



wwPDB X-ray Structure Validation Summary Report (i)

Feb 28, 2014 – 12:33 AM GMT

PDB ID : 3PYU
Title : Crystal structure of a complex containing domain 3 of CrPV IGR IRES RNA bound to the 70S ribosome. This file contains the 30S subunit of the second 70S ribosome.
Authors : Zhu, J.; Korostelev, A.; Costantino, D.; Noller, H.F.; Kieft, J.S.
Deposited on : 2010-12-13
Resolution : 3.40 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

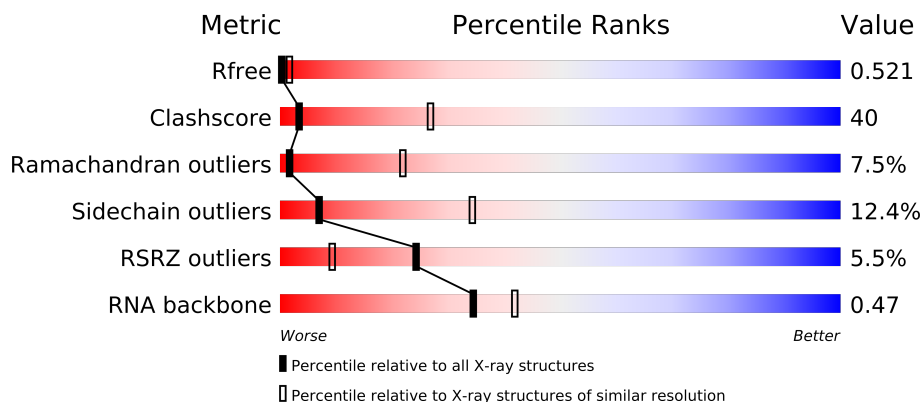
The following versions of software and data (see [references](#)) were used in the production of this report:

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

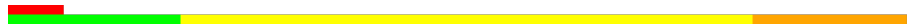


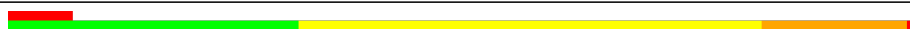
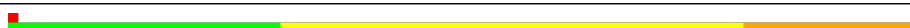
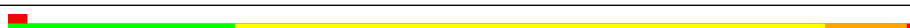

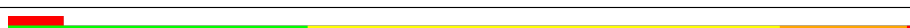

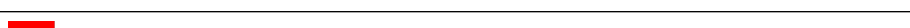


The reported resolution of this entry is 3.40 Å.

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.



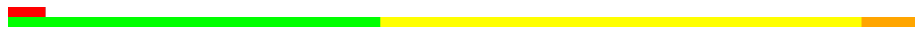
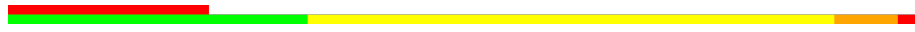


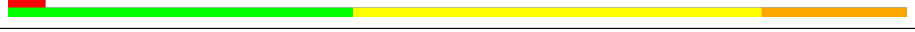
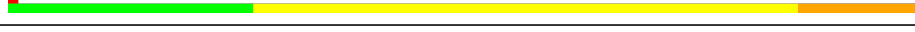
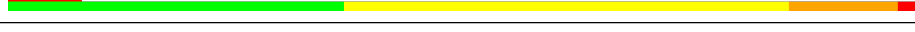


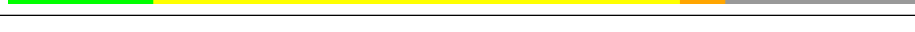
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1017 (3.52-3.28)
Clashscore	79885	1214 (3.50-3.30)
Ramachandran outliers	78287	1177 (3.50-3.30)
Sidechain outliers	78261	1177 (3.50-3.30)
RSRZ outliers	66119	1017 (3.52-3.28)
RNA backbone	1838	1002 (4.02-2.76)

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 ≥ 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	A	1506	
2	B	234	
3	C	206	
4	D	208	
5	E	151	
6	F	101	
7	G	155	
8	H	138	
9	I	127	
10	J	98	
11	K	119	
12	L	124	

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Mol	Chain	Length	Quality of chain
13	M	116	
14	N	60	
15	O	88	
16	P	83	
17	Q	99	
18	R	70	
19	S	78	
20	T	99	
21	U	24	
22	V	43	

2 Entry composition

There are 23 unique types of molecules in this entry. The entry contains 52199 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 ribosomal RNA 16S.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1506	Total	C	N	O	P	0	0	0
			32372	14409	5999	10459	1505			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	234	Total	C	N	O	S	0	0	0
			1901	1213	341	342	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	206	Total	C	N	O	S	0	0	0
			1613	1016	314	282	1			

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	151	Total	C	N	O	S	0	0	0
			1156	729	218	205	4			

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

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

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

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

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

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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	I	127	Total	C	N	O	0	0	0
			1011	639	198	174			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	98	Total	C	N	O	S	0	0	0
			795	499	156	139	1			

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	124	Total	C	N	O	S	0	0	0
			971	611	195	164	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
13	M	116	Total	C	N	O	S	0	0	0
			929	574	191	162	2			

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

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

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

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

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
16	P	83	Total	C	N	O	S	0	0	0
			701	443	139	118	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
17	Q	99	Total	C	N	O	S	0	0	0
			824	528	152	142	2			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
18	R	70	Total	C	N	O	0	0	0
			574	367	112	95			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
19	S	78	Total	C	N	O	S	0	0	0
			630	403	114	111	2			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
20	T	99	Total	C	N	O	S	0	0	0
			762	469	162	129	2			

- Molecule 21 is a protein called domain 3 of CrPV IGR IRES RNA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
21	U	24	Total	C	N	O	0	0	0
			209	128	50	31			

- Molecule 22 is a RNA chain called RNA (34-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
22	V	34	Total	C	N	O	P	0	0	0
			719	323	125	238	33			

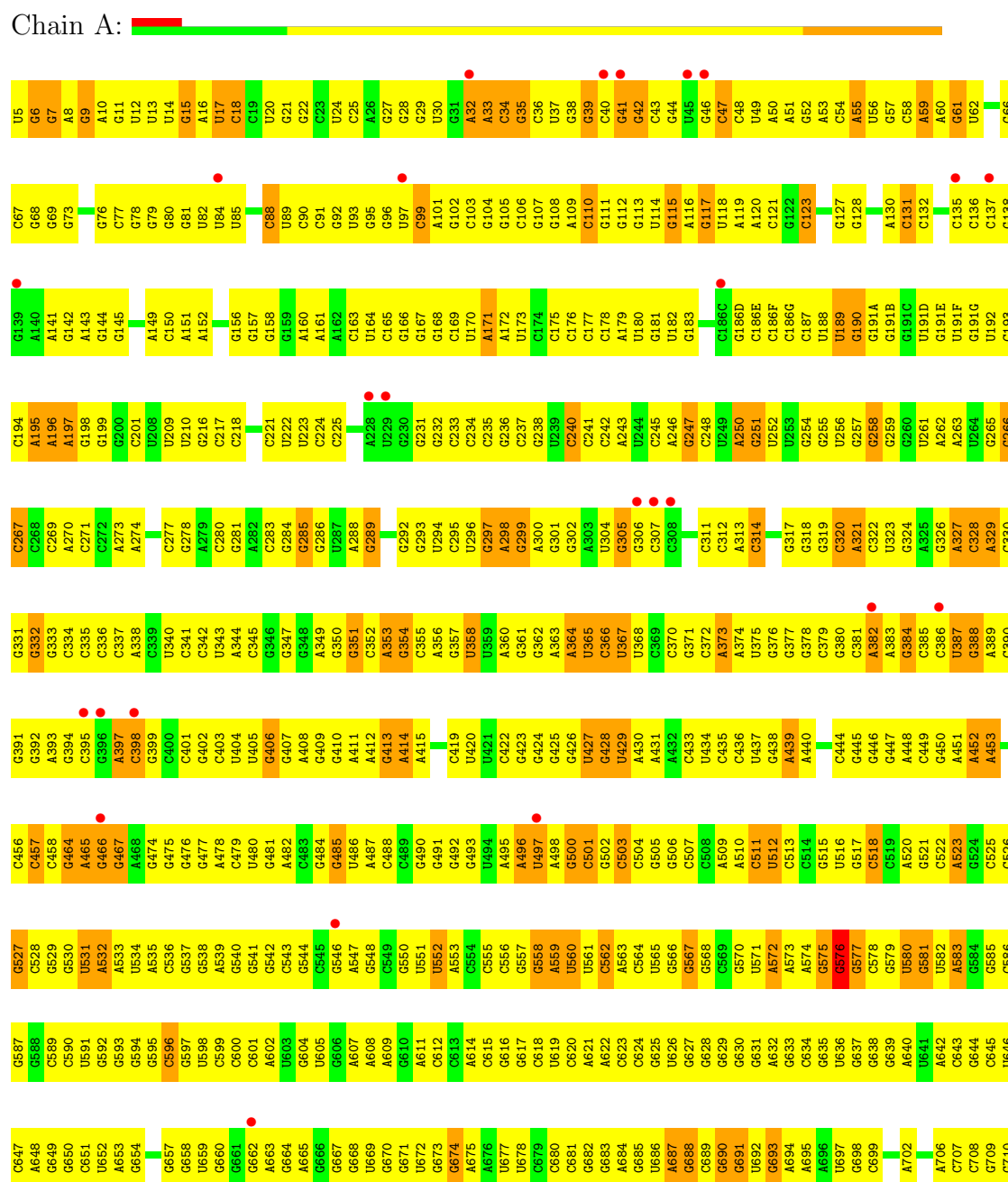
- Molecule 23 is ZINC ION (three-letter code: ZN) (formula: Zn).

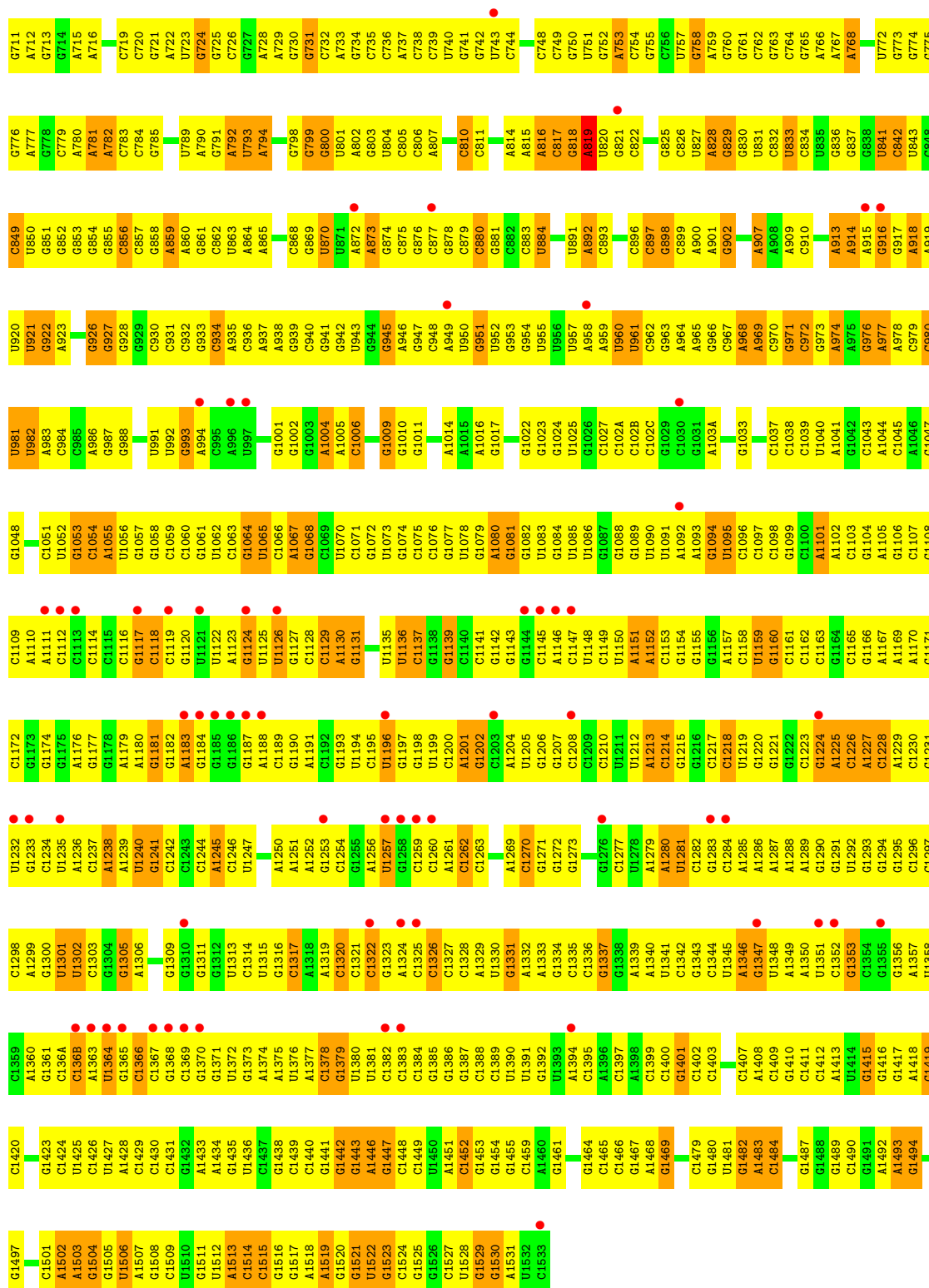
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
23	D	1	Total	Zn	0	0
			1	1		
23	N	1	Total	Zn	0	0
			1	1		

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 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: ribosomal RNA 16S

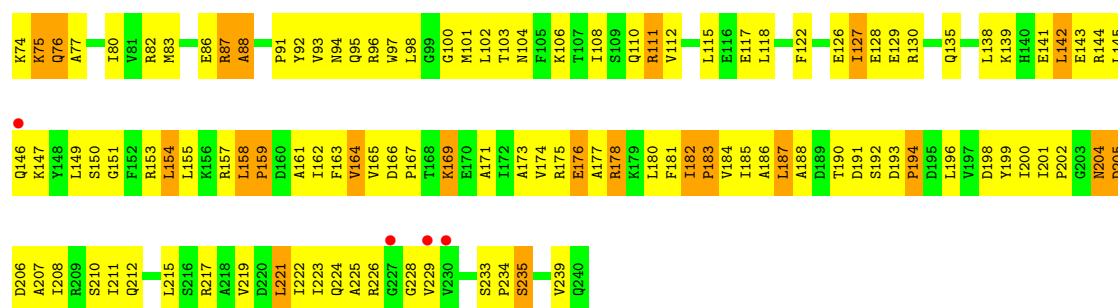




• Molecule 2: 30S ribosomal protein S2

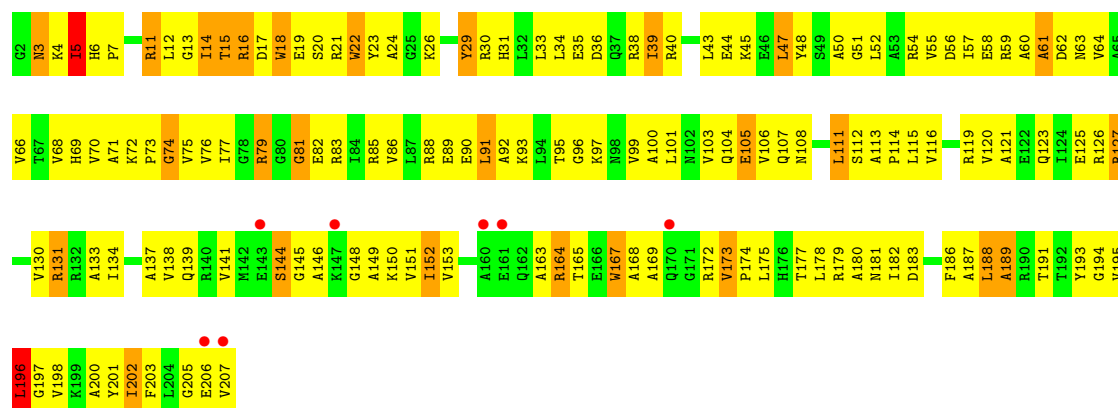
Chain B:





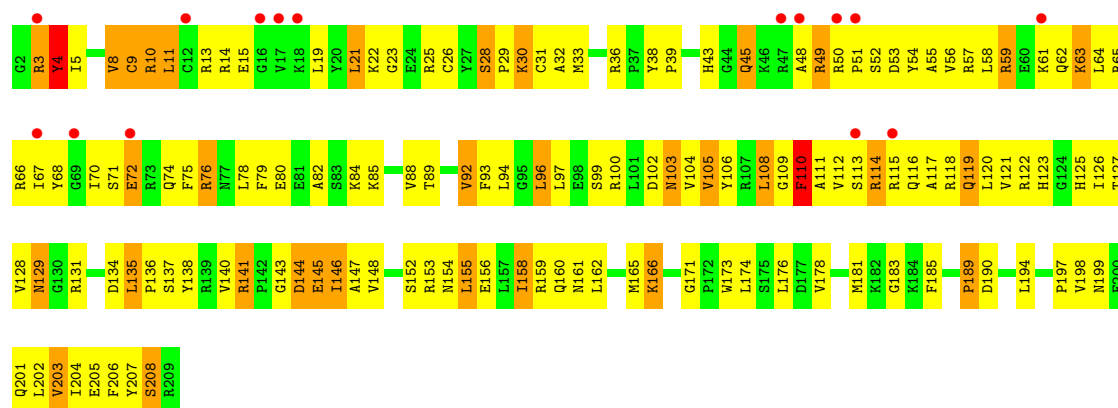
• Molecule 3: 30S ribosomal protein S3

Chain C:



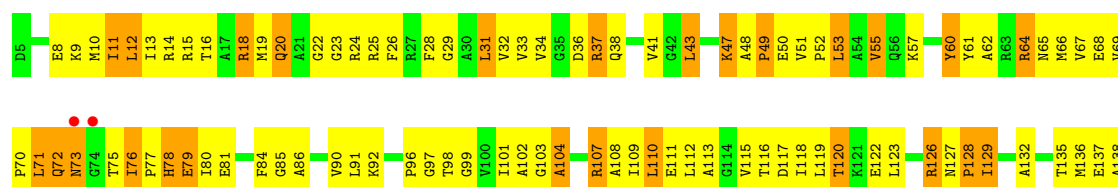
• Molecule 4: 30S ribosomal protein S4

Chain D:



• Molecule 5: 30S ribosomal protein S5

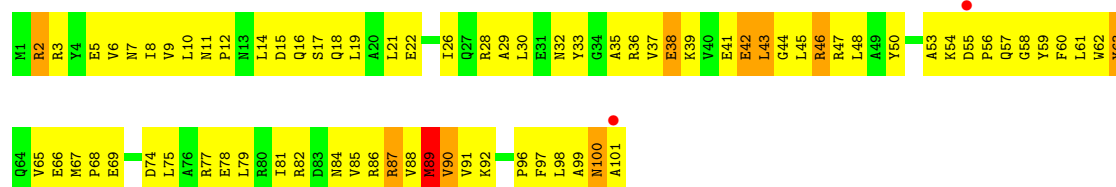
Chain E:





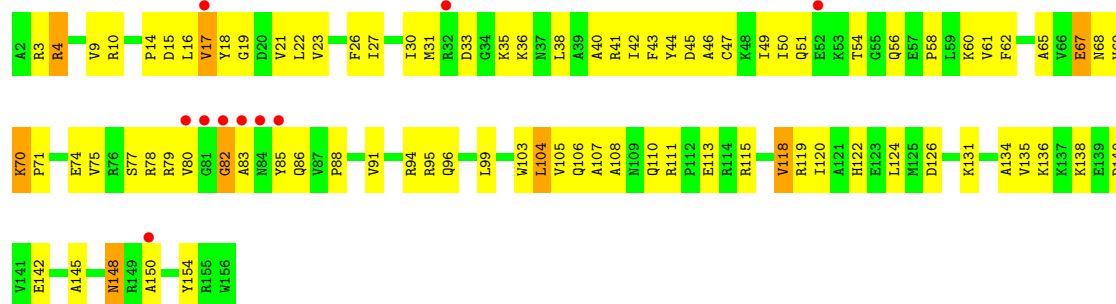
• Molecule 6: 30S ribosomal protein S6

Chain F:



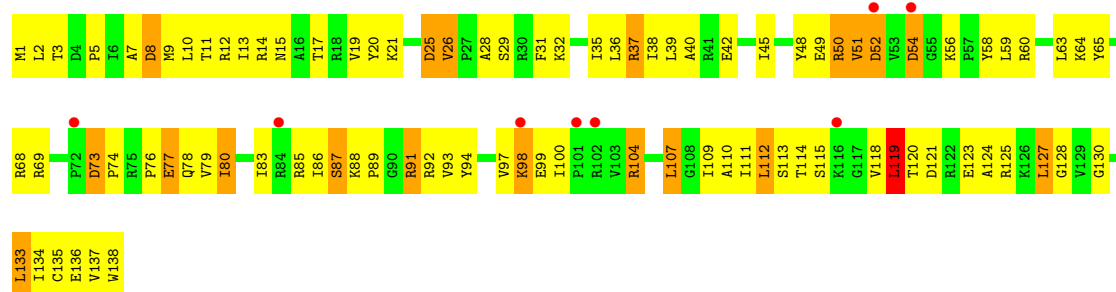
• Molecule 7: 30S ribosomal protein S7

Chain G:



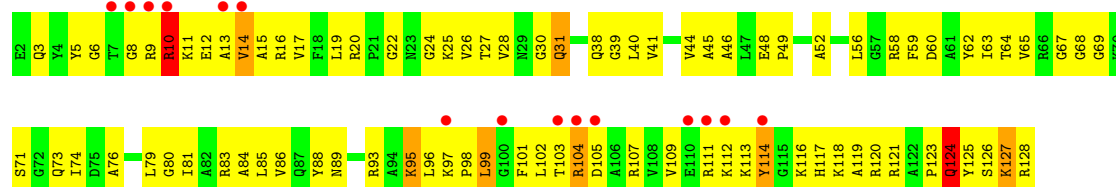
• Molecule 8: 30S ribosomal protein S8

Chain H:



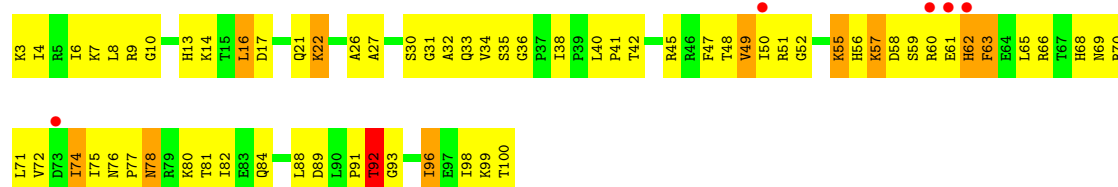
• Molecule 9: 30S ribosomal protein S9

Chain I:



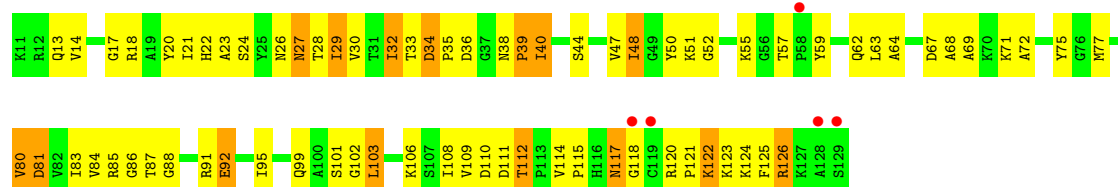
• Molecule 10: 30S ribosomal protein S10

Chain J:



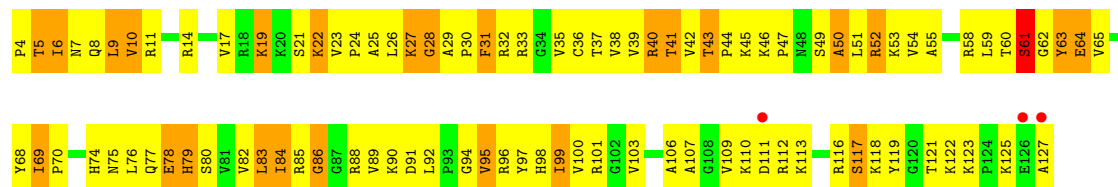
- Molecule 11: 30S ribosomal protein S11

Chain K:



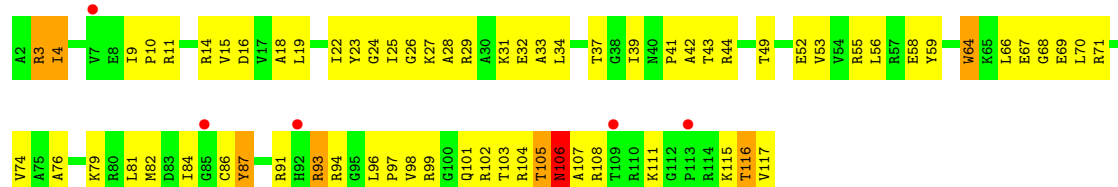
- Molecule 12: 30S ribosomal protein S12

Chain L:



- Molecule 13: 30S ribosomal protein S13

Chain M:



- Molecule 14: 30S ribosomal protein S14 type Z

Chain N:



- Molecule 15: 30S ribosomal protein S15

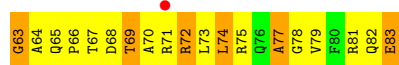
Chain O:





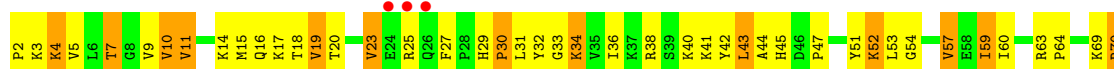
- Molecule 16: 30S ribosomal protein S16

Chain P:



- Molecule 17: 30S ribosomal protein S17

Chain Q:



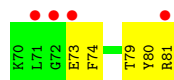
- Molecule 18: 30S ribosomal protein S18

Chain R:



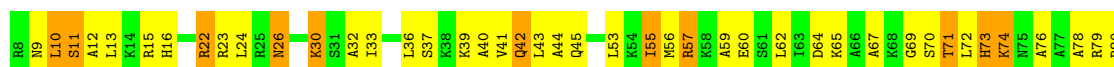
- Molecule 19: 30S ribosomal protein S19

Chain S:

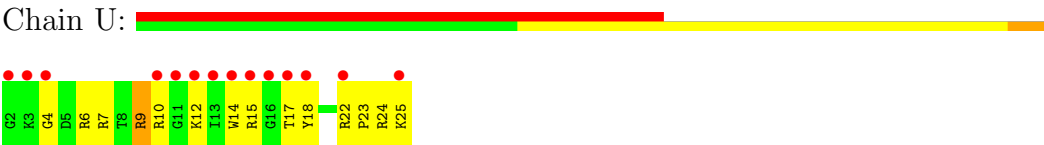


- Molecule 20: 30S ribosomal protein Thx

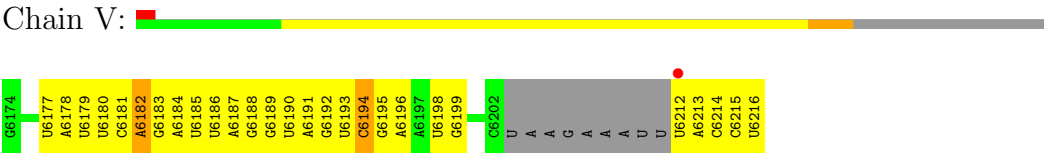
Chain T:



- Molecule 21: domain 3 of CrPV IGR IRES RNA



- Molecule 22: RNA (34-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	210.69Å 451.66Å 614.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.52 – 3.40 49.52 – 3.40	Depositor EDS
% Data completeness (in resolution range)	97.5 (49.52-3.40) 97.6 (49.52-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 3.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.228 , 0.266 0.517 , 0.521	Depositor DCC
R_{free} test set	7680 reflections (0.99%)	DCC
Wilson B-factor (Å ²)	86.0	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 76.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 775950 reflections	Xtriage
F_o, F_c correlation	0.48	EDS
Total number of atoms	52199	wwPDB-VP
Average B, all atoms (Å ²)	148.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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:
ZN

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	0/36238	0.96	75/56561 (0.1%)
2	B	0.28	0/1936	0.50	0/2609
3	C	0.29	0/1637	0.47	0/2205
4	D	0.34	0/1733	0.56	0/2318
5	E	0.36	0/1172	0.57	0/1576
6	F	0.37	0/856	0.59	0/1154
7	G	0.27	0/1276	0.46	0/1709
8	H	0.33	0/1136	0.58	0/1527
9	I	0.27	0/1029	0.45	0/1378
10	J	0.27	0/808	0.46	0/1085
11	K	0.41	0/900	0.61	0/1213
12	L	0.44	0/987	0.68	0/1320
13	M	0.24	0/939	0.44	0/1258
14	N	0.31	0/501	0.52	0/664
15	O	0.37	0/745	0.56	0/992
16	P	0.34	0/717	0.59	0/963
17	Q	0.37	0/837	0.56	0/1117
18	R	0.37	0/579	0.60	0/768
19	S	0.25	0/643	0.42	0/865
20	T	0.33	0/764	0.54	0/1006
21	U	0.24	0/213	0.42	0/277
22	V	0.43	0/802	0.69	0/1245
All	All	0.50	0/56448	0.85	75/83810 (0.1%)

There are no bond length outliers.

The worst 5 of 75 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	C	C6-N1-C2	10.59	124.53	120.30
1	A	285	G	C8-N9-C4	7.94	109.58	106.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1053	G	C4-N9-C1'	-7.85	116.30	126.50
1	A	1415	G	N1-C6-O6	7.64	124.48	119.90
1	A	117	G	N1-C6-O6	7.61	124.47	119.90

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	A	32372	0	16339	1760	0
2	B	1901	0	1951	180	0
3	C	1613	0	1677	186	0
4	D	1703	0	1764	182	0
5	E	1156	0	1213	141	0
6	F	843	0	857	93	0
7	G	1257	0	1296	92	0
8	H	1116	0	1177	140	0
9	I	1011	0	1043	112	0
10	J	795	0	840	92	0
11	K	885	0	904	72	0
12	L	971	0	1057	139	0
13	M	929	0	987	83	0
14	N	492	0	532	61	0
15	O	734	0	771	59	0
16	P	701	0	720	90	0
17	Q	824	0	893	77	0
18	R	574	0	644	70	0
19	S	630	0	652	60	0
20	T	762	0	859	70	0
21	U	209	0	221	17	0
22	V	719	0	366	57	0
23	D	1	0	0	0	0
23	N	1	0	0	0	0
All	All	52199	0	36763	3506	0

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 40.

The worst 5 of 3506 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
7:G:113:GLU:HB2	7:G:119:ARG:HG2	1.32	1.10
1:A:1347:G:C8	9:I:107:ARG:HB3	1.87	1.09
2:B:111:ARG:HH11	2:B:111:ARG:HG2	1.10	1.08
10:J:48:THR:HA	10:J:62:HIS:HB3	1.32	1.08
1:A:82:U:H2'	1:A:85:U:H5	1.18	1.07

There are no symmetry-related clashes.

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	
2	B	232/234 (99%)	173 (75%)	38 (16%)	21 (9%)	1	15
3	C	204/206 (99%)	134 (66%)	45 (22%)	25 (12%)	1	8
4	D	206/208 (99%)	151 (73%)	40 (19%)	15 (7%)	2	22
5	E	149/151 (99%)	104 (70%)	34 (23%)	11 (7%)	2	22
6	F	99/101 (98%)	71 (72%)	18 (18%)	10 (10%)	1	12
7	G	153/155 (99%)	121 (79%)	27 (18%)	5 (3%)	6	50
8	H	136/138 (99%)	98 (72%)	28 (21%)	10 (7%)	2	22
9	I	125/127 (98%)	89 (71%)	32 (26%)	4 (3%)	6	51
10	J	96/98 (98%)	74 (77%)	18 (19%)	4 (4%)	4	41
11	K	117/119 (98%)	82 (70%)	30 (26%)	5 (4%)	4	40
12	L	122/124 (98%)	80 (66%)	27 (22%)	15 (12%)	1	8
13	M	114/116 (98%)	93 (82%)	17 (15%)	4 (4%)	6	48
14	N	58/60 (97%)	46 (79%)	9 (16%)	3 (5%)	3	32
15	O	86/88 (98%)	61 (71%)	19 (22%)	6 (7%)	2	23
16	P	81/83 (98%)	46 (57%)	25 (31%)	10 (12%)	1	8

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
17	Q	97/99 (98%)	75 (77%)	16 (16%)	6 (6%)	2	27
18	R	68/70 (97%)	41 (60%)	18 (26%)	9 (13%)	0	7
19	S	76/78 (97%)	50 (66%)	21 (28%)	5 (7%)	2	25
20	T	97/99 (98%)	67 (69%)	23 (24%)	7 (7%)	2	23
21	U	22/24 (92%)	13 (59%)	8 (36%)	1 (4%)	4	38
All	All	2338/2378 (98%)	1669 (71%)	493 (21%)	176 (8%)	2	22

5 of 176 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	204	ASN
3	C	189	ALA
3	C	196	LEU
4	D	30	LYS
5	E	37	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	
2	B	202/202 (100%)	179 (89%)	23 (11%)	8	38
3	C	160/160 (100%)	145 (91%)	15 (9%)	13	50
4	D	180/180 (100%)	150 (83%)	30 (17%)	3	19
5	E	116/116 (100%)	94 (81%)	22 (19%)	2	11
6	F	90/90 (100%)	83 (92%)	7 (8%)	18	62
7	G	126/126 (100%)	121 (96%)	5 (4%)	42	84
8	H	119/119 (100%)	104 (87%)	15 (13%)	7	33
9	I	98/98 (100%)	88 (90%)	10 (10%)	11	46
10	J	88/88 (100%)	78 (89%)	10 (11%)	8	38
11	K	90/90 (100%)	76 (84%)	14 (16%)	4	23
12	L	104/104 (100%)	83 (80%)	21 (20%)	2	9
13	M	94/94 (100%)	87 (93%)	7 (7%)	20	65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
14	N	49/49 (100%)	45 (92%)	4 (8%)	17 60
15	O	79/79 (100%)	69 (87%)	10 (13%)	6 33
16	P	72/72 (100%)	56 (78%)	16 (22%)	1 7
17	Q	94/94 (100%)	79 (84%)	15 (16%)	3 21
18	R	61/61 (100%)	58 (95%)	3 (5%)	35 79
19	S	69/69 (100%)	60 (87%)	9 (13%)	6 31
20	T	76/76 (100%)	65 (86%)	11 (14%)	5 26
21	U	19/19 (100%)	19 (100%)	0	100 100
All	All	1986/1986 (100%)	1739 (88%)	247 (12%)	7 34

5 of 247 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
8	H	91	ARG
11	K	29	ILE
19	S	7	LYS
8	H	127	LEU
9	I	121	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 36 such sidechains are listed below:

Mol	Chain	Res	Type
8	H	82	HIS
10	J	56	HIS
19	S	65	ASN
9	I	117	HIS
10	J	78	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1505/1506 (99%)	294 (19%)	14 (0%)
22	V	32/43 (74%)	3 (9%)	0
All	All	1537/1549 (99%)	297 (19%)	14 (0%)

5 of 297 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	6	G
1	A	7	G
1	A	9	G
1	A	13	U
1	A	14	U

5 of 14 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	793	U
1	A	913	A
1	A	1201	A
1	A	687	A
1	A	1129	C

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 2 ligands modelled in this entry, 2 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, 95th 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 ²)	Q<0.9
1	A	1506/1506 (100%)	0.29	91 (6%) 21 8	51, 126, 251, 414	0
2	B	234/234 (100%)	0.09	6 (2%) 53 21	111, 177, 259, 325	0
3	C	206/206 (100%)	0.16	7 (3%) 43 17	105, 161, 226, 271	0
4	D	208/208 (100%)	0.32	15 (7%) 15 6	94, 146, 220, 300	0
5	E	151/151 (100%)	-0.17	2 (1%) 74 37	73, 117, 188, 252	0
6	F	101/101 (100%)	-0.18	2 (1%) 62 28	79, 131, 184, 246	0
7	G	155/155 (100%)	0.17	10 (6%) 18 7	119, 187, 237, 286	0
8	H	138/138 (100%)	0.29	8 (5%) 22 8	81, 123, 167, 219	0
9	I	127/127 (100%)	0.59	15 (11%) 5 3	121, 225, 286, 354	0
10	J	98/98 (100%)	0.27	5 (5%) 27 10	122, 197, 264, 351	0
11	K	119/119 (100%)	0.07	5 (4%) 35 13	74, 111, 178, 264	0
12	L	124/124 (100%)	0.12	3 (2%) 56 24	70, 109, 178, 252	0
13	M	116/116 (100%)	0.21	5 (4%) 34 13	135, 214, 309, 362	0
14	N	60/60 (100%)	1.12	13 (21%) 1 2	116, 167, 227, 281	0
15	O	88/88 (100%)	-0.13	2 (2%) 57 24	67, 110, 166, 241	0
16	P	83/83 (100%)	0.19	1 (1%) 75 39	87, 123, 177, 210	0
17	Q	99/99 (100%)	0.02	4 (4%) 36 14	79, 116, 166, 215	0
18	R	70/70 (100%)	-0.12	1 (1%) 72 35	82, 128, 192, 232	0
19	S	78/78 (100%)	0.70	6 (7%) 13 6	151, 216, 291, 350	0
20	T	99/99 (100%)	0.03	0 100 100	92, 136, 212, 269	0
21	U	24/24 (100%)	2.58	14 (58%) 0 0	163, 218, 265, 364	0
22	V	34/43 (79%)	-0.03	1 (2%) 49 21	92, 198, 333, 339	0
All	All	3918/3927 (99%)	0.23	216 (5%) 24 9	51, 144, 253, 414	0

The worst 5 of 216 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1257	U	9.4
1	A	1187	G	9.4
12	L	127	ALA	8.7
1	A	1112	C	6.8
1	A	1364	U	6.2

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, 95th 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 ²)	Q<0.9
23	ZN	N	62	1/1	0.04	-	144,144,144,144	0
23	ZN	D	210	1/1	0.10	-	131,131,131,131	0

6.5 Other polymers ⓘ

There are no such residues in this entry.