



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:24 PM GMT

PDB ID : 5A6T
Title : 1.65 Å resolution Sulfite inhibited *Sporosarcina pasteurii* urease
Authors : Mazzei, L.; Cianci, M.; Benini, S.; Bertini, L.; Musiani, F.; Ciurli, S.
Deposited on : 2015-07-01
Resolution : 1.65 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
<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-20026982
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20026982

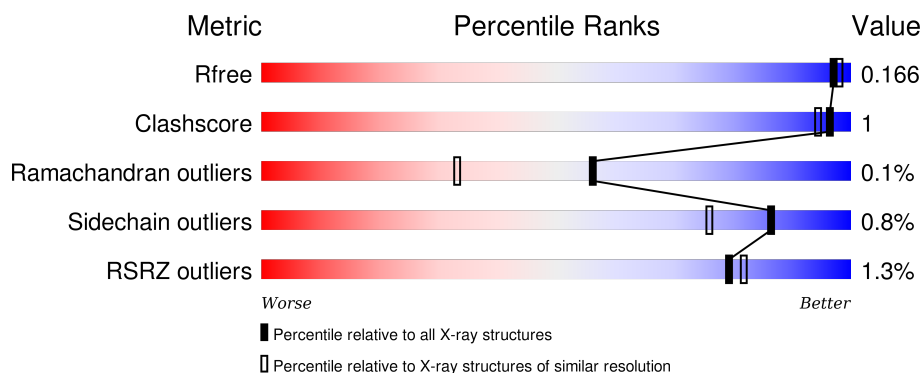
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 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}	91344	1226 (1.66-1.66)
Clashscore	102246	1323 (1.66-1.66)
Ramachandran outliers	100387	1295 (1.66-1.66)
Sidechain outliers	100360	1295 (1.66-1.66)
RSRZ outliers	91569	1227 (1.66-1.66)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	100	<div> <div style="width: 98%;"></div> <div>98%</div> </div>
2	B	126	<div> <div style="width: 89%;"></div> <div>89%</div> <div style="width: 8%;"></div> <div>8%</div> </div>
3	C	570	<div> <div style="width: 92%;"></div> <div>92%</div> <div style="width: 7%;"></div> <div>7%</div> </div>

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
4	EDO	A	1102	-	-	-	X
4	EDO	B	1127	-	-	-	X
4	EDO	C	1574	-	-	-	X
4	EDO	C	1575	-	-	-	X
4	EDO	C	1577	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6863 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 UREASE SUBUNIT GAMMA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	100	Total	C	N	O	S	0	2	0
			791	499	133	152	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	ALA	LEU	CONFLICT	UNP P41022
A	22	LYS	ARG	CONFLICT	UNP P41022

- Molecule 2 is a protein called UREASE SUBUNIT BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	122	Total	C	N	O	S	0	1	0
			960	594	172	193	1			

- Molecule 3 is a protein called UREASE SUBUNIT ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	570	Total	C	N	O	S	0	17	0
			4438	2786	762	864	26			

There are 10 discrepancies between the modelled and reference sequences:

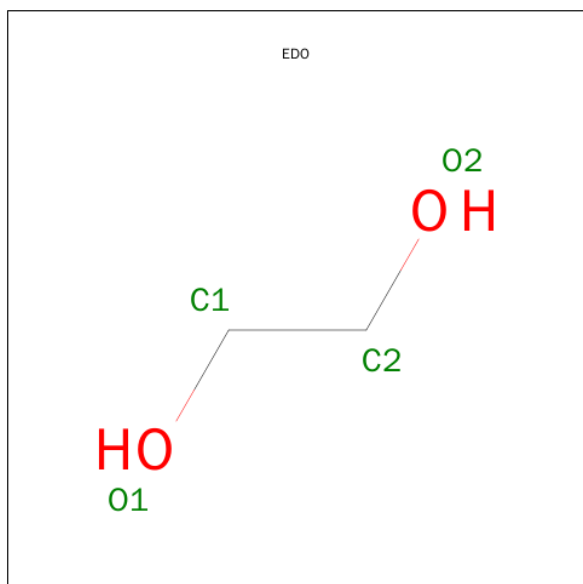
Chain	Residue	Modelled	Actual	Comment	Reference
C	19	GLN	ARG	CONFLICT	UNP P41020
C	28	TRP	GLY	CONFLICT	UNP P41020
C	29	ILE	-	INSERTION	UNP P41020
C	36	THR	TYR	CONFLICT	UNP P41020
C	37	THR	TYR	CONFLICT	UNP P41020
C	38	TYR	LEU	CONFLICT	UNP P41020
C	42	ALA	VAL	CONFLICT	UNP P41020
C	263	LEU	VAL	CONFLICT	UNP P41020

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Chain	Residue	Modelled	Actual	Comment	Reference
C	403	ALA	LEU	CONFLICT	UNP P41020
C	420	ILE	MET	CONFLICT	UNP P41020

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



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

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).

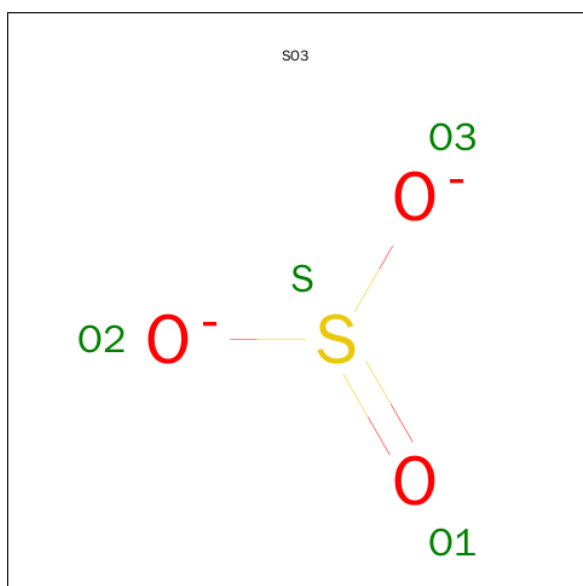


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	2	Total	Ni	0	0
			2	2		

- Molecule 7 is SULFITE ION (three-letter code: SO3) (formula: O₃S).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	97	Total	O		0	0
			97	97			
8	B	139	Total	O		0	0
			139	139			
8	C	399	Total	O		0	0
			399	399			

4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	131.26Å 131.26Å 189.07Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	113.68 – 1.65 55.12 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.7 (113.68-1.65) 99.7 (55.12-1.65)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.87 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.129 , 0.152 0.143 , 0.166	Depositor DCC
R_{free} test set	5779 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	17.2	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 114913 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6863	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NI, EDO, KCX, SO3, SO4, CXM

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	1.05	1/794 (0.1%)	1.04	1/1068 (0.1%)
2	B	1.17	4/972 (0.4%)	1.16	2/1308 (0.2%)
3	C	1.04	6/4531 (0.1%)	1.08	22/6136 (0.4%)
All	All	1.06	11/6297 (0.2%)	1.09	25/8512 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	C	0	1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	520	CYS	CB-SG	-18.73	1.50	1.82
2	B	33	SER	CB-OG	-8.03	1.31	1.42
3	C	394	GLU	CD-OE1	7.33	1.33	1.25
2	B	82	GLU	CD-OE1	-5.70	1.19	1.25
3	C	520	CYS	CA-CB	5.61	1.66	1.53
3	C	386	LYS	CD-CE	5.60	1.65	1.51
3	C	560	GLU	CD-OE2	-5.47	1.19	1.25
1	A	58	GLU	CG-CD	5.33	1.59	1.51
2	B	111	GLU	CD-OE1	5.28	1.31	1.25
3	C	386	LYS	CE-NZ	5.12	1.61	1.49
2	B	12	TYR	CG-CD2	-5.01	1.32	1.39

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	5	ARG	NE-CZ-NH2	-14.46	113.07	120.30
3	C	388	ARG	NE-CZ-NH1	9.31	124.95	120.30
3	C	5	ARG	NE-CZ-NH1	9.02	124.81	120.30
3	C	21	ARG	NE-CZ-NH1	7.96	124.28	120.30
2	B	65	ARG	NE-CZ-NH2	-7.36	116.62	120.30
3	C	1	MET	CG-SD-CE	-6.66	89.54	100.20
3	C	386	LYS	CD-CE-NZ	6.56	126.79	111.70
3	C	566	ARG	NE-CZ-NH1	6.33	123.46	120.30
3	C	363	ASP	CB-CG-OD1	6.14	123.82	118.30
3	C	341	ARG	NE-CZ-NH1	6.08	123.34	120.30
3	C	107	ASP	CB-CG-OD1	5.88	123.59	118.30
3	C	274	PHE	CB-CG-CD2	-5.86	116.70	120.80
1	A	66	ARG	NE-CZ-NH2	-5.80	117.40	120.30
3	C	251	ASP	CB-CG-OD2	-5.51	113.34	118.30
2	B	38	ARG	NE-CZ-NH2	-5.51	117.54	120.30
3	C	388	ARG	NE-CZ-NH2	-5.51	117.55	120.30
3	C	107	ASP	CB-CG-OD2	-5.44	113.40	118.30
3	C	566	ARG	NE-CZ-NH2	-5.26	117.67	120.30
3	C	526	LYS	CD-CE-NZ	5.25	123.77	111.70
3	C	449	ARG	NE-CZ-NH2	-5.15	117.72	120.30
3	C	240	ASP	CB-CG-OD1	5.14	122.93	118.30
3	C	405	ARG	NE-CZ-NH2	-5.09	117.75	120.30
3	C	234	ARG	NE-CZ-NH2	-5.08	117.76	120.30
3	C	369	ARG	NE-CZ-NH1	5.05	122.82	120.30
3	C	224	ASP	CB-CG-OD1	5.03	122.83	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	234	ARG	Sidechain

5.2 Too-close contacts ⓘ

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	791	0	811	0	0
2	B	960	0	942	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4438	0	4419	18	0
4	A	8	0	12	0	0
4	B	4	0	6	0	0
4	C	16	0	24	0	0
5	B	5	0	0	0	0
6	C	2	0	0	0	0
7	C	4	0	0	0	0
8	A	97	0	0	0	1
8	B	139	0	0	0	1
8	C	399	0	0	3	0
All	All	6863	0	6214	18	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 1.

All (18) 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)
3:C:272:HIS:HE1	3:C:299:SER:OG	1.75	0.70
3:C:507[A]:LYS:HD2	8:C:2170:HOH:O	1.91	0.69
3:C:270[B]:VAL:HG12	8:C:2359:HOH:O	2.00	0.61
3:C:270[B]:VAL:HG11	3:C:523:ILE:O	2.01	0.61
3:C:272:HIS:HD2	3:C:418[B]:GLN:NE2	1.98	0.60
3:C:293:HIS:HE1	8:C:2245:HOH:O	1.86	0.59
3:C:522[B]:ASN:OD1	3:C:522[B]:ASN:N	2.41	0.53
2:B:95:GLU:O	3:C:104:ASP:HB3	2.11	0.51
3:C:132:GLY:HA3	3:C:155:THR:OG1	2.14	0.48
3:C:272:HIS:CE1	3:C:299:SER:OG	2.63	0.47
3:C:272:HIS:HD2	3:C:418[B]:GLN:HE21	1.62	0.47
3:C:316:LEU:O	3:C:320:MET:HG2	2.16	0.46
3:C:70:LEU:HD11	3:C:86:ASP:HB3	1.98	0.46
2:B:71:SER:OG	3:C:49:VAL:HG21	2.19	0.43
3:C:507[A]:LYS:HE2	3:C:507[A]:LYS:HB3	1.80	0.43
3:C:137:HIS:CE1	3:C:274:PHE:CD2	3.08	0.42
3:C:362:THR:O	3:C:368:GLY:HA3	2.21	0.40
2:B:15:ALA:O	3:C:5:ARG:HD2	2.22	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)
8:A:2015:HOH:O	8:B:2049:HOH:O[8_676]	1.96	0.24

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	100/100 (100%)	100 (100%)	0	0	100	100
2	B	121/126 (96%)	118 (98%)	2 (2%)	1 (1%)	24	5
3	C	584/570 (102%)	561 (96%)	23 (4%)	0	100	100
All	All	805/796 (101%)	779 (97%)	25 (3%)	1 (0%)	56	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	99	ILE

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	85/83 (102%)	85 (100%)	0	100	100
2	B	102/105 (97%)	102 (100%)	0	100	100
3	C	475/458 (104%)	470 (99%)	5 (1%)	80	64
All	All	662/646 (102%)	657 (99%)	5 (1%)	86	76

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	1	MET
3	C	158	PHE
3	C	249	HIS
3	C	272	HIS
3	C	395	LYS

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

Mol	Chain	Res	Type
3	C	272	HIS
3	C	293	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
1	CXM	A	1	1	6,10,11	0.80	0	5,11,13	0.96	0
3	KCX	C	220	3,6	6,11,12	0.80	0	7,12,14	1.17	1 (14%)

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
1	CXM	A	1	1	-	0/6/10/12	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	KCX	C	220	3,6	-	0/6/10/12	0/0/0/0

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	220	KCX	CE-NZ-CX	-2.46	120.94	123.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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	$\# Z > 2$
4	EDO	A	1101	-	3,3,3	0.76	0	2,2,2	1.02	0
4	EDO	A	1102	-	3,3,3	1.53	0	2,2,2	0.58	0
4	EDO	B	1127	-	3,3,3	0.23	0	2,2,2	0.99	0
5	SO4	B	1128	-	4,4,4	0.58	0	6,6,6	1.26	1 (16%)
7	SO3	C	1573	6	1,3,3	5.88	1 (100%)	0,3,3	0.00	-
4	EDO	C	1574	-	3,3,3	0.53	0	2,2,2	0.38	0
4	EDO	C	1575	-	3,3,3	0.75	0	2,2,2	0.82	0
4	EDO	C	1576	-	3,3,3	0.53	0	2,2,2	0.26	0
4	EDO	C	1577	-	3,3,3	0.38	0	2,2,2	0.31	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
4	EDO	A	1101	-	-	0/1/1/1	0/0/0/0
4	EDO	A	1102	-	-	0/1/1/1	0/0/0/0
4	EDO	B	1127	-	-	0/1/1/1	0/0/0/0
5	SO4	B	1128	-	-	0/0/0/0	0/0/0/0
7	SO3	C	1573	6	-	0/0/0/0	0/0/0/0
4	EDO	C	1574	-	-	0/1/1/1	0/0/0/0
4	EDO	C	1575	-	-	0/1/1/1	0/0/0/0
4	EDO	C	1576	-	-	0/1/1/1	0/0/0/0
4	EDO	C	1577	-	-	0/1/1/1	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	1573	SO3	O1-S	5.88	1.67	1.44

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1128	SO4	O2-S-O1	2.35	117.44	109.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	99/100 (99%)	-0.82	0 100 100	13, 16, 23, 29	0
2	B	122/126 (96%)	-0.44	1 (0%) 87 88	14, 18, 33, 52	0
3	C	569/570 (99%)	-0.58	9 (1%) 74 78	12, 16, 28, 80	0
All	All	790/796 (99%)	-0.59	10 (1%) 79 82	12, 16, 28, 80	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	396	ASN	6.7
3	C	390	PRO	6.1
3	C	395	LYS	5.8
3	C	391	LEU	5.1
3	C	392	ALA	4.3
2	B	126	GLU	3.1
3	C	397	GLY	3.0
3	C	393	GLU	2.2
3	C	550	GLY	2.1
3	C	389	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CXM	A	1	11/12	0.99	0.05	-	14,16,19,22	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	KCX	C	220	12/13	0.97	0.05	-	11,13,15,15	0

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(\AA^2)	Q<0.9
4	EDO	C	1577	4/4	0.95	0.20	13.92	32,35,36,46	0
4	EDO	A	1102	4/4	0.92	0.10	4.09	23,24,26,28	0
4	EDO	C	1575	4/4	0.97	0.14	3.58	26,29,32,33	0
4	EDO	C	1574	4/4	0.96	0.09	3.26	27,30,30,31	0
4	EDO	B	1127	4/4	0.95	0.12	2.94	33,40,43,52	0
7	SO3	C	1573	4/4	0.97	0.10	1.87	22,23,25,28	0
4	EDO	A	1101	4/4	0.96	0.08	0.95	20,24,24,25	0
6	NI	C	1572	1/1	1.00	0.02	-2.69	16,16,16,16	0
6	NI	C	1571	1/1	1.00	0.01	-7.62	18,18,18,18	0
4	EDO	C	1576	4/4	0.89	0.23	-	41,41,41,43	0
5	SO4	B	1128	5/5	0.94	0.26	-	42,48,57,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.