Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 03:49 am BST

PDB ID	:	2WTP
Title	:	Crystal Structure of Cu-form Czce from C. metallidurans CH34
Authors	:	Haertlein, I.; Girard, E.; Sarret, G.; Hazemann, J.; Gourhant, P.; Kahn, R.; Coves, J.
Deposited on	:	2009-09-18
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 following versions of software and data (see references ⓘ) were used in the production of this report:

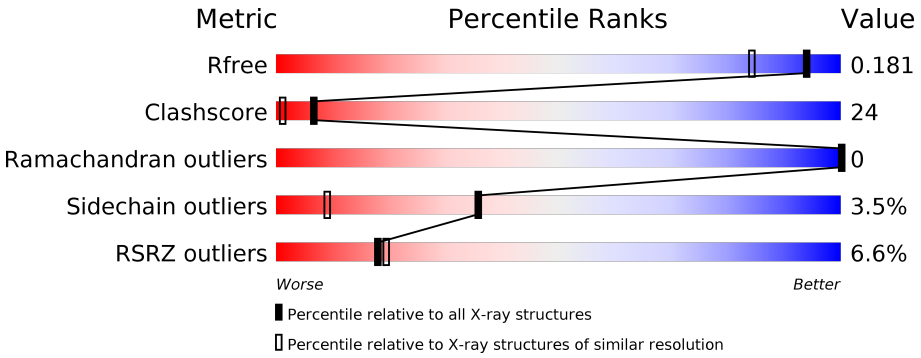
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.50 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 on 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	131	5% 49% 18% • • 29%
1	B	131	5% 53% 12% • 32%

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
2	PEG	A	1107	-	-	X	-
2	PEG	B	1106	-	-	X	-
2	PEG	B	1107	-	-	X	-
3	EDO	A	1108	-	-	X	-

2 Entry composition

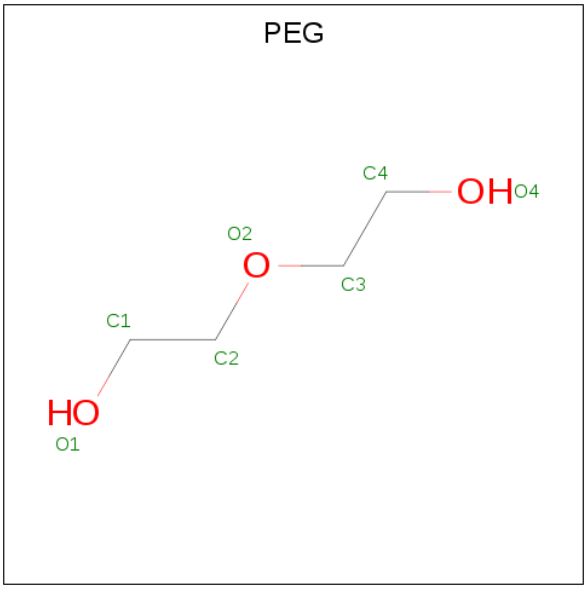
There are 8 unique types of molecules in this entry. The entry contains 1675 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ORF131 PROTEIN.

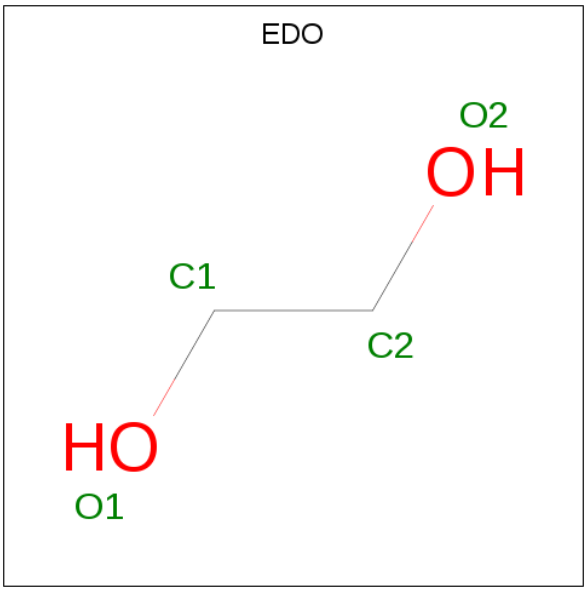
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	93	Total	C	N	O	S	0	6	0
			735	462	131	137	5			
1	B	89	Total	C	N	O	S	0	3	0
			685	430	121	130	4			

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



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

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

- Molecule 4 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Cu 2 2	0	0
4	A	1	Total Cu 1 1	0	0

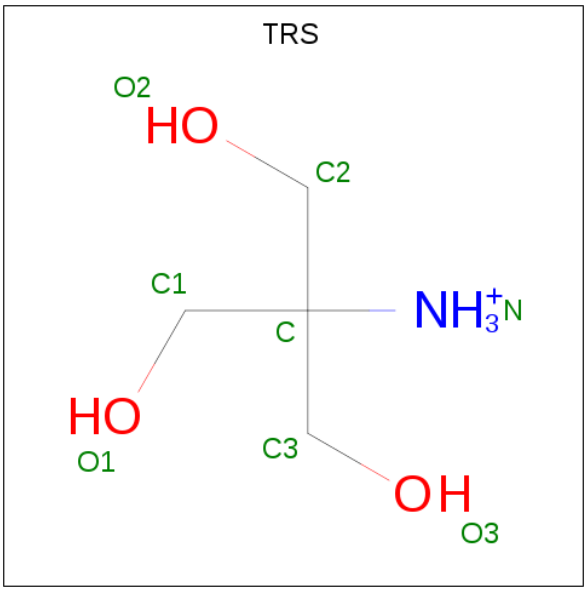
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Mg 1 1	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	5	Total Cl 5 5	0	0
6	A	5	Total Cl 5 5	0	0

- Molecule 7 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
7	B	1	Total	C	N	O	0	1
			16	8	2	6		

- Molecule 8 is water.

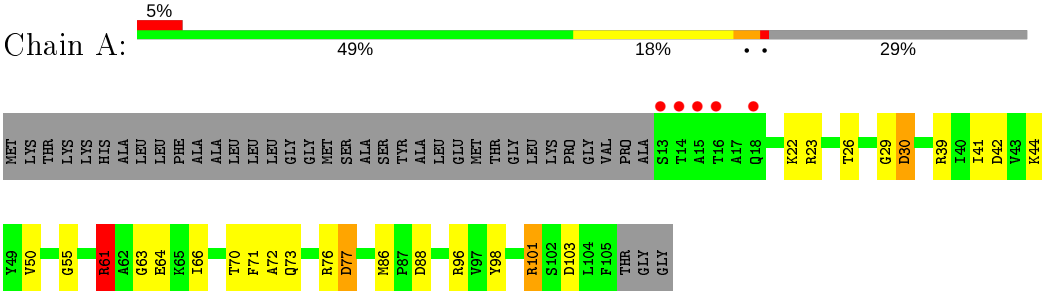
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	106	Total	O	0	0
			106	106		
8	B	94	Total	O	0	0
			94	94		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.

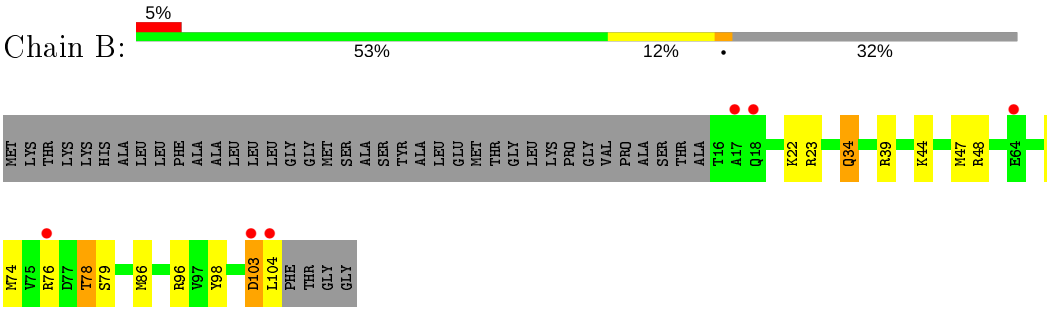
- Molecule 1: ORF131 PROTEIN

Chain A:



- Molecule 1: ORF131 PROTEIN

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	105.00Å 29.53Å 71.17Å 90.00° 113.41° 90.00°	Depositor
Resolution (Å)	19.14 – 1.50 19.14 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.14-1.50) 99.4 (19.14-1.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.82 (at 1.50Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.160 , 0.189 0.155 , 0.181	Depositor DCC
R_{free} test set	1645 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	14.2	Xtriage
Anisotropy	0.157	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.43 , 59.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	1675	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.58% 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: MG, CL, EDO, TRS, PEG, CU

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	1.23	2/765 (0.3%)	1.38	12/1034 (1.2%)
1	B	1.11	1/705 (0.1%)	1.40	9/955 (0.9%)
All	All	1.17	3/1470 (0.2%)	1.39	21/1989 (1.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	34	GLN	CB-CG	-6.93	1.33	1.52
1	A	61[A]	ARG	CG-CD	5.82	1.66	1.51
1	A	61[B]	ARG	CG-CD	5.82	1.66	1.51

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	74	MET	CG-SD-CE	-14.77	76.58	100.20
1	A	48	ARG	NE-CZ-NH2	-11.85	114.38	120.30
1	B	34	GLN	N-CA-CB	-9.78	92.99	110.60
1	A	96	ARG	NE-CZ-NH1	-9.69	115.46	120.30
1	A	48	ARG	NE-CZ-NH1	9.04	124.82	120.30
1	B	39	ARG	NE-CZ-NH2	-8.75	115.92	120.30
1	A	77	ASP	CB-CG-OD1	7.60	125.14	118.30
1	A	23	ARG	NE-CZ-NH1	7.50	124.05	120.30
1	A	30	ASP	CB-CG-OD1	-6.85	112.14	118.30
1	B	48	ARG	NE-CZ-NH2	-6.74	116.93	120.30
1	B	48	ARG	NE-CZ-NH1	6.73	123.67	120.30
1	A	23	ARG	NE-CZ-NH2	-6.56	117.02	120.30
1	A	103	ASP	CB-CG-OD2	6.12	123.81	118.30
1	A	96	ARG	NE-CZ-NH2	6.12	123.36	120.30
1	A	39	ARG	NE-CZ-NH2	-6.06	117.27	120.30
1	B	23	ARG	NE-CZ-NH2	-5.85	117.38	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	98	TYR	CZ-CE2-CD2	-5.79	114.59	119.80
1	B	86	MET	CG-SD-CE	5.62	109.19	100.20
1	A	86	MET	CG-SD-CE	5.58	109.13	100.20
1	A	98	TYR	CZ-CE2-CD2	-5.08	115.23	119.80
1	B	23	ARG	NE-CZ-NH1	5.08	122.84	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	735	0	751	48	2
1	B	685	0	695	19	0
2	A	7	0	9	10	0
2	B	14	0	20	12	0
3	A	4	0	6	12	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
5	A	1	0	0	0	0
6	A	5	0	0	1	0
6	B	5	0	0	1	0
7	B	16	0	21	0	0
8	A	106	0	0	9	5
8	B	94	0	0	6	2
All	All	1675	0	1502	70	6

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 24.

All (70) 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:61[A]:ARG:NH1	1:A:66:ILE:HD11	1.52	1.23
1:A:61[A]:ARG:NH1	8:A:2063:HOH:O	1.74	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ASP:OD1	3:A:1108:EDO:H12	1.58	1.04
1:A:30:ASP:OD1	3:A:1108:EDO:C1	2.07	1.02
1:A:61[A]:ARG:HH12	1:A:66:ILE:HD11	1.22	0.95
1:A:61[A]:ARG:NH2	1:A:64:GLU:OE1	2.01	0.92
1:B:72:ALA:HB1	2:B:1106:PEG:H31	1.50	0.90
2:B:1107:PEG:H42	8:B:2093:HOH:O	1.75	0.85
1:B:22:LYS:HG3	2:B:1107:PEG:H12	1.60	0.82
1:A:101[A]:ARG:CG	1:A:101[A]:ARG:HH11	1.93	0.81
1:A:101[A]:ARG:HH11	1:A:101[A]:ARG:HG3	1.44	0.81
1:A:30:ASP:OD1	3:A:1108:EDO:H11	1.81	0.79
1:B:73:GLN:H	2:B:1106:PEG:H32	1.47	0.79
1:A:41:ILE:HD13	1:A:50[B]:VAL:HG21	1.63	0.78
1:A:70:THR:HG22	2:A:1107:PEG:H32	1.68	0.76
6:B:1114:CL:CL	8:B:2045:HOH:O	2.41	0.74
1:B:73:GLN:N	2:B:1106:PEG:H32	2.03	0.74
1:B:73:GLN:O	2:B:1106:PEG:H21	1.88	0.73
1:A:55:GLY:HA2	2:A:1107:PEG:H22	1.71	0.72
1:A:73:GLN:HG2	2:A:1107:PEG:H21	1.73	0.70
1:A:26:THR:HG23	3:A:1108:EDO:H21	1.74	0.69
1:A:30:ASP:OD2	2:A:1107:PEG:C4	2.41	0.69
1:B:72:ALA:HB1	2:B:1106:PEG:C3	2.23	0.67
1:A:72:ALA:HA	2:A:1107:PEG:H31	1.76	0.66
1:A:29:GLY:O	3:A:1108:EDO:H22	1.96	0.65
1:B:76:ARG:O	1:B:78[B]:THR:HG23	1.96	0.64
1:A:30:ASP:HA	3:A:1108:EDO:H12	1.78	0.64
1:A:30:ASP:OD2	2:A:1107:PEG:H41	1.97	0.64
1:A:101[A]:ARG:CG	1:A:101[A]:ARG:NH1	2.55	0.64
1:B:104:LEU:O	8:B:2090:HOH:O	2.16	0.62
1:A:76:ARG:HD3	8:A:2071:HOH:O	1.98	0.62
1:A:42[A]:ASP:OD2	1:A:44:LYS:HE3	2.00	0.60
1:B:69:TRP:HZ3	2:B:1107:PEG:H21	1.66	0.60
1:A:76:ARG:NH1	8:A:2073:HOH:O	2.35	0.60
1:A:71:PHE:O	2:A:1107:PEG:H22	2.06	0.56
1:B:34:GLN:HB2	8:B:2022:HOH:O	2.06	0.56
1:A:26:THR:HA	3:A:1108:EDO:H22	1.86	0.56
1:A:61[A]:ARG:HH21	1:A:64:GLU:HA	1.70	0.56
1:B:69:TRP:CZ3	2:B:1107:PEG:H21	2.41	0.55
1:B:103:ASP:N	1:B:103:ASP:OD1	2.40	0.53
1:A:71:PHE:O	2:A:1107:PEG:C3	2.57	0.52
1:A:61[A]:ARG:NH2	1:A:64:GLU:HA	2.25	0.52
1:A:72:ALA:CA	2:A:1107:PEG:H31	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:GLN:O	2:B:1106:PEG:C2	2.59	0.51
1:A:26:THR:HG23	3:A:1108:EDO:C2	2.40	0.49
1:A:61[A]:ARG:CZ	1:A:66:ILE:HD11	2.36	0.49
1:A:26:THR:HA	3:A:1108:EDO:C2	2.42	0.48
1:B:78[B]:THR:HG21	8:B:2072:HOH:O	2.12	0.48
1:B:44:LYS:CG	1:B:47:MET:HB3	2.44	0.48
2:B:1107:PEG:C4	8:B:2093:HOH:O	2.49	0.47
1:A:101[A]:ARG:NH2	8:A:2097:HOH:O	2.47	0.47
1:B:44:LYS:HG2	1:B:47:MET:HB3	1.98	0.46
1:A:71:PHE:C	2:A:1107:PEG:H31	2.36	0.46
1:A:44:LYS:H	1:A:47[A]:MET:HE3	1.81	0.45
1:A:41:ILE:CD1	1:A:50[B]:VAL:HG21	2.42	0.45
1:A:77:ASP:OD2	8:A:2075:HOH:O	2.21	0.45
1:A:22:LYS:HG2	8:A:2008:HOH:O	2.16	0.44
1:B:22:LYS:HG3	2:B:1107:PEG:C1	2.40	0.43
1:B:78[A]:THR:HG22	1:B:79:SER:H	1.82	0.43
1:A:44:LYS:NZ	8:A:2043:HOH:O	2.51	0.43
1:B:22:LYS:HE3	1:B:22:LYS:HB2	1.68	0.42
1:A:30:ASP:CA	3:A:1108:EDO:H12	2.46	0.42
1:A:29:GLY:O	3:A:1108:EDO:C2	2.67	0.42
1:A:30:ASP:CG	3:A:1108:EDO:H12	2.35	0.42
1:A:101[B]:ARG:NH1	8:A:2096:HOH:O	2.53	0.41
1:A:88:ASP:HB2	6:A:1115:CL:CL	2.58	0.41
1:A:48:ARG:HD3	8:A:2047:HOH:O	2.21	0.41
1:A:101[A]:ARG:HG2	1:A:101[A]:ARG:NH1	2.36	0.40
1:A:61[A]:ARG:NE	1:A:63:GLY:O	2.53	0.40
1:A:47[A]:MET:HB2	1:A:47[A]:MET:HE3	1.64	0.40

All (6) 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:61[A]:ARG:NH2	8:A:2063:HOH:O[2_656]	1.83	0.37
1:A:61[A]:ARG:CZ	8:A:2063:HOH:O[2_656]	2.02	0.18
8:A:2003:HOH:O	8:B:2052:HOH:O[4_556]	2.07	0.13
8:B:2028:HOH:O	8:B:2092:HOH:O[1_545]	2.14	0.06
8:A:2022:HOH:O	8:A:2023:HOH:O[2_656]	2.16	0.04
8:A:2019:HOH:O	8:A:2047:HOH:O[1_565]	2.16	0.04

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	97/131 (74%)	96 (99%)	1 (1%)	0	100	100
1	B	90/131 (69%)	89 (99%)	1 (1%)	0	100	100
All	All	187/262 (71%)	185 (99%)	2 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	A	79/99 (80%)	75 (95%)	4 (5%)	24	4
1	B	73/99 (74%)	69 (94%)	4 (6%)	21	3
All	All	152/198 (77%)	144 (95%)	8 (5%)	36	3

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61[A]	ARG
1	A	61[B]	ARG
1	A	101[A]	ARG
1	A	101[B]	ARG
1	B	78[A]	THR
1	B	78[B]	THR
1	B	96	ARG
1	B	103	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	34	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 20 ligands modelled in this entry, 14 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	$\#\|Z\| > 2$
2	PEG	B	1106	-	6,6,6	0.65	0	5,5,5	1.33	1 (20%)
2	PEG	B	1107	-	6,6,6	0.72	0	5,5,5	1.71	2 (40%)
2	PEG	A	1107	-	6,6,6	1.26	1 (16%)	5,5,5	3.29	2 (40%)
3	EDO	A	1108	-	3,3,3	0.54	0	2,2,2	0.53	0
7	TRS	B	1105[B]	-	7,7,7	0.39	0	9,9,9	1.04	1 (11%)
7	TRS	B	1105[A]	-	7,7,7	0.49	0	9,9,9	1.33	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
2	PEG	B	1106	-	-	1/4/4/4	-
2	PEG	B	1107	-	-	1/4/4/4	-
2	PEG	A	1107	-	-	1/4/4/4	-
3	EDO	A	1108	-	-	0/1/1/1	-
7	TRS	B	1105[B]	-	-	0/9/9/9	-
7	TRS	B	1105[A]	-	-	1/9/9/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1107	PEG	O1-C1	-2.14	1.31	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1107	PEG	O2-C2-C1	-6.63	80.92	110.07
2	A	1107	PEG	C3-O2-C2	3.05	126.51	113.29
2	B	1107	PEG	O2-C2-C1	2.80	122.38	110.07
7	B	1105[A]	TRS	O1-C1-C	-2.62	102.68	111.00
7	B	1105[B]	TRS	C2-C-N	-2.26	101.23	107.98
2	B	1107	PEG	C3-O2-C2	-2.06	104.37	113.29
2	B	1106	PEG	O2-C2-C1	-2.01	101.24	110.07

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1107	PEG	C1-C2-O2-C3
2	A	1107	PEG	C1-C2-O2-C3
2	B	1106	PEG	C1-C2-O2-C3
7	B	1105[A]	TRS	C2-C-C1-O1

There are no ring outliers.

4 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1106	PEG	6	0
2	B	1107	PEG	6	0
2	A	1107	PEG	10	0
3	A	1108	EDO	12	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

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	93/131 (70%)	-0.02	6 (6%)	18	20	8, 13, 34, 69	0
1	B	89/131 (67%)	0.07	6 (6%)	17	19	9, 19, 41, 53	0
All	All	182/262 (69%)	0.02	12 (6%)	18	19	8, 16, 40, 69	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	15	ALA	5.0
1	A	14	THR	4.3
1	A	13	SER	3.7
1	B	17	ALA	3.3
1	B	64	GLU	3.3
1	A	45	PRO	2.6
1	A	18	GLN	2.3
1	B	18	GLN	2.3
1	B	76	ARG	2.2
1	A	16	THR	2.1
1	B	103	ASP	2.1
1	B	104	LEU	2.1

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 carbohydrates 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, 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
2	PEG	B	1106	7/7	0.85	0.20	33,33,33,33	0
2	PEG	B	1107	7/7	0.85	0.20	39,39,39,39	0
6	CL	B	1114	1/1	0.89	0.11	26,26,26,26	1
3	EDO	A	1108	4/4	0.90	0.31	30,30,30,30	0
2	PEG	A	1107	7/7	0.90	0.20	26,26,26,26	0
7	TRS	B	1105[A]	8/8	0.91	0.23	12,18,24,37	8
7	TRS	B	1105[B]	8/8	0.91	0.23	12,18,24,37	8
6	CL	B	1112	1/1	0.94	0.13	32,32,32,32	1
4	CU	A	1109	1/1	0.94	0.15	21,21,21,21	1
6	CL	B	1111	1/1	0.96	0.10	23,23,23,23	1
4	CU	B	1109	1/1	0.97	0.05	37,37,37,37	1
6	CL	B	1113	1/1	0.98	0.04	28,28,28,28	1
6	CL	A	1115	1/1	0.99	0.08	24,24,24,24	1
6	CL	A	1114	1/1	0.99	0.16	23,23,23,23	1
6	CL	B	1110	1/1	0.99	0.08	23,23,23,23	1
5	MG	A	1110	1/1	1.00	0.11	15,15,15,15	1
4	CU	B	1108	1/1	1.00	0.04	13,13,13,13	1
6	CL	A	1111	1/1	1.00	0.04	9,9,9,9	0
6	CL	A	1113	1/1	1.00	0.03	11,11,11,11	0
6	CL	A	1112	1/1	1.00	0.04	22,22,22,22	1

6.5 Other polymers (i)

There are no such residues in this entry.