



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 02:36 am BST

PDB ID : 6P1J  
Title : The structure of condensation and adenylation domains of teixobactin-producing nonribosomal peptide synthetase Txo2 serine module  
Authors : Tan, K.; Zhou, M.; Jedrzejczak, R.; Babnigg, G.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2019-05-20  
Resolution : 2.95 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](mailto: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.

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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

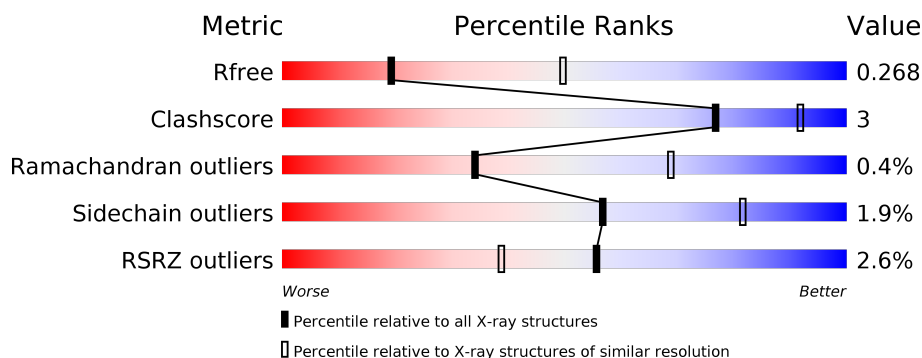
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.95 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	3104 (3.00-2.92)
Clashscore	141614	3462 (3.00-2.92)
Ramachandran outliers	138981	3340 (3.00-2.92)
Sidechain outliers	138945	3343 (3.00-2.92)
RSRZ outliers	127900	2986 (3.00-2.92)

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  $\geq 3$ , 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	964	<div> <div>3%</div> <div> <div></div> <div>90%</div> <div>9%</div> <div></div> </div> </div>
1	B	964	<div> <div>2%</div> <div> <div></div> <div>88%</div> <div>10%</div> <div></div> </div> </div>

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
5	CL	A	1107	-	-	-	X
6	GOL	A	1114	-	-	-	X

## 2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 15057 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.

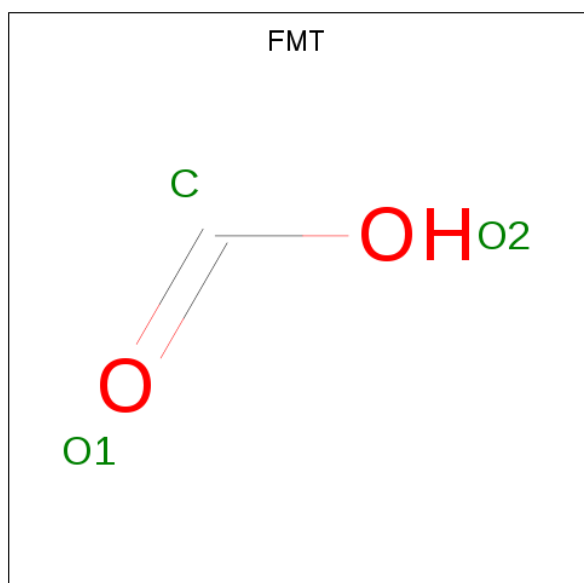
- Molecule 1 is a protein called Txo2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	958	Total	C	N	O	S	0	1	0
			7432	4692	1363	1354	23			
1	B	957	Total	C	N	O	S	0	2	0
			7444	4698	1364	1359	23			

There are 6 discrepancies between the modelled and reference sequences:

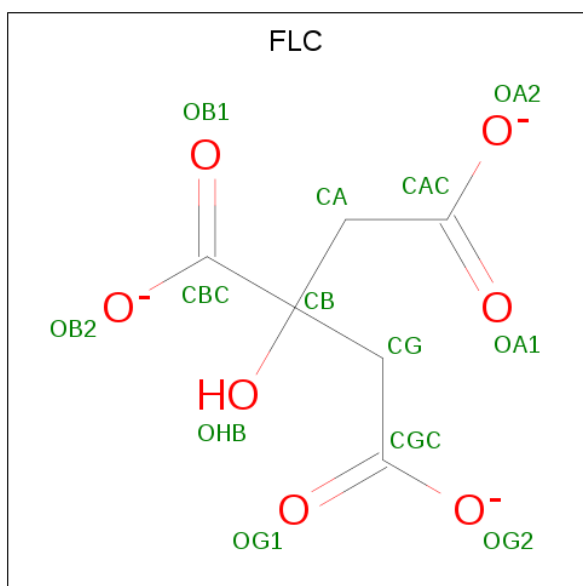
Chain	Residue	Modelled	Actual	Comment	Reference
A	42	SER	-	expression tag	UNP A0A0B5H0S3
A	43	ASN	-	expression tag	UNP A0A0B5H0S3
A	44	ALA	-	expression tag	UNP A0A0B5H0S3
B	42	SER	-	expression tag	UNP A0A0B5H0S3
B	43	ASN	-	expression tag	UNP A0A0B5H0S3
B	44	ALA	-	expression tag	UNP A0A0B5H0S3

- Molecule 2 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



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

- Molecule 3 is CITRATE ANION (three-letter code: FLC) (formula:  $C_6H_5O_7$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		
3	B	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).

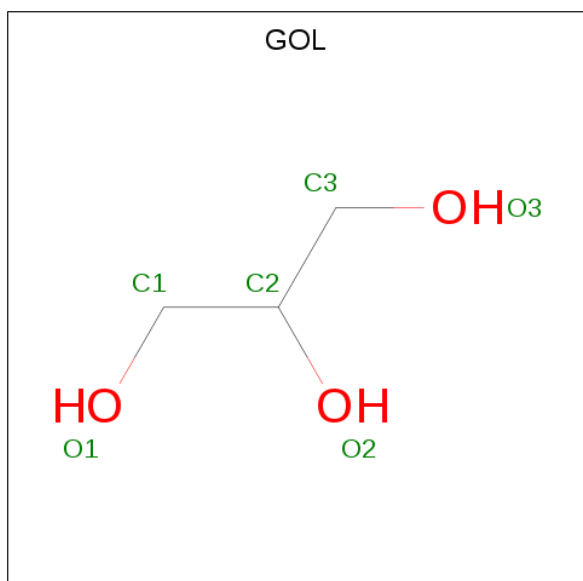


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

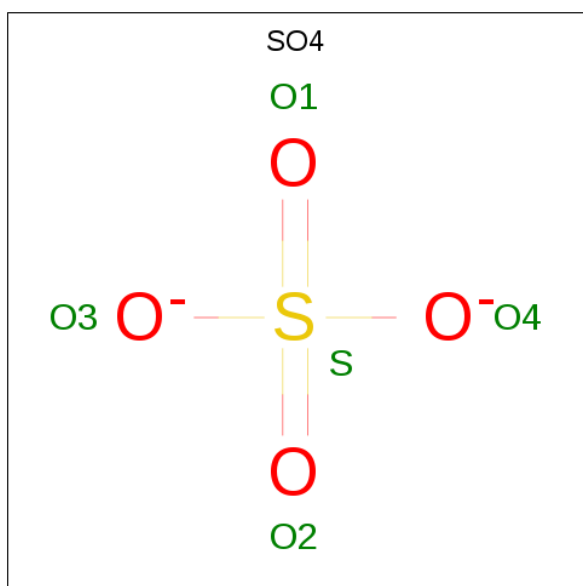
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	13	Total	Cl	0	0
			13	13		
5	A	6	Total	Cl	0	0
			6	6		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

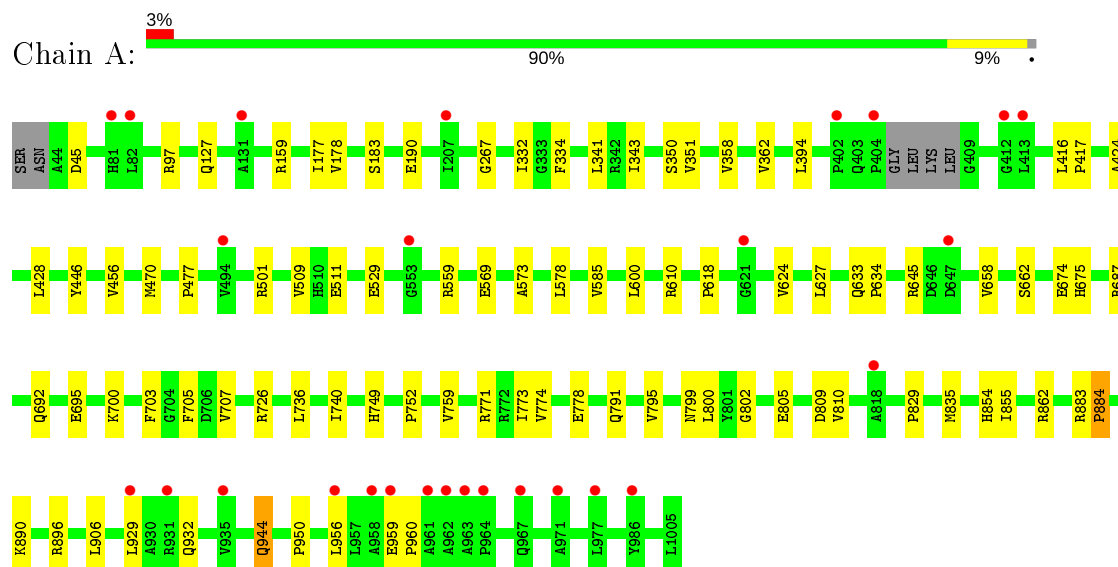
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	37	Total O 37 37	0	0
8	B	40	Total O 40 40	0	0

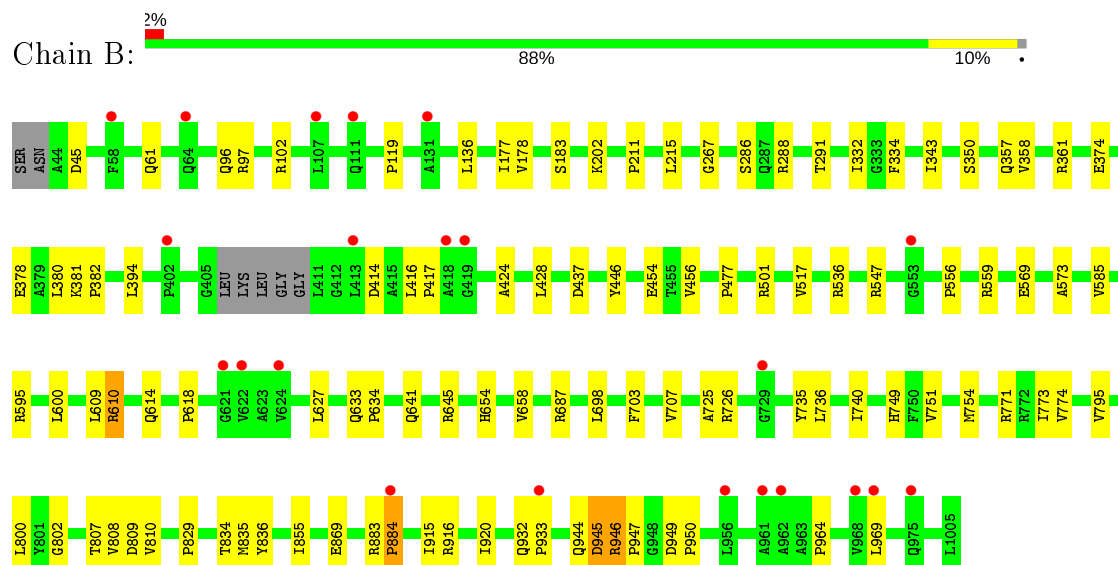
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Txo2



#### • Molecule 1: Txo2



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.86Å 399.88Å 144.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.20 – 2.95 48.48 – 2.95	Depositor EDS
% Data completeness (in resolution range)	97.5 (47.20-2.95) 97.5 (48.48-2.95)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 2.96Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, $R_{free}$	0.226 , 0.267 0.226 , 0.268	Depositor DCC
$R_{free}$ test set	3241 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.8	Xtriage
Anisotropy	0.717	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 45.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	15057	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.17 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1893e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, CL, FMT, SO4, ACT, FLC

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	0.23	0/7599	0.41	0/10345
1	B	0.23	0/7611	0.41	0/10361
All	All	0.23	0/15210	0.41	0/20706

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

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	7432	0	7410	43	0
1	B	7444	0	7424	45	0
2	A	3	0	1	0	0
2	B	12	0	4	0	0
3	A	13	0	5	0	0
3	B	13	0	5	1	0
4	A	4	0	3	0	0
5	A	6	0	0	0	0
5	B	13	0	0	0	0
6	A	30	0	40	0	0
7	A	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	A	37	0	0	0	0
8	B	40	0	0	0	0
All	All	15057	0	14892	88	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 3.

The worst 5 of 88 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:B:547:ARG:NH1	1:B:641:GLN:O	2.29	0.66
1:A:127:GLN:OE1	1:A:159:ARG:NH1	2.34	0.61
1:B:610:ARG:HD3	1:B:614:GLN:HB2	1.82	0.60
1:B:97:ARG:NH2	1:B:211:PRO:O	2.36	0.58
1:B:177:ILE:HG23	1:B:178:VAL:HG13	1.85	0.58

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	955/964 (99%)	927 (97%)	27 (3%)	1 (0%)	51 83
1	B	955/964 (99%)	924 (97%)	25 (3%)	6 (1%)	25 60
All	All	1910/1928 (99%)	1851 (97%)	52 (3%)	7 (0%)	34 69

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	884	PRO
1	B	807	THR

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Mol	Chain	Res	Type
1	B	808	VAL
1	B	884	PRO
1	B	119	PRO

### 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	755/762 (99%)	745 (99%)	10 (1%)	69	87
1	B	758/762 (100%)	739 (98%)	19 (2%)	47	76
All	All	1513/1524 (99%)	1484 (98%)	29 (2%)	57	81

5 of 29 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	202	LYS
1	B	334	PHE
1	B	916	ARG
1	B	215	LEU
1	B	381	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 34 ligands modelled in this entry, 19 are monoatomic - leaving 15 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	GOL	A	1110	-	5,5,5	0.89	0	5,5,5	1.01	0
6	GOL	A	1113	-	5,5,5	0.91	0	5,5,5	0.98	0
6	GOL	A	1114	-	5,5,5	0.91	0	5,5,5	1.01	0
7	SO4	A	1115	-	4,4,4	0.14	0	6,6,6	0.05	0
2	FMT	B	1101	-	0,2,2	0.00	-	0,1,1	0.00	-
6	GOL	A	1112	-	5,5,5	0.91	0	5,5,5	1.00	0
2	FMT	B	1102	-	0,2,2	0.00	-	0,1,1	0.00	-
6	GOL	A	1111	-	5,5,5	0.93	0	5,5,5	0.99	0
2	FMT	A	1101	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FMT	B	1103	-	0,2,2	0.00	-	0,1,1	0.00	-
2	FMT	B	1104	-	0,2,2	0.00	-	0,1,1	0.00	-
4	ACT	A	1103	-	1,3,3	6.50	1 (100%)	0,3,3	0.00	-
3	FLC	B	1105	-	3,12,12	1.56	0	3,17,17	2.77	2 (66%)
7	SO4	A	1116	-	4,4,4	0.14	0	6,6,6	0.06	0
3	FLC	A	1102	-	3,12,12	1.53	0	3,17,17	2.69	2 (66%)

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	GOL	A	1110	-	-	1/4/4/4	-
6	GOL	A	1113	-	-	0/4/4/4	-
3	FLC	B	1105	-	-	3/6/16/16	-
6	GOL	A	1112	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	A	1111	-	-	0/4/4/4	-
6	GOL	A	1114	-	-	2/4/4/4	-
3	FLC	A	1102	-	-	2/6/16/16	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1103	ACT	CH3-C	6.50	1.57	1.48

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1105	FLC	CB-CA-CAC	-3.56	109.28	114.98
3	A	1102	FLC	CB-CA-CAC	-3.36	109.60	114.98
3	A	1102	FLC	CB-CG-CGC	-3.22	109.83	114.98
3	B	1105	FLC	CB-CG-CGC	-3.21	109.84	114.98

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1114	GOL	C1-C2-C3-O3
3	B	1105	FLC	CA-CB-CG-CGC
6	A	1114	GOL	O2-C2-C3-O3
3	A	1102	FLC	CA-CB-CG-CGC
3	B	1105	FLC	CBC-CB-CG-CGC

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1105	FLC	1	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	958/964 (99%)	0.06	27 (2%)	53 36	36, 68, 116, 166	0
1	B	957/964 (99%)	0.02	22 (2%)	60 43	35, 68, 115, 168	0
All	All	1915/1928 (99%)	0.04	49 (2%)	56 39	35, 68, 116, 168	0

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	404	PRO	5.0
1	A	962	ALA	4.7
1	A	131	ALA	4.7
1	B	131	ALA	3.9
1	B	962	ALA	3.9

### 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 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<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
5	CL	A	1107	1/1	0.70	0.49	83,83,83,83	0
2	FMT	A	1101	3/3	0.73	0.35	75,75,82,82	0
6	GOL	A	1114	6/6	0.74	0.43	46,76,85,95	0
7	SO4	A	1115	5/5	0.76	0.34	121,124,128,132	0
2	FMT	B	1103	3/3	0.80	0.25	51,51,55,56	0
6	GOL	A	1111	6/6	0.80	0.28	47,77,89,90	0
2	FMT	B	1102	3/3	0.81	0.20	74,74,76,79	0
7	SO4	A	1116	5/5	0.82	0.31	109,111,115,117	5
5	CL	B	1113	1/1	0.82	0.50	89,89,89,89	0
5	CL	A	1108	1/1	0.83	0.38	105,105,105,105	0
6	GOL	A	1112	6/6	0.85	0.24	62,75,81,85	0
2	FMT	B	1101	3/3	0.85	0.28	51,51,60,63	0
5	CL	B	1111	1/1	0.86	0.14	70,70,70,70	1
5	CL	B	1112	1/1	0.86	0.12	83,83,83,83	0
5	CL	B	1106	1/1	0.86	0.10	77,77,77,77	0
5	CL	A	1106	1/1	0.87	0.19	99,99,99,99	0
5	CL	B	1115	1/1	0.87	0.23	72,72,72,72	0
5	CL	B	1108	1/1	0.88	0.13	91,91,91,91	0
4	ACT	A	1103	4/4	0.88	0.22	51,57,57,62	0
5	CL	B	1117	1/1	0.89	0.16	74,74,74,74	0
5	CL	A	1104	1/1	0.90	0.08	69,69,69,69	0
5	CL	B	1107	1/1	0.90	0.34	75,75,75,75	0
6	GOL	A	1110	6/6	0.90	0.19	41,58,68,70	0
2	FMT	B	1104	3/3	0.91	0.34	64,64,72,73	0
5	CL	A	1105	1/1	0.92	0.20	76,76,76,76	0
5	CL	B	1110	1/1	0.92	0.20	91,91,91,91	0
5	CL	B	1116	1/1	0.93	0.12	72,72,72,72	0
5	CL	A	1109	1/1	0.93	0.11	53,53,53,53	0
5	CL	B	1118	1/1	0.93	0.12	85,85,85,85	0
3	FLC	B	1105	13/13	0.93	0.16	46,53,70,73	0
6	GOL	A	1113	6/6	0.93	0.21	55,62,76,77	0
5	CL	B	1114	1/1	0.94	0.14	83,83,83,83	0
3	FLC	A	1102	13/13	0.94	0.19	44,52,71,75	0
5	CL	B	1109	1/1	0.95	0.09	64,64,64,64	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.