



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:26 am GMT

PDB ID : 1EFT  
Title : THE CRYSTAL STRUCTURE OF ELONGATION FACTOR EF-TU FROM THERMUS AQUATICUS IN THE GTP CONFORMATION  
Authors : Kjeldgaard, M.; Nissen, P.; Thirup, S.; Nyborg, J.  
Deposited on : 1993-08-24  
Resolution : 2.50 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

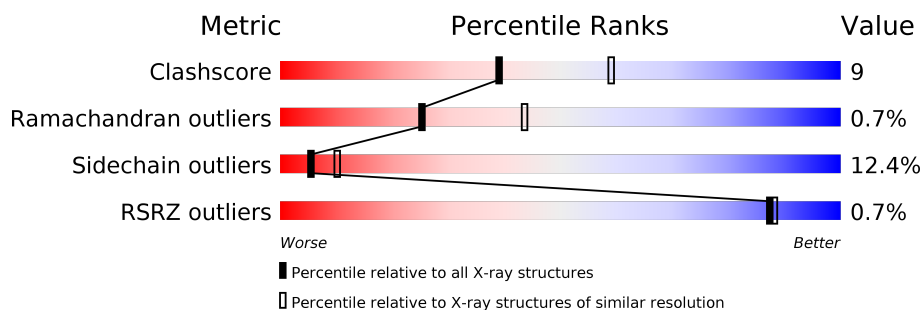
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	405	<div> <div></div> <div>73%</div> <div>22%</div> <div>.</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3302 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 ELONGATION FACTOR TU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3143	1985	548	598	12			

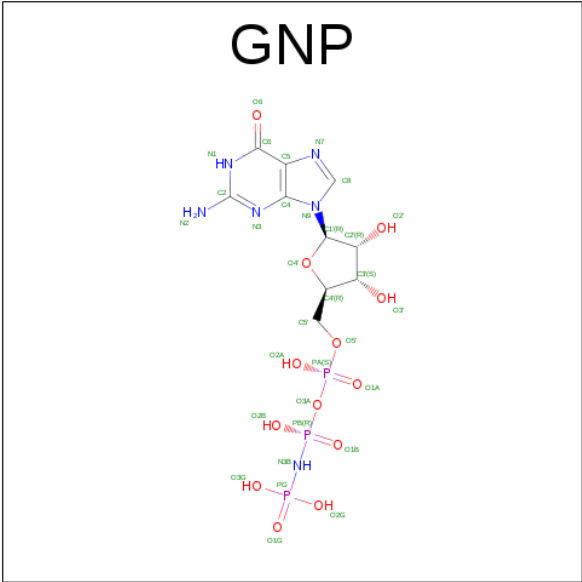
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	PHE	TYR	CONFLICT	UNP Q01698
A	35	THR	ALA	CONFLICT	UNP Q01698
A	322	VAL	ILE	CONFLICT	UNP Q01698
A	336	SER	THR	CONFLICT	UNP Q01698

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>6</sub>O<sub>13</sub>P<sub>3</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		

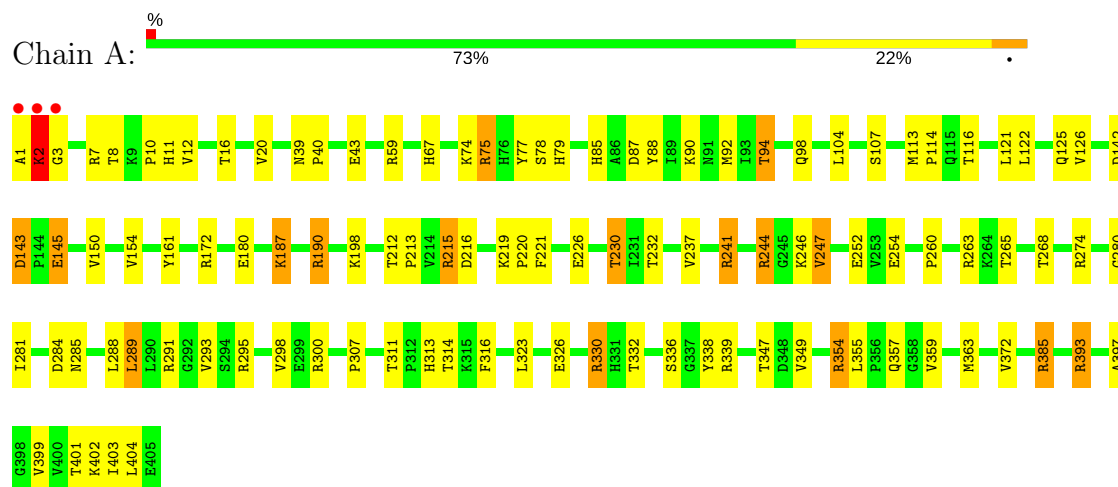
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	126	Total	O	0	0
			126	126		

### 3 Residue-property plots

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: ELONGATION FACTOR TU



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	150.70Å 99.70Å 40.20Å 90.00° 95.30° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50 15.07 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.50) 88.5 (15.07-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.25 (at 2.20Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.200 , (Not available) 0.193 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	33.6	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 68.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3302	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.28% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GNP, MG

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	0.57	0/3203	0.65	0/4343

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	A	0	5

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	ALA	Mainchain
1	A	161	TYR	Sidechain
1	A	338	TYR	Sidechain
1	A	339	ARG	Sidechain
1	A	357	GLN	Mainchain

### 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	3143	0	3158	54	0
2	A	1	0	0	0	0
3	A	32	0	13	1	0
4	A	126	0	0	1	0
All	All	3302	0	3171	54	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.

All (54) 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:2:LYS:HG2	1:A:274:ARG:HB3	1.52	0.91
1:A:220:PRO:O	1:A:244:ARG:HD3	1.84	0.77
1:A:85:HIS:HD2	1:A:87:ASP:H	1.38	0.71
1:A:190:ARG:HG3	1:A:190:ARG:HH11	1.57	0.70
1:A:237:VAL:HG22	1:A:289:LEU:HD22	1.73	0.69
1:A:2:LYS:HG3	1:A:3:GLY:N	2.08	0.69
1:A:254:GLU:HG3	1:A:307:PRO:HA	1.78	0.64
1:A:92:MET:HE1	1:A:122:LEU:HD12	1.81	0.62
1:A:121:LEU:HD21	1:A:323:LEU:HD11	1.82	0.61
1:A:2:LYS:HG3	1:A:3:GLY:H	1.64	0.61
1:A:190:ARG:NH1	1:A:190:ARG:HG3	2.16	0.60
1:A:2:LYS:HB3	1:A:274:ARG:O	2.05	0.57
1:A:150:VAL:O	1:A:154:VAL:HG23	2.05	0.56
1:A:316:PHE:CE1	1:A:372:VAL:HG22	2.42	0.55
1:A:90:LYS:O	1:A:94:THR:HG23	2.06	0.55
1:A:10:PRO:O	1:A:75:ARG:HB3	2.06	0.54
1:A:20:VAL:HA	3:A:406:GNP:O2G	2.07	0.54
1:A:94:THR:HG22	4:A:492:HOH:O	2.08	0.53
1:A:2:LYS:CG	1:A:274:ARG:HB3	2.34	0.52
1:A:143:ASP:OD1	1:A:145:GLU:HB2	2.08	0.52
1:A:221:PHE:CE1	1:A:247:VAL:HG22	2.45	0.51
1:A:254:GLU:HA	1:A:263:ARG:O	2.09	0.51
1:A:172:ARG:O	1:A:198:LYS:HD2	2.11	0.51
1:A:219:LYS:HB2	1:A:244:ARG:HG3	1.92	0.50
1:A:75:ARG:HD2	1:A:77:TYR:OH	2.11	0.50
1:A:11:HIS:O	1:A:215:ARG:NH2	2.45	0.49
1:A:246:LYS:HA	1:A:280:GLY:O	2.12	0.49
1:A:7:ARG:NH2	1:A:284:ASP:OD2	2.47	0.48
1:A:241:ARG:O	1:A:241:ARG:HG3	2.12	0.48
1:A:402:LYS:HG2	1:A:404:LEU:CD2	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:GLU:HG2	1:A:226:GLU:O	2.14	0.48
1:A:402:LYS:HG2	1:A:404:LEU:HD21	1.97	0.47
1:A:104:LEU:HD11	1:A:116:THR:HG23	1.96	0.47
1:A:187:LYS:HB2	1:A:187:LYS:HE2	1.65	0.47
1:A:354:ARG:HG3	1:A:354:ARG:NH1	2.30	0.46
1:A:212:THR:HA	1:A:213:PRO:HD3	1.74	0.45
1:A:355:LEU:HD13	1:A:359:VAL:HG12	1.97	0.45
1:A:11:HIS:HE1	1:A:78:SER:OG	2.00	0.45
1:A:2:LYS:HB2	1:A:2:LYS:HZ3	1.82	0.44
1:A:59:ARG:NH1	1:A:88:TYR:OH	2.51	0.43
1:A:2:LYS:CG	1:A:3:GLY:N	2.80	0.43
1:A:16:THR:HG23	1:A:79:HIS:CE1	2.54	0.43
1:A:330:ARG:NH1	1:A:332:THR:O	2.52	0.43
1:A:313:HIS:CD2	1:A:403:ILE:HG21	2.54	0.43
1:A:113:MET:HB3	1:A:114:PRO:CD	2.49	0.42
1:A:230:THR:OG1	1:A:295:ARG:HD3	2.19	0.42
1:A:12:VAL:HG21	1:A:75:ARG:CZ	2.50	0.42
1:A:354:ARG:NH1	1:A:354:ARG:CG	2.82	0.42
1:A:252:GLU:HA	1:A:265:THR:O	2.20	0.42
1:A:125:GLN:O	1:A:385:ARG:HB3	2.20	0.41
1:A:393:ARG:HH11	1:A:393:ARG:HG2	1.84	0.41
1:A:39:ASN:HA	1:A:40:PRO:HD2	1.78	0.41
1:A:67:HIS:HE1	1:A:98:GLN:OE1	2.04	0.41
1:A:125:GLN:CD	1:A:397:ALA:HB1	2.41	0.41

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	403/405 (100%)	380 (94%)	20 (5%)	3 (1%)	25 43

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	260	PRO
1	A	2	LYS
1	A	143	ASP

### 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	339/339 (100%)	297 (88%)	42 (12%)	5 10

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	8	THR
1	A	43	GLU
1	A	74	LYS
1	A	75	ARG
1	A	94	THR
1	A	107	SER
1	A	126	VAL
1	A	142	ASP
1	A	145	GLU
1	A	180	GLU
1	A	187	LYS
1	A	190	ARG
1	A	215	ARG
1	A	216	ASP
1	A	230	THR
1	A	232	THR
1	A	241	ARG
1	A	244	ARG
1	A	247	VAL
1	A	268	THR
1	A	281	ILE

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Mol	Chain	Res	Type
1	A	285	ASN
1	A	288	LEU
1	A	289	LEU
1	A	291	ARG
1	A	293	VAL
1	A	298	VAL
1	A	300	ARG
1	A	311	THR
1	A	314	THR
1	A	326	GLU
1	A	330	ARG
1	A	336	SER
1	A	347	THR
1	A	349	VAL
1	A	354	ARG
1	A	363	MET
1	A	385	ARG
1	A	393	ARG
1	A	399	VAL
1	A	401	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS
1	A	85	HIS
1	A	98	GLN
1	A	185	ASN
1	A	273	HIS
1	A	331	HIS
1	A	367	ASN

### 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 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	GNP	A	406	2	27,34,34	2.98	10 (37%)	26,54,54	1.76	6 (23%)

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	GNP	A	406	2	-	0/16/38/38	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	406	GNP	C4-N9	-10.44	1.33	1.47
3	A	406	GNP	C5-C6	-6.49	1.41	1.53
3	A	406	GNP	C8-N9	-3.76	1.35	1.46
3	A	406	GNP	C2'-C1'	-3.75	1.41	1.53
3	A	406	GNP	PB-N3B	-2.77	1.56	1.63
3	A	406	GNP	O4'-C4'	-2.23	1.40	1.45
3	A	406	GNP	PG-N3B	-2.22	1.57	1.63
3	A	406	GNP	PB-O3A	-2.15	1.56	1.59
3	A	406	GNP	PG-O1G	2.27	1.48	1.46
3	A	406	GNP	C6-N1	3.70	1.39	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	406	GNP	O6-C6-N1	-2.39	119.52	122.70
3	A	406	GNP	C2'-C3'-C4'	2.28	107.06	102.62
3	A	406	GNP	O2G-PG-O1G	2.36	119.39	113.41
3	A	406	GNP	O3A-PB-N3B	3.33	115.83	106.59
3	A	406	GNP	O6-C6-C5	3.87	127.10	119.69
3	A	406	GNP	O1B-PB-N3B	4.00	117.77	111.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	406	GNP	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, 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	405/405 (100%)	-0.73	3 (0%) 87 88	8, 22, 36, 43	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	LYS	4.2
1	A	3	GLY	3.0
1	A	1	ALA	2.9

### 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. 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
3	GNP	A	406	32/32	0.98	0.07	-1.07	2,14,24,26	0
2	MG	A	407	1/1	0.95	0.08	-2.39	2,2,2,2	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.