



# wwPDB X-ray Structure Validation Summary Report

Feb 26, 2014 – 10:00 PM GMT

PDB ID : 1FEU  
Title : CRYSTAL STRUCTURE OF RIBOSOMAL PROTEIN TL5, ONE OF THE CTC FAMILY PROTEINS, COMPLEXED WITH A FRAGMENT OF 5S RRNA.  
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Deposited on : 2000-07-23  
Resolution : 2.30 Å(reported)

This is a wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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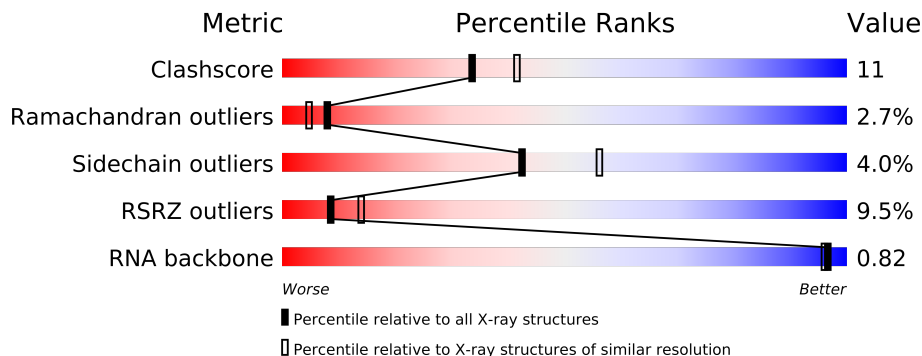
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.30 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	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)
RNA backbone	1838	1081 (3.00-1.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	B	19	
1	E	19	
2	C	21	
2	F	21	
3	A	206	
3	D	206	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	MG	C	309	-	X
5	MG	D	310	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
5	MG	E	315	-	X
5	MG	F	316	-	X

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5023 atoms, of which 0 are hydrogen and 0 are deuterium.

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 19 NT FRAGMENT OF 5S RRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	19	Total	C	N	O	P	0	0	0
			411	183	76	134	18			
1	E	19	Total	C	N	O	P	0	0	0
			411	183	76	134	18			

- Molecule 2 is a RNA chain called 21 NT FRAGMENT OF 5S RRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	21	Total	C	N	O	P	0	0	0
			450	202	87	141	20			
2	F	21	Total	C	N	O	P	0	0	0
			450	202	87	141	20			

- Molecule 3 is a protein called 50S RIBOSOMAL PROTEIN L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	185	Total	C	N	O	S	0	0	0
			1476	942	262	269	3			
3	D	189	Total	C	N	O	S	0	0	0
			1500	956	266	275	3			

- Molecule 4 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	Cd	0	0
			6	6		
4	D	7	Total	Cd	0	0
			7	7		

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total 2	Mg 2	0	0
5	D	4	Total 4	Mg 4	0	0
5	C	4	Total 4	Mg 4	0	0
5	F	5	Total 5	Mg 5	0	0
5	E	2	Total 2	Mg 2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	40	Total 40	O 40	0	0
6	B	26	Total 26	O 26	0	0
6	C	37	Total 37	O 37	0	0
6	D	121	Total 121	O 121	0	0
6	E	31	Total 31	O 31	0	0
6	F	40	Total 40	O 40	0	0

### 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 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: 19 NT FRAGMENT OF 5S RRNA

Chain B: 



- Molecule 1: 19 NT FRAGMENT OF 5S RRNA

Chain E: 



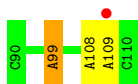
- Molecule 2: 21 NT FRAGMENT OF 5S RRNA

Chain C: 



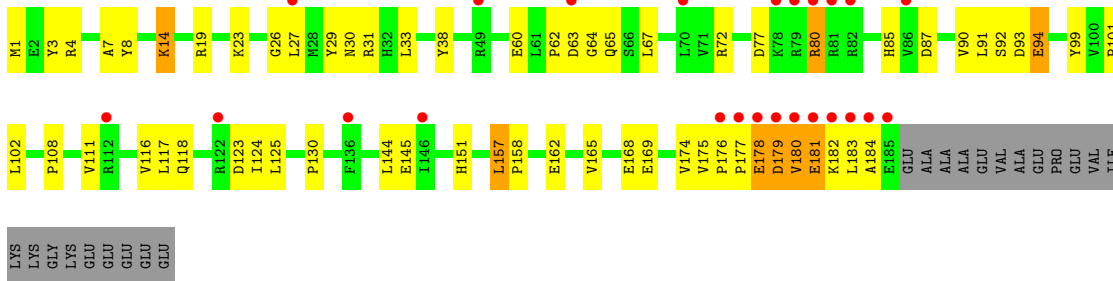
- Molecule 2: 21 NT FRAGMENT OF 5S RRNA

Chain F: 



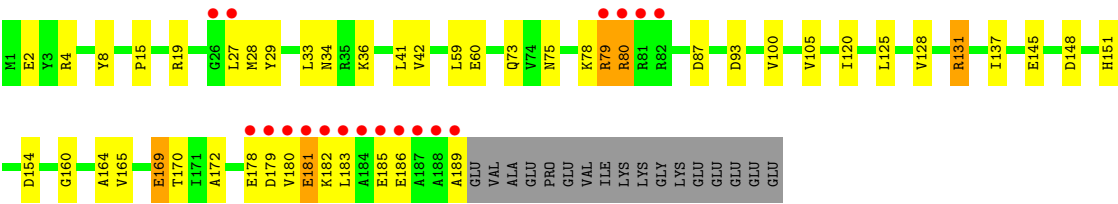
- Molecule 3: 50S RIBOSOMAL PROTEIN L25

Chain A: 



- Molecule 3: 50S RIBOSOMAL PROTEIN L25

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	109.90Å 109.90Å 137.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.30 19.88 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.4 (8.00-2.30) 95.3 (19.88-2.30)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.87 (at 2.30Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.208 , 0.245 0.221 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	31.5	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 36.4	EDS
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 40400 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5023	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% of the height of the origin peak. No significant pseudotranslation is detected.*

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



## 5 Model quality

### 5.1 Standard geometry

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

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	B	0.24	0/460	0.79	2/718 (0.3%)
1	E	0.25	0/460	0.66	0/718
2	C	0.23	0/504	0.81	5/785 (0.6%)
2	F	0.23	0/504	0.70	0/785
3	A	0.34	0/1508	0.60	0/2046
3	D	0.37	0/1532	0.65	0/2079
All	All	0.31	0/4968	0.68	7/7131 (0.1%)

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	81	G	P-O3'-C3'	6.88	127.96	119.70
1	B	81	G	OP1-P-O3'	6.17	118.77	105.20
2	C	103	U	P-O3'-C3'	-5.61	112.97	119.70
2	C	100	G	OP1-P-O3'	5.55	117.40	105.20
2	C	103	U	OP2-P-O3'	5.43	117.15	105.20

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	411	0	206	2	0
1	E	411	0	206	8	0
2	C	450	0	230	1	0
2	F	450	0	230	2	0
3	A	1476	0	1506	46	1
3	D	1500	0	1527	41	0
4	A	6	0	0	0	0
4	D	7	0	0	0	0
5	B	2	0	0	0	0
5	C	4	0	0	0	0
5	D	4	0	0	0	0
5	E	2	0	0	0	0
5	F	5	0	0	0	0
6	A	40	0	0	1	0
6	B	26	0	0	0	0
6	C	37	0	0	0	0
6	D	121	0	0	5	0
6	E	31	0	0	0	0
6	F	40	0	0	1	0
All	All	5023	0	3905	95	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

The worst 5 of 95 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:D:93:ASP:HB2	6:D:425:HOH:O	1.73	0.89
1:E:69:G:H3'	1:E:69:G:N3	1.91	0.85
3:D:4:ARG:HD3	3:D:60:GLU:OE1	1.78	0.83
1:E:75:G:H1'	3:D:27:LEU:HD22	1.59	0.83
3:A:30:ASN:HD22	3:A:90:VAL:HB	1.50	0.76

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	Distance(Å)	Clash(Å)
3:A:80:ARG:NH2	3:A:80:ARG:NH2[6_555]	2.14	0.06

## 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	
3	A	183/206 (89%)	162 (88%)	14 (8%)	7 (4%)	5	2
3	D	187/206 (91%)	170 (91%)	14 (8%)	3 (2%)	14	12
All	All	370/412 (90%)	332 (90%)	28 (8%)	10 (3%)	8	5

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	177	PRO
3	A	181	GLU
3	D	80	ARG
3	A	80	ARG
3	D	79	ARG

### 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	
3	A	163/179 (91%)	155 (95%)	8 (5%)	35	45
3	D	164/179 (92%)	159 (97%)	5 (3%)	53	70
All	All	327/358 (91%)	314 (96%)	13 (4%)	42	56

5 of 13 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	157	LEU
3	A	169	GLU
3	D	131	ARG
3	A	94	GLU

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Mol	Chain	Res	Type
3	D	105	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	75	ASN
3	A	151	HIS
3	D	75	ASN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	B	19/19 (100%)	1 (5%)	1 (5%)
1	E	18/19 (94%)	0	0
2	C	20/21 (95%)	0	0
2	F	20/21 (95%)	1 (5%)	0
All	All	77/80 (96%)	2 (2%)	1 (1%)

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	B	70	C
2	F	99	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	B	69	G

## 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 30 ligands modelled in this entry, 30 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

### 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	B	19/19 (100%)	0.08	0 <span>100</span> <span>100</span>	21, 30, 46, 56	0
1	E	19/19 (100%)	0.24	1 (5%) <span>25</span> <span>35</span>	19, 23, 55, 69	0
2	C	21/21 (100%)	-0.15	0 <span>100</span> <span>100</span>	19, 32, 46, 48	0
2	F	21/21 (100%)	0.10	1 (4%) <span>29</span> <span>39</span>	18, 26, 44, 47	0
3	A	185/206 (89%)	0.76	24 (12%) <span>4</span> <span>7</span>	25, 41, 68, 70	0
3	D	189/206 (91%)	0.55	18 (9%) <span>8</span> <span>14</span>	16, 28, 69, 70	0
All	All	454/492 (92%)	0.55	44 (9%) <span>8</span> <span>13</span>	16, 34, 69, 70	0

The worst 5 of 44 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	183	LEU	13.2
3	D	189	ALA	11.4
3	D	188	ALA	10.7
3	D	187	ALA	10.4
3	A	184	ALA	10.1

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

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	MG	D	310	1/1	0.70	21.95	47,47,47,47	1
5	MG	C	309	1/1	0.68	15.76	40,40,40,40	1
5	MG	E	315	1/1	0.19	3.30	43,43,43,43	0
5	MG	F	316	1/1	0.15	2.47	31,31,31,31	0
5	MG	D	319	1/1	0.19	1.82	48,48,48,48	0
5	MG	C	312	1/1	0.13	1.12	33,33,33,33	0
5	MG	F	325	1/1	0.15	-0.22	46,46,46,46	0
4	CD	D	306	1/1	0.10	-0.39	30,30,30,30	1
5	MG	F	317	1/1	0.13	-0.50	52,52,52,52	0
4	CD	A	296	1/1	0.09	-0.64	47,47,47,47	0
5	MG	B	313	1/1	0.09	-0.77	26,26,26,26	0
4	CD	D	304	1/1	0.11	-0.84	28,28,28,28	0
4	CD	D	308	1/1	0.12	-0.87	35,35,35,35	1
4	CD	A	302	1/1	0.08	-0.96	41,41,41,41	1
4	CD	A	301	1/1	0.09	-1.14	43,43,43,43	1
5	MG	B	311	1/1	0.11	-1.24	42,42,42,42	0
4	CD	D	303	1/1	0.09	-1.44	33,33,33,33	0
5	MG	D	322	1/1	0.13	-1.44	43,43,43,43	0
5	MG	F	324	1/1	0.07	-1.88	32,32,32,32	0
4	CD	A	297	1/1	0.04	-2.41	40,40,40,40	0
4	CD	A	298	1/1	0.03	-2.72	53,53,53,53	0
4	CD	D	305	1/1	0.06	-3.11	40,40,40,40	0
5	MG	C	318	1/1	0.13	-3.13	43,43,43,43	0
5	MG	E	321	1/1	0.07	-3.29	38,38,38,38	0
5	MG	D	320	1/1	0.05	-4.66	28,28,28,28	0
4	CD	A	299	1/1	0.04	-5.12	47,47,47,47	1
5	MG	F	323	1/1	0.07	-5.25	38,38,38,38	0
4	CD	D	307	1/1	0.05	-5.79	46,46,46,46	1
4	CD	D	300	1/1	0.04	-6.48	42,42,42,42	1
5	MG	C	314	1/1	0.07	-8.80	32,32,32,32	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.