



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 25, 2020 – 06:33 am BST

PDB ID : 1G1X  
Title : STRUCTURE OF RIBOSOMAL PROTEINS S15, S6, S18, AND 16S RIBOSOMAL RNA  
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Deposited on : 2000-10-13  
Resolution : 2.60 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
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

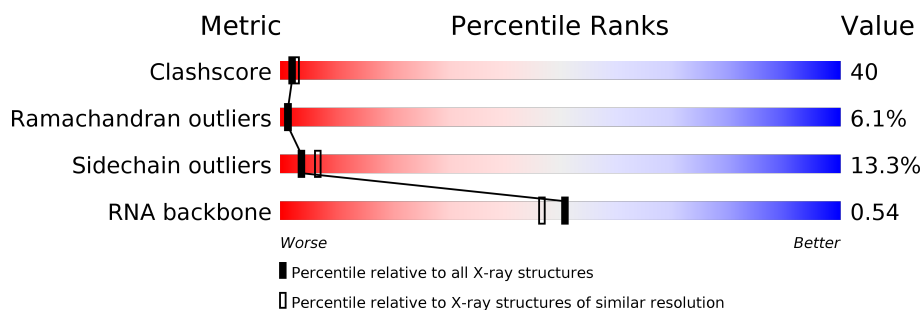
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RNA backbone	3102	1040 (2.90-2.30)


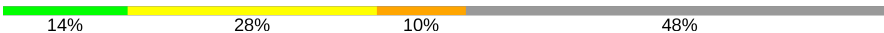

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria 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  $\leq 5\%$ .

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	D	41	41% 44% 10% ..
1	I	41	29% 41% 29%
2	E	44	27% 43% 25% 5%
2	J	44	23% 55% 16% 7%
3	A	98	32% 57% 10% .
3	F	98	16% 64% 18% .
4	B	88	44% 41% 15%

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Mol	Chain	Length	Quality of chain
4	G	88	 35% 51% 11% .
5	C	88	 14% 28% 10% 48%
5	H	88	 13% 36% 6% 45%

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7543 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 16S RIBOSOMAL RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	40	Total	C	N	O	P	0	0	0
			880	391	178	272	39			
1	I	41	Total	C	N	O	P	0	0	0
			902	401	183	278	40			

- Molecule 2 is a RNA chain called 16S RIBOSOMAL RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	44	Total	C	N	O	P	0	0	0
			932	417	167	305	43			
2	J	44	Total	C	N	O	P	0	0	0
			932	417	167	305	43			

- Molecule 3 is a protein called 30S RIBOSOMAL PROTEIN S6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	98	Total	C	N	O	S	0	0	0
			825	521	151	150	3			
3	F	98	Total	C	N	O	S	0	0	0
			825	521	151	150	3			

- Molecule 4 is a protein called 30S RIBOSOMAL PROTEIN S15.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	B	88	Total	C	N	O	S	0	0	0
			737	460	147	128	2			
4	G	88	Total	C	N	O	S	0	0	0
			737	460	147	128	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	79	GLU	ALA	CONFLICT	UNP Q5SJ76
B	80	ILE	LEU	CONFLICT	UNP Q5SJ76
B	81	VAL	ILE	CONFLICT	UNP Q5SJ76
B	86	LEU	ILE	CONFLICT	UNP Q5SJ76
G	79	GLU	ALA	CONFLICT	UNP Q5SJ76
G	80	ILE	LEU	CONFLICT	UNP Q5SJ76
G	81	VAL	ILE	CONFLICT	UNP Q5SJ76
G	86	LEU	ILE	CONFLICT	UNP Q5SJ76

- Molecule 5 is a protein called 30S RIBOSOMAL PROTEIN S18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	C	46	Total	C	N	O	0	0	0
			379	245	73	61			
5	H	48	Total	C	N	O	0	0	0
			394	254	76	64			

### 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. 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. 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.

Note EDS was not executed.

#### • Molecule 1: 16S RIBOSOMAL RNA

Chain D: 



#### • Molecule 1: 16S RIBOSOMAL RNA

Chain I: 



#### • Molecule 2: 16S RIBOSOMAL RNA

Chain E: 



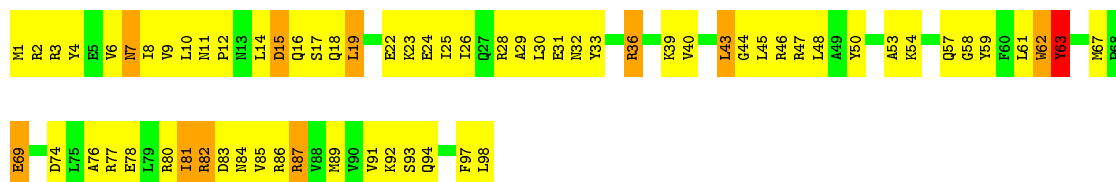
#### • Molecule 2: 16S RIBOSOMAL RNA

Chain J: 

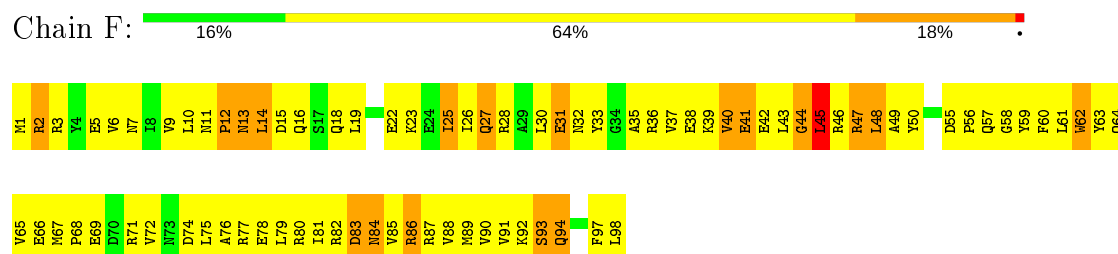


#### • Molecule 3: 30S RIBOSOMAL PROTEIN S6

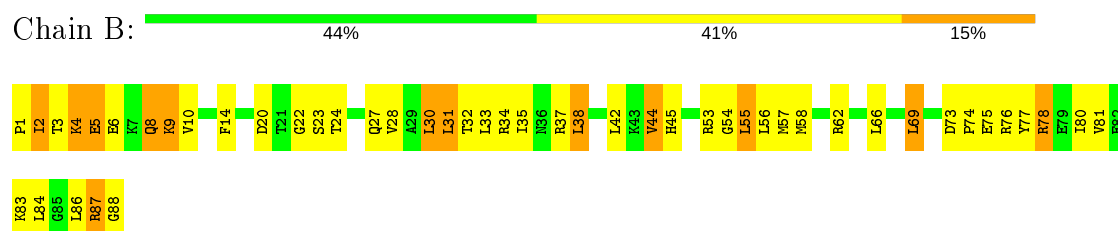
Chain A: 



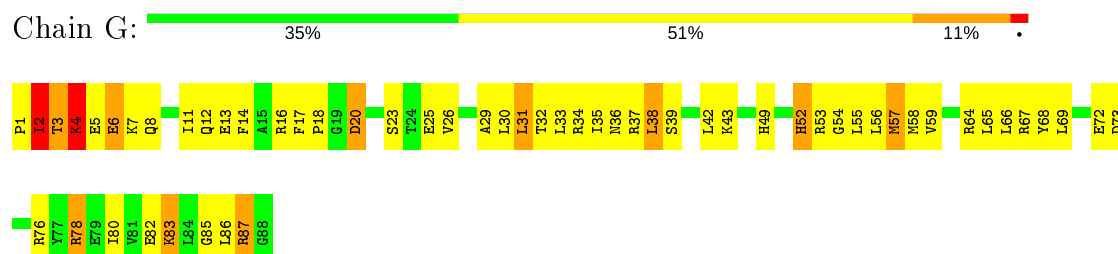
- Molecule 3: 30S RIBOSOMAL PROTEIN S6



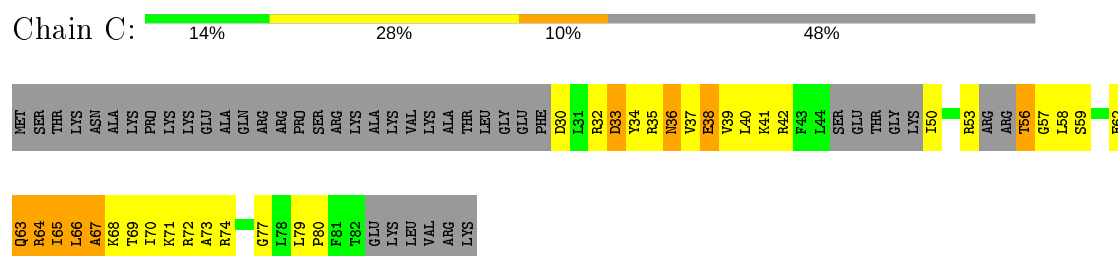
- Molecule 4: 30S RIBOSOMAL PROTEIN S15



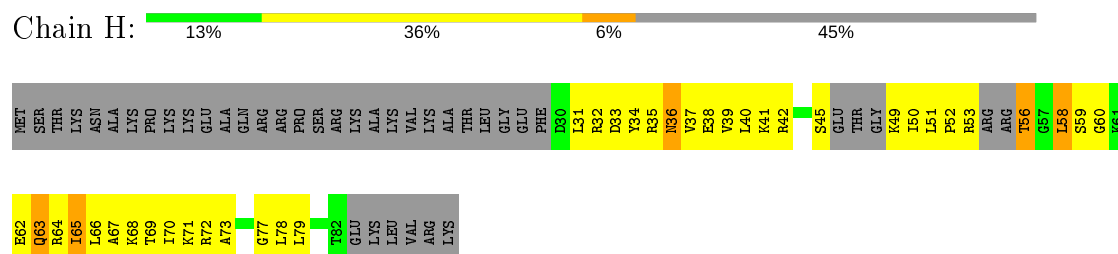
- Molecule 4: 30S RIBOSOMAL PROTEIN S15



- Molecule 5: 30S RIBOSOMAL PROTEIN S18



- Molecule 5: 30S RIBOSOMAL PROTEIN S18



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	169.50Å 169.50Å 113.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.60	Depositor
% Data completeness (in resolution range)	99.9 (50.00-2.60)	Depositor
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.28	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.266 , 0.298	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7543	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP



## 5 Model quality [i](#)

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

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	D	0.71	0/990	0.80	0/1549
1	I	0.75	0/1015	0.86	1/1588 (0.1%)
2	E	0.73	0/1040	0.87	3/1619 (0.2%)
2	J	0.74	0/1040	0.86	3/1619 (0.2%)
3	A	0.57	0/838	0.78	0/1129
3	F	0.67	1/838 (0.1%)	0.85	0/1129
4	B	0.56	0/748	0.75	1/996 (0.1%)
4	G	0.57	0/748	0.74	0/996
5	C	0.54	0/381	0.80	0/506
5	H	0.66	0/396	0.95	0/525
All	All	0.67	1/8034 (0.0%)	0.83	8/11656 (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	D	0	2
1	I	0	3
2	J	0	2
All	All	0	7

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	62	TRP	CE3-CZ3	8.24	1.52	1.38

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	717	C	N1-C1'-C2'	6.55	122.51	114.00
4	B	44	VAL	N-CA-C	-6.28	94.03	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	753	A	N9-C1'-C2'	6.08	121.91	114.00
1	I	593	A	N9-C1'-C2'	-6.06	105.33	112.00
2	J	748	C	N1-C1'-C2'	6.01	121.81	114.00

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	652	U	Sidechain
1	D	674	G	Sidechain
1	I	659	U	Sidechain
1	I	672	U	Sidechain
1	I	674	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	D	880	0	438	30	0
1	I	902	0	449	41	0
2	E	932	0	479	36	0
2	J	932	0	479	40	0
3	A	825	0	841	96	0
3	F	825	0	841	112	0
4	B	737	0	772	58	0
4	G	737	0	772	64	0
5	C	379	0	421	50	0
5	H	394	0	437	59	0
All	All	7543	0	5929	531	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 40.

The worst 5 of 531 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:I:652:U:O2'	1:I:653:A:H5''	1.37	1.21
1:D:659:U:H5''	3:F:36:ARG:HH22	1.11	1.10
4:G:4:LYS:HD3	4:G:5:GLU:H	1.31	0.96
5:H:53:ARG:HA	5:H:56:THR:HG23	1.47	0.95
1:D:659:U:H5''	3:F:36:ARG:NH2	1.82	0.94

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	
3	A	96/98 (98%)	74 (77%)	18 (19%)	4 (4%)	3	3
3	F	96/98 (98%)	68 (71%)	17 (18%)	11 (12%)	0	0
4	B	86/88 (98%)	75 (87%)	9 (10%)	2 (2%)	6	11
4	G	86/88 (98%)	71 (83%)	11 (13%)	4 (5%)	2	2
5	C	40/88 (46%)	26 (65%)	9 (22%)	5 (12%)	0	0
5	H	42/88 (48%)	31 (74%)	10 (24%)	1 (2%)	6	10
All	All	446/548 (81%)	345 (77%)	74 (17%)	27 (6%)	1	1

5 of 27 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	C	63	GLN
3	F	12	PRO
3	F	43	LEU
3	F	47	ARG
3	F	93	SER

### 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	89/89 (100%)	80 (90%)	9 (10%)	7	14
3	F	89/89 (100%)	78 (88%)	11 (12%)	4	8
4	B	80/80 (100%)	69 (86%)	11 (14%)	3	6
4	G	80/80 (100%)	66 (82%)	14 (18%)	2	3
5	C	41/77 (53%)	35 (85%)	6 (15%)	3	5
5	H	43/77 (56%)	38 (88%)	5 (12%)	5	10
All	All	422/492 (86%)	366 (87%)	56 (13%)	4	7

5 of 56 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	C	80	PRO
3	F	31	GLU
5	H	35	ARG
3	F	2	ARG
3	F	25	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 23 such sidechains are listed below:

Mol	Chain	Res	Type
4	B	70	GLN
3	F	16	GLN
4	G	70	GLN
5	C	36	ASN
3	F	27	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	D	39/41 (95%)	5 (12%)	2 (5%)
1	I	40/41 (97%)	9 (22%)	2 (5%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	E	43/44 (97%)	13 (30%)	6 (13%)
2	J	43/44 (97%)	12 (27%)	4 (9%)
All	All	165/170 (97%)	39 (23%)	14 (8%)

5 of 39 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	D	583	A
1	D	652	U
1	D	653	A
1	D	654	G
1	D	665	A

5 of 14 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	E	752	G
2	E	753	A
2	J	733	A
2	E	748	C
2	J	721	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](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

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 ⓘ

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.