



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 06:02 PM UTC

PDB ID : 6RKG / pdb_00006rkg
Title : 1.32 Å RESOLUTION OF SPOROSARCINA PASTEURII UREASE INHIBITED IN THE PRESENCE OF NBPTO AT pH 7.5
Authors : Mazzei, L.; Ciani, M.; Benini, S.; Ciurli, S.
Deposited on : 2019-04-30
Resolution : 1.32 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

<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

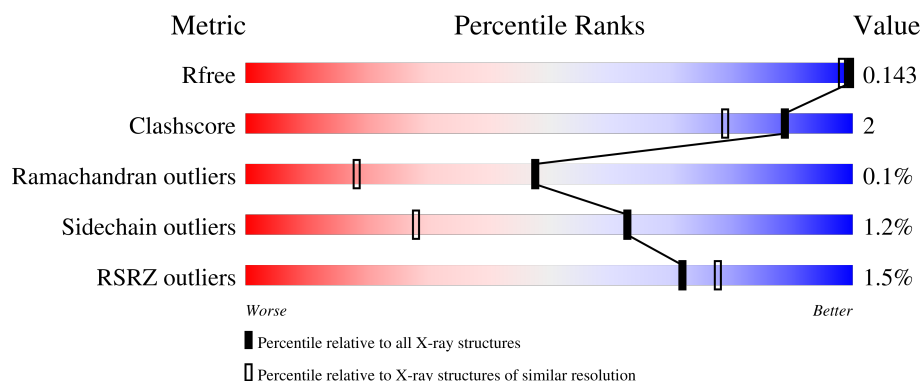
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.32 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	2531 (1.34-1.30)
Clashscore	190562	2585 (1.34-1.30)
Ramachandran outliers	187476	2528 (1.34-1.30)
Sidechain outliers	187428	2528 (1.34-1.30)
RSRZ outliers	180081	2528 (1.34-1.30)

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	% 99% .
2	B	122	2% 96% .
3	C	570	% 95% 5%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7127 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	5	0
			820	517	139	156	8			

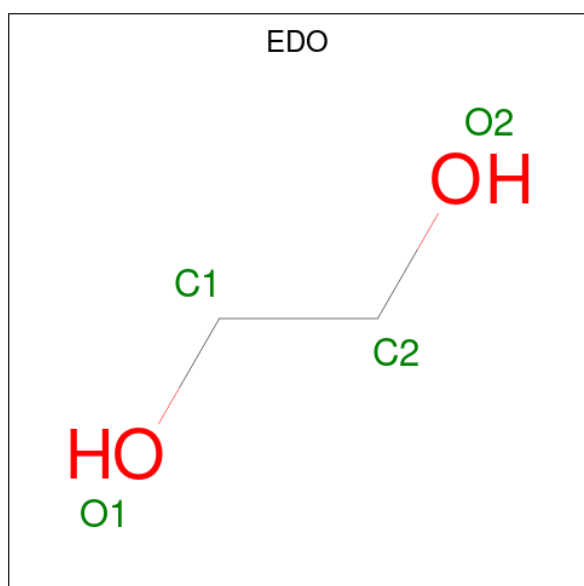
- Molecule 2 is a protein called Urease subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	122	Total	C	N	O	S	0	5	0
			990	610	179	200	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	18	0
			4469	2797	772	874	26			

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: 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	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
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 (CCD ID: SO4) (formula: O₄S).

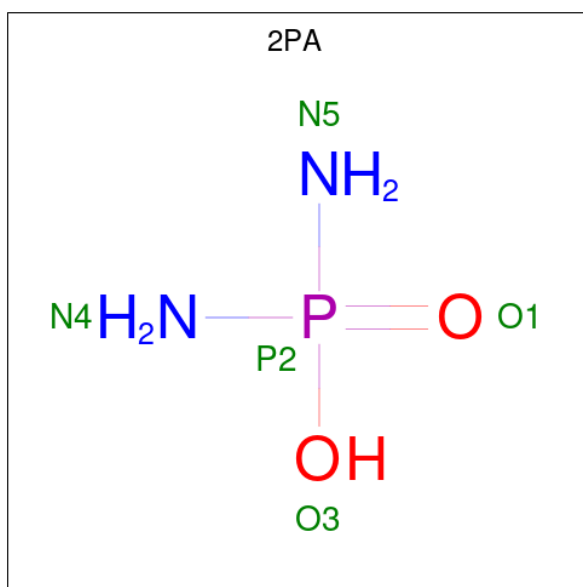


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

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

- Molecule 7 is DIAMIDOPHOSPHATE (CCD ID: 2PA) (formula: H₅N₂O₂P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	N	O	P	0	0
			5	2	2	1		

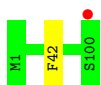
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	111	Total	O	0	0
			111	111		
8	B	146	Total	O	0	0
			146	146		
8	C	482	Total	O	0	0
			482	482		

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.

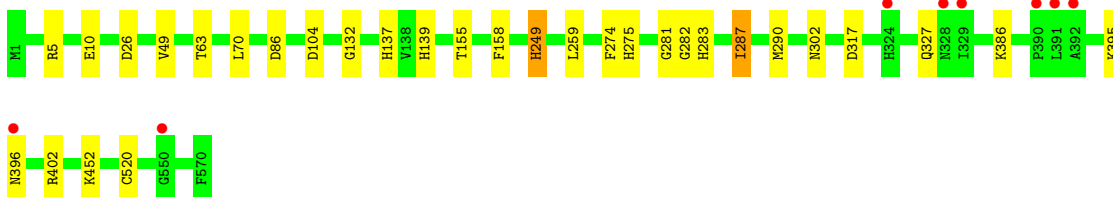
- Molecule 1: Urease subunit gamma



- Molecule 2: Urease subunit beta



- Molecule 3: Urease subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	131.78Å 131.78Å 188.77Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	97.67 – 1.32 97.67 – 1.32	Depositor EDS
% Data completeness (in resolution range)	98.6 (97.67-1.32) 98.6 (97.67-1.32)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 1.32Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.122 , 0.143 (Not available) , 0.143	Depositor DCC
R_{free} test set	11101 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	13.0	Xtriage
Anisotropy	0.588	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7127	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 2.69% 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: KCX, CXM, EDO, 2PA, NI, SO4

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	0.47	0/820	0.66	0/1100
2	B	0.43	0/1002	0.65	0/1347
3	C	0.49	0/4538	0.74	4/6147 (0.1%)
All	All	0.48	0/6360	0.72	4/8594 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	302	ASN	CA-C-N	-6.44	113.06	119.56
3	C	302	ASN	C-N-CA	-6.44	113.06	119.56
3	C	520	CYS	CB-CA-C	-5.74	98.84	109.29
3	C	302	ASN	N-CA-C	5.07	121.01	109.81

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	820	0	842	2	0
2	B	990	0	965	6	0
3	C	4469	0	4416	22	0
4	A	20	0	30	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	4	0	6	0	0
4	C	48	0	72	2	0
5	B	5	0	0	0	0
5	C	25	0	0	0	0
6	C	2	0	0	0	0
7	C	5	0	4	1	0
8	A	111	0	0	0	0
8	B	146	0	0	0	0
8	C	482	0	0	3	0
All	All	7127	0	6335	24	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 24 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)
2:B:17[B]:GLY:O	3:C:5[B]:ARG:CG	2.06	1.04
2:B:17[B]:GLY:O	3:C:5[B]:ARG:HG2	1.61	0.99
2:B:17[B]:GLY:O	3:C:5[B]:ARG:HG3	1.89	0.73
3:C:287[A]:ILE:O	3:C:287[A]:ILE:HD13	1.97	0.63
3:C:317:ASP:HB3	4:C:614:EDO:H11	1.84	0.60

There are no symmetry-related clashes.

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	103/100 (103%)	103 (100%)	0	0	100 100
2	B	125/122 (102%)	121 (97%)	3 (2%)	1 (1%)	16 2

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	584/570 (102%)	560 (96%)	24 (4%)	0	100	100
All	All	812/792 (102%)	784 (97%)	27 (3%)	1 (0%)	48	18

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	99	ILE

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	88/83 (106%)	88 (100%)	0	100	100
2	B	105/101 (104%)	105 (100%)	0	100	100
3	C	476/458 (104%)	466 (98%)	10 (2%)	47	11
All	All	669/642 (104%)	659 (98%)	10 (2%)	63	20

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	396	ASN
3	C	402	ARG
3	C	452	LYS
3	C	249	HIS
3	C	287[A]	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	HIS
3	C	179	ASN
3	C	327	GLN

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	9,10,11	0.74	0	9,11,13	1.04	0
3	KCX	C	220	6,3	10,11,12	0.60	0	6,12,14	0.65	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
1	CXM	A	1	1	-	3/9/10/12	-
3	KCX	C	220	6,3	-	0/9/10/12	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

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

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 2 are monoatomic - leaving 25 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	605	-	3,3,3	0.43	0	2,2,2	0.13	0
4	EDO	B	201	-	3,3,3	0.47	0	2,2,2	0.29	0
4	EDO	C	604	-	3,3,3	0.46	0	2,2,2	0.21	0
4	EDO	A	204	-	3,3,3	0.51	0	2,2,2	0.19	0
5	SO4	C	619	-	4,4,4	0.42	0	6,6,6	0.27	0
5	SO4	C	620	-	4,4,4	0.82	0	6,6,6	0.43	0
4	EDO	A	203	-	3,3,3	0.46	0	2,2,2	0.29	0
5	SO4	C	615	-	4,4,4	0.43	0	6,6,6	0.05	0
5	SO4	C	616	-	4,4,4	0.47	0	6,6,6	0.36	0
4	EDO	A	205	-	3,3,3	0.53	0	2,2,2	0.35	0
5	SO4	C	618	-	4,4,4	0.42	0	6,6,6	0.06	0
4	EDO	C	609	-	3,3,3	0.59	0	2,2,2	0.25	0
4	EDO	C	608	-	3,3,3	0.42	0	2,2,2	0.24	0
7	2PA	C	617	6	1,4,4	5.96	1 (100%)	0,6,6	-	-
4	EDO	A	201	-	3,3,3	0.47	0	2,2,2	0.44	0
4	EDO	C	614	-	3,3,3	0.36	0	2,2,2	0.29	0
5	SO4	B	202	-	4,4,4	0.44	0	6,6,6	0.10	0
4	EDO	C	607	-	3,3,3	0.50	0	2,2,2	0.18	0
4	EDO	C	610	-	3,3,3	0.44	0	2,2,2	0.34	0
4	EDO	C	613	-	3,3,3	0.39	0	2,2,2	0.34	0
4	EDO	C	603	-	3,3,3	0.39	0	2,2,2	0.33	0
4	EDO	C	606	-	3,3,3	0.53	0	2,2,2	0.12	0
4	EDO	A	202	-	3,3,3	0.54	0	2,2,2	0.65	0
4	EDO	C	612	-	3,3,3	0.44	0	2,2,2	0.50	0
4	EDO	C	611	-	3,3,3	0.40	0	2,2,2	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
4	EDO	C	607	-	-	0/1/1/1	-
4	EDO	C	605	-	-	0/1/1/1	-
4	EDO	C	614	-	-	1/1/1/1	-
4	EDO	C	610	-	-	0/1/1/1	-
4	EDO	B	201	-	-	0/1/1/1	-
4	EDO	C	604	-	-	0/1/1/1	-
4	EDO	A	204	-	-	0/1/1/1	-
4	EDO	C	609	-	-	1/1/1/1	-
4	EDO	C	613	-	-	1/1/1/1	-
4	EDO	A	203	-	-	0/1/1/1	-
4	EDO	C	603	-	-	0/1/1/1	-
4	EDO	C	606	-	-	0/1/1/1	-
4	EDO	A	201	-	-	0/1/1/1	-
4	EDO	C	608	-	-	0/1/1/1	-
4	EDO	A	202	-	-	1/1/1/1	-
4	EDO	C	612	-	-	0/1/1/1	-
4	EDO	A	205	-	-	0/1/1/1	-
4	EDO	C	611	-	-	1/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	617	2PA	P2-O1	5.96	1.56	1.46

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	613	EDO	O1-C1-C2-O2
4	C	614	EDO	O1-C1-C2-O2
4	A	202	EDO	O1-C1-C2-O2
4	C	611	EDO	O1-C1-C2-O2
4	C	609	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	205	EDO	2	0
7	C	617	2PA	1	0
4	C	614	EDO	2	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	99/100 (99%)	-0.46	1 (1%) 79 84	6, 13, 22, 26	5 (5%)
2	B	122/122 (100%)	-0.19	3 (2%) 58 65	8, 15, 29, 56	5 (4%)
3	C	569/570 (99%)	-0.37	8 (1%) 73 79	6, 13, 22, 84	18 (3%)
All	All	790/792 (99%)	-0.35	12 (1%) 72 78	6, 13, 23, 84	28 (3%)

The worst 5 of 12 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	391	LEU	4.1
3	C	396	ASN	3.5
3	C	324	HIS	2.9
2	B	126	GLU	2.8
3	C	390	PRO	2.8

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	0.98	0.06	14,16,18,20	0
3	KCX	C	220	12/13	0.98	0.04	10,11,11,12	0

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(Å ²)	Q<0.9
4	EDO	C	613	4/4	0.78	0.21	36,37,40,43	4
4	EDO	C	607	4/4	0.80	0.18	39,47,48,52	0
4	EDO	A	202	4/4	0.82	0.17	31,38,44,48	0
4	EDO	A	204	4/4	0.83	0.21	32,33,48,81	0
4	EDO	C	608	4/4	0.84	0.15	54,56,64,66	0
4	EDO	A	205	4/4	0.85	0.15	27,31,34,35	0
5	SO4	C	615	5/5	0.85	0.15	21,61,65,75	5
4	EDO	C	614	4/4	0.86	0.16	21,21,24,24	4
5	SO4	B	202	5/5	0.87	0.12	33,68,71,76	5
4	EDO	A	203	4/4	0.87	0.14	33,38,43,44	0
4	EDO	C	612	4/4	0.89	0.16	23,24,31,35	4
5	SO4	C	618	5/5	0.89	0.18	36,46,54,55	5
5	SO4	C	616	5/5	0.91	0.12	18,19,34,36	5
4	EDO	C	611	4/4	0.92	0.14	28,43,53,58	0
4	EDO	B	201	4/4	0.92	0.10	31,32,32,40	0
4	EDO	C	606	4/4	0.92	0.13	38,43,53,78	0
4	EDO	C	610	4/4	0.92	0.10	34,35,35,38	0
4	EDO	C	609	4/4	0.94	0.10	20,27,28,29	0
4	EDO	C	604	4/4	0.94	0.09	25,26,28,35	0
5	SO4	C	619	5/5	0.94	0.10	34,34,34,35	5
5	SO4	C	620	5/5	0.94	0.09	34,43,45,47	5
4	EDO	C	605	4/4	0.96	0.07	22,23,26,29	0
4	EDO	C	603	4/4	0.96	0.07	21,23,28,28	0
4	EDO	A	201	4/4	0.99	0.05	14,16,19,21	0
7	2PA	C	617	5/5	0.99	0.04	11,11,12,13	0
6	NI	C	602	1/1	1.00	0.02	11,11,11,11	0
6	NI	C	601	1/1	1.00	0.02	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.