



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2018 – 11:36 pm GMT

PDB ID : 2CV2  
Title : Glutamyl-tRNA synthetase from *Thermus thermophilus* in complex with tRNA(Glu) and an enzyme inhibitor, Glu-AMS  
Authors : Sekine, S.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)  
Deposited on : 2005-05-31  
Resolution : 2.69 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	:	1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix)	:	1.13
EDS	:	trunk30967
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk30967

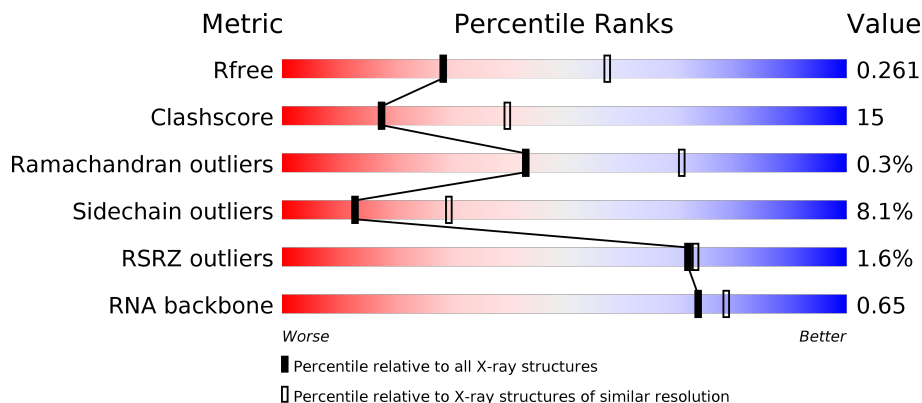
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.69 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))
$R_{free}$	111664	2449 (2.70-2.70)
Clashscore	122126	2756 (2.70-2.70)
Ramachandran outliers	120053	2716 (2.70-2.70)
Sidechain outliers	120020	2716 (2.70-2.70)
RSRZ outliers	108989	2376 (2.70-2.70)
RNA backbone	2636	1009 (3.00-2.40)

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 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	C	75	<div> <div>0%</div> <div> <div></div> <div>52%</div> <div>32%</div> <div>12%</div> <div>•</div> </div> </div>
1	D	75	<div> <div>0%</div> <div> <div></div> <div>53%</div> <div>31%</div> <div>12%</div> <div>•</div> </div> </div>
2	A	468	<div> <div>0%</div> <div> <div></div> <div>68%</div> <div>29%</div> <div>•</div> </div> </div>
2	B	468	<div> <div>2%</div> <div> <div></div> <div>61%</div> <div>34%</div> <div>5%</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11215 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 RNA chain called tRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	75	Total	C	N	O	P	0	0	0
			1597	711	284	527	75			
1	D	75	Total	C	N	O	P	0	0	0
			1597	711	284	527	75			

- Molecule 2 is a protein called glutamyl-tRNA synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	468	Total	C	N	O	S	0	0	0
			3814	2443	676	687	8			
2	B	468	Total	C	N	O	S	0	0	0
			3814	2443	676	687	8			

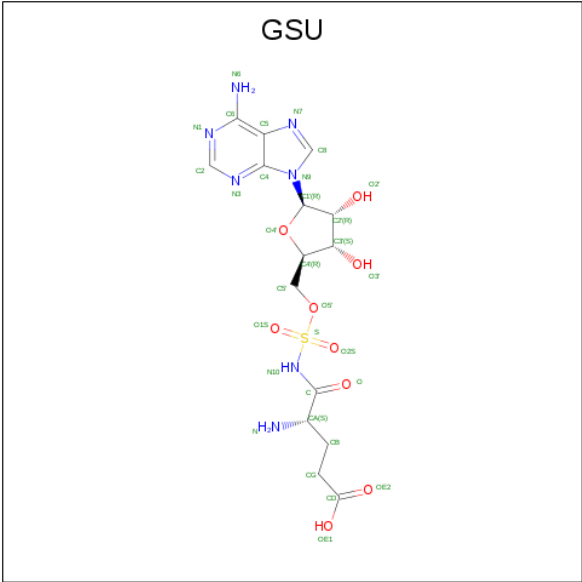
- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Cl	0	0
			1	1		
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is O5'-(L-GLUTAMYL-SULFAMOYL)-ADENOSINE (three-letter code: GSU) (formula: C<sub>15</sub>H<sub>21</sub>N<sub>7</sub>O<sub>9</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			32	15	7	9	1		
5	B	1	Total	C	N	O	S	0	0
			32	15	7	9	1		

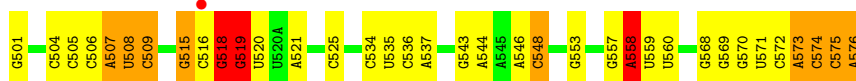
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	48	Total	O	0	0
			48	48		
6	D	36	Total	O	0	0
			36	36		
6	A	125	Total	O	0	0
			125	125		
6	B	116	Total	O	0	0
			116	116		

### 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.

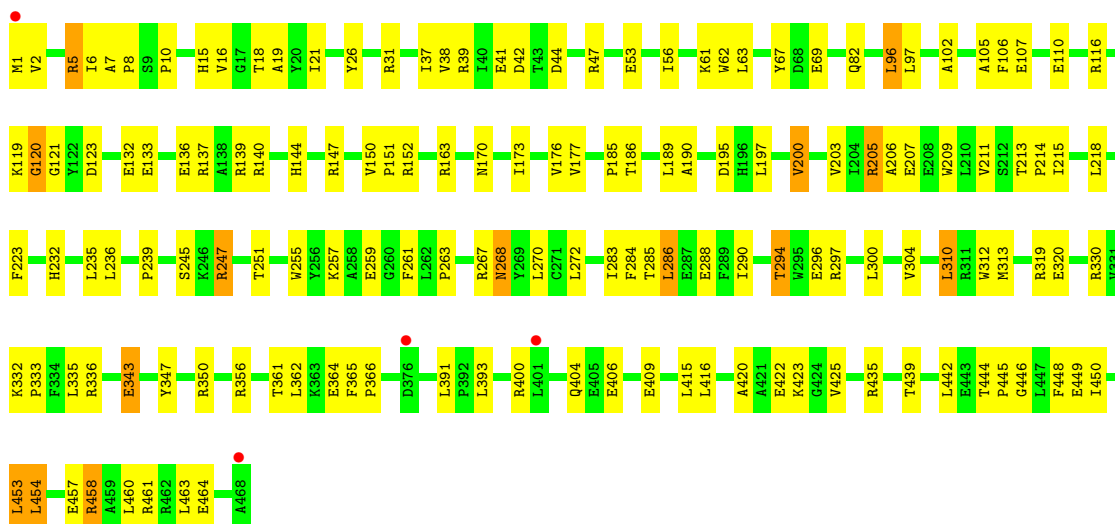
#### • Molecule 1: tRNA



#### • Molecule 1: tRNA

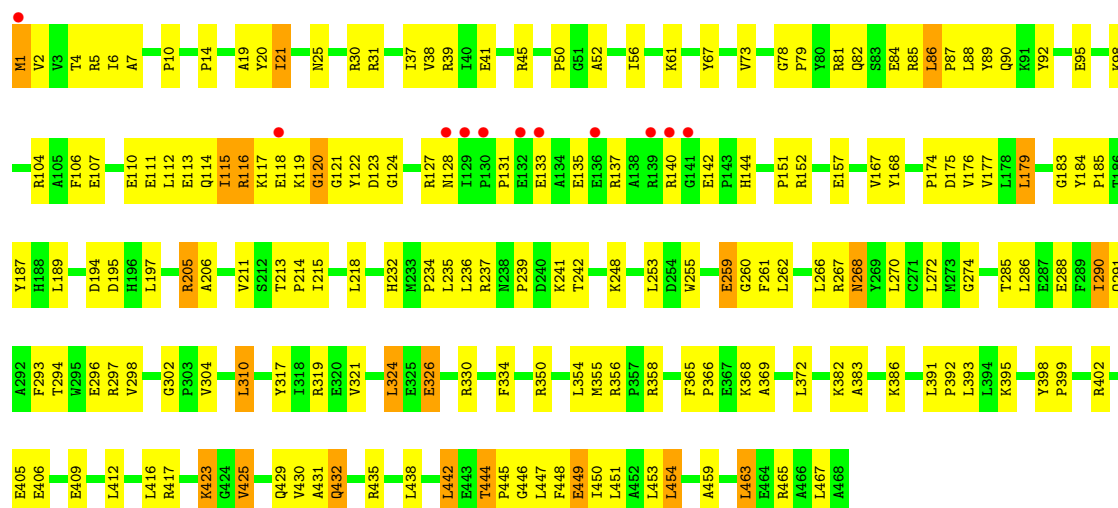


#### • Molecule 2: glutamyl-tRNA synthetase



#### • Molecule 2: glutamyl-tRNA synthetase





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	110.32Å 219.82Å 135.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.34 – 2.69 46.34 – 2.63	Depositor EDS
% Data completeness (in resolution range)	99.4 (46.34-2.69) 96.4 (46.34-2.63)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.14 (at 2.65Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.200 , 0.260 0.201 , 0.261	Depositor DCC
$R_{free}$ test set	2378 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.0	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 53.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	11215	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.61% 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: MG, GSU, 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	C	0.55	1/1782 (0.1%)	0.86	9/2774 (0.3%)
1	D	0.54	1/1782 (0.1%)	0.87	9/2774 (0.3%)
2	A	0.52	0/3910	0.70	0/5293
2	B	0.52	0/3910	0.68	0/5293
All	All	0.53	2/11384 (0.0%)	0.75	18/16134 (0.1%)

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	C	1	3
1	D	1	3
All	All	2	6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	501	G	OP3-P	-7.12	1.52	1.61
1	D	501	G	OP3-P	-5.85	1.54	1.61

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	558	A	C2'-C3'-O3'	8.16	127.46	109.50
1	C	558	A	C2'-C3'-O3'	7.83	126.72	109.50
1	D	507	A	C2'-C3'-O3'	7.76	126.56	109.50
1	C	573	A	C2'-C3'-O3'	7.31	125.58	109.50
1	D	573	A	C2'-C3'-O3'	7.30	125.57	109.50



All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	558	A	C3'
1	D	558	A	C3'

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	515	G	Sidechain
1	C	535	U	Sidechain
1	C	568	G	Sidechain
1	D	515	G	Sidechain
1	D	522	G	Sidechain

## 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	C	1597	0	813	19	0
1	D	1597	0	813	19	0
2	A	3814	0	3818	115	1
2	B	3814	0	3818	157	1
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	32	0	20	2	0
5	B	32	0	20	2	0
6	A	125	0	0	6	0
6	B	116	0	0	6	0
6	C	48	0	0	0	0
6	D	36	0	0	0	0
All	All	11215	0	9302	297	1

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 15.

The worst 5 of 297 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)
2:A:247:ARG:H	2:A:247:ARG:HD3	1.29	0.96
2:A:177:VAL:HG23	2:A:186:THR:HG21	1.49	0.95
2:A:163:ARG:HH12	2:A:300:LEU:HD22	1.31	0.92
2:B:454:LEU:H	2:B:454:LEU:HD22	1.38	0.86
2:B:435:ARG:HH11	2:B:444:THR:CG2	1.89	0.85

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:343:GLU:O	2:B:50:PRO:O[6_654]	2.16	0.04

## 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	
2	A	466/468 (100%)	453 (97%)	12 (3%)	1 (0%)	49	77
2	B	466/468 (100%)	448 (96%)	16 (3%)	2 (0%)	36	64
All	All	932/936 (100%)	901 (97%)	28 (3%)	3 (0%)	43	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	120	GLY
2	B	120	GLY
2	B	151	PRO

### 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	
2	A	393/393 (100%)	366 (93%)	27 (7%)	17	39
2	B	393/393 (100%)	356 (91%)	37 (9%)	9	22
All	All	786/786 (100%)	722 (92%)	64 (8%)	13	29

5 of 64 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	5	ARG
2	B	123	ASP
2	B	442	LEU
2	B	21	ILE
2	B	115	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	25	ASN
2	B	432	GLN
2	B	191	ASN
2	A	404	GLN
2	B	90	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	74/75 (98%)	14 (18%)	6 (8%)
1	D	74/75 (98%)	13 (17%)	6 (8%)
All	All	148/150 (98%)	27 (18%)	12 (8%)

5 of 27 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	C	507	A
1	C	508	U
1	C	509	C
1	C	516	C
1	C	518	G

5 of 12 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	C	573	A
1	D	507	A
1	D	558	A
1	C	558	A
1	D	519	G

## 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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$
5	GSU	A	601	-	27,34,34	2.04	4 (14%)	27,50,50	2.43	4 (14%)
5	GSU	B	602	-	27,34,34	1.56	3 (11%)	27,50,50	2.30	6 (22%)

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
5	GSU	A	601	-	-	0/16/40/40	0/3/3/3
5	GSU	B	602	-	-	0/16/40/40	0/3/3/3

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	602	GSU	O2S-S	2.60	1.44	1.42
5	A	601	GSU	C2-N3	2.65	1.36	1.32
5	B	602	GSU	S-N10	4.62	1.64	1.59
5	B	602	GSU	O1S-S	4.75	1.46	1.42
5	A	601	GSU	O1S-S	4.76	1.46	1.42

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	601	GSU	O2S-S-O1S	-9.92	111.16	121.23
5	B	602	GSU	O2S-S-O1S	-9.30	111.80	121.23
5	A	601	GSU	CB-CA-N	-3.08	102.62	110.50
5	B	602	GSU	CB-CA-N	-2.08	105.17	110.50
5	B	602	GSU	C4-C5-N7	2.09	111.43	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	601	GSU	2	0
5	B	602	GSU	2	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, 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	C	75/75 (100%)	-0.26	1 (1%) 77 78	25, 43, 76, 83	0
1	D	75/75 (100%)	-0.24	1 (1%) 77 78	21, 41, 75, 96	0
2	A	468/468 (100%)	-0.32	4 (0%) 84 85	13, 29, 61, 75	0
2	B	468/468 (100%)	-0.21	11 (2%) 59 59	13, 32, 76, 95	0
All	All	1086/1086 (100%)	-0.26	17 (1%) 72 73	13, 32, 70, 96	0

The worst 5 of 17 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	128	ASN	5.3
1	D	516	C	4.4
2	B	140	ARG	3.9
2	A	468	ALA	3.5
2	B	141	GLY	3.2

### 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
3	MG	C	902	1/1	0.88	0.20	36,36,36,36	0
3	MG	D	901	1/1	0.94	0.27	33,33,33,33	0
4	CL	B	903	1/1	0.98	0.25	27,27,27,27	0
5	GSU	B	602	32/32	0.98	0.14	16,25,31,32	0
5	GSU	A	601	32/32	0.98	0.14	16,22,28,31	0
4	CL	A	904	1/1	0.99	0.30	33,33,33,33	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.