



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 17, 2017 – 02:37 PM EST

PDB ID : 5IFU  
Title : Crystal Structure of Prolyl-tRNA synthetase (ProRS, Proline-tRNA ligase) from Plasmodium falciparum in complex with Glyburide  
Authors : Dranow, D.M.; Hewitt, S.N.; Abendroth, J.; Structural Genomics Consortium (SGC)  
Deposited on : 2016-02-26  
Resolution : 2.45 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](mailto: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.

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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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20028442  
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) : rb-20028442

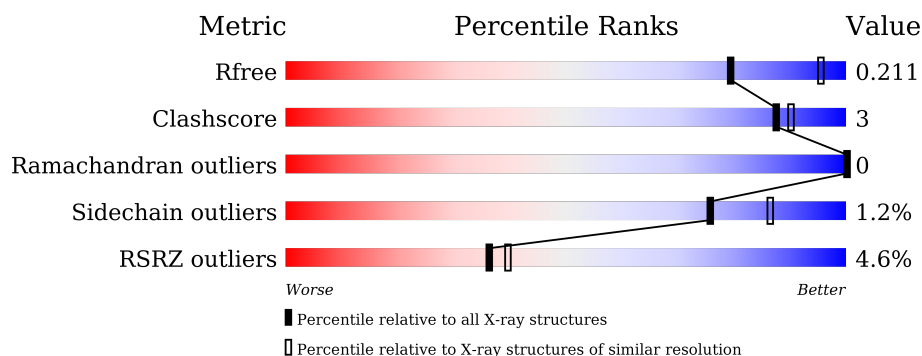
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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. 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	506	<div> <div>2%</div> <div>81%</div> <div>5%</div> <div>13%</div> </div>
1	B	506	<div> <div>6%</div> <div>79%</div> <div>8%</div> <div>12%</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
2	GBM	A	801	-	-	-	X
4	FMT	A	803	-	-	-	X
4	FMT	A	804	-	-	-	X
4	FMT	A	806	-	-	-	X
4	FMT	A	808	-	-	-	X
4	FMT	A	809	-	-	-	X
4	FMT	A	811	-	-	-	X
4	FMT	A	812	-	-	-	X
4	FMT	B	802	-	-	-	X
4	FMT	B	804	-	-	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7313 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.

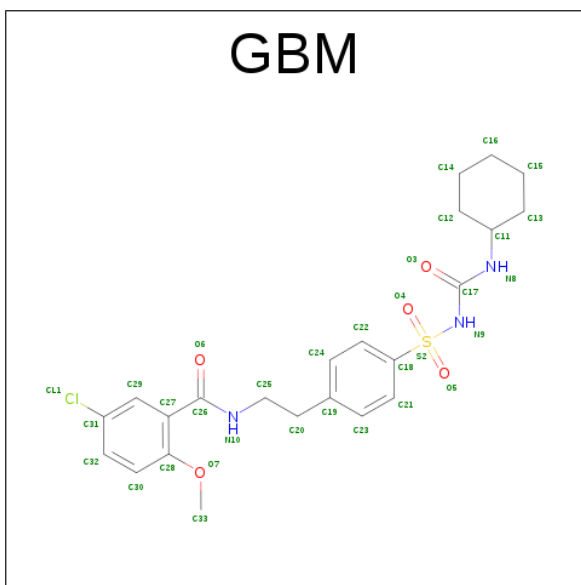
- Molecule 1 is a protein called Proline-tRNA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	438	Total	C	N	O	S	0	4	0
			3528	2273	585	648	22			
1	B	445	Total	C	N	O	S	0	0	0
			3438	2214	573	631	20			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	MET	-	expression tag	UNP Q8I5R7
A	242	ALA	-	expression tag	UNP Q8I5R7
A	243	HIS	-	expression tag	UNP Q8I5R7
A	244	HIS	-	expression tag	UNP Q8I5R7
A	245	HIS	-	expression tag	UNP Q8I5R7
A	246	HIS	-	expression tag	UNP Q8I5R7
A	247	HIS	-	expression tag	UNP Q8I5R7
A	248	HIS	-	expression tag	UNP Q8I5R7
B	241	MET	-	expression tag	UNP Q8I5R7
B	242	ALA	-	expression tag	UNP Q8I5R7
B	243	HIS	-	expression tag	UNP Q8I5R7
B	244	HIS	-	expression tag	UNP Q8I5R7
B	245	HIS	-	expression tag	UNP Q8I5R7
B	246	HIS	-	expression tag	UNP Q8I5R7
B	247	HIS	-	expression tag	UNP Q8I5R7
B	248	HIS	-	expression tag	UNP Q8I5R7

- Molecule 2 is 5-chloro-N-(2-{4-[(cyclohexylcarbamoyl)sulfamoyl]phenyl}ethyl)-2-methoxybenzamide (three-letter code: GBM) (formula: C<sub>23</sub>H<sub>28</sub>ClN<sub>3</sub>O<sub>5</sub>S).

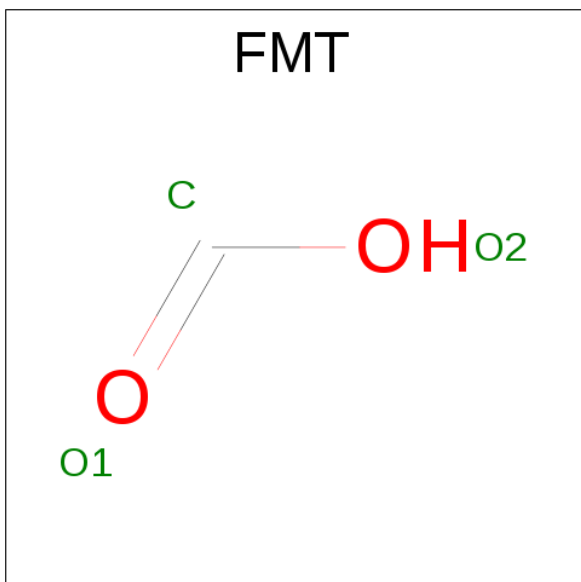


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	S	0	0
			33	23	1	3	5	1		
2	B	1	Total	C	N	O	S	0	0	
			20	14	2	3	1			

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

- Molecule 4 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

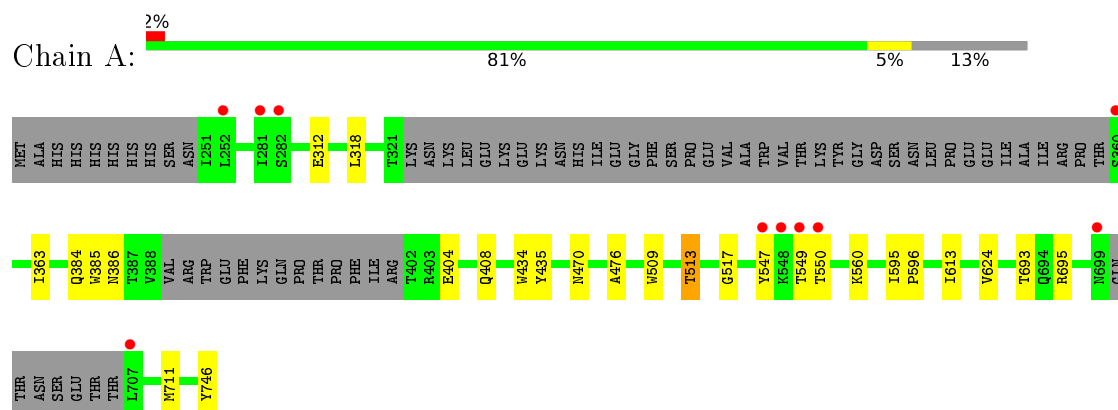
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	191	Total	O	0	0
			191	191		
6	B	52	Total	O	0	0
			52	52		

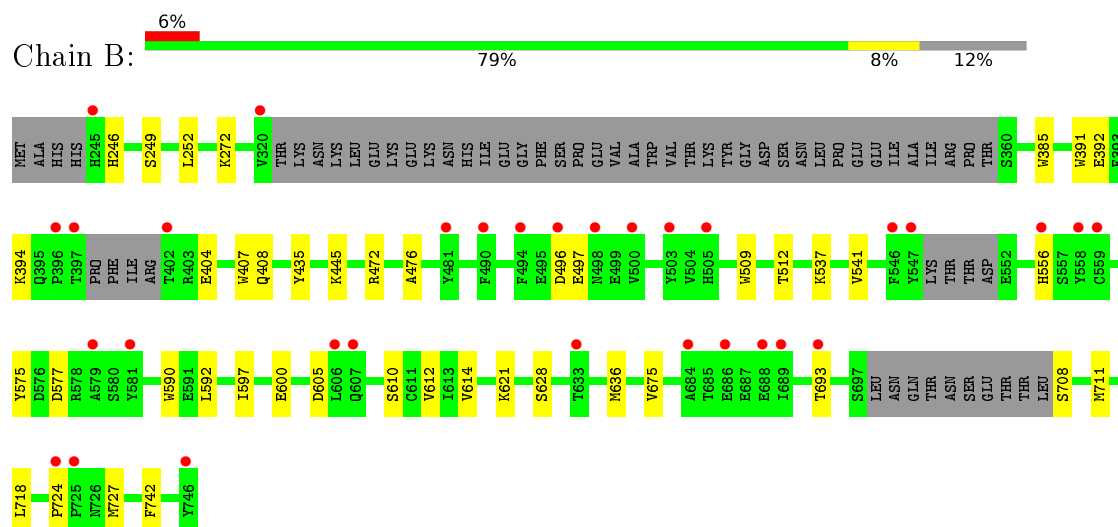
### 3 Residue-property plots [i](#)

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: Proline-tRNA ligase



#### • Molecule 1: Proline-tRNA ligase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	138.31Å 138.31Å 156.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.74 – 2.45 43.74 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.74-2.45) 96.5 (43.74-2.45)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.03 (at 2.45Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.185 , 0.213 0.182 , 0.211	Depositor DCC
$R_{free}$ test set	2687 reflections (4.94%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.2	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 47.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7313	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FMT, GBM, EDO, 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	0.24	0/3623	0.40	0/4906
1	B	0.22	0/3525	0.38	0/4795
All	All	0.23	0/7148	0.39	0/9701

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	3528	0	3429	13	0
1	B	3438	0	3188	23	0
2	A	33	0	28	1	0
2	B	20	0	17	1	0
3	A	1	0	0	0	0
4	A	30	0	10	0	0
4	B	12	0	4	0	0
5	A	4	0	6	0	0
5	B	4	0	6	0	0
6	A	191	0	0	1	0
6	B	52	0	0	1	0
All	All	7313	0	6688	35	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.

All (35) 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:272:LYS:HB3	1:B:592:LEU:HD21	1.76	0.66
1:B:392:GLU:O	6:B:901:HOH:O	2.14	0.65
1:B:693:THR:HG21	1:B:711:MET:HG2	1.84	0.60
1:B:496:ASP:OD1	1:B:497:GLU:N	2.36	0.59
1:B:435:TYR:CZ	1:B:476:ALA:HB1	2.39	0.58
1:B:407:TRP:HB2	1:B:512:THR:HG22	1.87	0.56
1:B:404:GLU:OE2	2:B:801:GBM:N9	2.33	0.56
1:A:312:GLU:OE2	1:B:537:LYS:NZ	2.32	0.55
1:A:385:TRP:CZ3	1:A:408:GLN:HB2	2.44	0.53
1:A:693:THR:HG21	1:A:711:MET:HG2	1.93	0.51
1:B:600:GLU:HB2	1:B:612:VAL:HB	1.92	0.51
1:B:249:SER:HB2	1:B:252:LEU:HG	1.94	0.49
1:A:318:LEU:HG	1:A:363:ILE:HD12	1.94	0.48
1:A:695:ARG:NH2	6:A:904:HOH:O	2.47	0.48
1:B:614:VAL:HG22	1:B:621:LYS:HG2	1.97	0.47
1:A:435:TYR:CZ	1:A:476:ALA:HB1	2.50	0.47
1:A:517:GLY:HA3	2:A:801:GBM:O3	2.15	0.47
1:B:249:SER:H	1:B:252:LEU:HD12	1.79	0.46
1:B:541:VAL:HG21	1:B:590:TRP:CD1	2.51	0.46
1:A:404:GLU:HG3	1:A:513:THR:OG1	2.16	0.46
1:B:385:TRP:CZ3	1:B:408:GLN:HB2	2.52	0.45
1:B:445:LYS:HB3	1:B:718:LEU:HD12	1.98	0.44
1:B:472:ARG:NH2	1:B:708:SER:O	2.50	0.44
1:A:384[B]:GLN:NE2	1:A:386:ASN:OD1	2.45	0.44
1:A:547:TYR:C	1:A:549:THR:H	2.22	0.43
1:B:391:TRP:HB3	1:B:394:LYS:HE3	2.00	0.43
1:B:724:PRO:HB2	1:B:727:MET:HB2	2.00	0.42
1:B:675:VAL:HG23	1:B:742:PHE:HB2	2.01	0.42
1:A:613:ILE:HD12	1:A:624:VAL:HG21	2.01	0.42
1:A:595:ILE:HA	1:A:596:PRO:HD3	1.85	0.42
1:B:249:SER:N	1:B:252:LEU:HD12	2.35	0.41
1:B:597:ILE:HG23	1:B:636:MET:HE2	2.03	0.41
1:B:575:TYR:CE2	1:B:577:ASP:HB3	2.56	0.41
1:A:470:ASN:ND2	1:A:746:TYR:OH	2.49	0.40
1:B:605:ASP:HB3	1:B:610:SER:O	2.20	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	434/506 (86%)	421 (97%)	13 (3%)	0	100	100
1	B	435/506 (86%)	425 (98%)	10 (2%)	0	100	100
All	All	869/1012 (86%)	846 (97%)	23 (3%)	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	379/460 (82%)	374 (99%)	5 (1%)	76	86
1	B	345/460 (75%)	341 (99%)	4 (1%)	78	87
All	All	724/920 (79%)	715 (99%)	9 (1%)	76	87

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	434	TRP
1	A	509	TRP
1	A	513	THR
1	A	550	THR
1	A	560	LYS
1	B	246	HIS
1	B	509	TRP
1	B	556	HIS

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Mol	Chain	Res	Type
1	B	628	SER

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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 19 ligands modelled in this entry, 1 is monoatomic - leaving 18 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	GBM	A	801	-	35,35,35	1.96	10 (28%)	48,48,48	2.63	10 (20%)
4	FMT	A	803	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	804	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	805	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	806	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	807	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	808	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	809	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	810	-	0,2,2	0.00	-	0,1,1	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	FMT	A	811	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	A	812	-	0,2,2	0.00	-	0,1,1	0.00	-
5	EDO	A	813	-	3,3,3	0.46	0	2,2,2	0.30	0
2	GBM	B	801	-	21,21,35	1.95	7 (33%)	29,29,48	3.03	6 (20%)
4	FMT	B	802	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	803	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	804	-	0,2,2	0.00	-	0,1,1	0.00	-
4	FMT	B	805	-	0,2,2	0.00	-	0,1,1	0.00	-
5	EDO	B	806	-	3,3,3	0.46	0	2,2,2	0.26	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	GBM	A	801	-	-	0/27/35/35	0/3/3/3
4	FMT	A	803	-	-	0/0/0/0	0/0/0/0
4	FMT	A	804	-	-	0/0/0/0	0/0/0/0
4	FMT	A	805	-	-	0/0/0/0	0/0/0/0
4	FMT	A	806	-	-	0/0/0/0	0/0/0/0
4	FMT	A	807	-	-	0/0/0/0	0/0/0/0
4	FMT	A	808	-	-	0/0/0/0	0/0/0/0
4	FMT	A	809	-	-	0/0/0/0	0/0/0/0
4	FMT	A	810	-	-	0/0/0/0	0/0/0/0
4	FMT	A	811	-	-	0/0/0/0	0/0/0/0
4	FMT	A	812	-	-	0/0/0/0	0/0/0/0
5	EDO	A	813	-	-	0/1/1/1	0/0/0/0
2	GBM	B	801	-	-	0/15/23/35	0/2/2/3
4	FMT	B	802	-	-	0/0/0/0	0/0/0/0
4	FMT	B	803	-	-	0/0/0/0	0/0/0/0
4	FMT	B	804	-	-	0/0/0/0	0/0/0/0
4	FMT	B	805	-	-	0/0/0/0	0/0/0/0
5	EDO	B	806	-	-	0/1/1/1	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	GBM	O3-C17	-2.42	1.18	1.23
2	A	801	GBM	O6-C26	-2.33	1.18	1.23
2	B	801	GBM	O3-C17	-2.26	1.18	1.23
2	A	801	GBM	O7-C28	2.19	1.40	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	GBM	O4-S2	2.86	1.46	1.43
2	B	801	GBM	O4-S2	2.97	1.46	1.43
2	A	801	GBM	O5-S2	3.10	1.46	1.43
2	B	801	GBM	O5-S2	3.10	1.46	1.43
2	B	801	GBM	S2-N9	3.54	1.71	1.64
2	B	801	GBM	C17-N9	3.65	1.45	1.39
2	B	801	GBM	C17-N8	3.68	1.45	1.35
2	A	801	GBM	C17-N8	3.74	1.45	1.35
2	A	801	GBM	C18-S2	3.76	1.82	1.76
2	B	801	GBM	C18-S2	3.91	1.82	1.76
2	A	801	GBM	S2-N9	3.92	1.72	1.64
2	A	801	GBM	C17-N9	4.04	1.46	1.39
2	A	801	GBM	C26-N10	5.57	1.45	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	GBM	O4-S2-O5	-14.54	100.26	119.54
2	B	801	GBM	O4-S2-O5	-13.98	101.00	119.54
2	A	801	GBM	C28-C27-C26	-3.29	120.44	126.13
2	A	801	GBM	O7-C28-C30	-2.62	119.96	124.35
2	A	801	GBM	C11-N8-C17	-2.27	118.05	123.38
2	A	801	GBM	C17-N9-S2	-2.10	119.09	123.66
2	B	801	GBM	C17-N9-S2	-2.07	119.16	123.66
2	A	801	GBM	C33-O7-C28	-2.02	114.57	117.53
2	B	801	GBM	O4-S2-C18	2.19	110.73	107.94
2	A	801	GBM	O5-S2-C18	2.99	111.75	107.94
2	A	801	GBM	O7-C28-C27	3.02	120.26	116.52
2	B	801	GBM	O5-S2-C18	3.14	111.94	107.94
2	B	801	GBM	N9-C17-N8	3.47	120.90	114.48
2	A	801	GBM	C18-S2-N9	3.59	110.33	105.97
2	B	801	GBM	C18-S2-N9	4.71	111.69	105.97
2	A	801	GBM	N9-C17-N8	4.72	123.21	114.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	GBM	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	GBM	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

### 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<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	438/506 (86%)	-0.24	10 (2%) 64 66	28, 43, 79, 141	0
1	B	445/506 (87%)	0.22	31 (6%) 19 21	35, 74, 116, 135	0
All	All	883/1012 (87%)	-0.01	41 (4%) 36 40	28, 55, 111, 141	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	547	TYR	5.6
1	B	558	TYR	5.0
1	B	684	ALA	4.1
1	B	320	VAL	4.0
1	B	397	THR	3.9
1	A	281	ILE	3.9
1	B	606	LEU	3.8
1	B	503	TYR	3.5
1	B	689	ILE	3.4
1	A	707	LEU	3.3
1	A	282	SER	3.2
1	B	725	PRO	3.2
1	B	245	HIS	3.1
1	B	498	ASN	2.8
1	B	693	THR	2.8
1	A	548	LYS	2.8
1	A	550	THR	2.7
1	B	559	CYS	2.7
1	A	549	THR	2.6
1	B	496	ASP	2.6
1	B	546	PHE	2.6
1	B	396	PRO	2.5
1	B	686	GLU	2.5
1	B	746	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	607	GLN	2.5
1	B	402	THR	2.4
1	B	494	PHE	2.4
1	B	581	TYR	2.4
1	B	579	ALA	2.4
1	B	481	TYR	2.3
1	B	688	GLU	2.2
1	B	556	HIS	2.2
1	B	724	PRO	2.2
1	B	500	VAL	2.2
1	B	633	THR	2.2
1	A	699	ASN	2.1
1	B	505	HIS	2.1
1	A	360	SER	2.1
1	B	490	PHE	2.0
1	B	547	TYR	2.0
1	A	252	LEU	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 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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	FMT	A	811	3/3	0.72	0.33	25.17	67,67,70,70	0
4	FMT	A	806	3/3	0.71	0.33	9.25	76,76,77,78	0
4	FMT	B	804	3/3	0.84	0.46	7.92	82,82,84,84	0
4	FMT	A	804	3/3	0.84	0.22	6.51	80,80,80,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	FMT	A	812	3/3	0.89	0.24	5.77	84,84,85,86	0
4	FMT	A	803	3/3	0.73	0.33	4.58	55,55,58,59	0
4	FMT	B	802	3/3	0.85	0.25	4.25	64,64,66,67	0
2	GBM	A	801	33/33	0.82	0.29	4.13	53,84,128,129	0
4	FMT	A	809	3/3	0.90	0.18	2.59	57,57,61,64	0
4	FMT	A	808	3/3	0.65	0.18	2.12	81,81,82,82	0
2	GBM	B	801	20/33	0.94	0.17	0.72	44,57,66,67	0
3	CL	A	802	1/1	0.99	0.12	-0.83	39,39,39,39	0
4	FMT	B	803	3/3	0.91	0.21	-	67,67,68,68	0
4	FMT	A	805	3/3	0.92	0.17	-	70,70,71,71	0
4	FMT	B	805	3/3	0.76	0.19	-	74,74,75,76	0
5	EDO	B	806	4/4	0.74	0.44	-	75,79,80,83	0
4	FMT	A	807	3/3	0.87	0.61	-	74,74,79,80	0
5	EDO	A	813	4/4	0.71	0.41	-	78,78,78,79	0
4	FMT	A	810	3/3	0.91	0.22	-	78,78,81,82	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.