



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 24, 2021 – 08:08 PM JST

PDB ID : 7F6Q
Title : Crystal structure of metal-citrate-binding mutant (S79A) protein (MctA) of ABC transporter in apo state
Authors : Kanaujia, S.P.; Mandal, S.K.
Deposited on : 2021-06-25
Resolution : 1.63 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 : 2.23.2
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) : 2.23.2

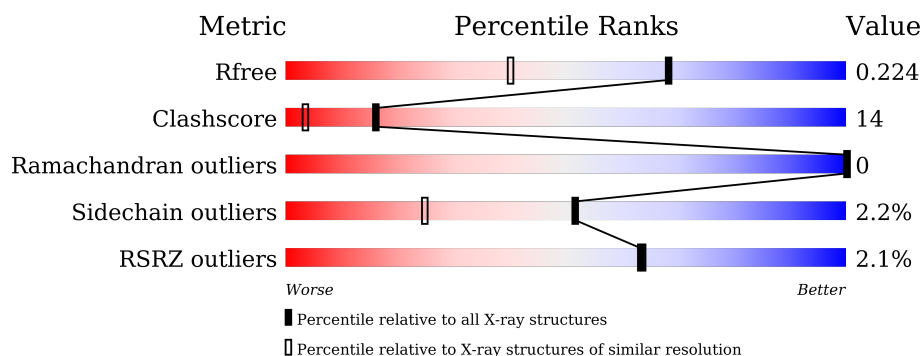
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.63 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	3122 (1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	342	

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	CO2	A	406	-	-	X	-
5	ACT	A	407	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	EDO	A	409	-	-	X	-
6	EDO	A	410	-	-	X	-
6	EDO	A	411	-	-	X	-
7	PDO	A	412	-	-	X	-
8	PEG	A	414	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 3128 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 Iron ABC transporter, periplasmic iron-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	11	0
			2726	1752	472	498	4			

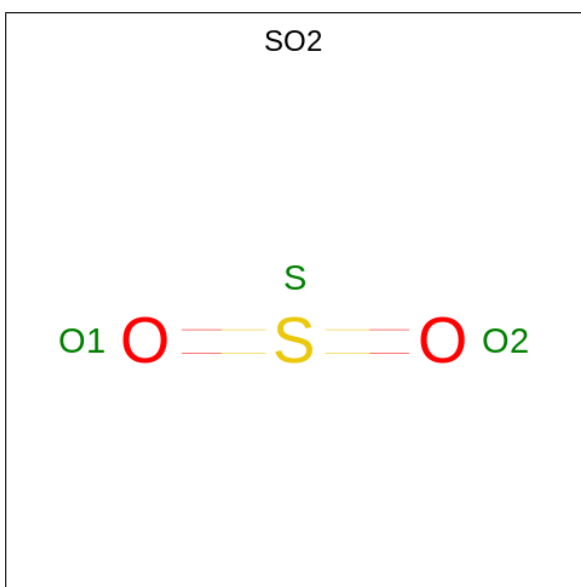
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP Q53VZ2
A	0	MET	-	expression tag	UNP Q53VZ2
A	79	ALA	SER	engineered mutation	UNP Q53VZ2

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

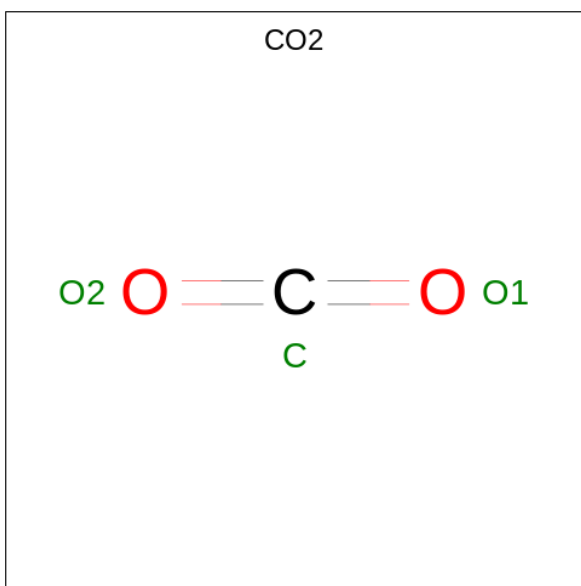
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Cl	0	0
			4	4		

- Molecule 3 is SULFUR DIOXIDE (three-letter code: SO2) (formula: O₂S).



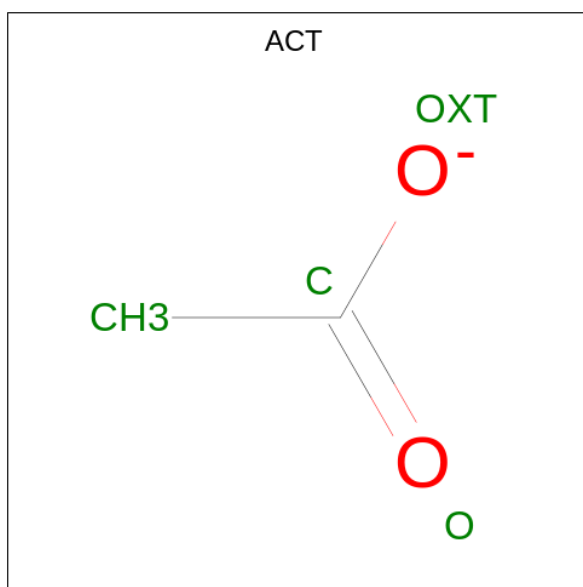
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			3	2	1		

- Molecule 4 is CARBON DIOXIDE (three-letter code: CO2) (formula: CO₂).



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

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



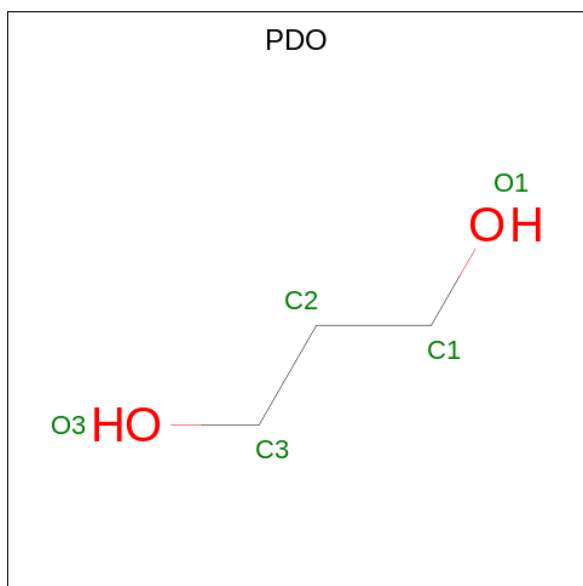
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 7 is 1,3-PROPANDIOL (three-letter code: PDO) (formula: $C_3H_8O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			5	3	2		
7	A	1	Total	C	O	0	0
			5	3	2		

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



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

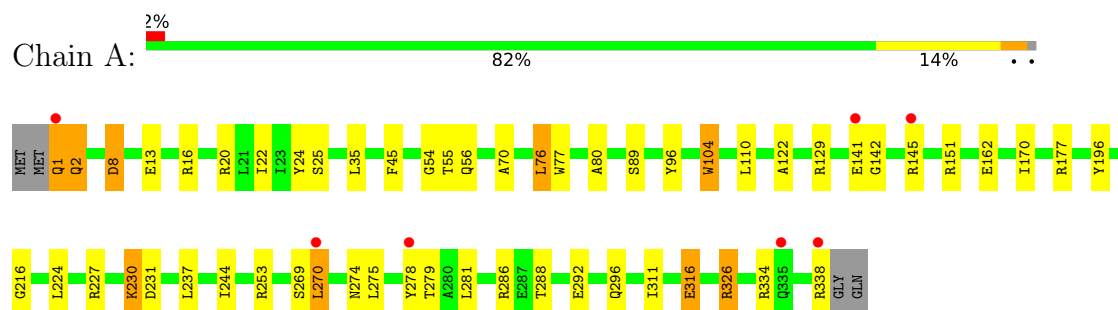
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	355	Total	O	0	0
			355	355		

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: Iron ABC transporter, periplasmic iron-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	72.17Å 72.17Å 114.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.21 – 1.63 42.18 – 1.63	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.21-1.63) 99.9 (42.18-1.63)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.30 (at 1.63Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.175 , 0.213 0.186 , 0.224	Depositor DCC
R_{free} test set	2174 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	19.7	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3128	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.70% 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: SO2, CL, PDO, PEG, EDO, CO2, ACT

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	1.00	5/2815 (0.2%)	1.19	12/3817 (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 (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	316	GLU	CD-OE1	7.33	1.33	1.25
1	A	2	GLN	C-O	6.66	1.36	1.23
1	A	13	GLU	CD-OE2	6.51	1.32	1.25
1	A	279	THR	C-O	5.36	1.33	1.23
1	A	292	GLU	CD-OE1	-5.16	1.20	1.25

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	ARG	NE-CZ-NH2	-10.61	115.00	120.30
1	A	253	ARG	NE-CZ-NH1	9.61	125.11	120.30
1	A	286	ARG	NE-CZ-NH1	7.91	124.26	120.30
1	A	326	ARG	NE-CZ-NH2	6.99	123.79	120.30
1	A	129	ARG	NE-CZ-NH1	6.90	123.75	120.30
1	A	270	LEU	CB-CG-CD1	6.76	122.49	111.00
1	A	196	TYR	CB-CG-CD1	6.35	124.81	121.00
1	A	129	ARG	NE-CZ-NH2	-5.47	117.57	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	286	ARG	CD-NE-CZ	5.47	131.25	123.60
1	A	104	TRP	CE3-CZ3-CH2	5.43	127.17	121.20
1	A	104	TRP	CH2-CZ2-CE2	-5.42	111.98	117.40
1	A	8	ASP	CB-CG-OD2	-5.29	113.54	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	269	SER	Mainchain

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	2726	0	2769	71	0
2	A	4	0	0	1	0
3	A	3	0	0	0	0
4	A	3	0	0	3	0
5	A	4	0	3	3	0
6	A	16	0	23	27	0
7	A	10	0	16	12	0
8	A	7	0	9	4	0
9	A	355	0	0	18	0
All	All	3128	0	2820	81	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 14.

All (81) 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:274:ASN:HB2	6:A:409:EDO:C1	1.78	1.11
1:A:122:ALA:CB	1:A:170[B]:ILE:HD11	1.84	1.06
1:A:122:ALA:HB2	1:A:170[B]:ILE:HD11	1.40	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:96:TYR:OH	7:A:412:PDO:H31	1.69	0.93
1:A:274:ASN:HB2	6:A:409:EDO:H12	1.49	0.92
1:A:2:GLN:HB3	6:A:410:EDO:H21	1.54	0.88
1:A:270:LEU:CD2	6:A:411:EDO:O2	2.24	0.85
6:A:410:EDO:H11	9:A:762:HOH:O	1.77	0.84
1:A:35:LEU:HD23	1:A:275[A]:LEU:HD22	1.62	0.81
1:A:122:ALA:HB1	1:A:170[B]:ILE:HD11	1.62	0.81
1:A:270:LEU:HD21	6:A:411:EDO:O2	1.83	0.77
1:A:316:GLU:HG2	9:A:757:HOH:O	1.84	0.77
2:A:403:CL:CL	9:A:670:HOH:O	2.40	0.77
1:A:274:ASN:HB2	6:A:409:EDO:H11	1.68	0.75
1:A:2:GLN:CB	6:A:410:EDO:H21	2.16	0.74
1:A:270:LEU:HD23	6:A:411:EDO:O2	1.87	0.74
1:A:122:ALA:HB2	1:A:170[B]:ILE:CD1	2.18	0.73
1:A:56:GLN:HB3	8:A:414:PEG:H32	1.71	0.73
1:A:145:ARG:NH2	6:A:408:EDO:O1	2.22	0.72
7:A:412:PDO:H11	9:A:714:HOH:O	1.88	0.72
1:A:274:ASN:CB	6:A:409:EDO:H12	2.18	0.72
1:A:20:ARG:HD2	9:A:603:HOH:O	1.89	0.72
1:A:2:GLN:HB3	6:A:410:EDO:C2	2.21	0.70
1:A:96:TYR:OH	7:A:412:PDO:C3	2.40	0.70
1:A:270:LEU:HB3	6:A:409:EDO:O2	1.94	0.68
1:A:54:GLY:HA3	8:A:414:PEG:H41	1.75	0.67
1:A:278:TYR:HB2	5:A:407:ACT:H2	1.77	0.66
1:A:8:ASP:HB2	7:A:412:PDO:H21	1.78	0.65
1:A:278:TYR:HD2	5:A:407:ACT:H2	1.62	0.65
1:A:326:ARG:HD2	9:A:533:HOH:O	1.96	0.65
1:A:141:GLU:HG3	9:A:706:HOH:O	1.97	0.63
1:A:2:GLN:CB	6:A:410:EDO:C2	2.76	0.62
1:A:2:GLN:CA	6:A:410:EDO:H21	2.29	0.62
1:A:244[B]:ILE:HD13	1:A:311:ILE:HD12	1.82	0.62
1:A:145:ARG:HD3	9:A:585:HOH:O	2.00	0.62
1:A:170[B]:ILE:HB	1:A:244[B]:ILE:HD11	1.82	0.62
1:A:170[B]:ILE:CD1	1:A:216:GLY:O	2.48	0.61
1:A:2:GLN:HA	6:A:410:EDO:O2	2.01	0.61
1:A:270:LEU:HD12	9:A:596:HOH:O	2.01	0.61
1:A:274:ASN:CB	6:A:409:EDO:C1	2.68	0.60
1:A:89[B]:SER:HB3	9:A:502:HOH:O	2.02	0.59
1:A:338:ARG:HA	9:A:756:HOH:O	2.05	0.57
1:A:1:GLN:N	9:A:503:HOH:O	2.16	0.57
6:A:410:EDO:C1	9:A:762:HOH:O	2.43	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:TYR:CD2	5:A:407:ACT:H2	2.40	0.56
1:A:270:LEU:HB3	6:A:409:EDO:C2	2.36	0.55
1:A:56:GLN:CB	8:A:414:PEG:H32	2.39	0.52
4:A:406:CO2:O2	7:A:412:PDO:C2	2.58	0.52
4:A:406:CO2:O2	7:A:412:PDO:H22	2.10	0.51
1:A:2:GLN:HA	6:A:410:EDO:C2	2.41	0.51
1:A:8:ASP:CB	7:A:412:PDO:H21	2.42	0.49
1:A:35:LEU:CD2	1:A:275[A]:LEU:HD22	2.37	0.49
1:A:230:LYS:HD2	1:A:230:LYS:C	2.34	0.49
1:A:224:LEU:HA	1:A:281:LEU:HD12	1.94	0.48
1:A:162:GLU:OE1	1:A:334:ARG:NH1	2.46	0.48
1:A:244[B]:ILE:HD13	1:A:311:ILE:CD1	2.42	0.48
7:A:413:PDO:H31	9:A:772:HOH:O	2.14	0.48
1:A:25:SER:HA	1:A:77:TRP:O	2.15	0.47
1:A:270:LEU:HB3	6:A:409:EDO:H21	1.96	0.47
1:A:1:GLN:O	6:A:410:EDO:H11	2.15	0.47
1:A:142:GLY:HA3	6:A:408:EDO:H11	1.97	0.46
1:A:16:ARG:HA	1:A:45:PHE:CD1	2.50	0.46
1:A:170[B]:ILE:HD13	1:A:216:GLY:O	2.16	0.46
8:A:414:PEG:H12	9:A:718:HOH:O	2.16	0.46
1:A:177[B]:ARG:HE	1:A:177[B]:ARG:HB3	1.57	0.44
1:A:70:ALA:HB1	7:A:413:PDO:H21	1.98	0.44
1:A:296:GLN:OE1	9:A:505:HOH:O	2.21	0.44
1:A:227:ARG:HD3	1:A:281:LEU:HD11	2.01	0.43
1:A:274:ASN:HB2	6:A:409:EDO:O1	2.15	0.42
1:A:55:THR:HB	1:A:80:ALA:HB2	2.00	0.42
1:A:270:LEU:O	6:A:409:EDO:H12	2.19	0.42
1:A:237:LEU:O	9:A:506:HOH:O	2.22	0.41
1:A:24:TYR:HB2	1:A:76[A]:LEU:HD22	2.02	0.41
1:A:96:TYR:HE1	7:A:412:PDO:H12	1.85	0.41
1:A:1:GLN:O	6:A:410:EDO:C1	2.68	0.41
1:A:20:ARG:HH11	1:A:22:ILE:HD11	1.86	0.40
1:A:288:THR:HG23	6:A:411:EDO:H11	2.04	0.40
1:A:230:LYS:HD2	1:A:231:ASP:OD1	2.22	0.40
4:A:406:CO2:O2	7:A:412:PDO:H21	2.22	0.40
7:A:412:PDO:C1	9:A:714:HOH:O	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	347/342 (102%)	343 (99%)	4 (1%)	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	281/273 (103%)	274 (98%)	7 (2%)	47	20

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLN
1	A	76[A]	LEU
1	A	76[B]	LEU
1	A	104	TRP
1	A	110	LEU
1	A	151	ARG
1	A	230	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](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	$\# Z > 2$
6	EDO	A	409	-	3,3,3	0.35	0	2,2,2	0.59	0
3	SO2	A	405	-	2,2,2	0.52	0	1,1,1	0.15	0
6	EDO	A	410	-	3,3,3	0.62	0	2,2,2	0.27	0
6	EDO	A	411	-	3,3,3	0.85	0	2,2,2	0.67	0
7	PDO	A	413	-	4,4,4	0.15	0	3,3,3	0.43	0
6	EDO	A	408	-	3,3,3	0.16	0	2,2,2	0.14	0
5	ACT	A	407	-	1,3,3	5.15	1 (100%)	0,3,3	-	-
4	CO2	A	406	-	2,2,2	0.19	0	1,1,1	0.81	0
7	PDO	A	412	-	4,4,4	0.35	0	3,3,3	0.95	0
8	PEG	A	414	-	6,6,6	0.55	0	5,5,5	0.17	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
6	EDO	A	410	-	-	1/1/1/1	-
6	EDO	A	409	-	-	1/1/1/1	-
6	EDO	A	411	-	-	1/1/1/1	-
7	PDO	A	413	-	-	0/2/2/2	-
6	EDO	A	408	-	-	1/1/1/1	-
7	PDO	A	412	-	-	2/2/2/2	-
8	PEG	A	414	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	407	ACT	CH3-C	5.15	1.55	1.48

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	412	PDO	O1-C1-C2-C3
7	A	412	PDO	C1-C2-C3-O3
8	A	414	PEG	O2-C3-C4-O4
6	A	408	EDO	O1-C1-C2-O2
8	A	414	PEG	O1-C1-C2-O2
6	A	409	EDO	O1-C1-C2-O2
6	A	411	EDO	O1-C1-C2-O2
6	A	410	EDO	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 46 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	409	EDO	10	0
6	A	410	EDO	11	0
6	A	411	EDO	4	0
7	A	413	PDO	2	0
6	A	408	EDO	2	0
5	A	407	ACT	3	0
4	A	406	CO2	3	0
7	A	412	PDO	10	0
8	A	414	PEG	4	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	A	338/342 (98%)	-0.13	7 (2%) 63 64	13, 20, 33, 67	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	GLN	4.5
1	A	338	ARG	4.1
1	A	145	ARG	2.5
1	A	278	TYR	2.3
1	A	141	GLU	2.1
1	A	270	LEU	2.0
1	A	335	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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
6	EDO	A	408	4/4	0.64	0.17	40,44,45,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	ACT	A	407	4/4	0.72	0.28	26,29,31,32	0
2	CL	A	401	1/1	0.84	0.07	48,48,48,48	0
4	CO2	A	406	3/3	0.84	0.12	43,43,47,52	0
6	EDO	A	409	4/4	0.88	0.23	34,38,38,42	0
6	EDO	A	411	4/4	0.89	0.33	34,34,34,35	0
7	PDO	A	412	5/5	0.89	0.20	28,29,34,34	0
8	PEG	A	414	7/7	0.89	0.13	30,33,37,37	0
2	CL	A	403	1/1	0.92	0.09	47,47,47,47	0
7	PDO	A	413	5/5	0.92	0.16	36,38,41,41	0
2	CL	A	402	1/1	0.92	0.05	51,51,51,51	0
6	EDO	A	410	4/4	0.93	0.27	24,28,32,33	0
2	CL	A	404	1/1	0.95	0.13	49,49,49,49	0
3	SO2	A	405	3/3	0.96	0.05	34,34,48,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.