



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 20, 2024 – 08:38 pm GMT

PDB ID : 6ZNY
Title : 1.50 Å resolution 3-methylcatechol (3-methylbenzene-1,2-diol) inhibited Sporosarcina pasteurii urease
Authors : Mazzei, L.; Cianci, M.; Musiani, F.; Ciurli, S.
Deposited on : 2020-07-07
Resolution : 1.50 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

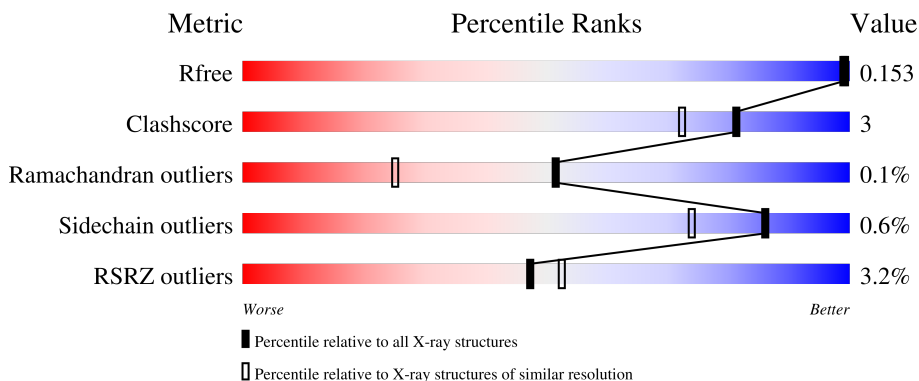
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

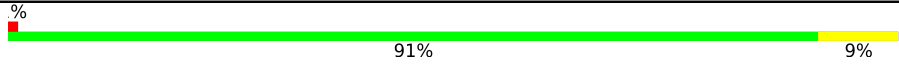
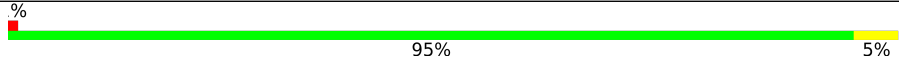
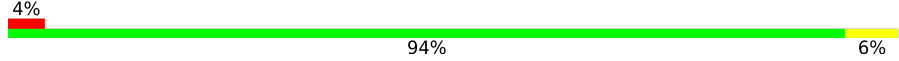
The reported resolution of this entry is 1.50 Å.

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	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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	
2	BBB	122	
3	CCC	570	

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	605	-	-	X	-
4	EDO	CCC	606	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 7184 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	AAA	100	851	534	146	163	8	0	9	0

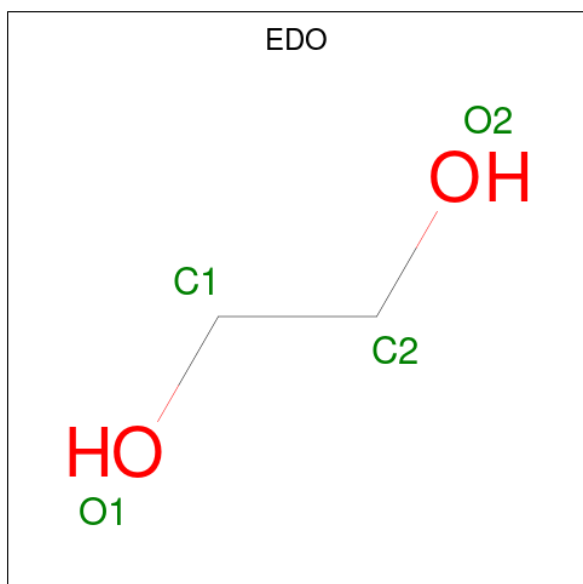
- Molecule 2 is a protein called Urease subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	BBB	122	967	597	173	196	1	0	2	0

- Molecule 3 is a protein called Urease subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	CCC	570	4497	2824	774	870	29	0	23	0

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



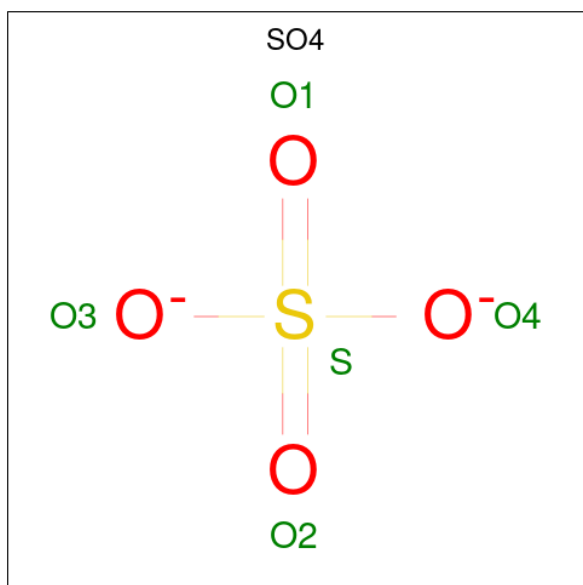
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	AAA	1	Total C O 4 2 2	0	0
4	BBB	1	Total C O 4 2 2	0	0
4	BBB	1	Total C O 4 2 2	0	0
4	CCC	1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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
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 (three-letter code: SO4) (formula: O₄S).



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

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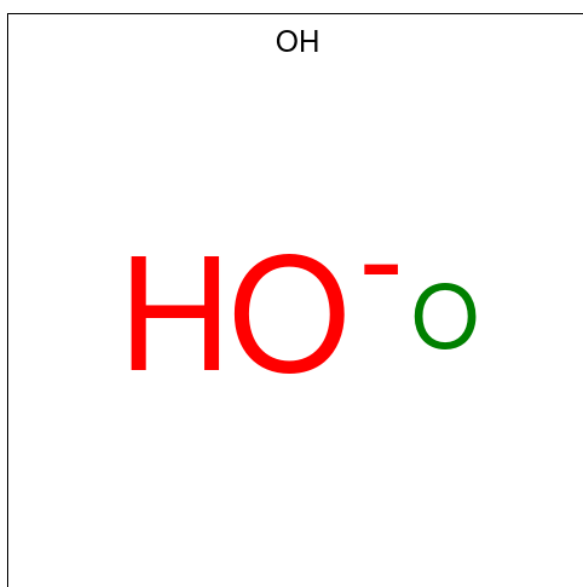
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	CCC	1	Total O S 5 4 1	0	0
5	CCC	1	Total O S 5 4 1	0	0

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

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

- Molecule 7 is HYDROXIDE ION (three-letter code: OH) (formula: HO).



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

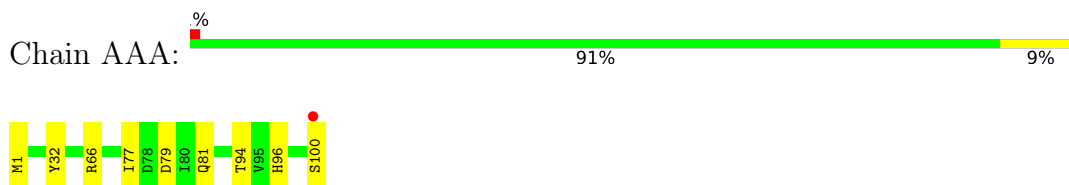
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	AAA	89	Total O 89 89	0	0
8	BBB	140	Total O 140 140	0	0
8	CCC	486	Total O 486 486	0	0

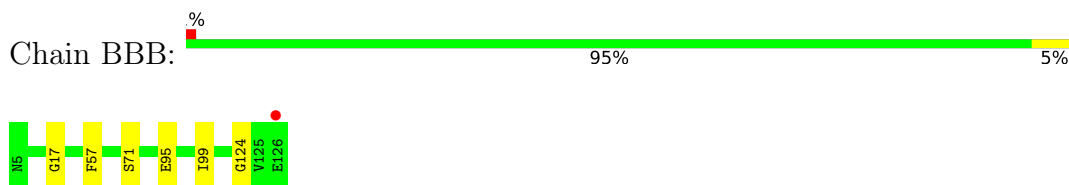
3 Residue-property plots [i](#)

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.

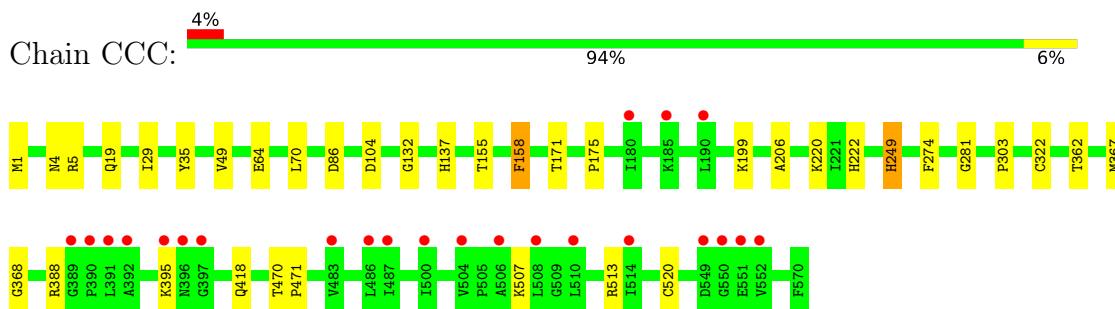
- Molecule 1: Urease subunit gamma



- 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	131.58Å 131.58Å 189.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	94.86 – 1.50 94.68 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (94.86-1.50) 99.6 (94.68-1.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 1.50Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.133 , 0.152 0.134 , 0.153	Depositor DCC
R_{free} test set	7683 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	16.5	Xtrriage
Anisotropy	0.659	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7184	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% 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: SO4, KCX, EDO, QNW, CXM, NI, 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	0.70	0/858	0.82	0/1152
2	BBB	0.73	0/979	0.81	0/1316
3	CCC	0.72	0/4574	0.85	3/6190 (0.0%)
All	All	0.72	0/6411	0.84	3/8658 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	CCC	388	ARG	NE-CZ-NH1	6.17	123.39	120.30
3	CCC	520[A]	CYS	CB-CA-C	-5.94	98.52	110.40
3	CCC	520[B]	CYS	CB-CA-C	-5.94	98.52	110.40

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	AAA	851	0	864	6	0
2	BBB	967	0	946	4	0
3	CCC	4497	0	4477	27	0
4	AAA	28	0	42	2	0
4	BBB	8	0	12	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	CCC	60	0	90	8	0
5	AAA	10	0	0	0	0
5	BBB	15	0	0	0	0
5	CCC	30	0	0	0	0
6	CCC	2	0	0	0	0
7	CCC	1	0	0	0	0
8	AAA	89	0	0	0	0
8	BBB	140	0	0	1	0
8	CCC	486	0	0	3	1
All	All	7184	0	6431	34	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 3.

All (34) 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:64[B]:GLU:HG2	4:CCC:605:EDO:C1	2.14	0.76
3:CCC:64[B]:GLU:HG2	4:CCC:605:EDO:H11	1.71	0.73
1:AAA:32:TYR:OH	4:AAA:205:EDO:H22	1.90	0.71
8:BBB:301:HOH:O	3:CCC:4[A]:ASN:HB2	1.93	0.69
3:CCC:206:ALA:HB2	4:CCC:607:EDO:H12	1.76	0.67
3:CCC:513[A]:ARG:NH1	8:CCC:704:HOH:O	2.30	0.65
2:BBB:71:SER:OG	3:CCC:49[A]:VAL:HG21	2.01	0.60
3:CCC:64[B]:GLU:CG	4:CCC:605:EDO:H11	2.32	0.59
3:CCC:64[B]:GLU:HG2	4:CCC:605:EDO:H12	1.85	0.58
3:CCC:367[B]:MET:HG2	3:CCC:367[B]:MET:O	2.02	0.58
2:BBB:17:GLY:O	3:CCC:5[A]:ARG:HG2	2.06	0.56
1:AAA:77:ILE:HA	4:AAA:205:EDO:H21	1.91	0.53
1:AAA:66[B]:ARG:NH1	1:AAA:100:SER:OXT	2.44	0.51
3:CCC:64[B]:GLU:CG	4:CCC:605:EDO:C1	2.88	0.50
3:CCC:35:TYR:CD2	4:CCC:611:EDO:H22	2.47	0.49
2:BBB:95:GLU:O	3:CCC:104:ASP:HB3	2.14	0.48
1:AAA:79[B]:ASP:OD1	1:AAA:96[B]:HIS:ND1	2.41	0.47
3:CCC:132:GLY:HA3	3:CCC:155:THR:OG1	2.15	0.47
3:CCC:158:PHE:CE2	3:CCC:418[A]:GLN:CG	2.99	0.46
3:CCC:249:HIS:CE1	3:CCC:281:GLY:HA3	2.50	0.45
1:AAA:79[B]:ASP:CG	1:AAA:96[B]:HIS:HD1	2.18	0.44
1:AAA:81[B]:GLN:HG2	1:AAA:94:THR:OG1	2.18	0.44
3:CCC:206:ALA:CB	4:CCC:607:EDO:H12	2.46	0.43
3:CCC:470:THR:N	3:CCC:471:PRO:CD	2.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:CCC:362:THR:O	3:CCC:368:GLY:HA3	2.19	0.43
3:CCC:303:PRO:HG3	3:CCC:368:GLY:HA2	2.01	0.42
3:CCC:64[B]:GLU:CD	8:CCC:707:HOH:O	2.56	0.42
3:CCC:137:HIS:CE1	3:CCC:274:PHE:CD2	3.08	0.41
2:BBB:57:PHE:O	2:BBB:124:GLY:HA3	2.20	0.41
3:CCC:175:PRO:HB3	3:CCC:199:LYS:HE3	2.03	0.41
3:CCC:70:LEU:HD11	3:CCC:86:ASP:HB3	2.03	0.41
3:CCC:19:GLN:HA	3:CCC:29:ILE:O	2.21	0.41
3:CCC:507[A]:LYS:HD3	8:CCC:1087:HOH:O	2.20	0.41
3:CCC:171:THR:HG22	3:CCC:222:HIS:CG	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:CCC:1123:HOH:O	8:CCC:1123:HOH:O[12_565]	2.17	0.03

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	107/100 (107%)	107 (100%)	0	0	100	100
2	BBB	122/122 (100%)	117 (96%)	4 (3%)	1 (1%)	19	5
3	CCC	588/570 (103%)	569 (97%)	19 (3%)	0	100	100
All	All	817/792 (103%)	793 (97%)	23 (3%)	1 (0%)	51	25

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	BBB	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	AAA	92/83 (111%)	92 (100%)	0	100	100
2	BBB	103/101 (102%)	103 (100%)	0	100	100
3	CCC	480/457 (105%)	475 (99%)	5 (1%)	76	57
All	All	675/641 (105%)	670 (99%)	5 (1%)	86	69

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

Mol	Chain	Res	Type
3	CCC	1[A]	MET
3	CCC	1[B]	MET
3	CCC	158	PHE
3	CCC	249	HIS
3	CCC	395	LYS

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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	KCX	CCC	220	6,3	9,11,12	0.52	0	5,12,14	1.64	2 (40%)
1	CXM	AAA	1	1	8,10,11	2.04	1 (12%)	7,11,13	1.60	1 (14%)
3	QNW	CCC	322	3	14,15,16	1.12	1 (7%)	13,20,22	1.55	2 (15%)

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	KCX	CCC	220	6,3	-	0/9/10/12	-
1	CXM	AAA	1	1	-	4/9/10/12	-
3	QNW	CCC	322	3	-	0/5/7/9	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	1	CXM	ON1-CN	5.24	1.31	1.21
3	CCC	322	QNW	C2-SG	-3.10	1.72	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	CCC	322	QNW	C6-C5-C4	3.85	121.47	117.82
1	AAA	1	CXM	ON1-CN-N	-3.65	118.86	124.85
3	CCC	322	QNW	C1-C6-C5	-3.16	117.17	121.97
3	CCC	220	KCX	CE-NZ-CX	2.57	126.02	121.89
3	CCC	220	KCX	OQ1-CX-NZ	-2.54	121.02	124.96

There are no chirality outliers.

All (4) torsion outliers are listed below:

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

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 38 ligands modelled in this entry, 2 are monoatomic and 1 is modelled with single atom - leaving 35 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
5	SO4	AAA	202	-	4,4,4	0.34	0	6,6,6	0.25	0
4	EDO	CCC	609	-	3,3,3	0.50	0	2,2,2	0.74	0
4	EDO	BBB	201	-	3,3,3	0.66	0	2,2,2	0.14	0
4	EDO	CCC	617	-	3,3,3	0.48	0	2,2,2	0.33	0
4	EDO	CCC	610	-	3,3,3	0.80	0	2,2,2	0.10	0
4	EDO	CCC	608	-	3,3,3	0.53	0	2,2,2	0.44	0
5	SO4	CCC	613	-	4,4,4	0.33	0	6,6,6	0.11	0
4	EDO	AAA	206	-	3,3,3	0.89	0	2,2,2	0.51	0
4	EDO	AAA	204	-	3,3,3	0.67	0	2,2,2	0.39	0
5	SO4	BBB	204	-	4,4,4	0.42	0	6,6,6	0.10	0
5	SO4	CCC	612	-	4,4,4	0.33	0	6,6,6	0.72	0
4	EDO	BBB	202	-	3,3,3	0.31	0	2,2,2	0.40	0
5	SO4	AAA	203	-	4,4,4	0.35	0	6,6,6	0.06	0
4	EDO	CCC	619	-	3,3,3	1.01	0	2,2,2	0.61	0
4	EDO	AAA	201	-	3,3,3	0.50	0	2,2,2	0.18	0
4	EDO	CCC	606	-	3,3,3	0.65	0	2,2,2	0.39	0
4	EDO	CCC	616	-	3,3,3	0.33	0	2,2,2	0.49	0
4	EDO	AAA	208	-	3,3,3	0.74	0	2,2,2	0.18	0
5	SO4	CCC	614	-	4,4,4	0.37	0	6,6,6	0.06	0
4	EDO	CCC	618	-	3,3,3	0.31	0	2,2,2	0.41	0
4	EDO	AAA	205	-	3,3,3	1.29	0	2,2,2	1.05	0
5	SO4	BBB	205	-	4,4,4	0.47	0	6,6,6	0.10	0
5	SO4	CCC	624	-	4,4,4	0.46	0	6,6,6	0.24	0
4	EDO	CCC	607	-	3,3,3	0.43	0	2,2,2	0.87	0
5	SO4	CCC	623	-	4,4,4	0.31	0	6,6,6	0.38	0
4	EDO	AAA	209	-	3,3,3	0.51	0	2,2,2	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	EDO	CCC	611	-	3,3,3	0.22	0	2,2,2	1.05	0
4	EDO	CCC	605	-	3,3,3	0.27	0	2,2,2	0.61	0
5	SO4	CCC	622	-	4,4,4	0.61	0	6,6,6	0.17	0
4	EDO	CCC	604	-	3,3,3	0.55	0	2,2,2	0.24	0
4	EDO	CCC	620	-	3,3,3	0.44	0	2,2,2	0.98	0
4	EDO	CCC	621	-	3,3,3	0.50	0	2,2,2	0.70	0
4	EDO	AAA	207	-	3,3,3	0.72	0	2,2,2	1.17	0
5	SO4	BBB	203	-	4,4,4	0.31	0	6,6,6	0.14	0
4	EDO	CCC	615	-	3,3,3	0.58	0	2,2,2	0.80	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	CCC	609	-	-	0/1/1/1	-
4	EDO	BBB	201	-	-	1/1/1/1	-
4	EDO	CCC	617	-	-	0/1/1/1	-
4	EDO	CCC	610	-	-	0/1/1/1	-
4	EDO	CCC	608	-	-	0/1/1/1	-
4	EDO	AAA	206	-	-	0/1/1/1	-
4	EDO	AAA	204	-	-	0/1/1/1	-
4	EDO	BBB	202	-	-	0/1/1/1	-
4	EDO	CCC	619	-	-	0/1/1/1	-
4	EDO	AAA	201	-	-	1/1/1/1	-
4	EDO	CCC	606	-	-	0/1/1/1	-
4	EDO	CCC	616	-	-	0/1/1/1	-
4	EDO	AAA	208	-	-	1/1/1/1	-
4	EDO	CCC	618	-	-	0/1/1/1	-
4	EDO	AAA	205	-	-	1/1/1/1	-
4	EDO	CCC	607	-	-	1/1/1/1	-
4	EDO	AAA	209	-	-	0/1/1/1	-
4	EDO	CCC	611	-	-	0/1/1/1	-
4	EDO	CCC	605	-	-	0/1/1/1	-
4	EDO	CCC	604	-	-	1/1/1/1	-
4	EDO	CCC	620	-	-	1/1/1/1	-
4	EDO	CCC	621	-	-	0/1/1/1	-
4	EDO	AAA	207	-	-	1/1/1/1	-
4	EDO	CCC	615	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	AAA	205	EDO	O1-C1-C2-O2
4	AAA	208	EDO	O1-C1-C2-O2
4	BBB	201	EDO	O1-C1-C2-O2
4	AAA	201	EDO	O1-C1-C2-O2
4	CCC	604	EDO	O1-C1-C2-O2
4	CCC	607	EDO	O1-C1-C2-O2
4	AAA	207	EDO	O1-C1-C2-O2
4	CCC	620	EDO	O1-C1-C2-O2

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	AAA	205	EDO	2	0
4	CCC	607	EDO	2	0
4	CCC	611	EDO	1	0
4	CCC	605	EDO	5	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

6.1 Protein, DNA and RNA chains

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%)	-0.56	1 (1%) 82 85	16, 20, 27, 46	0
2	BBB	122/122 (100%)	-0.39	1 (0%) 86 89	17, 22, 38, 66	0
3	CCC	568/570 (99%)	-0.08	23 (4%) 38 42	15, 19, 32, 70	0
All	All	789/792 (99%)	-0.19	25 (3%) 47 52	15, 19, 32, 70	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	CCC	396	ASN	5.1
3	CCC	391	LEU	5.0
2	BBB	126	GLU	4.3
3	CCC	392	ALA	4.1
3	CCC	390	PRO	4.1
3	CCC	508	LEU	3.7
3	CCC	550	GLY	3.6
3	CCC	510	LEU	3.5
3	CCC	483	VAL	3.5
3	CCC	395	LYS	3.5
3	CCC	190	LEU	3.3
3	CCC	486	LEU	3.3
3	CCC	487	ILE	3.2
3	CCC	514	ILE	3.0
3	CCC	551	GLU	2.7
3	CCC	389	GLY	2.6
3	CCC	506	ALA	2.6
3	CCC	549	ASP	2.6
3	CCC	185[A]	LYS	2.5
3	CCC	552	VAL	2.4
1	AAA	100	SER	2.4
3	CCC	397	GLY	2.4
3	CCC	504	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
3	CCC	500	ILE	2.2
3	CCC	180	ILE	2.1

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	AAA	1	11/12	0.97	0.07	19,22,27,27	0
3	QNW	CCC	322	15/16	0.97	0.07	23,27,30,31	0
3	KCX	CCC	220	12/13	0.98	0.06	15,16,18,20	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(Å ²)	Q<0.9
4	EDO	BBB	201	4/4	0.55	0.22	57,60,61,68	0
4	EDO	AAA	201	4/4	0.65	0.21	72,73,73,74	0
4	EDO	AAA	209	4/4	0.70	0.14	66,66,66,68	0
4	EDO	AAA	207	4/4	0.71	0.19	38,47,49,58	0
4	EDO	CCC	606	4/4	0.73	0.70	56,57,61,70	0
4	EDO	CCC	621	4/4	0.80	0.23	46,46,50,51	0
4	EDO	CCC	609	4/4	0.82	0.26	43,47,49,65	0
4	EDO	CCC	610	4/4	0.83	0.17	35,36,37,42	0
4	EDO	CCC	604	4/4	0.84	0.10	48,53,57,57	0
5	SO4	AAA	202	5/5	0.84	0.24	37,60,75,84	0
4	EDO	CCC	607	4/4	0.86	0.21	44,46,49,51	0
4	EDO	CCC	611	4/4	0.87	0.20	51,55,56,57	0
4	EDO	CCC	615	4/4	0.87	0.15	32,36,40,45	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	SO4	CCC	624	5/5	0.87	0.27	36,50,58,58	5
4	EDO	CCC	620	4/4	0.88	0.23	35,39,41,42	0
4	EDO	AAA	206	4/4	0.88	0.12	37,42,45,46	0
4	EDO	CCC	605	4/4	0.89	0.17	37,40,42,45	0
5	SO4	AAA	203	5/5	0.89	0.20	84,84,89,90	5
4	EDO	CCC	617	4/4	0.89	0.12	37,43,46,48	0
5	SO4	CCC	613	5/5	0.90	0.11	68,68,70,71	5
5	SO4	BBB	205	5/5	0.91	0.24	34,51,53,54	5
4	EDO	CCC	608	4/4	0.91	0.16	43,45,46,47	0
5	SO4	CCC	614	5/5	0.91	0.28	79,86,87,87	5
4	EDO	AAA	208	4/4	0.91	0.20	36,47,49,60	0
4	EDO	AAA	205	4/4	0.92	0.09	25,27,30,35	0
5	SO4	CCC	622	5/5	0.92	0.16	27,49,54,60	5
5	SO4	BBB	204	5/5	0.92	0.22	44,54,57,62	5
4	EDO	BBB	202	4/4	0.94	0.16	37,37,38,47	0
7	OH	CCC	603	1/1	0.94	0.07	18,18,18,18	0
5	SO4	BBB	203	5/5	0.95	0.23	52,57,61,63	0
4	EDO	CCC	619	4/4	0.96	0.07	31,32,33,36	0
5	SO4	CCC	623	5/5	0.96	0.21	25,25,37,39	5
4	EDO	CCC	618	4/4	0.97	0.19	29,33,34,36	0
5	SO4	CCC	612	5/5	0.97	0.09	33,34,38,41	0
4	EDO	AAA	204	4/4	0.98	0.09	22,24,26,26	0
4	EDO	CCC	616	4/4	0.99	0.09	33,33,35,36	0
6	NI	CCC	602	1/1	1.00	0.03	18,18,18,18	0
6	NI	CCC	601	1/1	1.00	0.03	19,19,19,19	0

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.