



wwPDB X-ray Structure Validation Summary Report

Dec 12, 2023 – 12:36 PM EST

PDB ID : 8U5E
Title : Crystal Structure of C-terminal domain of Clostridium perfringens Enterotoxin
in Space Group P 21 21 21
Authors : Kapoor, S.; Vecchio, A.J.
Deposited on : 2023-09-12
Resolution : 1.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

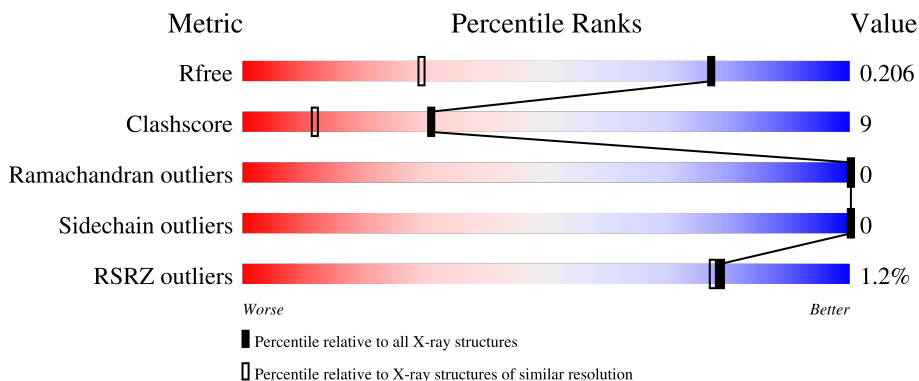
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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	1714 (1.40-1.40)
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763 (1.40-1.40)
Sidechain outliers	138945	1762 (1.40-1.40)
RSRZ outliers	127900	1674 (1.40-1.40)

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	145	
1	B	145	
1	C	145	
1	D	145	

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	GOL	A	405	-	-	-	X
2	GOL	B	402	-	-	X	-
3	ACT	B	401	-	-	X	-
3	ACT	C	403	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4956 atoms, of which 38 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 Heat-labile enterotoxin B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	119	1017	647	171	198	1	0	10	0
1	B	127	1051	672	171	207	1	0	7	0
1	C	127	1011	642	168	200	1	0	0	0
1	D	128	1055	677	171	206	1	0	8	0

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	175	GLY	-	expression tag	UNP P01558
A	176	GLY	-	expression tag	UNP P01558
A	177	GLY	-	expression tag	UNP P01558
A	178	GLY	-	expression tag	UNP P01558
A	179	SER	-	expression tag	UNP P01558
A	180	GLY	-	expression tag	UNP P01558
A	181	GLY	-	expression tag	UNP P01558
A	182	GLY	-	expression tag	UNP P01558
A	183	GLY	-	expression tag	UNP P01558
A	184	SER	-	expression tag	UNP P01558
A	185	GLY	-	expression tag	UNP P01558
A	186	GLY	-	expression tag	UNP P01558
A	187	GLY	-	expression tag	UNP P01558
A	188	LEU	-	expression tag	UNP P01558
A	189	VAL	-	expression tag	UNP P01558
A	190	PRO	-	expression tag	UNP P01558
A	191	ARG	-	expression tag	UNP P01558
A	192	GLY	-	expression tag	UNP P01558
A	193	SER	-	expression tag	UNP P01558
B	175	GLY	-	expression tag	UNP P01558
B	176	GLY	-	expression tag	UNP P01558

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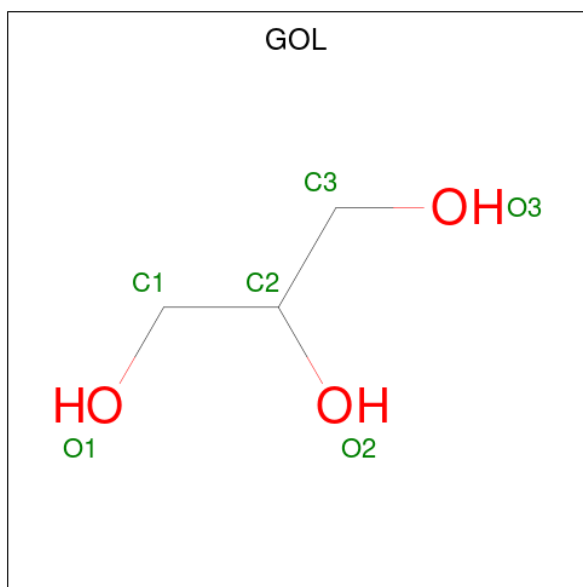
Chain	Residue	Modelled	Actual	Comment	Reference
B	177	GLY	-	expression tag	UNP P01558
B	178	GLY	-	expression tag	UNP P01558
B	179	SER	-	expression tag	UNP P01558
B	180	GLY	-	expression tag	UNP P01558
B	181	GLY	-	expression tag	UNP P01558
B	182	GLY	-	expression tag	UNP P01558
B	183	GLY	-	expression tag	UNP P01558
B	184	SER	-	expression tag	UNP P01558
B	185	GLY	-	expression tag	UNP P01558
B	186	GLY	-	expression tag	UNP P01558
B	187	GLY	-	expression tag	UNP P01558
B	188	LEU	-	expression tag	UNP P01558
B	189	VAL	-	expression tag	UNP P01558
B	190	PRO	-	expression tag	UNP P01558
B	191	ARG	-	expression tag	UNP P01558
B	192	GLY	-	expression tag	UNP P01558
B	193	SER	-	expression tag	UNP P01558
C	175	GLY	-	expression tag	UNP P01558
C	176	GLY	-	expression tag	UNP P01558
C	177	GLY	-	expression tag	UNP P01558
C	178	GLY	-	expression tag	UNP P01558
C	179	SER	-	expression tag	UNP P01558
C	180	GLY	-	expression tag	UNP P01558
C	181	GLY	-	expression tag	UNP P01558
C	182	GLY	-	expression tag	UNP P01558
C	183	GLY	-	expression tag	UNP P01558
C	184	SER	-	expression tag	UNP P01558
C	185	GLY	-	expression tag	UNP P01558
C	186	GLY	-	expression tag	UNP P01558
C	187	GLY	-	expression tag	UNP P01558
C	188	LEU	-	expression tag	UNP P01558
C	189	VAL	-	expression tag	UNP P01558
C	190	PRO	-	expression tag	UNP P01558
C	191	ARG	-	expression tag	UNP P01558
C	192	GLY	-	expression tag	UNP P01558
C	193	SER	-	expression tag	UNP P01558
D	175	GLY	-	expression tag	UNP P01558
D	176	GLY	-	expression tag	UNP P01558
D	177	GLY	-	expression tag	UNP P01558
D	178	GLY	-	expression tag	UNP P01558
D	179	SER	-	expression tag	UNP P01558
D	180	GLY	-	expression tag	UNP P01558

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Chain	Residue	Modelled	Actual	Comment	Reference
D	181	GLY	-	expression tag	UNP P01558
D	182	GLY	-	expression tag	UNP P01558
D	183	GLY	-	expression tag	UNP P01558
D	184	SER	-	expression tag	UNP P01558
D	185	GLY	-	expression tag	UNP P01558
D	186	GLY	-	expression tag	UNP P01558
D	187	GLY	-	expression tag	UNP P01558
D	188	LEU	-	expression tag	UNP P01558
D	189	VAL	-	expression tag	UNP P01558
D	190	PRO	-	expression tag	UNP P01558
D	191	ARG	-	expression tag	UNP P01558
D	192	GLY	-	expression tag	UNP P01558
D	193	SER	-	expression tag	UNP P01558

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



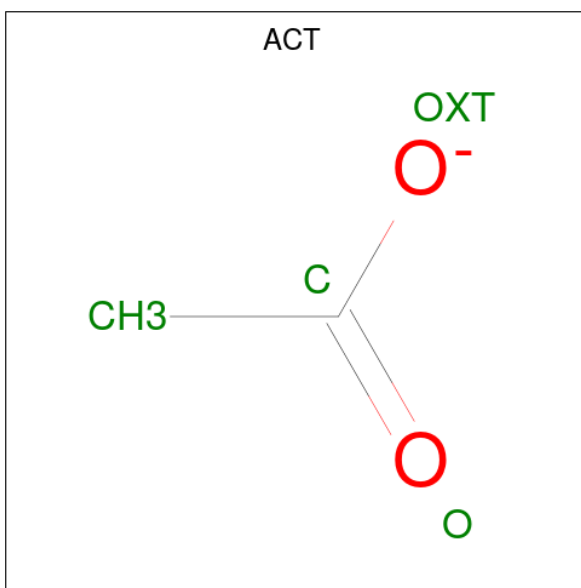
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O			
			6	3	3	0	0	
2	A	1	Total	C	O			
			6	3	3	0	0	
2	A	1	Total	C	H	O		
			14	3	8	3	0	
2	B	1	Total	C	O			
			6	3	3	0	0	
2	B	1	Total	C	H	O		
			14	3	8	3	0	

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C H O 14 3 8 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C H O 14 3 8 3	0	0

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



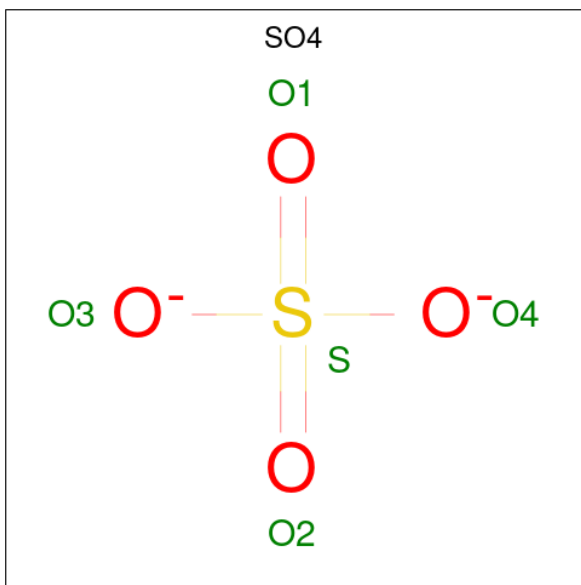
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C H O 7 2 3 2	0	0
3	A	1	Total C H O 7 2 3 2	0	0
3	B	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0

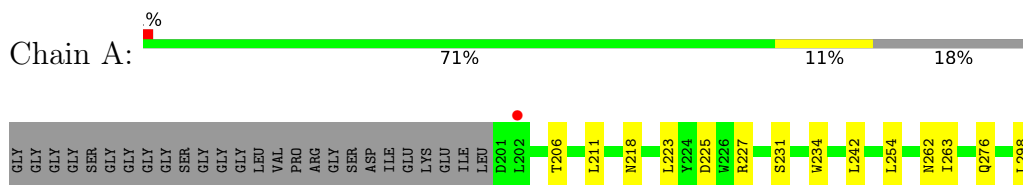
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	176	Total O 176 176	0	0
5	B	171	Total O 171 171	0	0
5	C	170	Total O 170 170	0	0
5	D	156	Total O 156 156	0	0

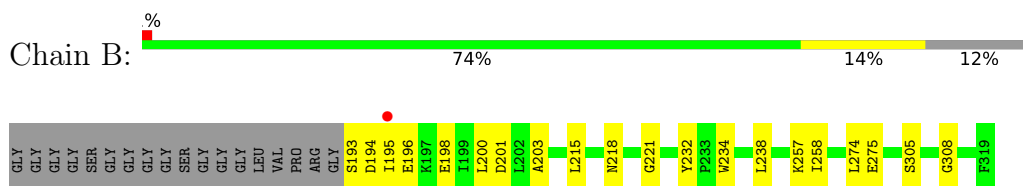
3 Residue-property plots [i](#)

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.

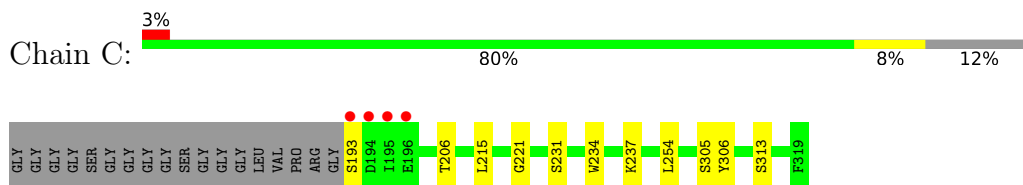
- Molecule 1: Heat-labile enterotoxin B chain



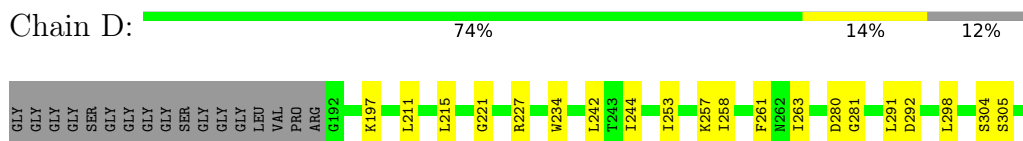
- Molecule 1: Heat-labile enterotoxin B chain



- Molecule 1: Heat-labile enterotoxin B chain



- Molecule 1: Heat-labile enterotoxin B chain



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.63Å 64.00Å 136.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.19 – 1.40 59.19 – 1.40	Depositor EDS
% Data completeness (in resolution range)	83.8 (59.19-1.40) 76.6 (59.19-1.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.31 (at 1.40Å)	Xtrriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.161 , 0.205 0.163 , 0.206	Depositor DCC
R_{free} test set	1951 reflections (1.95%)	wwPDB-VP
Wilson B-factor (Å ²)	15.5	Xtrriage
Anisotropy	0.430	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.037 for k,h,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4956	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% 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: GOL, ACT, SO4

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	0.28	0/1044	0.52	0/1417
1	B	0.28	0/1088	0.50	0/1477
1	C	0.28	0/1032	0.50	0/1400
1	D	0.27	0/1098	0.49	0/1491
All	All	0.27	0/4262	0.50	0/5785

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	1017	0	989	22	0
1	B	1051	0	1036	16	0
1	C	1011	0	979	10	0
1	D	1055	0	1053	26	0
2	A	18	8	24	6	0
2	B	12	8	16	4	0
2	C	12	8	16	2	0
2	D	30	8	40	7	0
3	A	8	6	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	4	0	3	2	0
3	C	8	0	6	3	0
3	D	4	0	3	0	0
4	B	5	0	0	1	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
5	A	176	0	0	5	0
5	B	171	0	0	2	0
5	C	170	0	0	4	0
5	D	156	0	0	5	0
All	All	4918	38	4171	77	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 9.

The worst 5 of 77 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:211:LEU:HB2	1:A:242[B]:LEU:HD11	1.52	0.91
1:B:308:GLY:H	3:B:401:ACT:H3	1.43	0.81
1:B:258[B]:ILE:HG22	5:B:502:HOH:O	1.88	0.74
1:B:196:GLU:HB3	1:B:203:ALA:HB2	1.71	0.73
1:A:263[B]:ILE:HG12	1:A:298:LEU:CD1	2.20	0.71

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	127/145 (88%)	126 (99%)	1 (1%)	0	100	100
1	B	132/145 (91%)	130 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	125/145 (86%)	123 (98%)	2 (2%)	0	100	100
1	D	134/145 (92%)	132 (98%)	2 (2%)	0	100	100
All	All	518/580 (89%)	511 (99%)	7 (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	113/118 (96%)	113 (100%)	0	100	100
1	B	119/118 (101%)	119 (100%)	0	100	100
1	C	111/118 (94%)	111 (100%)	0	100	100
1	D	120/118 (102%)	120 (100%)	0	100	100
All	All	463/472 (98%)	463 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

21 ligands 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	ACT	C	403	-	3,3,3	1.25	0	3,3,3	1.37	0
2	GOL	D	406	-	5,5,5	0.08	0	5,5,5	0.29	0
2	GOL	C	402	-	5,5,5	0.89	0	5,5,5	1.00	0
2	GOL	D	402	-	5,5,5	0.95	0	5,5,5	1.04	0
2	GOL	A	405	-	5,5,5	0.10	0	5,5,5	0.26	0
2	GOL	B	403	-	5,5,5	0.77	0	5,5,5	0.92	0
4	SO4	C	405	-	4,4,4	0.13	0	6,6,6	0.07	0
2	GOL	D	404	-	5,5,5	0.94	0	5,5,5	0.91	0
2	GOL	A	402	-	5,5,5	0.94	0	5,5,5	1.04	0
3	ACT	A	403	-	3,3,3	1.12	0	3,3,3	1.53	0
3	ACT	A	404	-	3,3,3	1.07	0	3,3,3	1.52	0
3	ACT	B	401	-	3,3,3	1.32	0	3,3,3	1.34	0
2	GOL	A	401	-	5,5,5	0.92	0	5,5,5	1.00	0
3	ACT	C	401	-	3,3,3	1.36	0	3,3,3	1.47	0
3	ACT	D	401	-	3,3,3	1.30	0	3,3,3	1.39	0
4	SO4	B	404	-	4,4,4	0.13	0	6,6,6	0.05	0
2	GOL	D	403	-	5,5,5	0.85	0	5,5,5	1.04	0
2	GOL	D	405	-	5,5,5	1.03	0	5,5,5	1.32	1 (20%)
4	SO4	D	407	-	4,4,4	0.15	0	6,6,6	0.05	0
2	GOL	B	402	-	5,5,5	1.14	0	5,5,5	0.82	0
2	GOL	C	404	-	5,5,5	0.74	0	5,5,5	0.98	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
2	GOL	D	405	-	-	2/4/4/4	-
2	GOL	D	406	-	-	0/4/4/4	-
2	GOL	C	402	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	D	402	-	-	0/4/4/4	-
2	GOL	D	404	-	-	0/4/4/4	-
2	GOL	A	401	-	-	0/4/4/4	-
2	GOL	A	405	-	-	2/4/4/4	-
2	GOL	B	402	-	-	4/4/4/4	-
2	GOL	B	403	-	-	0/4/4/4	-
2	GOL	C	404	-	-	0/4/4/4	-
2	GOL	A	402	-	-	0/4/4/4	-
2	GOL	D	403	-	-	0/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	405	GOL	C3-C2-C1	-2.34	102.61	111.70

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	402	GOL	O1-C1-C2-C3
2	B	402	GOL	C1-C2-C3-O3
2	D	405	GOL	C1-C2-C3-O3
2	B	402	GOL	O1-C1-C2-O2
2	B	402	GOL	O2-C2-C3-O3

There are no ring outliers.

13 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	403	ACT	3	0
2	D	406	GOL	2	0
2	C	402	GOL	1	0
2	D	402	GOL	1	0
2	A	405	GOL	3	0
2	D	404	GOL	1	0
2	A	402	GOL	2	0
3	B	401	ACT	2	0
2	A	401	GOL	1	0
4	B	404	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	405	GOL	3	0
2	B	402	GOL	4	0
2	C	404	GOL	1	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	119/145 (82%)	-0.67	1 (0%) 86 84	17, 21, 29, 108	0
1	B	127/145 (87%)	-0.64	1 (0%) 86 84	15, 22, 41, 79	0
1	C	127/145 (87%)	-0.56	4 (3%) 49 48	15, 22, 41, 120	0
1	D	128/145 (88%)	-0.55	0 100 100	17, 25, 37, 82	0
All	All	501/580 (86%)	-0.60	6 (1%) 79 77	15, 23, 40, 120	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	194	ASP	4.1
1	B	195	ILE	4.0
1	A	202	LEU	2.6
1	C	196	GLU	2.4
1	C	193	SER	2.3

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(\AA^2)	Q<0.9
2	GOL	C	404	6/6	0.59	0.30	71,110,135,145	0
2	GOL	A	405	6/6	0.69	0.48	20,20,20,20	0
3	ACT	B	401	4/4	0.71	0.22	49,51,54,56	0
2	GOL	B	403	6/6	0.74	0.24	38,62,116,140	0
3	ACT	A	403	4/4	0.76	0.14	58,71,100,100	0
2	GOL	D	404	6/6	0.84	0.26	62,63,66,70	0
2	GOL	B	402	6/6	0.84	0.26	32,36,37,39	0
2	GOL	D	403	6/6	0.84	0.18	52,53,54,57	0
2	GOL	D	406	6/6	0.85	0.26	47,87,149,179	0
2	GOL	D	402	6/6	0.87	0.18	49,62,68,71	0
3	ACT	D	401	4/4	0.87	0.14	45,51,52,53	0
4	SO4	C	405	5/5	0.87	0.19	70,75,75,77	0
2	GOL	C	402	6/6	0.90	0.15	49,55,58,60	0
3	ACT	C	401	4/4	0.90	0.09	31,34,36,42	0
3	ACT	C	403	4/4	0.91	0.17	50,50,52,53	0
2	GOL	A	402	6/6	0.93	0.21	31,36,43,51	0
4	SO4	B	404	5/5	0.93	0.12	91,92,92,92	0
2	GOL	A	401	6/6	0.93	0.15	43,46,47,51	0
2	GOL	D	405	6/6	0.94	0.20	23,33,39,49	0
3	ACT	A	404	4/4	0.95	0.09	31,44,53,80	0
4	SO4	D	407	5/5	0.96	0.12	64,66,67,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.