Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2024 – 03:19 PM JST

PDB ID	:	4ZK8
Title	:	Copper-containing nitrite reductase from thermophilic bacterium Geobacillus thermodenitrificans (Re-refined)
Authors	:	Fukuda, Y.; Inoue, T.
Deposited on	:	2015-04-30
Resolution	:	1.15 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:

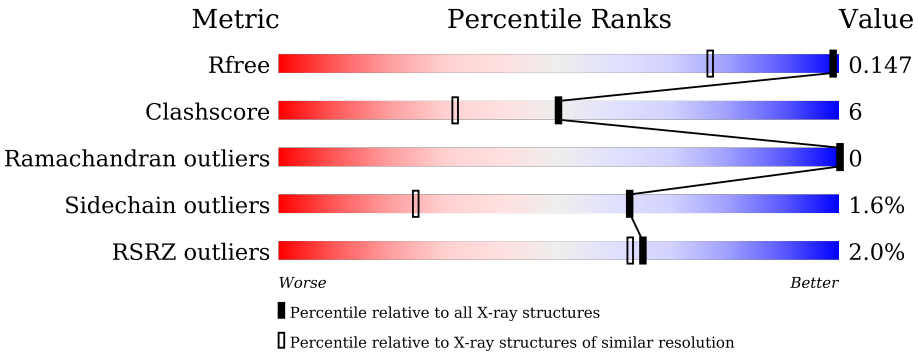
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.15 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	1492 (1.18-1.10)
Clashscore	141614	1537 (1.18-1.10)
Ramachandran outliers	138981	1483 (1.18-1.10)
Sidechain outliers	138945	1480 (1.18-1.10)
RSRZ outliers	127900	1464 (1.18-1.10)

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 <=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	323	2% 83% 7% • 9%

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	MPD	A	423	-	-	X	-
5	PEG	A	424[B]	-	X	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	TRS	A	429	-	X	-	-

2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 3011 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 Nitrite reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	294	Total	C	N	O	S	0	28	0
			2524	1612	428	465	19			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP A4IL26

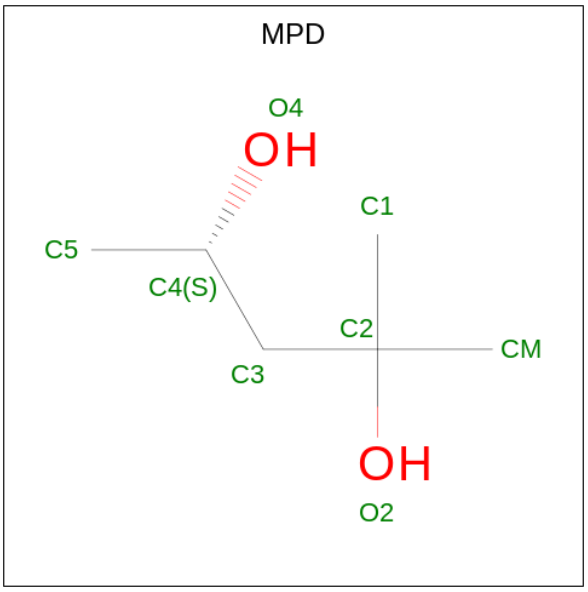
- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	18	Total	Cu	0	10
			25	25		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

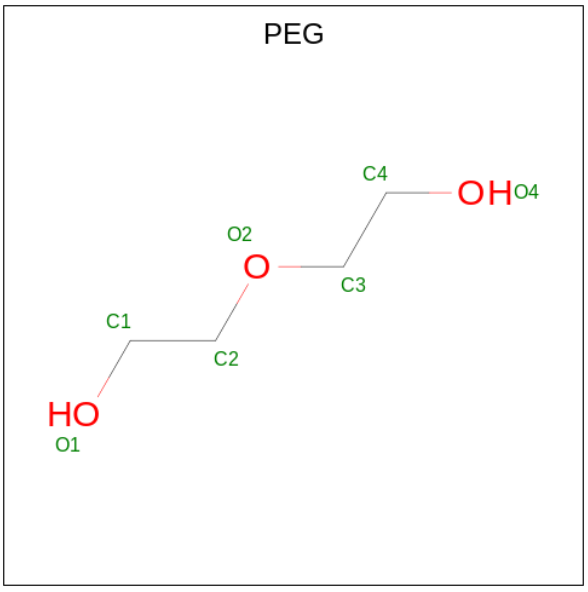
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	1
			3	3		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



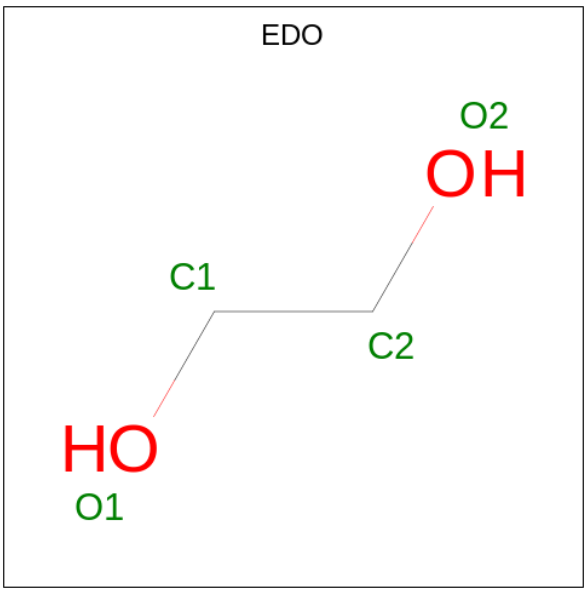
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 8 6 2	0	0
4	A	1	Total C O 8 6 2	0	1
4	A	1	Total C O 8 6 2	0	0

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



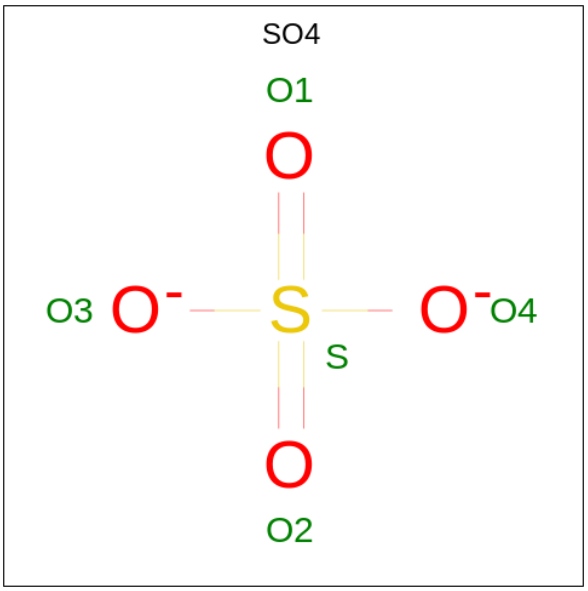
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 7 4 3	0	1

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



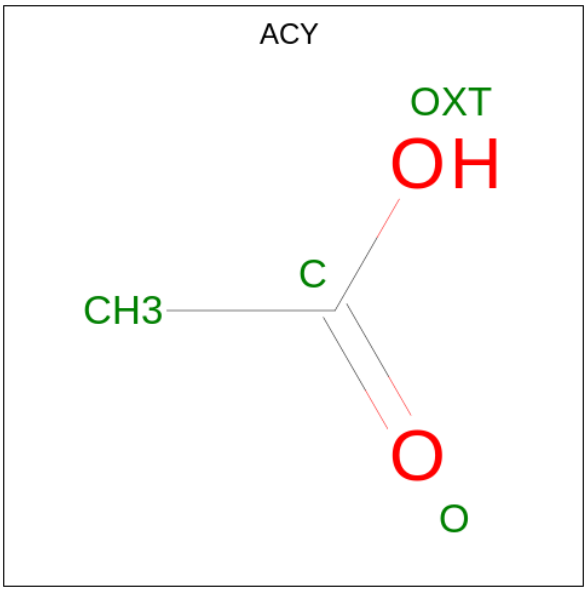
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
6	A	1	4	2	2	0	1

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



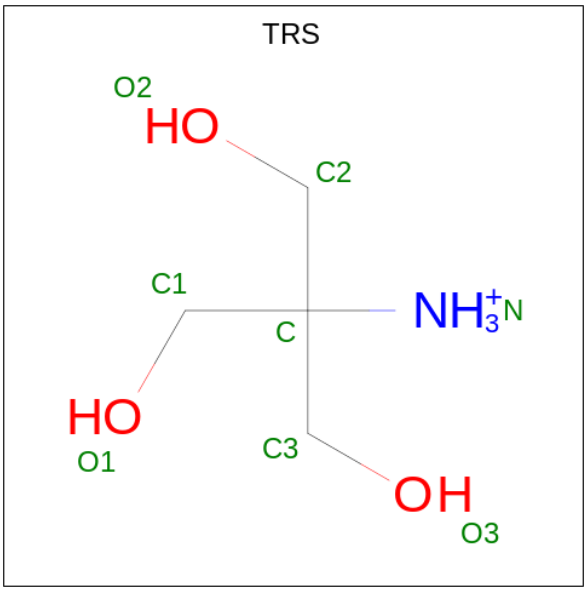
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
7	A	1	5	4	1	0	0

- Molecule 8 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C O 4 2 2	0	0
8	A	1	Total C O 4 2 2	0	0

- Molecule 9 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C N O 8 4 1 3	0	0

- Molecule 10 is water.

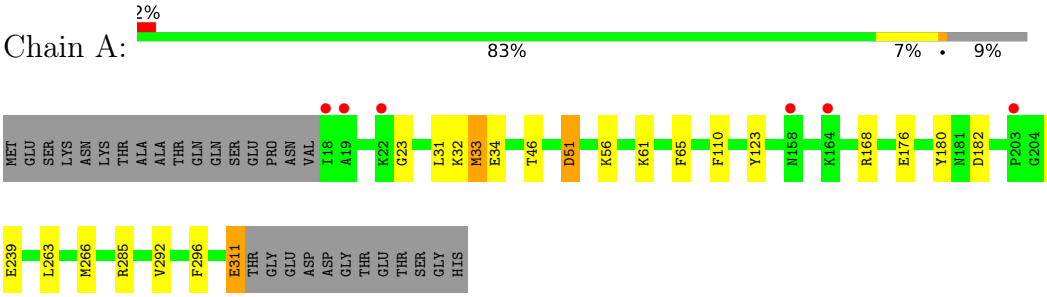
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	376	Total	O	0	40
			403	403		

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: Nitrite reductase

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	115.05Å 115.05Å 84.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	22.06 – 1.15 22.06 – 1.15	Depositor EDS
% Data completeness (in resolution range)	95.6 (22.06-1.15) 95.6 (22.06-1.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	1.66 (at 1.15Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.120 , 0.146 0.121 , 0.147	Depositor DCC
R_{free} test set	7102 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	9.8	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 78.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.036 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	3011	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.42% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: MPD, TRS, PEG, SO4, EDO, CU, ACY, CL

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	1.19	1/2592 (0.0%)	1.09	12/3518 (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
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	311	GLU	CD-OE2	5.16	1.31	1.25

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	229	ARG	NE-CZ-NH2	-7.72	116.44	120.30
1	A	182	ASP	CB-CG-OD1	7.62	125.15	118.30
1	A	180[A]	TYR	CB-CG-CD1	-6.77	116.94	121.00
1	A	180[B]	TYR	CB-CG-CD1	-6.77	116.94	121.00
1	A	168	ARG	NE-CZ-NH1	-6.47	117.06	120.30
1	A	110	PHE	CB-CG-CD1	6.30	125.21	120.80
1	A	51	ASP	CB-CG-OD1	6.25	123.93	118.30
1	A	168	ARG	NE-CZ-NH2	5.84	123.22	120.30
1	A	285	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	A	33[A]	MET	CG-SD-CE	5.44	108.90	100.20
1	A	33[B]	MET	CG-SD-CE	5.44	108.90	100.20
1	A	205	ASP	CB-CG-OD1	5.32	123.09	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	292	VAL	Peptide

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	2524	0	2473	23	0
2	A	25	0	0	1	0
3	A	3	0	0	0	0
4	A	24	0	41	10	0
5	A	7	0	9	10	0
6	A	4	0	6	0	0
7	A	5	0	0	0	0
8	A	8	0	6	0	0
9	A	8	0	12	0	0
10	A	403	0	0	7	0
All	All	3011	0	2547	30	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 6.

All (30) 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:A:23:GLY:H	4:A:423:MPD:H11	1.37	0.87
1:A:23:GLY:HA2	4:A:423:MPD:H32	1.64	0.78
1:A:263:LEU:HD13	1:A:266[B]:MET:SD	2.29	0.72
1:A:34:GLU:HG2	5:A:424[B]:PEG:C4	2.25	0.67
2:A:413[C]:CU:CU	10:A:501:HOH:O	1.45	0.67
1:A:61:LYS:NZ	4:A:423:MPD:H31	2.11	0.64
1:A:61:LYS:HZ2	4:A:423:MPD:H31	1.62	0.64
5:A:424[B]:PEG:C2	10:A:796[B]:HOH:O	2.49	0.61
1:A:33[B]:MET:H	5:A:424[B]:PEG:H11	1.67	0.59
5:A:424[B]:PEG:H22	10:A:796[B]:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239[B]:GLU:OE2	10:A:501:HOH:O	2.15	0.58
4:A:423:MPD:H4	10:A:731:HOH:O	2.04	0.57
1:A:34:GLU:HG2	5:A:424[B]:PEG:H41	1.87	0.57
1:A:34:GLU:HG2	5:A:424[B]:PEG:H42	1.86	0.55
1:A:33[B]:MET:H	5:A:424[B]:PEG:C1	2.22	0.53
1:A:31:LEU:HD21	1:A:33[B]:MET:CE	2.39	0.52
1:A:33[B]:MET:N	5:A:424[B]:PEG:H11	2.25	0.52
1:A:263:LEU:CD1	1:A:266[B]:MET:SD	2.98	0.51
1:A:32[B]:LYS:HB3	5:A:424[B]:PEG:H31	1.93	0.50
4:A:422[A]:MPD:O4	4:A:422[A]:MPD:H12	2.13	0.48
1:A:32[A]:LYS:HB2	1:A:32[A]:LYS:HE2	1.75	0.45
1:A:51:ASP:OD2	4:A:423:MPD:H51	2.17	0.45
4:A:422[A]:MPD:H52	10:A:866:HOH:O	2.17	0.44
1:A:56[A]:LYS:HG2	10:A:764:HOH:O	2.17	0.44
1:A:32[B]:LYS:HA	5:A:424[B]:PEG:H11	2.00	0.43
1:A:46:THR:O	1:A:65:PHE:HA	2.20	0.41
1:A:31:LEU:HD21	1:A:33[B]:MET:HE3	2.02	0.41
4:A:422[A]:MPD:O4	4:A:422[A]:MPD:C1	2.69	0.41
1:A:23:GLY:N	4:A:423:MPD:H11	2.19	0.41
1:A:33[B]:MET:HE2	1:A:33[B]:MET:HB2	1.91	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	A	320/323 (99%)	314 (98%)	6 (2%)	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	281/277 (101%)	277 (99%)	4 (1%)	67 29

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	TYR
1	A	176	GLU
1	A	296	PHE
1	A	311	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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 37 ligands modelled in this entry, 28 are monoatomic - leaving 9 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	MPD	A	422[A]	-	7,7,7	0.98	0	9,10,10	2.63	2 (22%)
6	EDO	A	425[A]	-	3,3,3	1.48	1 (33%)	2,2,2	1.61	1 (50%)
4	MPD	A	423	-	7,7,7	3.33	3 (42%)	9,10,10	1.89	3 (33%)
8	ACY	A	428	-	3,3,3	1.09	0	3,3,3	1.18	0
5	PEG	A	424[B]	-	6,6,6	1.66	1 (16%)	5,5,5	3.42	3 (60%)
8	ACY	A	427	2	3,3,3	0.57	0	3,3,3	0.96	0
9	TRS	A	429	-	7,7,7	1.02	0	9,9,9	3.13	1 (11%)
7	SO4	A	426	-	4,4,4	1.08	0	6,6,6	0.44	0
4	MPD	A	421	2	7,7,7	1.20	1 (14%)	9,10,10	2.22	2 (22%)

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	MPD	A	422[A]	-	-	2/5/5/5	-
6	EDO	A	425[A]	-	-	0/1/1/1	-
4	MPD	A	423	-	-	2/5/5/5	-
5	PEG	A	424[B]	-	-	4/4/4/4	-
9	TRS	A	429	-	-	9/9/9/9	-
4	MPD	A	421	2	-	4/5/5/5	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	423	MPD	O2-C2	-7.02	1.26	1.44
4	A	423	MPD	C3-C2	-4.05	1.42	1.53
5	A	424[B]	PEG	O2-C2	3.28	1.56	1.42
4	A	423	MPD	CM-C2	-2.88	1.43	1.52
6	A	425[A]	EDO	O2-C2	2.25	1.53	1.42
4	A	421	MPD	CM-C2	-2.09	1.45	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
9	A	429	TRS	O3-C3-C	8.85	139.04	111.00
5	A	424[B]	PEG	C3-O2-C2	6.69	142.28	113.29
4	A	422[A]	MPD	CM-C2-C1	-6.14	97.78	110.57
4	A	421	MPD	CM-C2-C1	-5.98	98.11	110.57
4	A	422[A]	MPD	O2-C2-C3	3.77	123.96	109.80
4	A	423	MPD	O2-C2-CM	3.31	118.69	108.08
4	A	423	MPD	CM-C2-C1	-3.04	104.23	110.57
4	A	421	MPD	C1-C2-C3	2.73	122.65	109.96
4	A	423	MPD	O2-C2-C1	-2.61	99.71	108.08
5	A	424[B]	PEG	O1-C1-C2	2.33	125.30	111.81
6	A	425[A]	EDO	O2-C2-C1	-2.16	96.38	111.91
5	A	424[B]	PEG	O2-C3-C4	2.11	119.33	110.07

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	422[A]	MPD	O2-C2-C3-C4
9	A	429	TRS	C3-C-C1-O1
9	A	429	TRS	N-C-C1-O1
9	A	429	TRS	C3-C-C2-O2
5	A	424[B]	PEG	C1-C2-O2-C3
5	A	424[B]	PEG	O1-C1-C2-O2
5	A	424[B]	PEG	O2-C3-C4-O4
9	A	429	TRS	C2-C-C1-O1
9	A	429	TRS	C1-C-C2-O2
9	A	429	TRS	N-C-C2-O2
9	A	429	TRS	C1-C-C3-O3
9	A	429	TRS	C2-C-C3-O3
5	A	424[B]	PEG	C4-C3-O2-C2
4	A	421	MPD	C1-C2-C3-C4
4	A	421	MPD	CM-C2-C3-C4
4	A	423	MPD	CM-C2-C3-C4
4	A	421	MPD	O2-C2-C3-C4
4	A	423	MPD	O2-C2-C3-C4
9	A	429	TRS	N-C-C3-O3
4	A	421	MPD	C2-C3-C4-O4
4	A	422[A]	MPD	C2-C3-C4-O4

There are no ring outliers.

3 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	422[A]	MPD	3	0
4	A	423	MPD	7	0
5	A	424[B]	PEG	10	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, 95^{th} 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	294/323 (91%)	-0.51	6 (2%) 65 63	7, 11, 22, 43	1 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	18	ILE	4.1
1	A	19	ALA	3.3
1	A	164	LYS	2.8
1	A	203	PRO	2.4
1	A	158[A]	ASN	2.3
1	A	22	LYS	2.3

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

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, 95^{th} 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
7	SO4	A	426	5/5	0.86	0.14	60,61,64,66	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
8	ACY	A	427	4/4	0.86	0.12	22,33,34,34	0
6	EDO	A	425[A]	4/4	0.87	0.12	20,21,24,34	4
2	CU	A	417[B]	1/1	0.90	0.10	33,33,33,33	1
2	CU	A	417[A]	1/1	0.90	0.10	31,31,31,31	1
4	MPD	A	422[A]	8/8	0.91	0.21	21,38,44,58	8
4	MPD	A	421	8/8	0.91	0.14	25,28,35,39	0
3	CL	A	420[B]	1/1	0.92	0.14	52,52,52,52	1
3	CL	A	420[A]	1/1	0.92	0.14	50,50,50,50	1
8	ACY	A	428	4/4	0.92	0.18	30,36,39,48	0
9	TRS	A	429	8/8	0.93	0.21	25,49,59,79	0
5	PEG	A	424[B]	7/7	0.95	0.11	15,18,22,22	7
4	MPD	A	423	8/8	0.95	0.17	22,31,45,98	0
2	CU	A	413[B]	1/1	0.97	0.09	12,12,12,12	1
2	CU	A	413[C]	1/1	0.97	0.09	19,19,19,19	1
2	CU	A	406[A]	1/1	0.97	0.10	31,31,31,31	1
2	CU	A	406[C]	1/1	0.97	0.10	17,17,17,17	1
2	CU	A	413[A]	1/1	0.97	0.09	23,23,23,23	1
2	CU	A	408[B]	1/1	0.98	0.12	18,18,18,18	1
2	CU	A	405	1/1	0.99	0.04	17,17,17,17	1
2	CU	A	418[A]	1/1	0.99	0.05	23,23,23,23	1
2	CU	A	418[B]	1/1	0.99	0.05	14,14,14,14	1
2	CU	A	407[A]	1/1	0.99	0.06	16,16,16,16	1
2	CU	A	404	1/1	0.99	0.05	18,18,18,18	1
2	CU	A	414	1/1	0.99	0.08	29,29,29,29	1
2	CU	A	412[A]	1/1	0.99	0.06	34,34,34,34	1
2	CU	A	402	1/1	1.00	0.01	9,9,9,9	0
2	CU	A	403	1/1	1.00	0.03	9,9,9,9	1
2	CU	A	409[A]	1/1	1.00	0.06	16,16,16,16	1
2	CU	A	409[B]	1/1	1.00	0.06	21,21,21,21	1
2	CU	A	415	1/1	1.00	0.05	23,23,23,23	1
2	CU	A	416	1/1	1.00	0.02	17,17,17,17	1
2	CU	A	410[A]	1/1	1.00	0.03	25,25,25,25	1
2	CU	A	411[A]	1/1	1.00	0.04	18,18,18,18	1
2	CU	A	411[B]	1/1	1.00	0.04	34,34,34,34	1
2	CU	A	401	1/1	1.00	0.02	9,9,9,9	0
3	CL	A	419	1/1	1.00	0.02	13,13,13,13	1

6.5 Other polymers (i)

There are no such residues in this entry.