



## wwPDB/EMDataBank EM Map/Model Validation Summary Report ⓘ

Jul 24, 2017 – 07:28 AM EDT

PDB ID : 5LMR  
EMDB ID: : EMD-4077  
Title : Structure of bacterial 30S-IF1-IF3-mRNA-tRNA translation pre-initiation complex(state-2B)  
Authors : Hussain, T.; Llacer, J.L.; Wimberly, B.T.; Ramakrishnan, V.  
Deposited on : unknown  
Resolution : 4.45 Å(reported)

This is a wwPDB/EMDataBank EM Map/Model Validation Summary Report  
for a publicly released PDB/EMDB 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/EMValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20029824

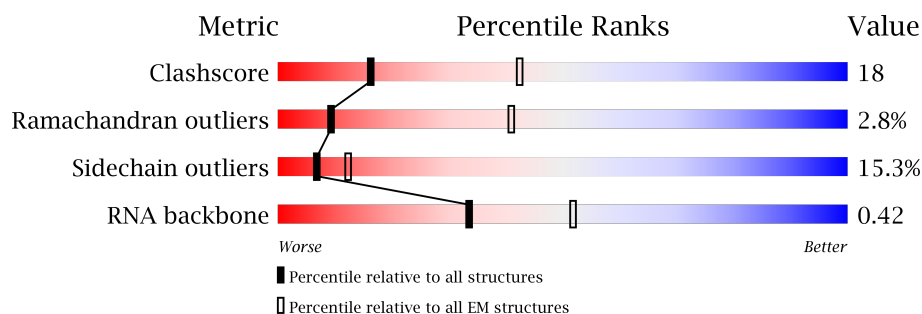
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 4.45 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.














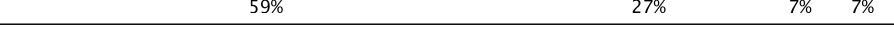





Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	125131	1336
Ramachandran outliers	121729	1120
Sidechain outliers	121581	1026
RNA backbone	3398	335

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	1522	27% 55% 17% .
2	B	256	56% 30% 5% . 9%
3	C	239	46% 33% 7% 14%
4	D	209	57% 38% .
5	E	162	59% 27% 6% . 7%
6	F	101	66% 31% ..
7	G	156	73% 26% .
8	H	138	63% 31% 6%

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Mol	Chain	Length	Quality of chain
9	I	128	
10	J	105	
11	K	129	
12	L	132	
13	M	126	
14	N	61	
15	O	89	
16	P	88	
17	Q	105	
18	R	88	
19	S	93	
20	T	106	
21	V	27	
22	W	72	
23	X	171	
24	Y	42	
25	Z	77	

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
26	ZN	D	300	-	-	X	-

## 2 Entry composition

There are 27 unique types of molecules in this entry. The entry contains 55662 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1515	Total	C	N	O	P	0	0
			32542	14490	6022	10519	1511		

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	234	Total	C	N	O	S	0	0
			1900	1213	341	341	5		

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	206	Total	C	N	O	S	0	0
			1612	1016	314	281	1		

- Molecule 4 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	208	Total	C	N	O	S	0	0
			1703	1066	339	291	7		

- Molecule 5 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	150	Total	C	N	O	S	0	0
			1146	724	217	201	4		

- Molecule 6 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	101	Total	C	N	O	S	0	0
			843	531	155	154	3		

- Molecule 7 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	155	Total	C	N	O	S	0	0
			1257	781	252	218	6		

- Molecule 8 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	138	Total	C	N	O	S	0	0
			1116	705	215	193	3		

- Molecule 9 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms				AltConf	Trace
9	I	127	Total	C	N	O	0	0
			1010	639	197	174		

- Molecule 10 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	98	Total	C	N	O	S	0	0
			792	498	156	137	1		

- Molecule 11 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	119	Total	C	N	O	S	0	0
			885	549	168	165	3		

- Molecule 12 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	124	Total	C	N	O	S	0	0
			970	611	195	163	1		

- Molecule 13 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	121	Total	C	N	O	S	0	0
			964	597	199	166	2		

- Molecule 14 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	60	Total	C	N	O	S	0	0
			492	312	104	72	4		

- Molecule 15 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	88	Total	C	N	O	S	0	0
			734	459	147	126	2		

- Molecule 16 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	83	Total	C	N	O	S	0	0
			700	443	139	117	1		

- Molecule 17 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Q	99	Total	C	N	O	S	0	0
			823	528	151	142	2		

- Molecule 18 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				AltConf	Trace
18	R	73	Total	C	N	O	0	0
			598	381	118	99		

- Molecule 19 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	S	82	Total	C	N	O	S	0	0
			655	419	120	114	2		

- Molecule 20 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	T	99	Total	C	N	O	S	0	0
			763	470	162	129	2		

- Molecule 21 is a protein called 30S ribosomal protein Thx.

Mol	Chain	Residues	Atoms				AltConf	Trace
21	V	24	Total	C	N	O	0	0
			208	128	50	30		

- Molecule 22 is a protein called Translation initiation factor IF-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	W	71	Total	C	N	O	S	0	0
			570	362	103	103	2		

- Molecule 23 is a protein called Translation initiation factor IF-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	X	168	Total	C	N	O	S	0	0
			1356	853	249	245	9		

- Molecule 24 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	Y	17	Total	C	N	O	P	0	0
			373	166	74	116	17		

- Molecule 25 is a RNA chain called tRNAi.

Mol	Chain	Residues	Atoms						AltConf	Trace
25	Z	77	Total	C	N	O	P	S	0	0
			1646	735	297	536	77	1		

- Molecule 26 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
26	D	1	Total	Zn	0
			1	1	
26	N	1	Total	Zn	0
			1	1	

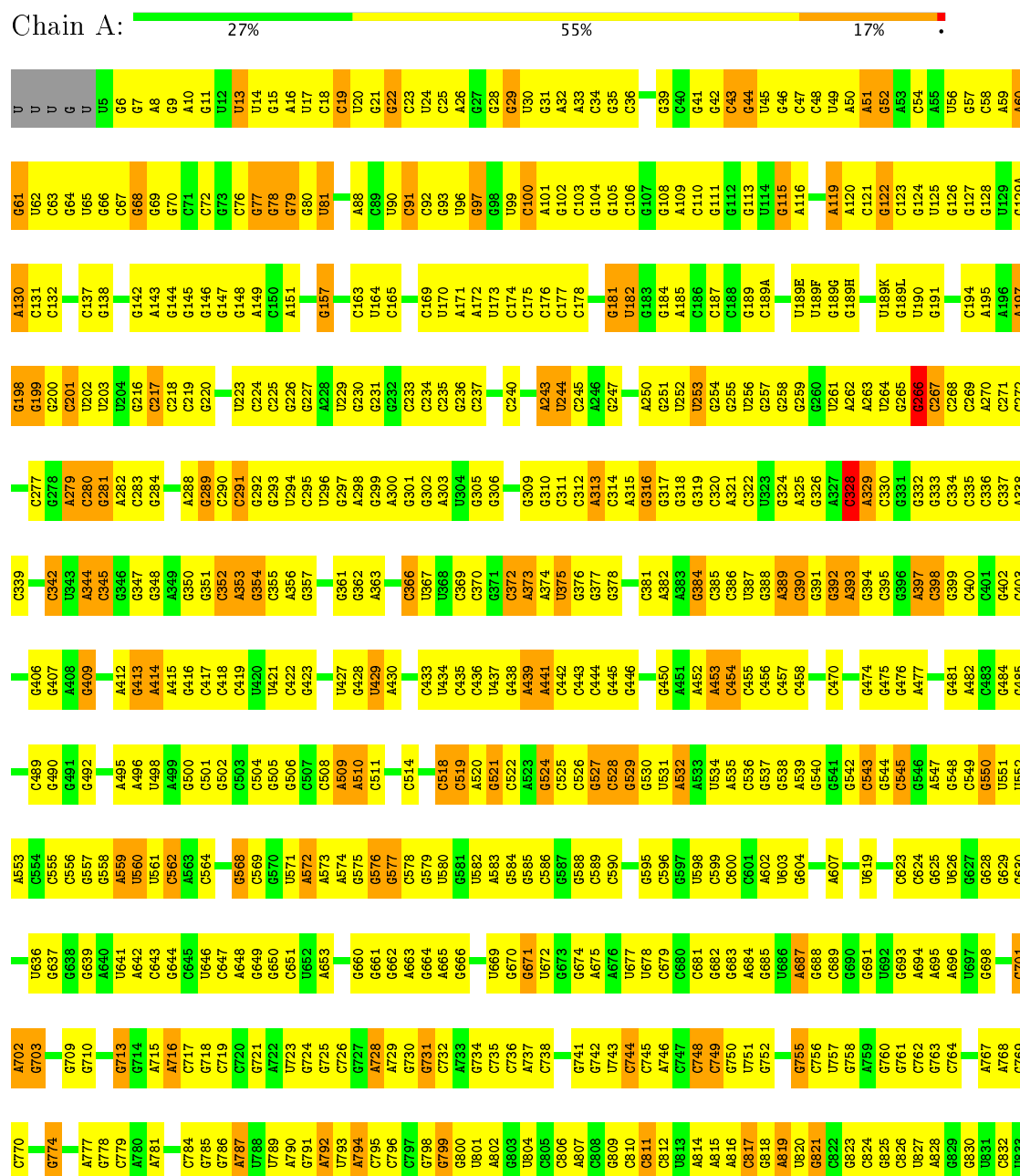
- Molecule 27 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
27	W	1	Total	Mg	0
			1	1	
27	Z	1	Total	Mg	0
			1	1	

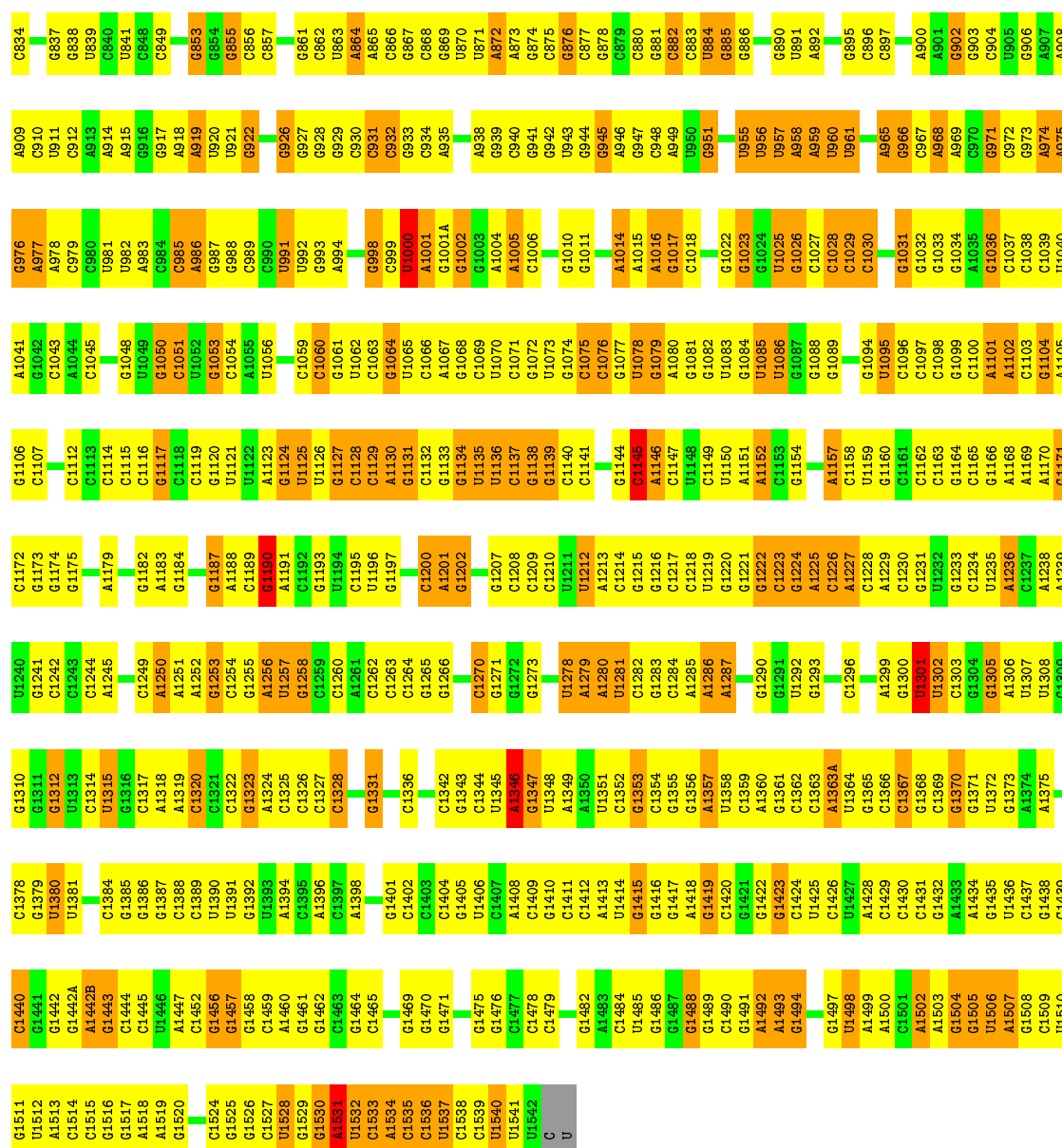
### 3 Residue-property plots

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. 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.

#### • Molecule 1: 16S rRNA

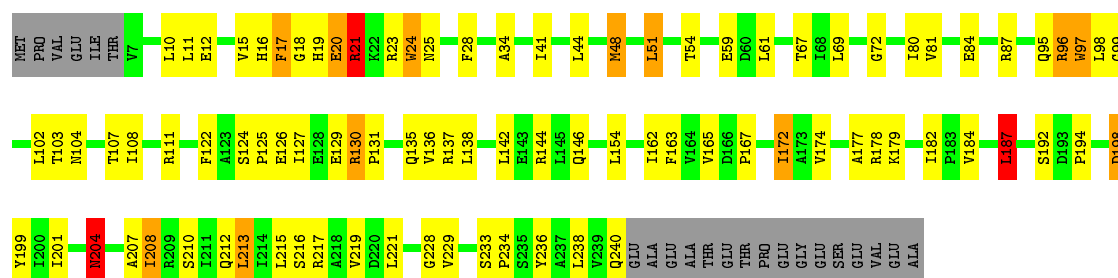






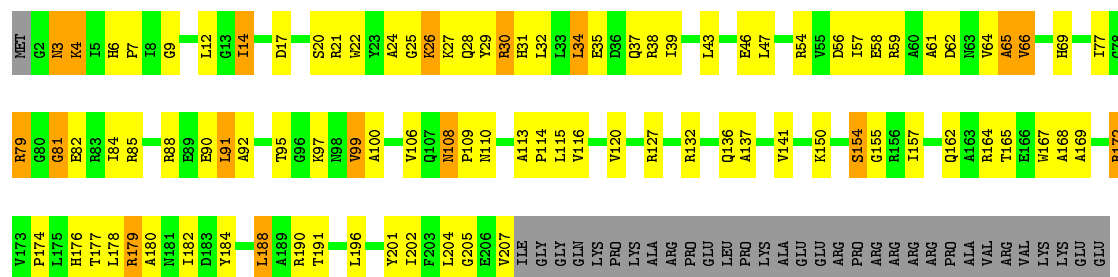
- Molecule 2: 30S ribosomal protein S2

Chain B:  56% 30% 5% 9%



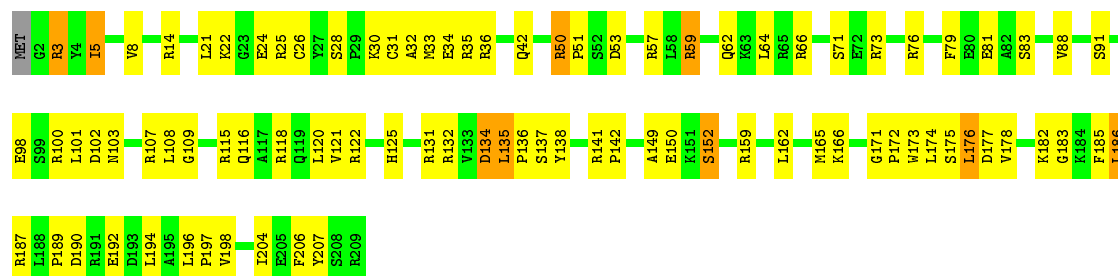
- Molecule 3: 30S ribosomal protein S3

Chain C: 



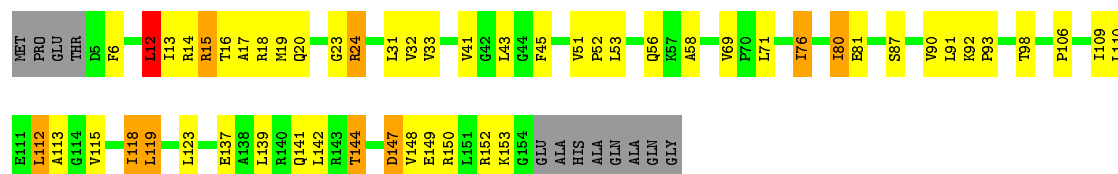
- Molecule 4: 30S ribosomal protein S4

Chain D: 



- Molecule 5: 30S ribosomal protein S5

Chain E: 



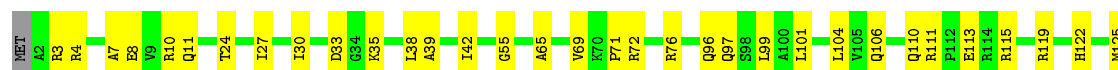
- Molecule 6: 30S ribosomal protein S6

Chain F: 



- Molecule 7: 30S ribosomal protein S7

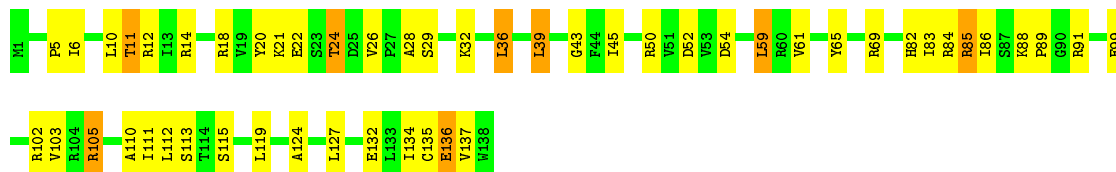
Chain G: 





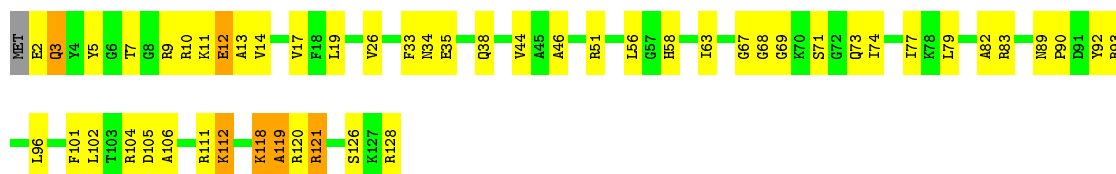
• Molecule 8: 30S ribosomal protein S8

Chain H: 63% 31% 6%



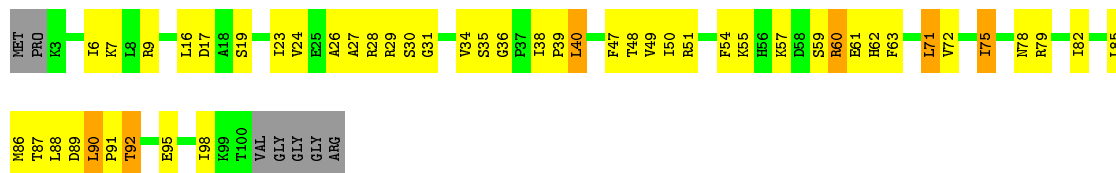
• Molecule 9: 30S ribosomal protein S9

Chain I: 59% 35% 5%



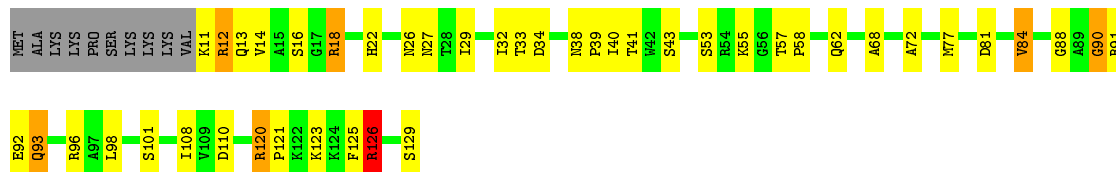
• Molecule 10: 30S ribosomal protein S10

Chain J: 47% 41% 6% 7%



• Molecule 11: 30S ribosomal protein S11

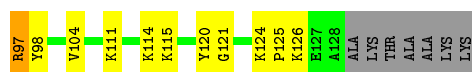
Chain K: 58% 29% 5% 8%



• Molecule 12: 30S ribosomal protein S12

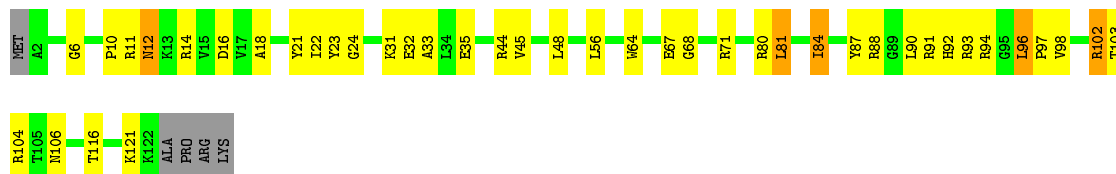
Chain L: 56% 33% 5% 6%





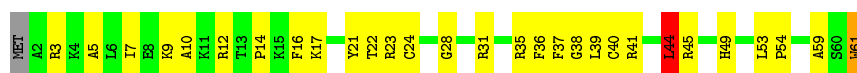
- Molecule 13: 30S ribosomal protein S13

Chain M: 63% 29%



- Molecule 14: 30S ribosomal protein S14 type Z

Chain N: 51% 44%



- Molecule 15: 30S ribosomal protein S15

Chain O: 63% 31%



- Molecule 16: 30S ribosomal protein S16

Chain P: 59% 34% 6%



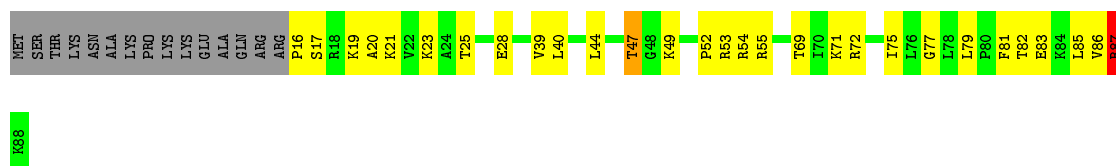
- Molecule 17: 30S ribosomal protein S17

Chain Q: 63% 25% 6% 6%



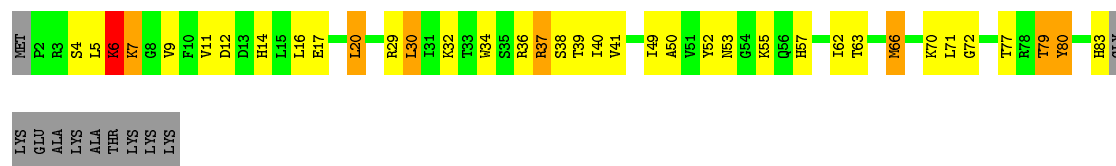
- Molecule 18: 30S ribosomal protein S18

Chain R: 50% 31% 17%



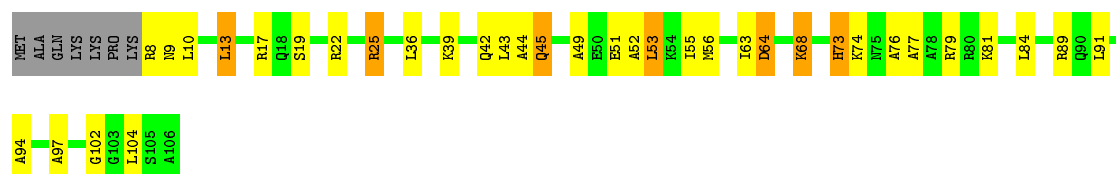
- Molecule 19: 30S ribosomal protein S19

Chain S: 



- Molecule 20: 30S ribosomal protein S20

Chain T: 



- Molecule 21: 30S ribosomal protein Thx

Chain V: 



- Molecule 22: Translation initiation factor IF-1

Chain W: 



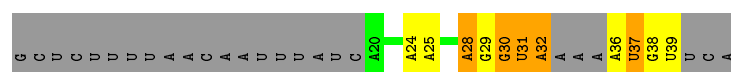
- Molecule 23: Translation initiation factor IF-3

Chain X: 




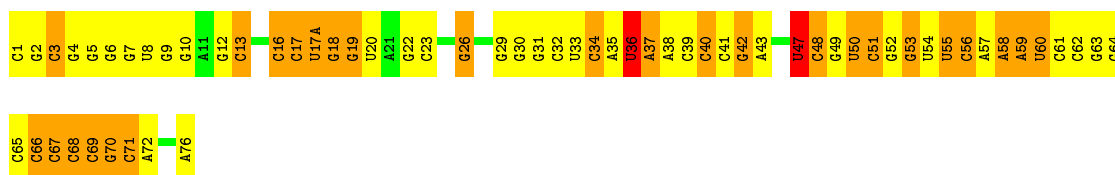
- Molecule 24: mRNA

Chain Y: 



- Molecule 25: tRNAi

Chain Z: 



## 4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of particles used	17176	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	30	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	104478	Depositor
Image detector	OTHER	Depositor

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 5MU, OMC, ZN, G7M, MG, 4SU, PSU

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	$\# Z  > 2$	RMSZ	$\# Z  > 2$
1	A	0.33	0/36416	0.73	13/56813 (0.0%)
10	J	0.48	0/805	0.80	2/1082 (0.2%)
11	K	0.47	0/900	0.82	0/1213
12	L	0.38	0/986	0.75	0/1320
13	M	0.45	0/974	0.78	1/1303 (0.1%)
14	N	0.41	0/501	0.82	1/664 (0.2%)
15	O	0.44	0/745	0.82	0/992
16	P	0.42	0/716	0.77	0/963
17	Q	0.40	0/836	0.75	0/1117
18	R	0.45	0/604	0.81	0/801
19	S	0.48	0/670	0.80	1/903 (0.1%)
2	B	0.51	0/1935	0.82	2/2609 (0.1%)
20	T	0.45	0/765	0.88	0/1007
21	V	0.48	0/212	0.75	0/277
22	W	0.53	0/580	0.90	1/782 (0.1%)
23	X	0.48	0/1375	0.78	0/1844
24	Y	0.43	0/418	0.69	0/649
25	Z	0.43	0/1719	0.86	2/2674 (0.1%)
3	C	0.44	0/1636	0.88	6/2205 (0.3%)
4	D	0.42	0/1733	0.78	0/2318
5	E	0.43	0/1162	0.89	3/1564 (0.2%)
6	F	0.43	0/856	0.78	1/1154 (0.1%)
7	G	0.44	0/1276	0.77	0/1709
8	H	0.40	0/1136	0.78	0/1527
9	I	0.42	0/1029	0.72	0/1379
All	All	0.38	0/59985	0.76	33/88869 (0.0%)

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
2	B	0	1
23	X	0	1
All	All	0	2

There are no bond length outliers.

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	65	ALA	CB-CA-C	-14.85	87.83	110.10
5	E	15	ARG	CB-CA-C	-12.38	85.65	110.40
22	W	23	ARG	N-CA-C	-11.78	79.19	111.00
5	E	15	ARG	N-CA-C	-8.75	87.37	111.00
3	C	66	VAL	N-CA-C	-8.72	87.46	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	16	HIS	Peptide
23	X	53	ASP	Peptide

## 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	32542	0	16447	1121	0
2	B	1900	0	1951	39	0
3	C	1612	0	1674	77	0
4	D	1703	0	1766	42	0
5	E	1146	0	1207	46	0
6	F	843	0	857	15	0
7	G	1257	0	1296	24	0
8	H	1116	0	1177	40	0
9	I	1010	0	1037	32	0
10	J	792	0	835	48	0
11	K	885	0	904	20	0
12	L	970	0	1057	32	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	964	0	1034	25	0
14	N	492	0	529	28	0
15	O	734	0	771	13	0
16	P	700	0	720	19	0
17	Q	823	0	891	20	0
18	R	598	0	670	17	0
19	S	655	0	672	35	0
20	T	763	0	861	17	0
21	V	208	0	221	3	0
22	W	570	0	599	25	0
23	X	1356	0	1401	15	0
24	Y	373	0	186	17	0
25	Z	1646	0	845	109	0
26	D	1	0	0	4	0
26	N	1	0	0	0	0
27	W	1	0	0	0	0
27	Z	1	0	0	0	0
All	All	55662	0	39608	1690	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 18.

The worst 5 of 1690 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)
22:W:21:THR:CG2	22:W:33:LEU:HD11	1.53	1.36
10:J:7:LYS:HG3	10:J:71:LEU:CD2	1.55	1.33
1:A:1358:U:H3	1:A:1363(A):A:N6	1.36	1.24
1:A:1345:U:N3	1:A:1375:A:N6	1.87	1.21
5:E:93:PRO:HG3	8:H:105:ARG:NH2	1.53	1.20

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 PDB entries followed by that with respect to all EM entries.

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	B	232/256 (91%)	195 (84%)	24 (10%)	13 (6%)	2	26
3	C	204/239 (85%)	177 (87%)	18 (9%)	9 (4%)	3	31
4	D	206/209 (99%)	191 (93%)	12 (6%)	3 (2%)	12	54
5	E	148/162 (91%)	140 (95%)	8 (5%)	0	100	100
6	F	99/101 (98%)	94 (95%)	5 (5%)	0	100	100
7	G	153/156 (98%)	144 (94%)	7 (5%)	2 (1%)	14	56
8	H	136/138 (99%)	129 (95%)	6 (4%)	1 (1%)	25	68
9	I	125/128 (98%)	107 (86%)	15 (12%)	3 (2%)	7	44
10	J	96/105 (91%)	72 (75%)	16 (17%)	8 (8%)	1	17
11	K	117/129 (91%)	99 (85%)	13 (11%)	5 (4%)	3	31
12	L	122/132 (92%)	101 (83%)	17 (14%)	4 (3%)	4	37
13	M	119/126 (94%)	102 (86%)	16 (13%)	1 (1%)	22	67
14	N	58/61 (95%)	51 (88%)	5 (9%)	2 (3%)	4	37
15	O	86/89 (97%)	84 (98%)	1 (1%)	1 (1%)	15	58
16	P	81/88 (92%)	73 (90%)	7 (9%)	1 (1%)	15	58
17	Q	97/105 (92%)	88 (91%)	6 (6%)	3 (3%)	5	39
18	R	71/88 (81%)	66 (93%)	4 (6%)	1 (1%)	13	54
19	S	80/93 (86%)	66 (82%)	11 (14%)	3 (4%)	4	34
20	T	97/106 (92%)	80 (82%)	12 (12%)	5 (5%)	2	27
21	V	22/27 (82%)	21 (96%)	1 (4%)	0	100	100
22	W	69/72 (96%)	58 (84%)	7 (10%)	4 (6%)	2	25
23	X	166/171 (97%)	150 (90%)	13 (8%)	3 (2%)	10	50
All	All	2584/2781 (93%)	2288 (88%)	224 (9%)	72 (3%)	9	41

5 of 72 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	21	ARG
2	B	24	TRP
9	I	119	ALA
11	K	101	SER
12	L	79	GLU

### 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 PDB entries followed by that with respect to all EM entries.

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	B	202/220 (92%)	166 (82%)	36 (18%)	2	15
3	C	160/188 (85%)	142 (89%)	18 (11%)	7	31
4	D	180/181 (99%)	148 (82%)	32 (18%)	2	15
5	E	115/123 (94%)	93 (81%)	22 (19%)	2	12
6	F	90/90 (100%)	75 (83%)	15 (17%)	2	18
7	G	126/127 (99%)	121 (96%)	5 (4%)	36	66
8	H	119/119 (100%)	98 (82%)	21 (18%)	2	15
9	I	98/99 (99%)	84 (86%)	14 (14%)	4	24
10	J	87/92 (95%)	76 (87%)	11 (13%)	5	28
11	K	90/99 (91%)	73 (81%)	17 (19%)	2	12
12	L	104/109 (95%)	89 (86%)	15 (14%)	4	24
13	M	97/101 (96%)	84 (87%)	13 (13%)	4	26
14	N	49/50 (98%)	43 (88%)	6 (12%)	6	29
15	O	79/80 (99%)	64 (81%)	15 (19%)	2	12
16	P	72/74 (97%)	59 (82%)	13 (18%)	2	14
17	Q	94/97 (97%)	80 (85%)	14 (15%)	3	22
18	R	64/77 (83%)	52 (81%)	12 (19%)	2	13
19	S	71/80 (89%)	58 (82%)	13 (18%)	2	14
20	T	76/82 (93%)	60 (79%)	16 (21%)	1	9
21	V	19/22 (86%)	19 (100%)	0	100	100
22	W	62/63 (98%)	52 (84%)	10 (16%)	3	19
23	X	145/150 (97%)	126 (87%)	19 (13%)	5	27
All	All	2199/2323 (95%)	1862 (85%)	337 (15%)	7	21

5 of 337 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	I	14	VAL
11	K	120	ARG
22	W	21	THR
9	I	71	SER
10	J	87	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
7	G	28	ASN
9	I	124	GLN
19	S	83	HIS
7	G	37	ASN
9	I	73	GLN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1505/1522 (98%)	420 (27%)	103 (6%)
24	Y	15/42 (35%)	7 (46%)	0
25	Z	74/77 (96%)	33 (44%)	9 (12%)
All	All	1594/1641 (97%)	460 (28%)	112 (7%)

5 of 460 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	8	A
1	A	9	G
1	A	13	U
1	A	19	C
1	A	22	G

5 of 112 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	864	A
1	A	992	U
25	Z	17	C
1	A	884	U
1	A	956	U

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

5 non-standard protein/DNA/RNA residues are modelled in this entry.

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$
25	OMC	Z	32	25	15,22,23	0.78	1 (6%)	19,31,34	1.30	3 (15%)
25	G7M	Z	46	25	19,26,27	2.77	4 (21%)	19,39,42	2.26	5 (26%)
25	5MU	Z	54	25	14,22,23	0.85	0	16,32,35	2.96	3 (18%)
25	PSU	Z	55	25	16,21,22	1.17	1 (6%)	20,30,33	3.71	6 (30%)
25	4SU	Z	8	25	14,21,22	1.41	2 (14%)	15,30,33	1.88	3 (20%)

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
25	OMC	Z	32	25	-	0/5/27/28	0/2/2/2
25	G7M	Z	46	25	-	0/3/25/26	0/3/3/3
25	5MU	Z	54	25	-	0/3/25/26	0/2/2/2
25	PSU	Z	55	25	-	0/7/25/26	0/2/2/2
25	4SU	Z	8	25	-	0/3/25/26	0/2/2/2

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
25	Z	8	4SU	C4-S4	-3.90	1.60	1.67
25	Z	55	PSU	C5-C1'	-2.77	1.49	1.52
25	Z	46	G7M	O4'-C1'	2.08	1.44	1.41
25	Z	8	4SU	O4'-C1'	2.20	1.44	1.41
25	Z	32	OMC	O4'-C1'	2.30	1.44	1.41

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	Z	55	PSU	N1-C2-N3	-9.76	121.38	128.40
25	Z	55	PSU	C5-C4-N3	-9.15	117.92	125.43
25	Z	54	5MU	C5-C4-N3	-7.21	117.29	125.24
25	Z	46	G7M	C5-C6-N1	-4.89	116.52	123.48
25	Z	8	4SU	C5-C4-N3	-3.54	119.27	123.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
25	Z	32	OMC	2	0
25	Z	54	5MU	2	0
25	Z	55	PSU	5	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	7
25	Z	2

The worst 5 of 9 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	841:U	O3'	848:C	P	5.77
1	A	84:U	O3'	88:A	P	5.45
1	A	93:G	O3'	96:U	P	5.02
1	A	204:U	O3'	216:G	P	3.93
1	A	1442(A):G	O3'	1442(B):A	P	3.82