



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 30, 2026 – 11:00 AM UTC

PDB ID : 6BIA / pdb_00006bia
Title : Crystal structure of Ps i-CgsB
Authors : Hettle, A.G.; Boraston, A.B.
Deposited on : 2017-11-01
Resolution : 2.80 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

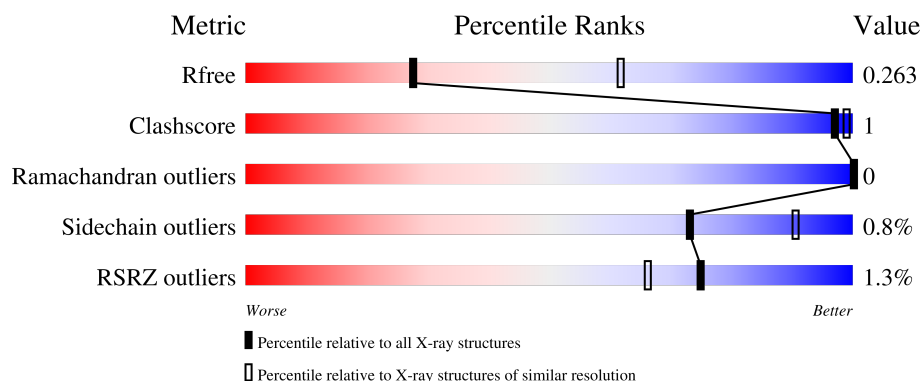
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	477	0% 91% 5%
1	B	477	0% 92% 5%
1	C	477	2% 92% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10786 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 Sulfatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	451	Total	C	N	O	S	0	0	0
			3567	2279	617	658	13			
1	B	452	Total	C	N	O	S	0	0	0
			3561	2272	617	659	13			
1	C	451	Total	C	N	O	S	0	0	0
			3543	2260	611	659	13			

There are 87 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	MET	-	initiating methionine	UNP A0A063KPH1
A	5	GLY	-	expression tag	UNP A0A063KPH1
A	6	SER	-	expression tag	UNP A0A063KPH1
A	7	SER	-	expression tag	UNP A0A063KPH1
A	8	HIS	-	expression tag	UNP A0A063KPH1
A	9	HIS	-	expression tag	UNP A0A063KPH1
A	10	HIS	-	expression tag	UNP A0A063KPH1
A	11	HIS	-	expression tag	UNP A0A063KPH1
A	12	HIS	-	expression tag	UNP A0A063KPH1
A	13	HIS	-	expression tag	UNP A0A063KPH1
A	14	SER	-	expression tag	UNP A0A063KPH1
A	15	SER	-	expression tag	UNP A0A063KPH1
A	16	GLY	-	expression tag	UNP A0A063KPH1
A	17	LEU	-	expression tag	UNP A0A063KPH1
A	18	VAL	-	expression tag	UNP A0A063KPH1
A	19	PRO	-	expression tag	UNP A0A063KPH1
A	20	ARG	-	expression tag	UNP A0A063KPH1
A	21	GLY	-	expression tag	UNP A0A063KPH1
A	22	SER	-	expression tag	UNP A0A063KPH1
A	23	HIS	-	expression tag	UNP A0A063KPH1
A	24	MET	-	expression tag	UNP A0A063KPH1
A	25	ALA	-	expression tag	UNP A0A063KPH1
A	29	LYS	ARG	conflict	UNP A0A063KPH1

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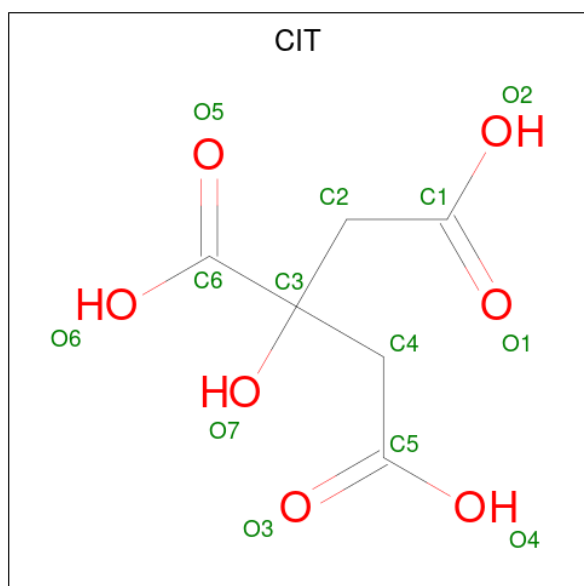
Chain	Residue	Modelled	Actual	Comment	Reference
A	138	THR	ALA	conflict	UNP A0A063KPH1
A	141	HIS	TYR	conflict	UNP A0A063KPH1
A	174	LYS	GLN	conflict	UNP A0A063KPH1
A	221	HIS	ASN	conflict	UNP A0A063KPH1
A	234	THR	VAL	conflict	UNP A0A063KPH1
A	274	ALA	THR	conflict	UNP A0A063KPH1
B	4	MET	-	initiating methionine	UNP A0A063KPH1
B	5	GLY	-	expression tag	UNP A0A063KPH1
B	6	SER	-	expression tag	UNP A0A063KPH1
B	7	SER	-	expression tag	UNP A0A063KPH1
B	8	HIS	-	expression tag	UNP A0A063KPH1
B	9	HIS	-	expression tag	UNP A0A063KPH1
B	10	HIS	-	expression tag	UNP A0A063KPH1
B	11	HIS	-	expression tag	UNP A0A063KPH1
B	12	HIS	-	expression tag	UNP A0A063KPH1
B	13	HIS	-	expression tag	UNP A0A063KPH1
B	14	SER	-	expression tag	UNP A0A063KPH1
B	15	SER	-	expression tag	UNP A0A063KPH1
B	16	GLY	-	expression tag	UNP A0A063KPH1
B	17	LEU	-	expression tag	UNP A0A063KPH1
B	18	VAL	-	expression tag	UNP A0A063KPH1
B	19	PRO	-	expression tag	UNP A0A063KPH1
B	20	ARG	-	expression tag	UNP A0A063KPH1
B	21	GLY	-	expression tag	UNP A0A063KPH1
B	22	SER	-	expression tag	UNP A0A063KPH1
B	23	HIS	-	expression tag	UNP A0A063KPH1
B	24	MET	-	expression tag	UNP A0A063KPH1
B	25	ALA	-	expression tag	UNP A0A063KPH1
B	29	LYS	ARG	conflict	UNP A0A063KPH1
B	138	THR	ALA	conflict	UNP A0A063KPH1
B	141	HIS	TYR	conflict	UNP A0A063KPH1
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B	234	THR	VAL	conflict	UNP A0A063KPH1
B	274	ALA	THR	conflict	UNP A0A063KPH1
C	4	MET	-	initiating methionine	UNP A0A063KPH1
C	5	GLY	-	expression tag	UNP A0A063KPH1
C	6	SER	-	expression tag	UNP A0A063KPH1
C	7	SER	-	expression tag	UNP A0A063KPH1
C	8	HIS	-	expression tag	UNP A0A063KPH1
C	9	HIS	-	expression tag	UNP A0A063KPH1
C	10	HIS	-	expression tag	UNP A0A063KPH1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	11	HIS	-	expression tag	UNP A0A063KPH1
C	12	HIS	-	expression tag	UNP A0A063KPH1
C	13	HIS	-	expression tag	UNP A0A063KPH1
C	14	SER	-	expression tag	UNP A0A063KPH1
C	15	SER	-	expression tag	UNP A0A063KPH1
C	16	GLY	-	expression tag	UNP A0A063KPH1
C	17	LEU	-	expression tag	UNP A0A063KPH1
C	18	VAL	-	expression tag	UNP A0A063KPH1
C	19	PRO	-	expression tag	UNP A0A063KPH1
C	20	ARG	-	expression tag	UNP A0A063KPH1
C	21	GLY	-	expression tag	UNP A0A063KPH1
C	22	SER	-	expression tag	UNP A0A063KPH1
C	23	HIS	-	expression tag	UNP A0A063KPH1
C	24	MET	-	expression tag	UNP A0A063KPH1
C	25	ALA	-	expression tag	UNP A0A063KPH1
C	29	LYS	ARG	conflict	UNP A0A063KPH1
C	138	THR	ALA	conflict	UNP A0A063KPH1
C	141	HIS	TYR	conflict	UNP A0A063KPH1
C	174	LYS	GLN	conflict	UNP A0A063KPH1
C	221	HIS	ASN	conflict	UNP A0A063KPH1
C	234	THR	VAL	conflict	UNP A0A063KPH1
C	274	ALA	THR	conflict	UNP A0A063KPH1

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).

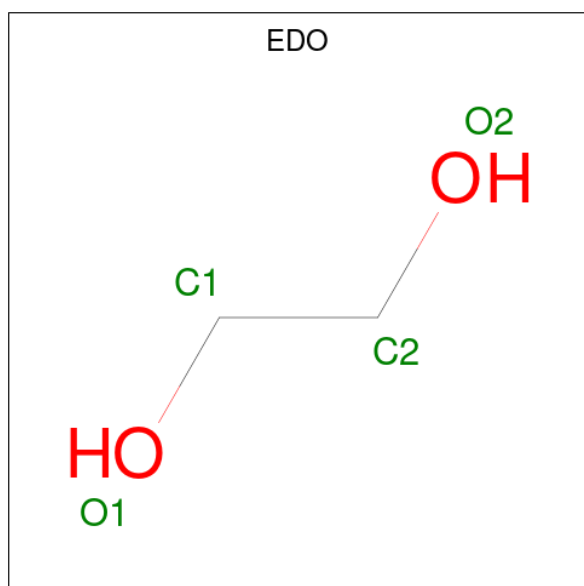


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 13 6 7	0	0
2	B	1	Total C O 13 6 7	0	0
2	C	1	Total C O 13 6 7	0	0

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Ca 1 1	0	0
3	B	1	Total Ca 1 1	0	0
3	C	1	Total Ca 1 1	0	0

- 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	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0

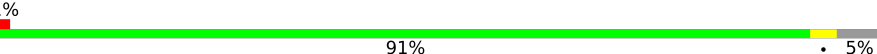
- Molecule 5 is water.

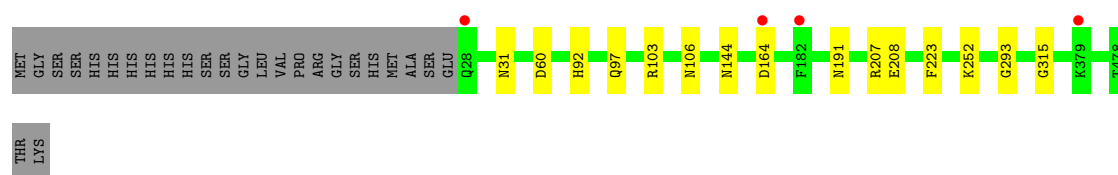
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	24	Total 24	O 24	0	0
5	B	21	Total 21	O 21	0	0
5	C	16	Total 16	O 16	0	0

3 Residue-property plots [i](#)

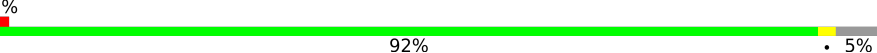
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Sulfatase

Chain A: 

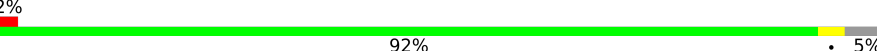


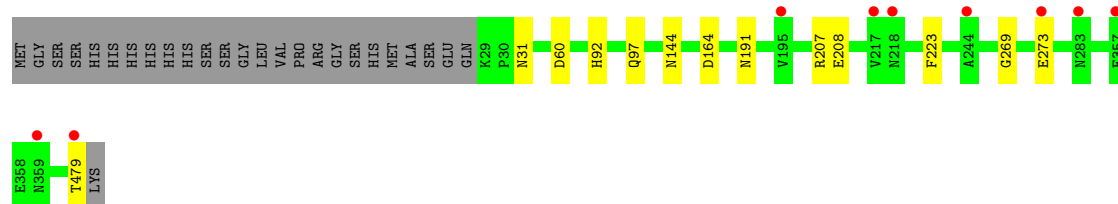
- Molecule 1: Sulfatase

Chain B: 



- Molecule 1: Sulfatase

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	133.85Å 133.85Å 223.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	114.86 – 2.80 114.86 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.1 (114.86-2.80) 97.1 (114.86-2.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.76 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.238 , 0.265 0.240 , 0.263	Depositor DCC
R_{free} test set	2464 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	51.8	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.0	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	0.91	EDS
Total number of atoms	10786	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.56% 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: EDO, CIT, CA

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	0.59	0/3673	0.78	0/4986
1	B	0.58	0/3667	0.78	0/4983
1	C	0.59	0/3649	0.78	0/4960
All	All	0.59	0/10989	0.78	0/14929

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	3567	0	3366	5	1
1	B	3561	0	3342	8	1
1	C	3543	0	3297	5	0
2	A	13	0	5	0	0
2	B	13	0	5	0	0
2	C	13	0	5	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	8	0	12	1	0
5	A	24	0	0	0	0
5	B	21	0	0	0	0
5	C	16	0	0	0	0
All	All	10786	0	10038	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)
1:C:207:ARG:NH1	1:C:208:GLU:OE2	2.23	0.71
1:B:245:MET:HE2	1:B:245:MET:HA	1.86	0.57
1:A:207:ARG:NH1	1:A:208:GLU:OE2	2.37	0.56
1:B:197:GLU:OE2	1:B:207:ARG:NH2	2.45	0.50
1:B:31:ASN:HD22	1:B:223:PHE:H	1.61	0.49
1:B:197:GLU:CD	1:B:207:ARG:HH22	2.21	0.48
1:C:31:ASN:HD22	1:C:223:PHE:H	1.60	0.48
1:A:31:ASN:HD22	1:A:223:PHE:H	1.61	0.47
1:A:144:ASN:HD21	1:A:191:ASN:HA	1.79	0.47
1:B:144:ASN:HD21	1:B:191:ASN:HA	1.79	0.47
1:B:245:MET:HE2	1:B:245:MET:CA	2.45	0.46
1:C:144:ASN:HD21	1:C:191:ASN:HA	1.80	0.45
1:B:99:ASN:HD22	4:B:503:EDO:H11	1.84	0.43
1:B:92:HIS:HA	1:B:97:GLN:HG3	2.01	0.43
1:A:92:HIS:HA	1:A:97:GLN:HG3	2.02	0.42
1:C:92:HIS:HA	1:C:97:GLN:HG3	2.02	0.41
1:C:269:GLY:O	1:C:273:GLU:HG2	2.21	0.41
1:A:293:GLY:HA3	1:A:315:GLY:O	2.21	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:60:ASP:OD2	1:B:241:GLU:OE1[3_744]	2.11	0.09

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	449/477 (94%)	434 (97%)	15 (3%)	0	100	100
1	B	450/477 (94%)	435 (97%)	15 (3%)	0	100	100
1	C	449/477 (94%)	434 (97%)	15 (3%)	0	100	100
All	All	1348/1431 (94%)	1303 (97%)	45 (3%)	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	364/404 (90%)	360 (99%)	4 (1%)	65	87
1	B	363/404 (90%)	361 (99%)	2 (1%)	78	92
1	C	357/404 (88%)	354 (99%)	3 (1%)	73	90
All	All	1084/1212 (89%)	1075 (99%)	9 (1%)	73	90

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	ARG
1	A	106	ASN
1	A	164	ASP
1	A	252	LYS
1	B	241	GLU

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Mol	Chain	Res	Type
1	B	252	LYS
1	C	60	ASP
1	C	164	ASP
1	C	479	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	A	57	ASN
1	A	121	ASN
1	A	141	HIS
1	A	144	ASN
1	A	221	HIS
1	A	278	ASN
1	A	410	GLN
1	B	31	ASN
1	B	106	ASN
1	B	121	ASN
1	B	141	HIS
1	B	144	ASN
1	B	169	GLN
1	B	180	ASN
1	B	221	HIS
1	B	278	ASN
1	B	313	GLN
1	C	31	ASN
1	C	121	ASN
1	C	141	HIS
1	C	144	ASN
1	C	173	GLN
1	C	180	ASN
1	C	221	HIS
1	C	278	ASN
1	C	279	ASN
1	C	410	GLN

5.3.3 RNA ⓘ

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 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	B	503	-	3,3,3	0.47	0	2,2,2	0.44	0
4	EDO	A	503	-	3,3,3	0.53	0	2,2,2	0.28	0
2	CIT	C	501	-	12,12,12	1.00	0	17,17,17	1.31	2 (11%)
2	CIT	A	501	-	12,12,12	1.10	0	17,17,17	1.29	3 (17%)
4	EDO	B	504	-	3,3,3	0.53	0	2,2,2	0.11	0
2	CIT	B	501	-	12,12,12	1.09	0	17,17,17	0.99	1 (5%)

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	B	503	-	-	1/1/1/1	-
4	EDO	A	503	-	-	0/1/1/1	-
2	CIT	C	501	-	-	5/16/16/16	-
2	CIT	A	501	-	-	11/16/16/16	-
4	EDO	B	504	-	-	1/1/1/1	-
2	CIT	B	501	-	-	0/16/16/16	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	CIT	O6-C6-C3	3.26	119.39	113.14
2	B	501	CIT	O6-C6-C3	2.81	118.54	113.14
2	A	501	CIT	O6-C6-C3	2.66	118.24	113.14
2	A	501	CIT	O1-C1-C2	-2.40	116.16	122.95
2	A	501	CIT	O2-C1-C2	2.27	121.55	114.35
2	C	501	CIT	O3-C5-C4	-2.21	116.69	122.95

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	501	CIT	C1-C2-C3-O7
2	C	501	CIT	C1-C2-C3-C4
2	C	501	CIT	C1-C2-C3-C6
4	B	503	EDO	O1-C1-C2-O2
2	A	501	CIT	O7-C3-C6-O5
2	A	501	CIT	C4-C3-C6-O6
2	A	501	CIT	C2-C3-C6-O5
2	A	501	CIT	C2-C3-C6-O6
2	A	501	CIT	O7-C3-C4-C5
2	A	501	CIT	C6-C3-C4-C5
2	A	501	CIT	C2-C3-C4-C5
2	A	501	CIT	O7-C3-C6-O6
2	A	501	CIT	C4-C3-C6-O5
4	B	504	EDO	O1-C1-C2-O2
2	C	501	CIT	C3-C4-C5-O3
2	C	501	CIT	C3-C4-C5-O4
2	A	501	CIT	O1-C1-C2-C3
2	A	501	CIT	O2-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	EDO	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	451/477 (94%)	-0.13	4 (0%) 81 74	33, 42, 60, 70	0
1	B	452/477 (94%)	-0.10	4 (0%) 81 74	31, 42, 57, 80	2 (0%)
1	C	451/477 (94%)	0.34	9 (1%) 65 56	33, 55, 71, 87	1 (0%)
All	All	1354/1431 (94%)	0.04	17 (1%) 75 66	31, 45, 67, 87	3 (0%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	28	GLN	5.1
1	B	479	THR	4.4
1	C	479	THR	3.7
1	C	283	ASN	3.2
1	C	359	ASN	2.7
1	B	408	ASN	2.6
1	A	182	PHE	2.3
1	C	218	ASN	2.3
1	C	357	GLU	2.3
1	B	220	LYS	2.2
1	A	379	LYS	2.2
1	C	195	VAL	2.1
1	B	118	LYS	2.1
1	A	164	ASP	2.1
1	C	217	VAL	2.0
1	C	244	ALA	2.0
1	C	273	GLU	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
4	EDO	A	503	4/4	0.57	0.32	57,58,58,59	0
4	EDO	B	504	4/4	0.69	0.30	63,64,66,67	0
4	EDO	B	503	4/4	0.70	0.24	55,59,62,65	0
2	CIT	A	501	13/13	0.82	0.13	55,64,72,74	0
2	CIT	C	501	13/13	0.84	0.12	63,68,78,79	0
2	CIT	B	501	13/13	0.85	0.11	57,62,66,67	0
3	CA	B	502	1/1	0.97	0.04	45,45,45,45	0
3	CA	A	502	1/1	0.98	0.03	43,43,43,43	0
3	CA	C	502	1/1	0.99	0.03	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.