



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:15 PM GMT

PDB ID : 4HPE
Title : Crystal structure of a putative cell wall hydrolase (CD630_03720) from Clostridium difficile 630 at 2.38 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2012-10-23
Resolution : 2.38 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

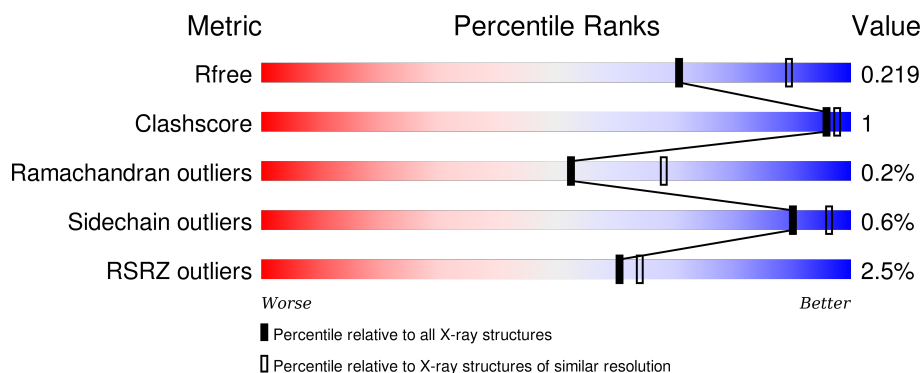
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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}	91344	4019 (2.40-2.36)
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (2.40-2.36)
RSRZ outliers	91569	4034 (2.40-2.36)

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. 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	308	<div> <div></div> <div>90%</div> <div>6%</div> </div>
1	B	308	<div> <div></div> <div>87%</div> <div>6%</div> <div>6%</div> </div>
1	C	308	<div> <div>3%</div> <div>91%</div> <div>6%</div> </div>
1	D	308	<div> <div>3%</div> <div>92%</div> <div>6%</div> </div>
1	E	308	<div> <div>2%</div> <div>89%</div> <div>5%</div> <div>6%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	308	

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	CL	A	404	-	-	-	X
2	CL	F	401	-	-	-	X
3	GOL	A	405	-	-	-	X
3	GOL	B	402	-	-	-	X
3	GOL	D	402	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13970 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 Putative cell wall hydrolase Tn916-like,CTn1-Orf17.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	Se	0	1	0
			2230	1408	378	435	3	6			
1	B	289	Total	C	N	O	S	Se	0	1	0
			2222	1403	377	433	3	6			
1	C	289	Total	C	N	O	S	Se	0	1	0
			2222	1403	377	433	3	6			
1	D	290	Total	C	N	O	S	Se	0	1	0
			2225	1406	376	434	3	6			
1	E	289	Total	C	N	O	S	Se	0	1	0
			2222	1403	375	435	3	6			
1	F	289	Total	C	N	O	S	Se	0	1	0
			2226	1405	377	435	3	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	leader sequence	UNP Q18DB2
B	0	GLY	-	leader sequence	UNP Q18DB2
C	0	GLY	-	leader sequence	UNP Q18DB2
D	0	GLY	-	leader sequence	UNP Q18DB2
E	0	GLY	-	leader sequence	UNP Q18DB2
F	0	GLY	-	leader sequence	UNP Q18DB2

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	1	Total	Cl	0	0
			1	1		
2	E	1	Total	Cl	0	0
			1	1		
2	B	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	1	Total	Cl	0	0
			1	1		
2	A	4	Total	Cl	0	0
			4	4		
2	F	1	Total	Cl	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	90	Total	O	0	0
			90	90		
4	B	103	Total	O	0	0
			103	103		

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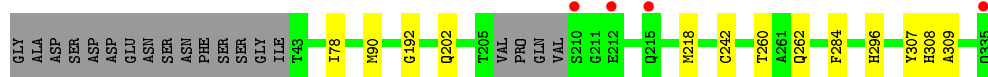
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	87	Total 87	O 87	0	0
4	D	108	Total 108	O 108	0	0
4	E	109	Total 109	O 109	0	0
4	F	93	Total 93	O 93	0	0

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 errors 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.

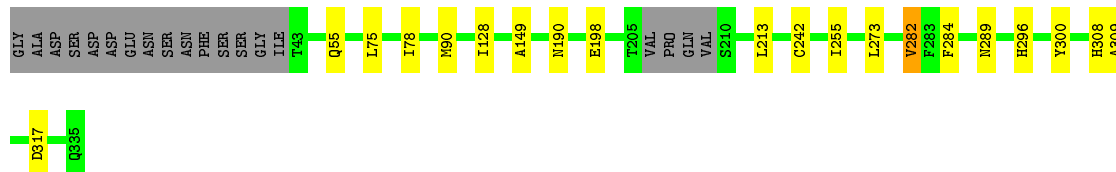
- Molecule 1: Putative cell wall hydrolase Tn916-like,CTn1-Orf17

Chain A: 




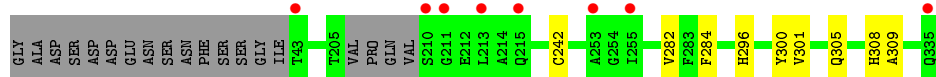
- Molecule 1: Putative cell wall hydrolase Tn916-like,CTn1-Orf17

Chain B: 

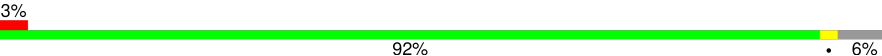


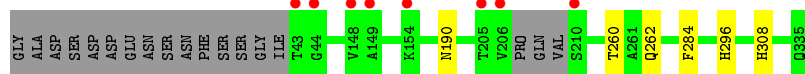
- Molecule 1: Putative cell wall hydrolase Tn916-like,CTn1-Orf17

Chain C: 




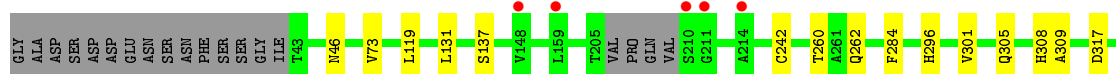
- Molecule 1: Putative cell wall hydrolase Tn916-like,CTn1-Orf17

Chain D: 



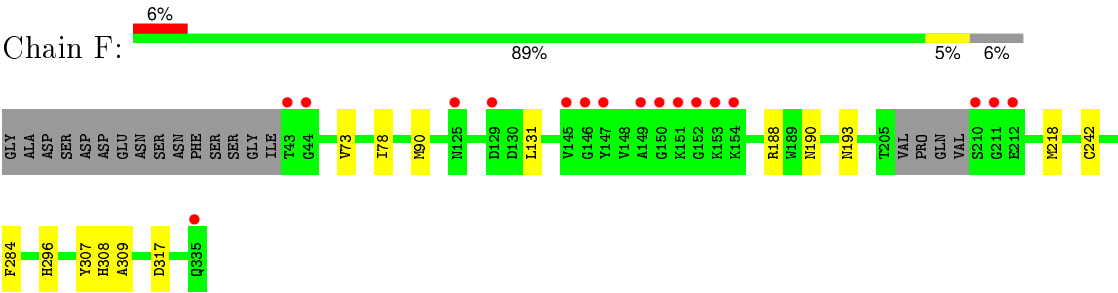
- Molecule 1: Putative cell wall hydrolase Tn916-like,CTn1-Orf17

Chain E: 



Q335

● Molecule 1: Putative cell wall hydrolase Tn916-like,CTn1-Orf17



F284
H296
Y307
H308
A309
D317
Q335

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	65.52Å 101.21Å 101.44Å 109.71° 108.86° 101.36°	Depositor
Resolution (Å)	49.40 – 2.38 49.40 – 2.38	Depositor EDS
% Data completeness (in resolution range)	87.6 (49.40-2.38) 74.6 (49.40-2.38)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.37Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.175 , 0.211 0.182 , 0.219	Depositor DCC
R_{free} test set	3897 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 76631 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13970	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: GOL, OCS, 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	0.51	0/2269	0.63	0/3060
1	B	0.51	0/2261	0.63	0/3051
1	C	0.51	0/2261	0.63	0/3051
1	D	0.52	0/2264	0.62	0/3055
1	E	0.51	0/2261	0.63	0/3050
1	F	0.50	0/2265	0.62	0/3056
All	All	0.51	0/13581	0.63	0/18323

There are no bond length outliers.

There are no bond angle outliers.

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	A	2230	0	2126	7	0
1	B	2222	0	2111	11	0
1	C	2222	0	2111	5	0
1	D	2225	0	2112	2	0
1	E	2222	0	2107	7	0
1	F	2226	0	2115	7	0
2	A	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	12	0	16	0	0
3	B	6	0	8	0	0
3	D	6	0	8	0	0
4	A	90	0	0	1	0
4	B	103	0	0	2	0
4	C	87	0	0	0	0
4	D	108	0	0	0	0
4	E	109	0	0	1	0
4	F	93	0	0	1	0
All	All	13970	0	12714	38	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 1.

All (38) 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:E:317:ASP:HB2	4:E:547:HOH:O	1.90	0.70
1:F:242:OCS:HA	1:F:309:ALA:HB3	1.85	0.59
1:C:282:VAL:CG2	1:C:300:TYR:HB2	2.34	0.56
1:B:55:GLN:HG3	1:B:75:LEU:HD11	1.87	0.56
1:A:242:OCS:HA	1:A:309:ALA:HB3	1.88	0.55
1:B:242:OCS:HA	1:B:309:ALA:HB3	1.90	0.54
1:D:284:PHE:HB2	1:D:296:HIS:HB3	1.90	0.52
1:B:78:ILE:HG12	1:B:90:MSE:HG2	1.91	0.52
1:C:242:OCS:HA	1:C:309:ALA:HB3	1.92	0.52
1:C:282:VAL:HG23	1:C:300:TYR:HB2	1.93	0.50
1:B:282:VAL:HG22	1:B:300:TYR:HB2	1.93	0.49
1:B:213:LEU:HD13	1:B:255:ILE:HD12	1.94	0.49
1:F:317:ASP:HB2	4:F:556:HOH:O	2.13	0.49
1:E:260:THR:HG22	1:E:262:GLN:H	1.78	0.48
1:B:317:ASP:HB2	4:B:581:HOH:O	2.13	0.48
1:F:78:ILE:HG12	1:F:90:MSE:HG2	1.94	0.48
1:E:284:PHE:HB2	1:E:296:HIS:HB3	1.97	0.46
1:A:260:THR:HG22	1:A:262:GLN:H	1.79	0.46
1:C:284:PHE:HB2	1:C:296:HIS:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:242:OCS:HA	1:E:309:ALA:HB3	1.98	0.45
1:E:301:VAL:HG23	1:E:305:GLN:HG3	1.98	0.45
1:B:198:GLU:HB3	4:B:585:HOH:O	2.16	0.44
1:B:284:PHE:HB2	1:B:296:HIS:HB3	1.99	0.44
1:A:78:ILE:HG12	1:A:90:MSE:HG2	1.99	0.44
1:D:260:THR:HG22	1:D:262:GLN:H	1.83	0.44
1:A:202:GLN:NE2	4:A:558:HOH:O	2.52	0.43
1:F:188:ARG:HD2	1:F:193:ASN:HA	2.00	0.42
1:E:73:VAL:HG21	1:E:131:LEU:HD21	2.02	0.42
1:F:73:VAL:HG21	1:F:131:LEU:HD21	2.01	0.42
1:F:284:PHE:HB2	1:F:296:HIS:HB3	2.02	0.42
1:E:119:LEU:HD22	1:E:137:SER:HB3	2.02	0.42
1:A:192:GLY:HA2	1:B:289:ASN:OD1	2.20	0.41
1:F:218:MSE:HE1	1:F:307:TYR:CD2	2.56	0.41
1:B:128:ILE:HD11	1:B:149:ALA:HB2	2.02	0.41
1:B:273:LEU:HD13	1:B:282:VAL:HG13	2.03	0.41
1:A:284:PHE:HB2	1:A:296:HIS:HB3	2.03	0.40
1:A:218:MSE:HE1	1:A:307:TYR:CD2	2.57	0.40
1:C:301:VAL:HG23	1:C:305:GLN:HG3	2.03	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	285/308 (92%)	277 (97%)	8 (3%)	0	100	100
1	B	285/308 (92%)	277 (97%)	7 (2%)	1 (0%)	39	53
1	C	285/308 (92%)	277 (97%)	8 (3%)	0	100	100
1	D	286/308 (93%)	278 (97%)	7 (2%)	1 (0%)	46	61
1	E	285/308 (92%)	277 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	285/308 (92%)	279 (98%)	5 (2%)	1 (0%)	39	53
All	All	1711/1848 (93%)	1665 (97%)	43 (2%)	3 (0%)	52	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	190	ASN
1	D	190	ASN
1	F	190	ASN

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	230/242 (95%)	229 (100%)	1 (0%)	93	98
1	B	228/242 (94%)	226 (99%)	2 (1%)	84	93
1	C	228/242 (94%)	227 (100%)	1 (0%)	93	98
1	D	228/242 (94%)	227 (100%)	1 (0%)	93	98
1	E	228/242 (94%)	226 (99%)	2 (1%)	84	93
1	F	229/242 (95%)	228 (100%)	1 (0%)	93	98
All	All	1371/1452 (94%)	1363 (99%)	8 (1%)	90	96

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

Mol	Chain	Res	Type
1	A	308	HIS
1	B	282	VAL
1	B	308	HIS
1	C	308	HIS
1	D	308	HIS
1	E	46	ASN
1	E	308	HIS
1	F	308	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	303	ASN
1	E	325	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 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	$\# Z > 2$
1	OCS	A	242	1	7,8,9	1.09	0	7,11,13	6.29	5 (71%)
1	OCS	B	242	1	7,8,9	0.85	0	7,11,13	5.20	3 (42%)
1	OCS	C	242	1	7,8,9	1.22	1 (14%)	7,11,13	3.97	4 (57%)
1	OCS	D	242	1	7,8,9	1.87	1 (14%)	7,11,13	6.30	4 (57%)
1	OCS	E	242	1	7,8,9	1.30	1 (14%)	7,11,13	9.90	6 (85%)
1	OCS	F	242	1	7,8,9	0.81	0	7,11,13	6.66	4 (57%)

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
1	OCS	A	242	1	-	0/4/7/9	0/0/0/0
1	OCS	B	242	1	-	0/4/7/9	0/0/0/0
1	OCS	C	242	1	-	0/4/7/9	0/0/0/0
1	OCS	D	242	1	-	0/4/7/9	0/0/0/0
1	OCS	E	242	1	-	0/4/7/9	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	OCS	F	242	1	-	0/4/7/9	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	242	OCS	CB-SG	2.51	1.81	1.77
1	E	242	OCS	CB-SG	2.62	1.81	1.77
1	D	242	OCS	CB-SG	4.64	1.84	1.77

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	242	OCS	OD1-SG-CB	-12.69	96.24	106.94
1	D	242	OCS	OD2-SG-OD1	-7.55	94.04	111.61
1	E	242	OCS	OD2-SG-OD3	-5.68	98.40	111.61
1	F	242	OCS	OD2-SG-OD3	-3.80	102.77	111.61
1	A	242	OCS	OD2-SG-OD3	-3.38	103.74	111.61
1	E	242	OCS	CB-CA-C	-3.32	102.35	111.46
1	B	242	OCS	OD2-SG-OD3	-3.27	103.99	111.61
1	C	242	OCS	O-C-CA	-2.71	118.43	125.49
1	F	242	OCS	O-C-CA	-2.70	118.46	125.49
1	A	242	OCS	O-C-CA	-2.62	118.66	125.49
1	C	242	OCS	OD2-SG-OD1	-2.61	105.54	111.61
1	D	242	OCS	O-C-CA	-2.58	118.76	125.49
1	F	242	OCS	OD2-SG-OD1	-2.51	105.76	111.61
1	E	242	OCS	O-C-CA	-2.38	119.29	125.49
1	A	242	OCS	OD2-SG-OD1	-2.25	106.38	111.61
1	B	242	OCS	O-C-CA	-2.20	119.75	125.49
1	D	242	OCS	OD1-SG-CB	2.03	108.65	106.94
1	E	242	OCS	OD2-SG-OD1	3.61	120.02	111.61
1	C	242	OCS	OD1-SG-CB	5.79	111.82	106.94
1	C	242	OCS	OD3-SG-CB	7.48	113.25	106.94
1	A	242	OCS	OD1-SG-CB	10.00	115.37	106.94
1	A	242	OCS	OD3-SG-CB	12.35	117.36	106.94
1	B	242	OCS	OD3-SG-CB	12.95	117.86	106.94
1	D	242	OCS	OD3-SG-CB	14.26	118.97	106.94
1	F	242	OCS	OD3-SG-CB	16.59	120.93	106.94
1	E	242	OCS	OD3-SG-CB	21.50	125.07	106.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	242	OCS	1	0
1	B	242	OCS	1	0
1	C	242	OCS	1	0
1	E	242	OCS	1	0
1	F	242	OCS	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 9 are monoatomic - leaving 4 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	$\# Z > 2$
3	GOL	A	405	-	5,5,5	0.16	0	5,5,5	0.36	0
3	GOL	A	406	-	5,5,5	0.21	0	5,5,5	0.41	0
3	GOL	B	402	-	5,5,5	0.22	0	5,5,5	0.47	0
3	GOL	D	402	-	5,5,5	0.20	0	5,5,5	0.38	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
3	GOL	A	405	-	-	0/4/4/4	0/0/0/0
3	GOL	A	406	-	-	0/4/4/4	0/0/0/0
3	GOL	B	402	-	-	0/4/4/4	0/0/0/0
3	GOL	D	402	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	A	282/308 (91%)	-0.09	4 (1%) 78 80	33, 46, 71, 104	0
1	B	282/308 (91%)	-0.13	0 100 100	33, 47, 71, 99	0
1	C	282/308 (91%)	-0.08	8 (2%) 56 59	31, 46, 75, 113	0
1	D	283/308 (91%)	-0.06	8 (2%) 56 59	28, 46, 72, 98	0
1	E	282/308 (91%)	-0.09	6 (2%) 67 70	27, 48, 77, 97	0
1	F	282/308 (91%)	0.19	17 (6%) 25 29	31, 53, 89, 106	0
All	All	1693/1848 (91%)	-0.05	43 (2%) 61 64	27, 47, 78, 113	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	44	GLY	5.7
1	F	153	LYS	5.4
1	D	43	THR	5.3
1	F	149	ALA	4.3
1	C	211	GLY	4.2
1	E	211	GLY	4.0
1	C	335	GLN	4.0
1	F	43	THR	4.0
1	F	146	GLY	3.8
1	E	210	SER	3.8
1	D	206	VAL	3.6
1	D	149	ALA	3.4
1	A	335	GLN	3.4
1	D	205	THR	3.3
1	D	154	LYS	3.2
1	F	44	GLY	3.2
1	C	213	LEU	3.1
1	F	211	GLY	3.1
1	F	335	GLN	3.0

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Mol	Chain	Res	Type	RSRZ
1	F	147	TYR	2.8
1	F	154	LYS	2.8
1	F	125	ASN	2.7
1	E	335	GLN	2.6
1	A	215	GLN	2.6
1	C	210	SER	2.6
1	F	129	ASP	2.6
1	F	145	VAL	2.6
1	C	255	ILE	2.6
1	C	215	GLN	2.5
1	C	253	ALA	2.5
1	F	152	GLY	2.4
1	C	43	THR	2.4
1	F	150	GLY	2.4
1	A	210	SER	2.4
1	A	212	GLU	2.4
1	F	212	GLU	2.4
1	F	210	SER	2.3
1	E	159	LEU	2.2
1	D	148	VAL	2.2
1	E	148	VAL	2.1
1	D	210	SER	2.1
1	F	151	LYS	2.0
1	E	214	ALA	2.0

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

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
1	OCS	C	242	9/10	0.96	0.11	-	56,57,66,69	0
1	OCS	E	242	9/10	0.97	0.12	-	38,40,51,53	0
1	OCS	D	242	9/10	0.98	0.12	-	44,46,60,61	0
1	OCS	F	242	9/10	0.97	0.10	-	33,40,55,56	0
1	OCS	B	242	9/10	0.98	0.11	-	47,50,67,67	0
1	OCS	A	242	9/10	0.96	0.14	-	51,51,65,65	0

6.3 Carbohydrates [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	CL	A	404	1/1	0.96	0.21	7.59	75,75,75,75	0
3	GOL	B	402	6/6	0.93	0.20	6.72	49,54,55,59	0
3	GOL	A	405	6/6	0.95	0.18	4.82	53,56,57,57	0
3	GOL	D	402	6/6	0.94	0.18	3.43	50,52,53,54	0
2	CL	F	401	1/1	0.96	0.15	2.91	69,69,69,69	0
2	CL	B	401	1/1	0.95	0.16	1.89	81,81,81,81	0
2	CL	E	401	1/1	0.96	0.14	1.53	62,62,62,62	0
2	CL	C	401	1/1	0.97	0.16	1.05	70,70,70,70	0
3	GOL	A	406	6/6	0.96	0.14	0.12	39,48,50,52	0
2	CL	A	403	1/1	0.93	0.18	-0.65	87,87,87,87	0
2	CL	D	401	1/1	0.92	0.08	-1.56	60,60,60,60	0
2	CL	A	401	1/1	0.99	0.07	-2.85	41,41,41,41	0
2	CL	A	402	1/1	0.96	0.11	-	75,75,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.