



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 16, 2023 – 01:32 pm GMT

PDB ID : 4CEU
Title : 1.58 Å resolution native Sporosarcina pasteurii urease
Authors : Benini, S.; Cianci, M.; Ciurli, S.
Deposited on : 2013-11-12
Resolution : 1.58 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

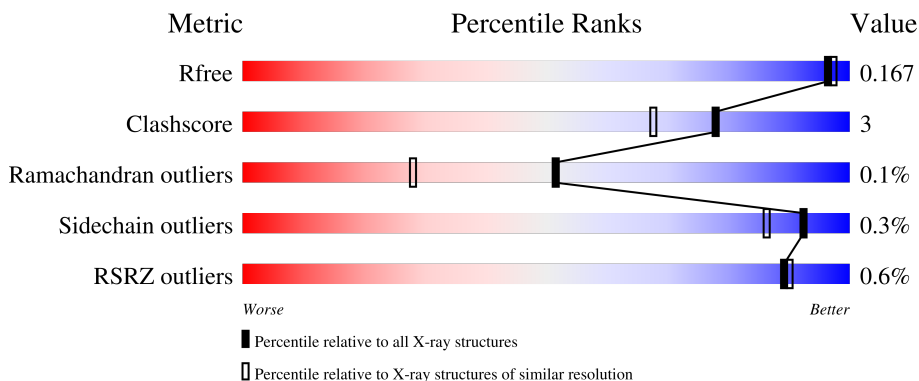
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.58 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}	130704	5534 (1.60-1.56)
Clashscore	141614	5861 (1.60-1.56)
Ramachandran outliers	138981	5708 (1.60-1.56)
Sidechain outliers	138945	5703 (1.60-1.56)
RSRZ outliers	127900	5431 (1.60-1.56)

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	100	97%
2	B	126	89% 8%
3	C	570	92% 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
4	EDO	C	1575	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7030 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
			Total	C	N	O	S			
1	A	100	790	498	134	151	7	0	2	0

- Molecule 2 is a protein called UREASE SUBUNIT BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	122	951	590	172	188	1	0	1	0

- Molecule 3 is a protein called UREASE SUBUNIT ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	570	4353	2737	745	846	25	0	9	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	19	GLU	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	263	LEU	VAL	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 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	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

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

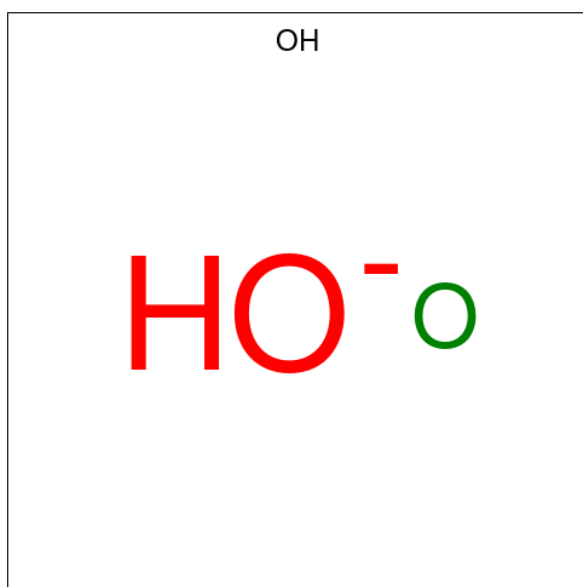


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		

- 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 HYDROXIDE ION (three-letter code: OH) (formula: HO).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	124	Total O 124 124	0	0
8	B	203	Total O 203 203	0	0
8	C	532	Total O 532 532	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.

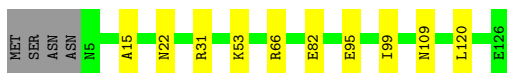
- Molecule 1: UREASE SUBUNIT GAMMA

Chain A:  97%



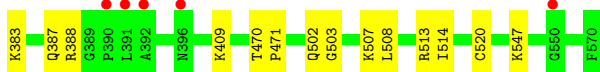
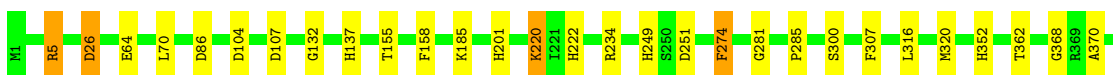
- Molecule 2: UREASE SUBUNIT BETA

Chain B:  89% 8%



- Molecule 3: UREASE SUBUNIT ALPHA

Chain C:  92% 7%



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	131.12Å 131.12Å 188.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	97.51 – 1.58 97.32 – 1.58	Depositor EDS
% Data completeness (in resolution range)	99.8 (97.51-1.58) 99.8 (97.32-1.58)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.21 (at 1.58Å)	Xtrriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.135 , 0.155 0.150 , 0.167	Depositor DCC
R_{free} test set	6570 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	12.1	Xtrriage
Anisotropy	0.557	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7030	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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: NI, SO4, CXM, OH, KCX, 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	0.58	0/796	0.77	0/1071
2	B	0.58	0/966	0.76	0/1300
3	C	0.59	1/4449 (0.0%)	0.83	11/6031 (0.2%)
All	All	0.58	1/6211 (0.0%)	0.81	11/8402 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	520	CYS	CB-SG	-9.55	1.66	1.82

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	5	ARG	NE-CZ-NH2	-9.30	115.65	120.30
3	C	388	ARG	NE-CZ-NH1	6.54	123.57	120.30
3	C	5	ARG	NE-CZ-NH1	6.32	123.46	120.30
3	C	274	PHE	CB-CG-CD2	-6.12	116.52	120.80
3	C	388	ARG	NE-CZ-NH2	-5.87	117.36	120.30
3	C	520	CYS	CB-CA-C	-5.80	98.79	110.40
3	C	274	PHE	CB-CG-CD1	5.44	124.61	120.80
3	C	26	ASP	CB-CG-OD1	5.42	123.18	118.30
3	C	107	ASP	CB-CG-OD1	5.16	122.95	118.30
3	C	251	ASP	CB-CG-OD2	-5.12	113.69	118.30
3	C	234	ARG	NE-CZ-NH1	5.10	122.85	120.30

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	790	0	809	1	0
2	B	951	0	935	7	0
3	C	4353	0	4321	30	0
4	A	12	0	18	0	0
4	B	4	0	6	0	0
4	C	28	0	42	4	1
5	A	5	0	0	0	0
5	B	5	0	0	0	0
5	C	20	0	0	0	0
6	C	2	0	0	0	0
7	C	1	0	0	0	0
8	A	124	0	0	0	0
8	B	203	0	0	2	0
8	C	532	0	0	5	2
All	All	7030	0	6131	36	2

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 3.

All (36) 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:201:HIS:HE1	3:C:222:HIS:H	1.26	0.83
3:C:64[A]:GLU:OE2	8:C:2113:HOH:O	2.02	0.76
3:C:201:HIS:CE1	3:C:222:HIS:H	2.03	0.76
3:C:300:SER:OG	3:C:352:HIS:HE1	1.79	0.65
3:C:513[A]:ARG:NE	4:C:1575:EDO:O2	2.30	0.63
3:C:370:ALA:N	8:C:2364:HOH:O	2.15	0.58
3:C:513[A]:ARG:NH1	8:C:2459:HOH:O	2.36	0.58
2:B:109:ASN:HB3	8:B:2184:HOH:O	2.06	0.54
3:C:383:LYS:NZ	3:C:387:GLN:HE22	2.07	0.51
3:C:249:HIS:CE1	3:C:281:GLY:HA3	2.46	0.50
3:C:507[A]:LYS:HD2	8:C:2225:HOH:O	2.11	0.50
2:B:22:ASN:HD21	2:B:66:ARG:HH22	1.60	0.49
3:C:513[A]:ARG:NH2	4:C:1575:EDO:O2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:132:GLY:HA3	3:C:155:THR:OG1	2.13	0.49
3:C:70:LEU:HD11	3:C:86:ASP:HB3	1.95	0.48
2:B:95:GLU:O	3:C:104:ASP:HB3	2.15	0.47
1:A:81[B]:GLN:HG2	1:A:94:THR:OG1	2.14	0.47
3:C:26:ASP:HB2	8:C:2372:HOH:O	2.15	0.47
2:B:31[B]:ARG:NH1	8:B:2082:HOH:O	2.46	0.47
3:C:502:GLN:O	3:C:507[B]:LYS:HE2	2.15	0.46
3:C:362:THR:O	3:C:368:GLY:HA3	2.15	0.46
3:C:383:LYS:HZ3	3:C:387:GLN:HE22	1.64	0.46
3:C:503:GLY:O	3:C:507[B]:LYS:HG3	2.15	0.46
2:B:31[A]:ARG:HD2	2:B:82:GLU:OE2	2.17	0.44
3:C:137:HIS:CE1	3:C:274:PHE:CD2	3.06	0.44
3:C:514:ILE:O	4:C:1575:EDO:H11	2.18	0.44
3:C:352:HIS:HD2	3:C:409:LYS:NZ	2.16	0.43
3:C:316:LEU:O	3:C:320:MET:HG2	2.19	0.43
3:C:470:THR:N	3:C:471:PRO:CD	2.82	0.42
3:C:307:PHE:O	3:C:383:LYS:HE3	2.20	0.41
3:C:547:LYS:O	4:C:1578:EDO:H21	2.20	0.41
2:B:15:ALA:O	3:C:5:ARG:HD2	2.20	0.41
3:C:220:KCX:CX	3:C:222:HIS:HD2	2.34	0.41
3:C:352:HIS:HD2	3:C:409:LYS:HZ2	1.69	0.41
3:C:185:LYS:HE3	3:C:508:LEU:HD22	2.03	0.40
2:B:53:LYS:NZ	2:B:120:LEU:O	2.49	0.40

All (2) 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:C:2364:HOH:O	8:C:2423:HOH:O[3_565]	1.72	0.48
4:C:1575:EDO:O1	8:C:2447:HOH:O[12_565]	2.04	0.16

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%)	117 (97%)	3 (2%)	1 (1%)	19	5
3	C	576/570 (101%)	556 (96%)	20 (4%)	0	100	100
All	All	797/796 (100%)	773 (97%)	23 (3%)	1 (0%)	51	28

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	85/84 (101%)	85 (100%)	0	100	100
2	B	100/105 (95%)	100 (100%)	0	100	100
3	C	462/460 (100%)	460 (100%)	2 (0%)	91	84
All	All	647/649 (100%)	645 (100%)	2 (0%)	92	86

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

Mol	Chain	Res	Type
3	C	158	PHE
3	C	285	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
2	B	22	ASN
3	C	201	HIS

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Mol	Chain	Res	Type
3	C	267	ASN
3	C	352	HIS
3	C	387	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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	8,10,11	2.21	1 (12%)	7,11,13	1.05	0
3	KCX	C	220	6,3	9,11,12	2.41	1 (11%)	5,12,14	1.47	1 (20%)

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	-	4/9/10/12	-
3	KCX	C	220	6,3	-	0/9/10/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	220	KCX	OQ1-CX	7.14	1.34	1.21
1	A	1	CXM	ON1-CN	5.96	1.32	1.21

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	220	KCX	OQ1-CX-NZ	-2.94	120.40	124.96

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	CXM	O-C-CA-CB
1	A	1	CXM	CB-CA-N-CN
1	A	1	CXM	ON1-CN-N-CA
1	A	1	CXM	C-CA-N-CN

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	220	KCX	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 2 are monoatomic and 1 is modelled with single atom - leaving 17 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	C	1574	-	3,3,3	0.45	0	2,2,2	0.44	0
5	SO4	B	1128	-	4,4,4	0.43	0	6,6,6	0.51	0
5	SO4	C	1580	-	4,4,4	0.38	0	6,6,6	0.94	0
4	EDO	C	1572	-	3,3,3	0.42	0	2,2,2	0.39	0
4	EDO	C	1573	-	3,3,3	0.53	0	2,2,2	0.50	0
4	EDO	C	1578	-	3,3,3	0.38	0	2,2,2	0.18	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	A	1101	-	3,3,3	0.56	0	2,2,2	0.74	0
5	SO4	C	1579	-	4,4,4	0.30	0	6,6,6	0.43	0
5	SO4	C	1582	-	4,4,4	0.60	0	6,6,6	0.53	0
5	SO4	A	1104	-	4,4,4	0.29	0	6,6,6	0.39	0
4	EDO	C	1575	-	3,3,3	0.50	0	2,2,2	0.08	0
4	EDO	A	1102	-	3,3,3	0.60	0	2,2,2	0.09	0
5	SO4	C	1581	-	4,4,4	0.35	0	6,6,6	0.05	0
4	EDO	C	1576	-	3,3,3	0.47	0	2,2,2	0.33	0
4	EDO	A	1103	-	3,3,3	1.07	0	2,2,2	0.54	0
4	EDO	B	1127	-	3,3,3	0.52	0	2,2,2	0.06	0
4	EDO	C	1577	-	3,3,3	0.35	0	2,2,2	0.55	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	C	1574	-	-	0/1/1/1	-
4	EDO	C	1572	-	-	0/1/1/1	-
4	EDO	C	1573	-	-	0/1/1/1	-
4	EDO	C	1578	-	-	0/1/1/1	-
4	EDO	A	1101	-	-	0/1/1/1	-
4	EDO	C	1575	-	-	0/1/1/1	-
4	EDO	A	1102	-	-	0/1/1/1	-
4	EDO	C	1576	-	-	0/1/1/1	-
4	EDO	A	1103	-	-	0/1/1/1	-
4	EDO	B	1127	-	-	0/1/1/1	-
4	EDO	C	1577	-	-	0/1/1/1	-

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.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1578	EDO	1	0
4	C	1575	EDO	3	1

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%)	-0.75	0	100 100	10, 13, 19, 27	0
2	B	122/126 (96%)	-0.72	0	100 100	11, 15, 24, 39	0
3	C	569/570 (99%)	-0.71	5 (0%)	84 85	10, 12, 24, 48	1 (0%)
All	All	790/796 (99%)	-0.72	5 (0%)	89 90	10, 13, 24, 48	1 (0%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	396	ASN	4.2
3	C	390	PRO	4.2
3	C	391	LEU	4.2
3	C	392	ALA	3.4
3	C	550	GLY	2.5

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. 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
1	CXM	A	1	11/12	0.97	0.06	12,14,16,16	0
3	KCX	C	220	12/13	0.98	0.05	9,10,11,11	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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(\AA^2)	Q<0.9
4	EDO	C	1575	4/4	0.67	0.29	33,36,36,40	0
4	EDO	C	1578	4/4	0.85	0.36	40,41,41,45	0
4	EDO	C	1577	4/4	0.86	0.21	46,48,48,49	0
4	EDO	A	1103	4/4	0.86	0.12	19,19,21,22	0
5	SO4	C	1580	5/5	0.88	0.18	19,19,28,28	5
4	EDO	C	1576	4/4	0.89	0.19	31,32,33,34	0
4	EDO	A	1102	4/4	0.89	0.12	33,36,37,39	0
4	EDO	B	1127	4/4	0.90	0.18	28,30,31,33	0
4	EDO	C	1574	4/4	0.91	0.10	21,24,25,28	0
5	SO4	B	1128	5/5	0.92	0.18	42,45,50,51	0
4	EDO	A	1101	4/4	0.92	0.14	16,21,21,26	0
5	SO4	A	1104	5/5	0.94	0.17	34,38,41,42	0
5	SO4	C	1579	5/5	0.95	0.10	24,30,34,36	5
5	SO4	C	1581	5/5	0.95	0.11	59,60,60,61	5
4	EDO	C	1573	4/4	0.96	0.18	23,25,26,27	0
4	EDO	C	1572	4/4	0.97	0.07	24,24,25,26	0
7	OH	C	1571	1/1	0.97	0.05	11,11,11,11	0
5	SO4	C	1582	5/5	0.99	0.09	25,26,28,33	0
6	NI	C	601	1/1	1.00	0.02	11,11,11,11	0
6	NI	C	600	1/1	1.00	0.02	12,12,12,12	0

6.5 Other polymers [i](#)

There are no such residues in this entry.