



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 5FV3 / pdb_00005fv3
Title : Crystal structure of human JARID1B construct c2 in complex with N- Oxalylglycine.
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Deposited on : 2016-02-02
Resolution : 2.37 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	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

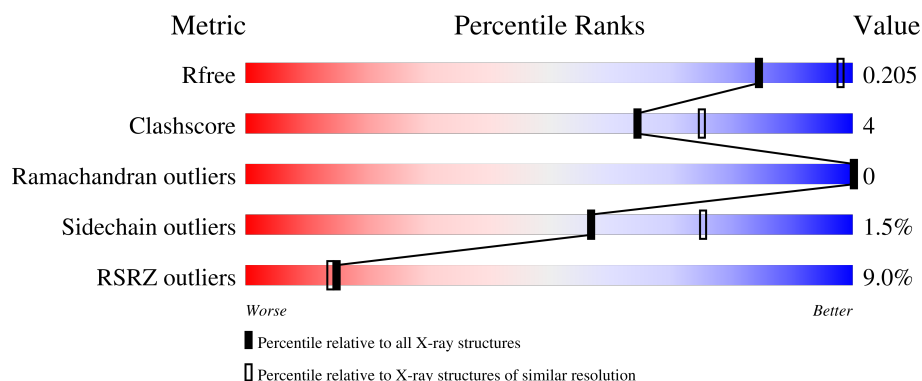
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.37 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	484	

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
7	CL	A	1769	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 3911 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 5B, LYSINE-SPECIFIC DEMETHYLASE 5B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total	C	N	O	S	0	8	0
			3718	2397	617	664	40			

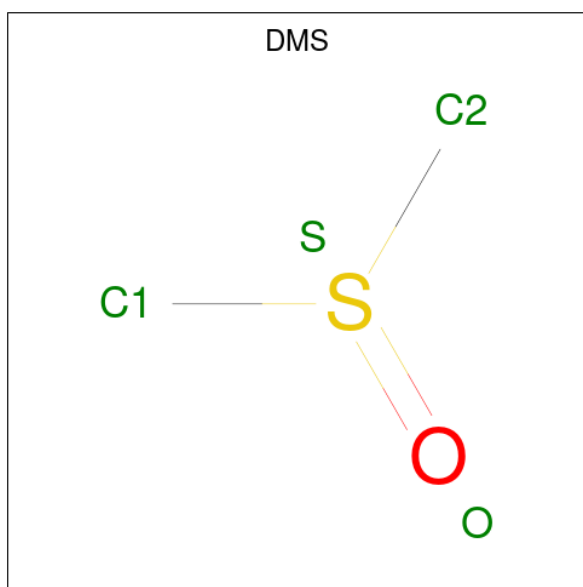
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP Q9UGL1
A	0	MET	-	expression tag	UNP Q9UGL1
A	102	GLY	-	linker	UNP Q9UGL1
A	103	GLY	-	linker	UNP Q9UGL1
A	104	GLY	-	linker	UNP Q9UGL1
A	105	GLY	-	linker	UNP Q9UGL1

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

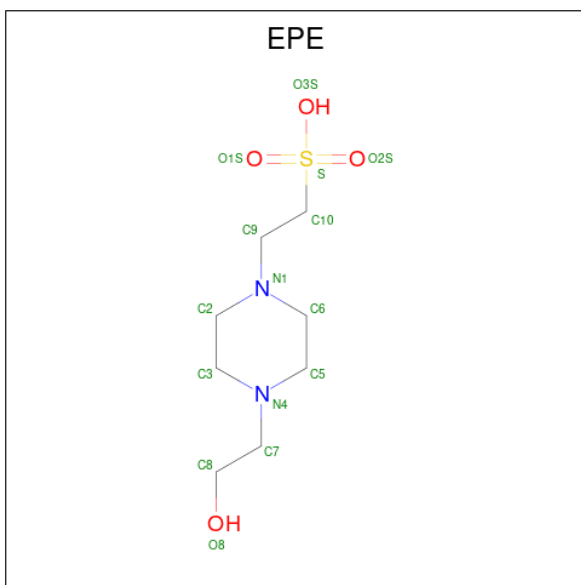


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	S	0	0
			4	2	1	1		
3	A	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

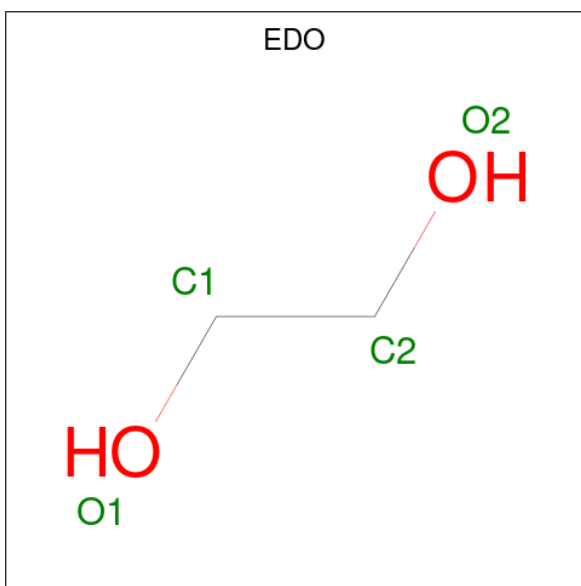
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	0
			2	2		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: C₈H₁₈N₂O₄S).



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

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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		

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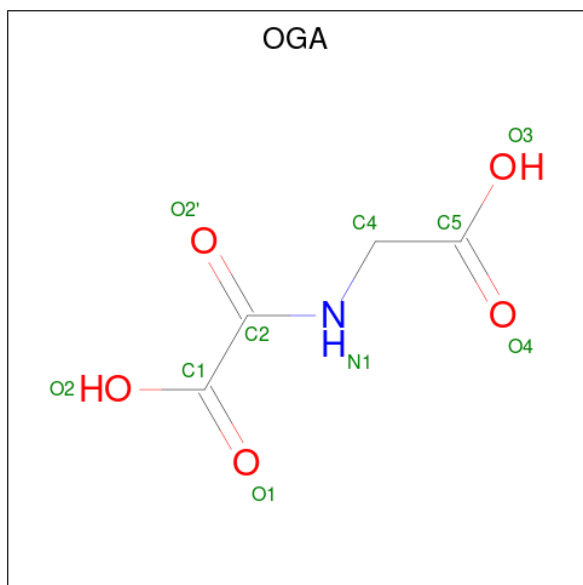
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 8 is N-OXALYLGLYCINE (CCD ID: OGA) (formula: C₄H₅NO₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			10	4	1	5		

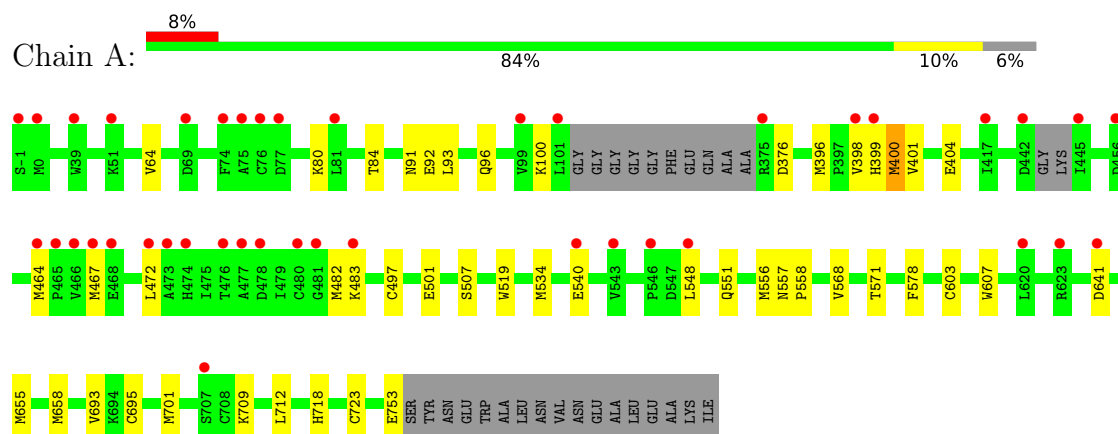
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	120	Total 120	O 120	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: LYSINE-SPECIFIC DEMETHYLASE 5B, LYSINE-SPECIFIC DEMETHYLASE 5B



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	142.35Å 142.35Å 152.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	123.28 – 2.37 123.28 – 2.37	Depositor EDS
% Data completeness (in resolution range)	99.9 (123.28-2.37) 100.0 (123.28-2.37)	Depositor EDS
R_{merge}	0.31	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.194 , 0.235 0.198 , 0.205	Depositor DCC
R_{free} test set	1823 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtrriage
Anisotropy	(Not available)	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 39.4	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3911	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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: ZN, OGA, MN, CL, EPE, DMS, 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	0.28	0/3828	0.69	0/5204

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	3718	0	3579	27	1
2	A	1	0	0	0	0
3	A	8	0	12	0	0
4	A	2	0	0	0	0
5	A	15	0	17	1	0
6	A	36	0	54	2	0
7	A	1	0	0	2	0
8	A	10	0	3	0	0
9	A	120	0	0	1	0
All	All	3911	0	3665	29	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 (29) 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:91:ASN:ND2	7:A:1769:CL:CL	2.33	0.98
7:A:1769:CL:CL	9:A:2020:HOH:O	2.41	0.74
1:A:548:LEU:O	1:A:551:GLN:NE2	2.23	0.70
1:A:96:GLN:NE2	1:A:497:CYS:SG	2.69	0.66
1:A:100:LYS:HE2	1:A:551:GLN:HE22	1.71	0.55
1:A:519:TRP:HB2	1:A:571:THR:HG22	1.88	0.54
1:A:92:GLU:HG2	1:A:93:LEU:HG	1.89	0.54
1:A:568:VAL:H	6:A:1763:EDO:H11	1.74	0.53
1:A:701[A]:MET:HB2	6:A:1760:EDO:H21	1.93	0.50
1:A:695:CYS:HB3	1:A:718:HIS:CE1	2.48	0.49
1:A:709:LYS:HG3	1:A:723:CYS:HB3	1.95	0.48
1:A:603:CYS:HB3	1:A:607:TRP:CE3	2.49	0.47
1:A:709:LYS:HB3	1:A:712:LEU:HD12	1.95	0.47
1:A:396:MET:HE3	1:A:401:VAL:HG22	1.96	0.47
5:A:1759:EPE:H81	5:A:1759:EPE:H52	1.77	0.46
1:A:472:LEU:HD11	1:A:482:MET:HE2	1.96	0.45
1:A:398:VAL:HG23	1:A:399:HIS:ND1	2.31	0.45
1:A:558:PRO:HB3	1:A:568:VAL:HG11	1.99	0.44
1:A:641:ASP:N	1:A:641:ASP:OD1	2.49	0.44
1:A:84:THR:OG1	1:A:376:ASP:OD1	2.33	0.43
1:A:534:MET:HA	1:A:556:MET:HE1	1.99	0.43
1:A:64:VAL:HB	1:A:578:PHE:HB2	1.99	0.43
1:A:501:GLU:HG3	1:A:507[A]:SER:HB3	2.02	0.42
1:A:464:MET:HE2	1:A:464:MET:HB2	1.94	0.42
1:A:400:MET:HE3	1:A:400:MET:HB2	1.86	0.41
1:A:557:ASN:HA	1:A:558:PRO:HD3	1.95	0.41
1:A:534:MET:HE3	1:A:534:MET:HB2	1.88	0.41
1:A:655:MET:HG3	1:A:658:MET:HE2	2.02	0.41
1:A:467:MET:HE2	1:A:467:MET:HB3	1.91	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:80:LYS:NZ	1:A:723:CYS:O[3_665]	2.18	0.02

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	457/484 (94%)	446 (98%)	11 (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	407/430 (95%)	401 (98%)	6 (2%)	57	75

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	400	MET
1	A	404	GLU
1	A	483	LYS
1	A	540	GLU
1	A	693	VAL
1	A	753	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	462	ASN

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Mol	Chain	Res	Type
1	A	463	ASN
1	A	551	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 17 ligands modelled in this entry, 4 are monoatomic - leaving 13 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	A	1762	-	3,3,3	0.46	0	2,2,2	0.35	0
6	EDO	A	1764	-	3,3,3	0.46	0	2,2,2	0.32	0
6	EDO	A	1763	-	3,3,3	0.46	0	2,2,2	0.27	0
3	DMS	A	1755	-	3,3,3	0.67	0	3,3,3	0.47	0
6	EDO	A	1766	-	3,3,3	0.45	0	2,2,2	0.35	0
6	EDO	A	1765	-	3,3,3	0.46	0	2,2,2	0.33	0
6	EDO	A	1768	-	3,3,3	0.46	0	2,2,2	0.37	0
5	EPE	A	1759	-	15,15,15	0.82	1 (6%)	18,20,20	1.71	5 (27%)
8	OGA	A	1770	4	9,9,9	1.31	2 (22%)	10,11,11	1.34	2 (20%)
6	EDO	A	1760	-	3,3,3	0.46	0	2,2,2	0.34	0
6	EDO	A	1767	-	3,3,3	0.46	0	2,2,2	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	A	1761	-	3,3,3	0.45	0	2,2,2	0.36	0
3	DMS	A	1756	-	3,3,3	0.68	0	3,3,3	0.51	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	A	1762	-	-	0/1/1/1	-
6	EDO	A	1764	-	-	0/1/1/1	-
6	EDO	A	1763	-	-	0/1/1/1	-
6	EDO	A	1766	-	-	0/1/1/1	-
6	EDO	A	1765	-	-	0/1/1/1	-
6	EDO	A	1768	-	-	1/1/1/1	-
5	EPE	A	1759	-	-	4/9/19/19	0/1/1/1
8	OGA	A	1770	4	-	0/8/9/9	-
6	EDO	A	1760	-	-	0/1/1/1	-
6	EDO	A	1767	-	-	0/1/1/1	-
6	EDO	A	1761	-	-	0/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	1770	OGA	O4-C5	2.78	1.31	1.22
5	A	1759	EPE	C10-S	2.69	1.81	1.77
8	A	1770	OGA	O3-C5	-2.59	1.22	1.30

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1759	EPE	C5-N4-C3	3.82	117.44	108.83
8	A	1770	OGA	O3-C5-C4	3.23	123.97	112.74
5	A	1759	EPE	C7-N4-C3	2.89	118.62	111.23
5	A	1759	EPE	C7-N4-C5	2.83	118.48	111.23
8	A	1770	OGA	O4-C5-C4	-2.59	112.20	122.68
5	A	1759	EPE	O1S-S-C10	2.51	109.94	106.92
5	A	1759	EPE	O3S-S-C10	2.14	109.22	105.77

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1759	EPE	C8-C7-N4-C5
5	A	1759	EPE	C9-C10-S-O1S
5	A	1759	EPE	C9-C10-S-O2S
6	A	1768	EDO	O1-C1-C2-O2
5	A	1759	EPE	C9-C10-S-O3S

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1763	EDO	1	0
5	A	1759	EPE	1	0
6	A	1760	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

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	455/484 (94%)	0.53	41 (9%) 15 14	19, 52, 84, 117	8 (1%)

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	474[A]	HIS	6.9
1	A	442	ASP	5.2
1	A	-1	SER	5.1
1	A	468	GLU	4.9
1	A	399	HIS	4.9
1	A	445	ILE	4.2
1	A	478	ASP	4.0
1	A	375	ARG	3.9
1	A	398	VAL	3.6
1	A	69	ASP	3.5
1	A	39	TRP	3.5
1	A	465	PRO	3.4
1	A	74	PHE	3.3
1	A	101	LEU	3.3
1	A	76	CYS	3.3
1	A	540	GLU	3.0
1	A	546	PRO	3.0
1	A	466	VAL	2.9
1	A	417	ILE	2.9
1	A	0	MET	2.9
1	A	473	ALA	2.8
1	A	543	VAL	2.8
1	A	548	LEU	2.8
1	A	620[A]	LEU	2.7
1	A	467	MET	2.7
1	A	75	ALA	2.5
1	A	456	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	480	CYS	2.5
1	A	464	MET	2.4
1	A	99	VAL	2.3
1	A	476	THR	2.2
1	A	707	SER	2.2
1	A	77	ASP	2.2
1	A	641	ASP	2.2
1	A	477	ALA	2.2
1	A	481	GLY	2.1
1	A	51	LYS	2.1
1	A	623	ARG	2.1
1	A	81	LEU	2.0
1	A	472	LEU	2.0
1	A	483	LYS	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($\text{\AA}^2$)	Q<0.9
3	DMS	A	1755	4/4	0.77	0.38	66,74,95,140	0
6	EDO	A	1767	4/4	0.83	0.23	69,79,81,87	0
3	DMS	A	1756	4/4	0.87	0.22	53,58,59,82	0
6	EDO	A	1765	4/4	0.88	0.19	65,70,75,86	0
6	EDO	A	1768	4/4	0.88	0.18	54,62,66,68	0
6	EDO	A	1763	4/4	0.90	0.18	62,64,67,76	0
5	EPE	A	1759	15/15	0.90	0.20	44,72,85,86	15
6	EDO	A	1761	4/4	0.91	0.21	53,63,76,80	0
6	EDO	A	1762	4/4	0.91	0.22	60,82,84,88	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MN	A	1758	1/1	0.91	0.13	103,103,103,103	0
6	EDO	A	1764	4/4	0.93	0.16	56,62,62,71	0
6	EDO	A	1760	4/4	0.95	0.15	40,44,58,67	0
7	CL	A	1769	1/1	0.95	0.10	77,77,77,77	0
6	EDO	A	1766	4/4	0.96	0.17	51,63,63,68	0
8	OGA	A	1770	10/10	0.96	0.09	36,43,48,50	0
4	MN	A	1757	1/1	1.00	0.04	30,30,30,30	0
2	ZN	A	1754	1/1	1.00	0.02	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.