



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:07 AM UTC

PDB ID : 6ZO0 / pdb_00006zo0
Title : 2.23 Å resolution 3,4-dimethylcatechol (3,4-dimethylbenzene-1,2-diol) inhibited *Sporosarcina pasteurii* urease
Authors : Mazzei, L.; Ciani, M.; Musiani, F.; Ciurli, S.
Deposited on : 2020-07-07
Resolution : 2.23 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

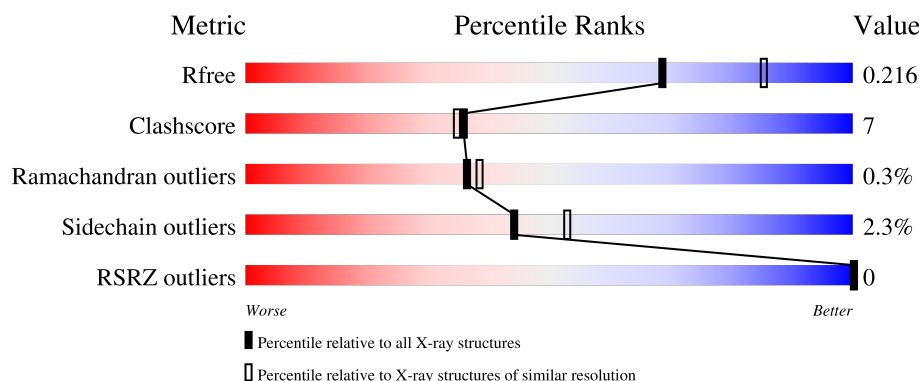
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.23 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	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	AAA	100	 83% 15% .
2	BBB	122	 84% 16% .
3	CCC	570	 82% 16% .

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	CCC	608	-	-	X	-
4	EDO	CCC	611	-	-	X	-
5	SO4	BBB	203	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6776 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	AAA	100	Total	C	N	O	S	0	5	0
			815	512	138	158	7			

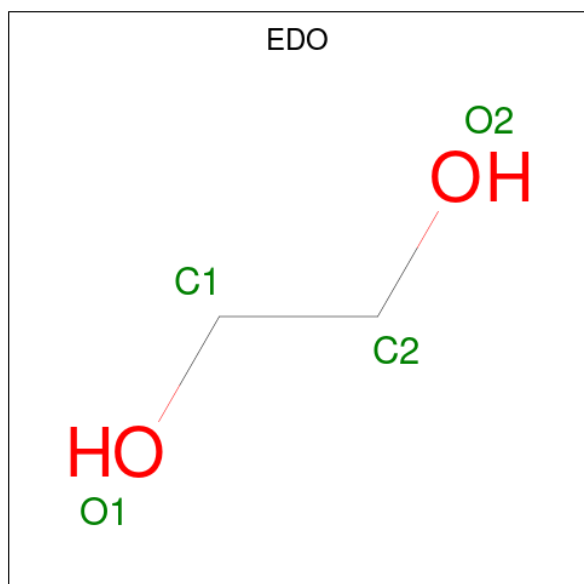
- Molecule 2 is a protein called Urease subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	BBB	122	Total	C	N	O	S	0	4	0
			984	607	177	199	1			

- Molecule 3 is a protein called Urease subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	CCC	570	Total	C	N	O	S	0	17	0
			4449	2794	765	864	26			

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	BBB	1	Total C O 4 2 2	0	0
4	BBB	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
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4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0
4	CCC	1	Total C O 4 2 2	0	0

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

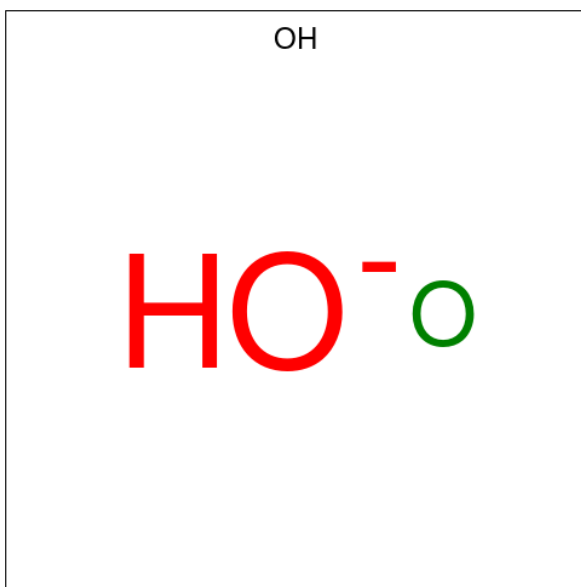


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	BBB	1	Total	O	S	0	0
			5	4	1		
5	BBB	1	Total	O	S	0	0
			5	4	1		
5	BBB	1	Total	O	S	0	0
			5	4	1		
5	CCC	1	Total	O	S	0	0
			5	4	1		
5	CCC	1	Total	O	S	0	0
			5	4	1		
5	CCC	1	Total	O	S	0	0
			5	4	1		
5	CCC	1	Total	O	S	0	0
			5	4	1		
5	CCC	1	Total	O	S	0	0
			5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

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

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

- Molecule 7 is HYDROXIDE ION (CCD ID: OH) (formula: HO).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	AAA	50	Total O 50 50	0	0
8	BBB	87	Total O 87 87	0	0
8	CCC	255	Total O 255 255	0	0

3 Residue-property plots

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 AAA:  83% 15% .



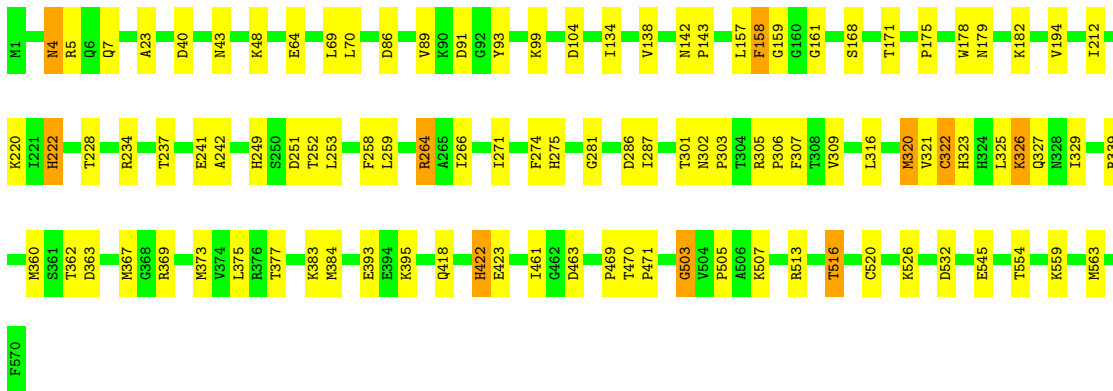
- Molecule 2: Urease subunit beta

Chain BBB:  84% 16% .



- Molecule 3: Urease subunit alpha

Chain CCC:  82% 16% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	131.50Å 131.50Å 188.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	113.88 – 2.23 113.88 – 2.23	Depositor EDS
% Data completeness (in resolution range)	100.0 (113.88-2.23) 100.0 (113.88-2.23)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.161 , 0.216 0.161 , 0.216	Depositor DCC
R_{free} test set	2279 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	27.9	Xtriage
Anisotropy	0.577	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6776	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage'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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, EDO, NI, CXM, QO8, KCX, OH

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	AAA	0.65	0/821	1.15	1/1104 (0.1%)
2	BBB	0.66	0/999	1.27	3/1342 (0.2%)
3	CCC	0.70	2/4525 (0.0%)	1.21	19/6125 (0.3%)
All	All	0.69	2/6345 (0.0%)	1.21	23/8571 (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
2	BBB	0	1
3	CCC	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	CCC	222	HIS	CE1-NE2	5.83	1.38	1.32
3	CCC	422	HIS	CE1-NE2	5.01	1.37	1.32

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	29	THR	CA-CB-OG1	-7.61	98.19	109.60
3	CCC	4	ASN	CB-CA-C	7.56	122.58	109.65
3	CCC	234	ARG	CG-CD-NE	-6.88	96.85	112.00
3	CCC	503	GLY	CA-C-N	-6.82	115.94	120.24
3	CCC	503	GLY	C-N-CA	-6.82	115.94	120.24
3	CCC	520	CYS	CB-CA-C	-6.79	96.94	109.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	CCC	554	THR	CA-CB-OG1	-6.74	99.49	109.60
3	CCC	264	ARG	CG-CD-NE	6.35	125.98	112.00
3	CCC	516	THR	CA-CB-OG1	-6.33	100.11	109.60
3	CCC	251	ASP	CA-CB-CG	6.07	118.67	112.60
3	CCC	228	THR	CA-CB-OG1	-6.06	100.51	109.60
3	CCC	40	ASP	CA-CB-CG	6.00	118.60	112.60
2	BBB	13	ARG	CG-CD-NE	5.92	125.03	112.00
3	CCC	369	ARG	CA-C-N	5.68	128.16	120.38
3	CCC	369	ARG	C-N-CA	5.68	128.16	120.38
1	AAA	74	PRO	N-CA-C	-5.46	105.94	113.53
2	BBB	5	ASN	CA-CB-CG	-5.39	107.21	112.60
3	CCC	393	GLU	CB-CG-CD	5.37	121.73	112.60
3	CCC	237	THR	CA-CB-OG1	-5.35	101.58	109.60
3	CCC	23	ALA	CB-CA-C	-5.29	110.45	116.54
3	CCC	326	LYS	CB-CA-C	5.19	118.38	109.51
3	CCC	463	ASP	CA-CB-CG	5.16	117.76	112.60
3	CCC	91	ASP	CB-CA-C	-5.04	103.55	111.51

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	BBB	125	VAL	Peptide
3	CCC	322	QO8	Mainchain

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	AAA	815	0	831	13	0
2	BBB	984	0	964	12	0
3	CCC	4449	0	4418	69	0
4	AAA	16	0	24	2	0
4	BBB	8	0	12	0	0
4	CCC	64	0	96	17	0
5	BBB	15	0	0	2	0
5	CCC	30	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	CCC	2	0	0	0	0
7	CCC	1	0	0	1	0
8	AAA	50	0	0	0	0
8	BBB	87	0	0	5	0
8	CCC	255	0	0	5	0
All	All	6776	0	6345	93	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 7.

All (93) 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:CCC:513[A]:ARG:HH21	4:CCC:611:EDO:C1	1.52	1.21
3:CCC:513[A]:ARG:NH2	4:CCC:611:EDO:H12	1.62	1.15
2:BBB:111[A]:GLU:OE2	8:BBB:301:HOH:O	1.70	1.09
3:CCC:513[A]:ARG:NH2	4:CCC:611:EDO:C1	2.19	1.02
2:BBB:111[A]:GLU:CD	8:BBB:301:HOH:O	2.10	0.94
3:CCC:513[A]:ARG:HH21	4:CCC:611:EDO:H12	0.79	0.93
2:BBB:31[A]:ARG:HG3	5:BBB:203:SO4:O3	1.74	0.87
3:CCC:505:PRO:HD2	4:CCC:608:EDO:H11	1.65	0.79
3:CCC:326:LYS:HB2	3:CCC:329:ILE:HD12	1.65	0.79
3:CCC:264:ARG:NH2	8:CCC:703:HOH:O	2.15	0.78
7:CCC:603:OH:O	8:CCC:702:HOH:O	2.03	0.77
2:BBB:111[A]:GLU:OE1	8:BBB:301:HOH:O	2.00	0.75
3:CCC:48[B]:LYS:NZ	8:CCC:701:HOH:O	1.94	0.74
3:CCC:64[B]:GLU:OE2	4:CCC:605:EDO:O2	2.06	0.74
3:CCC:505:PRO:HB2	4:CCC:608:EDO:H22	1.71	0.72
3:CCC:513[A]:ARG:NH2	4:CCC:611:EDO:H11	2.09	0.65
3:CCC:309:VAL:HG11	3:CCC:559:LYS:HD2	1.82	0.61
2:BBB:31[A]:ARG:NH1	5:BBB:203:SO4:O4	2.33	0.60
3:CCC:158:PHE:CE2	3:CCC:418[A]:GLN:CG	2.86	0.58
3:CCC:220:KCX:HE2	3:CCC:274:PHE:CD2	2.38	0.58
3:CCC:503:GLY:HA2	4:CCC:608:EDO:C1	2.33	0.58
3:CCC:301:THR:CG2	3:CCC:363:ASP:HB2	2.34	0.58
1:AAA:44:MET:HG2	1:AAA:84:ALA:HB2	1.85	0.57
3:CCC:303:PRO:HG3	3:CCC:367:MET:O	2.06	0.56
3:CCC:360[A]:MET:HE3	3:CCC:377:THR:HA	1.87	0.55
3:CCC:323:HIS:HB2	3:CCC:325:LEU:HG	1.89	0.55
3:CCC:134:ILE:HD11	8:CCC:826:HOH:O	2.07	0.54
3:CCC:70:LEU:HD11	3:CCC:86:ASP:HB3	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CCC:323:HIS:HE1	4:CCC:609:EDO:H22	1.72	0.54
3:CCC:220:KCX:HE2	3:CCC:274:PHE:HD2	1.73	0.54
3:CCC:220:KCX:CX	3:CCC:222:HIS:HD2	2.22	0.53
3:CCC:503:GLY:HA2	4:CCC:608:EDO:H12	1.90	0.53
2:BBB:19:ILE:HG12	3:CCC:5[A]:ARG:HD2	1.91	0.52
1:AAA:73:VAL:N	1:AAA:74:PRO:CD	2.74	0.51
3:CCC:302:ASN:N	3:CCC:303:PRO:HD2	2.26	0.51
2:BBB:13:ARG:HD2	8:BBB:355:HOH:O	2.12	0.50
3:CCC:43:ASN:HA	8:CCC:795:HOH:O	2.11	0.50
3:CCC:362:THR:HG22	3:CCC:373:MET:HB2	1.94	0.49
1:AAA:44:MET:HE3	1:AAA:82:ALA:HB1	1.95	0.48
3:CCC:307:PHE:O	3:CCC:383:LYS:HE3	2.14	0.48
2:BBB:103:THR:HG23	2:BBB:113:ILE:HD11	1.95	0.48
3:CCC:323:HIS:CE1	4:CCC:609:EDO:H22	2.49	0.48
1:AAA:8:LYS:HE3	1:AAA:8:LYS:HB3	1.61	0.47
3:CCC:158:PHE:CE2	3:CCC:418[A]:GLN:HG3	2.50	0.47
3:CCC:320:MET:HA	3:CCC:325:LEU:HB2	1.97	0.47
3:CCC:212:ILE:HG21	3:CCC:242:ALA:HB1	1.96	0.46
3:CCC:157:LEU:O	3:CCC:194:VAL:HA	2.16	0.46
3:CCC:175:PRO:O	3:CCC:179:ASN:HB2	2.16	0.46
3:CCC:69:LEU:HB3	3:CCC:89:VAL:HB	1.97	0.46
3:CCC:503:GLY:HA2	4:CCC:608:EDO:H11	1.96	0.46
1:AAA:44:MET:HE2	1:AAA:83:GLU:O	2.16	0.45
3:CCC:93:TYR:OH	4:CCC:611:EDO:H11	2.16	0.45
3:CCC:171:THR:HG22	3:CCC:222:HIS:CG	2.51	0.45
3:CCC:286:ASP:O	3:CCC:287:ILE:C	2.60	0.45
3:CCC:274:PHE:O	3:CCC:275:HIS:C	2.60	0.45
3:CCC:266:ILE:HD13	3:CCC:271:ILE:HG21	1.99	0.45
3:CCC:320:MET:HA	3:CCC:325:LEU:HD12	1.99	0.44
1:AAA:30:LEU:HD13	1:AAA:38:ILE:HD12	2.00	0.44
3:CCC:422:HIS:CD2	3:CCC:423:GLU:HG2	2.53	0.44
3:CCC:384:MET:HE2	3:CCC:384:MET:HA	2.00	0.44
3:CCC:323:HIS:CE1	4:CCC:618:EDO:H11	2.52	0.44
3:CCC:375:LEU:HD12	3:CCC:563:MET:HE2	2.00	0.44
3:CCC:142:ASN:HA	3:CCC:143:PRO:HD3	1.89	0.43
2:BBB:27:LYS:HG2	8:BBB:342:HOH:O	2.18	0.43
3:CCC:559:LYS:N	5:CCC:622:SO4:O2	2.52	0.43
3:CCC:253:LEU:HD21	4:CCC:609:EDO:H11	2.01	0.43
3:CCC:470:THR:N	3:CCC:471:PRO:CD	2.81	0.43
1:AAA:66[B]:ARG:HD2	1:AAA:100:SER:OXT	2.19	0.43
2:BBB:57:PHE:O	2:BBB:124:GLY:HA3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CCC:158:PHE:CD2	3:CCC:418[A]:GLN:CD	2.97	0.42
3:CCC:360[A]:MET:HG3	3:CCC:377:THR:OG1	2.18	0.42
3:CCC:161:GLY:HA2	3:CCC:168:SER:OG	2.18	0.42
3:CCC:138:VAL:O	3:CCC:159:GLY:HA3	2.19	0.42
3:CCC:93:TYR:OH	4:CCC:611:EDO:C1	2.67	0.42
3:CCC:362:THR:CG2	3:CCC:373:MET:HB2	2.50	0.42
2:BBB:95:GLU:O	3:CCC:104:ASP:HB3	2.19	0.42
1:AAA:42:PHE:HZ	1:AAA:59:GLU:OE1	2.03	0.42
3:CCC:249:HIS:CE1	3:CCC:281:GLY:HA3	2.55	0.42
3:CCC:305:ARG:HA	3:CCC:306:PRO:HA	1.76	0.42
3:CCC:422:HIS:CE1	3:CCC:516:THR:O	2.73	0.41
3:CCC:461:ILE:HD11	3:CCC:469:PRO:HB3	2.02	0.41
3:CCC:99:LYS:C	3:CCC:99:LYS:HD3	2.45	0.41
1:AAA:6:ALA:O	1:AAA:10:LYS:HG3	2.20	0.41
3:CCC:252:THR:HG21	3:CCC:339:ARG:HD2	2.03	0.41
1:AAA:10:LYS:HD3	1:AAA:44:MET:CE	2.51	0.41
3:CCC:532:ASP:OD1	3:CCC:532:ASP:N	2.46	0.41
3:CCC:178:TRP:O	3:CCC:182:LYS:HG2	2.22	0.40
3:CCC:316:LEU:HD23	3:CCC:316:LEU:HA	1.91	0.40
3:CCC:258:PHE:O	3:CCC:259:LEU:C	2.64	0.40
1:AAA:6:ALA:CB	4:AAA:202:EDO:H12	2.50	0.40
1:AAA:6:ALA:HB3	4:AAA:202:EDO:H12	2.04	0.40
2:BBB:44:SER:O	2:BBB:66:ARG:HD2	2.22	0.40
1:AAA:66[A]:ARG:NH2	1:AAA:75:GLU:OE2	2.54	0.40

There are no symmetry-related clashes.

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	AAA	103/100 (103%)	103 (100%)	0	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	BBB	124/122 (102%)	119 (96%)	3 (2%)	2 (2%)	7	4
3	CCC	583/570 (102%)	556 (95%)	27 (5%)	0	100	100
All	All	810/792 (102%)	778 (96%)	30 (4%)	2 (0%)	36	48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	BBB	99	ILE
2	BBB	52	ASN

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	AAA	88/83 (106%)	86 (98%)	2 (2%)	44	53
2	BBB	105/101 (104%)	101 (96%)	4 (4%)	29	34
3	CCC	473/457 (104%)	460 (97%)	13 (3%)	39	47
All	All	666/641 (104%)	647 (97%)	19 (3%)	44	44

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	66[A]	ARG
1	AAA	66[B]	ARG
2	BBB	26	GLU
2	BBB	33[A]	SER
2	BBB	33[B]	SER
2	BBB	115	GLN
3	CCC	4	ASN
3	CCC	7	GLN
3	CCC	158	PHE
3	CCC	241	GLU
3	CCC	320	MET
3	CCC	321	VAL

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Mol	Chain	Res	Type
3	CCC	327	GLN
3	CCC	395	LYS
3	CCC	507[A]	LYS
3	CCC	507[B]	LYS
3	CCC	526	LYS
3	CCC	545[A]	GLU
3	CCC	545[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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
3	QO8	CCC	322	3	16,16,17	1.67	3 (18%)	15,22,24	2.21	5 (33%)
3	KCX	CCC	220	6,3	10,11,12	0.64	0	6,12,14	0.97	0
1	CXM	AAA	1	1	9,10,11	2.22	1 (11%)	9,11,13	1.82	1 (11%)

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
3	QO8	CCC	322	3	-	3/7/7/9	0/1/1/1
3	KCX	CCC	220	6,3	-	0/9/10/12	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CXM	AAA	1	1	-	4/9/10/12	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	1	CXM	ON1-CN	6.29	1.33	1.21
3	CCC	322	QO8	O-C	-4.22	1.24	1.42
3	CCC	322	QO8	C2-SG	-2.79	1.73	1.77
3	CCC	322	QO8	CB-SG	-2.39	1.76	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	CCC	322	QO8	CB-SG-C2	5.75	114.67	102.42
1	AAA	1	CXM	ON1-CN-N	-4.72	117.12	124.86
3	CCC	322	QO8	O4-C4-C5	2.74	125.25	118.10
3	CCC	322	QO8	C7-C6-C5	2.58	124.83	121.18
3	CCC	322	QO8	C7-C6-C1	-2.19	115.72	119.57
3	CCC	322	QO8	C1-C2-C3	-2.19	118.19	120.09

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	AAA	1	CXM	O-C-CA-CB
3	CCC	322	QO8	O-C-CA-N
3	CCC	322	QO8	O-C-CA-CB
1	AAA	1	CXM	CB-CA-N-CN
3	CCC	322	QO8	CA-CB-SG-C2
1	AAA	1	CXM	C-CA-N-CN
1	AAA	1	CXM	ON1-CN-N-CA

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	CCC	220	KCX	3	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 34 ligands modelled in this entry, 2 are monoatomic and 1 is modelled with single atom - leaving 31 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	AAA	202	-	3,3,3	0.64	0	2,2,2	0.09	0
4	EDO	CCC	608	-	3,3,3	0.62	0	2,2,2	0.78	0
4	EDO	BBB	202	-	3,3,3	0.43	0	2,2,2	0.63	0
4	EDO	CCC	610	-	3,3,3	0.39	0	2,2,2	0.85	0
5	SO4	CCC	626	-	4,4,4	0.33	0	6,6,6	0.07	0
5	SO4	CCC	621	-	4,4,4	0.32	0	6,6,6	0.22	0
4	EDO	CCC	605	-	3,3,3	1.15	0	2,2,2	0.29	0
5	SO4	BBB	203	-	4,4,4	0.33	0	6,6,6	0.11	0
4	EDO	CCC	615	-	3,3,3	0.60	0	2,2,2	0.30	0
5	SO4	BBB	205	-	4,4,4	0.30	0	6,6,6	0.10	0
4	EDO	AAA	203	-	3,3,3	0.47	0	2,2,2	0.76	0
4	EDO	CCC	611	-	3,3,3	0.63	0	2,2,2	1.76	1 (50%)
4	EDO	CCC	607	-	3,3,3	0.45	0	2,2,2	1.01	0
4	EDO	CCC	613	-	3,3,3	0.46	0	2,2,2	0.46	0
4	EDO	CCC	618	-	3,3,3	0.63	0	2,2,2	0.66	0
4	EDO	CCC	609	-	3,3,3	0.69	0	2,2,2	0.08	0
4	EDO	CCC	612	-	3,3,3	0.72	0	2,2,2	0.20	0
5	SO4	CCC	623	-	4,4,4	0.32	0	6,6,6	0.16	0
4	EDO	CCC	614	-	3,3,3	0.56	0	2,2,2	0.18	0
4	EDO	BBB	201	-	3,3,3	0.50	0	2,2,2	0.93	0
4	EDO	CCC	619	-	3,3,3	0.84	0	2,2,2	0.43	0
4	EDO	CCC	620	-	3,3,3	0.60	0	2,2,2	0.25	0
4	EDO	CCC	604	-	3,3,3	0.30	0	2,2,2	1.40	0
5	SO4	CCC	625	-	4,4,4	0.29	0	6,6,6	0.07	0
4	EDO	AAA	201	-	3,3,3	1.00	0	2,2,2	0.40	0
4	EDO	CCC	606	-	3,3,3	0.64	0	2,2,2	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	CCC	622	-	4,4,4	0.29	0	6,6,6	0.14	0
5	SO4	BBB	204	-	4,4,4	0.29	0	6,6,6	0.18	0
5	SO4	CCC	624	-	4,4,4	0.28	0	6,6,6	0.12	0
4	EDO	CCC	616	-	3,3,3	0.54	0	2,2,2	0.55	0
4	EDO	AAA	204	-	3,3,3	0.45	0	2,2,2	0.52	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	AAA	202	-	-	0/1/1/1	-
4	EDO	CCC	608	-	-	0/1/1/1	-
4	EDO	BBB	202	-	-	1/1/1/1	-
4	EDO	CCC	610	-	-	0/1/1/1	-
4	EDO	CCC	605	-	-	1/1/1/1	-
4	EDO	CCC	615	-	-	1/1/1/1	-
4	EDO	AAA	203	-	-	0/1/1/1	-
4	EDO	CCC	611	-	-	1/1/1/1	-
4	EDO	CCC	607	-	-	1/1/1/1	-
4	EDO	CCC	613	-	-	1/1/1/1	-
4	EDO	CCC	618	-	-	1/1/1/1	-
4	EDO	CCC	609	-	-	1/1/1/1	-
4	EDO	CCC	612	-	-	1/1/1/1	-
4	EDO	CCC	614	-	-	1/1/1/1	-
4	EDO	BBB	201	-	-	1/1/1/1	-
4	EDO	CCC	619	-	-	1/1/1/1	-
4	EDO	CCC	620	-	-	1/1/1/1	-
4	EDO	CCC	604	-	-	1/1/1/1	-
4	EDO	AAA	201	-	-	0/1/1/1	-
4	EDO	CCC	606	-	-	1/1/1/1	-
4	EDO	CCC	616	-	-	0/1/1/1	-
4	EDO	AAA	204	-	-	1/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	CCC	611	EDO	O2-C2-C1	-2.47	93.59	112.39

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	CCC	609	EDO	O1-C1-C2-O2
4	CCC	605	EDO	O1-C1-C2-O2
4	CCC	606	EDO	O1-C1-C2-O2
4	CCC	612	EDO	O1-C1-C2-O2
4	CCC	614	EDO	O1-C1-C2-O2
4	CCC	619	EDO	O1-C1-C2-O2
4	CCC	620	EDO	O1-C1-C2-O2
4	BBB	201	EDO	O1-C1-C2-O2
4	CCC	604	EDO	O1-C1-C2-O2
4	CCC	607	EDO	O1-C1-C2-O2
4	AAA	204	EDO	O1-C1-C2-O2
4	CCC	613	EDO	O1-C1-C2-O2
4	CCC	618	EDO	O1-C1-C2-O2
4	BBB	202	EDO	O1-C1-C2-O2
4	CCC	611	EDO	O1-C1-C2-O2
4	CCC	615	EDO	O1-C1-C2-O2

There are no ring outliers.

8 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	AAA	202	EDO	2	0
4	CCC	608	EDO	5	0
4	CCC	605	EDO	1	0
5	BBB	203	SO4	2	0
4	CCC	611	EDO	7	0
4	CCC	618	EDO	1	0
4	CCC	609	EDO	3	0
5	CCC	622	SO4	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 [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	AAA	99/100 (99%)	-0.81	0 100 100	17, 29, 41, 56	5 (5%)
2	BBB	122/122 (100%)	-0.68	0 100 100	16, 31, 55, 99	4 (3%)
3	CCC	568/570 (99%)	-0.75	0 100 100	12, 29, 54, 98	17 (2%)
All	All	789/792 (99%)	-0.75	0 100 100	12, 29, 52, 99	26 (3%)

There are no RSRZ outliers to report.

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
3	QO8	CCC	322	16/17	0.87	0.16	71,85,96,97	0
3	KCX	CCC	220	12/13	0.98	0.06	24,32,47,48	0
1	CXM	AAA	1	11/12	0.98	0.05	26,29,35,39	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($\text{\AA}^2$)	Q<0.9
4	EDO	CCC	606	4/4	0.73	0.19	56,59,71,74	0
5	SO4	CCC	625	5/5	0.76	0.13	85,86,93,93	5
5	SO4	BBB	205	5/5	0.79	0.12	99,117,119,120	0
5	SO4	BBB	203	5/5	0.79	0.14	88,122,137,150	0
4	EDO	CCC	612	4/4	0.80	0.15	53,63,64,68	0
5	SO4	CCC	626	5/5	0.81	0.09	89,89,93,93	5
5	SO4	CCC	621	5/5	0.84	0.15	89,90,94,108	0
4	EDO	CCC	605	4/4	0.84	0.18	35,38,44,46	0
4	EDO	CCC	620	4/4	0.84	0.15	69,74,74,76	0
4	EDO	CCC	611	4/4	0.85	0.18	43,51,54,61	0
4	EDO	BBB	202	4/4	0.85	0.19	67,70,88,98	0
5	SO4	CCC	624	5/5	0.86	0.22	113,113,134,135	5
4	EDO	CCC	618	4/4	0.87	0.12	59,67,72,74	0
5	SO4	BBB	204	5/5	0.87	0.10	84,87,103,112	0
5	SO4	CCC	623	5/5	0.87	0.20	102,119,144,147	0
4	EDO	CCC	615	4/4	0.88	0.14	48,53,59,59	0
4	EDO	CCC	619	4/4	0.90	0.14	56,61,61,62	0
4	EDO	CCC	604	4/4	0.91	0.11	46,49,50,54	0
4	EDO	CCC	616	4/4	0.92	0.11	48,51,51,54	0
4	EDO	CCC	609	4/4	0.93	0.13	43,51,58,81	0
4	EDO	AAA	204	4/4	0.93	0.13	59,60,62,63	0
5	SO4	CCC	622	5/5	0.93	0.08	78,84,96,98	0
4	EDO	CCC	607	4/4	0.93	0.09	37,38,44,47	0
4	EDO	CCC	613	4/4	0.93	0.10	54,59,61,61	0
4	EDO	CCC	614	4/4	0.93	0.14	54,55,61,64	0
4	EDO	CCC	608	4/4	0.93	0.20	35,43,44,49	0
4	EDO	AAA	202	4/4	0.94	0.08	40,41,41,51	0
4	EDO	CCC	610	4/4	0.94	0.10	41,45,45,56	0
4	EDO	AAA	201	4/4	0.95	0.09	32,35,35,43	0
4	EDO	BBB	201	4/4	0.95	0.11	46,48,54,58	0
4	EDO	AAA	203	4/4	0.98	0.05	36,37,38,40	0
7	OH	CCC	603	1/1	0.99	0.04	35,35,35,35	0
6	NI	CCC	602	1/1	1.00	0.01	35,35,35,35	0
6	NI	CCC	601	1/1	1.00	0.01	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.