416[A]	FJ8	CAE-NAC-NAA	2.36	105.94	102.45
8	A	415[A]	FJ8	CAG-NAI-NAJ	-2.23	117.62	122.12
8	A	416[A]	FJ8	NAC-NAA-NAB	-2.13	109.07	113.52
9	A	417[A]	EPE	C7-N4-C3	2.08	116.79	111.24
9	A	417[A]	EPE	C6-C5-N4	2.00	114.69	110.65

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	415[A]	FJ8	NAC-CAE-CAF-CAG
8	A	415[A]	FJ8	NAD-CAE-CAF-CAG
8	A	416[A]	FJ8	NAC-CAE-CAF-CAG
9	A	417[A]	EPE	C9-C10-S-O3S
9	A	417[A]	EPE	C8-C7-N4-C3
5	A	404[A]	EDO	O1-C1-C2-O2
9	A	417[A]	EPE	C9-C10-S-O2S
5	A	404[B]	EDO	O1-C1-C2-O2
5	A	411[B]	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
8	A	416[A]	FJ8	NAD-CAE-CAF-CAG
5	A	409	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	408	EDO	2	0
5	A	410[A]	EDO	1	0
5	A	406	EDO	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 [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	330/342 (96%)	0.05	17 (5%) 33 36	5, 11, 24, 64	59 (17%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319[A]	PHE	7.9
1	A	315	ALA	6.5
1	A	318	THR	6.1
1	A	317	VAL	5.8
1	A	316	ARG	3.9
1	A	65	ASN	3.6
1	A	339	GLY	3.3
1	A	168[A]	CYS	3.3
1	A	170	VAL	2.5
1	A	334	ASP	2.4
1	A	336	TRP	2.4
1	A	335	LEU	2.3
1	A	179	TYR	2.2
1	A	313	GLY	2.2
1	A	338	ARG	2.2
1	A	340	GLN	2.2
1	A	337	LYS	2.2

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

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
5	EDO	A	404[A]	4/4	0.74	0.25	35,35,36,36	4
5	EDO	A	404[B]	4/4	0.74	0.25	35,35,35,36	4
5	EDO	A	409	4/4	0.80	0.17	31,31,33,35	0
5	EDO	A	408	4/4	0.87	0.14	38,39,39,40	0
9	EPE	A	417[A]	15/15	0.88	0.16	13,17,21,23	15
5	EDO	A	407	4/4	0.89	0.11	31,32,33,33	0
5	EDO	A	410[A]	4/4	0.90	0.12	40,40,40,41	0
8	FJ8	A	416[A]	18/18	0.91	0.11	13,17,21,24	18
5	EDO	A	412	4/4	0.91	0.14	31,31,33,34	4
5	EDO	A	405	4/4	0.92	0.11	17,20,22,22	0
5	EDO	A	411[A]	4/4	0.92	0.15	35,35,35,36	4
5	EDO	A	411[B]	4/4	0.92	0.15	35,35,35,35	4
5	EDO	A	406	4/4	0.93	0.11	31,33,34,35	0
6	SO4	A	413[A]	5/5	0.95	0.09	18,20,20,21	5
8	FJ8	A	415[A]	10/18	0.97	0.06	10,12,14,14	10
3	CL	A	402	1/1	0.98	0.06	33,33,33,33	0
4	NA	A	403	1/1	0.99	0.10	12,12,12,12	1
7	NI	A	414	1/1	1.00	0.03	9,9,9,9	1
2	ZN	A	401	1/1	1.00	0.01	17,17,17,17	0

6.5 Other polymers

There are no such residues in this entry.