



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 26, 2026 – 02:54 AM UTC

PDB ID : 4LFC / pdb\_00004lfc  
Title : Crystal Structure of 30S ribosomal subunit from *Thermus thermophilus*  
Authors : Demirci, H.; Belardinelli, R.; Carr, J.; Murphy IV, F.; Jogl, G.; Dahlberg, A.E.; Gregory, S.T.  
Deposited on : 2013-06-26  
Resolution : 3.60 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

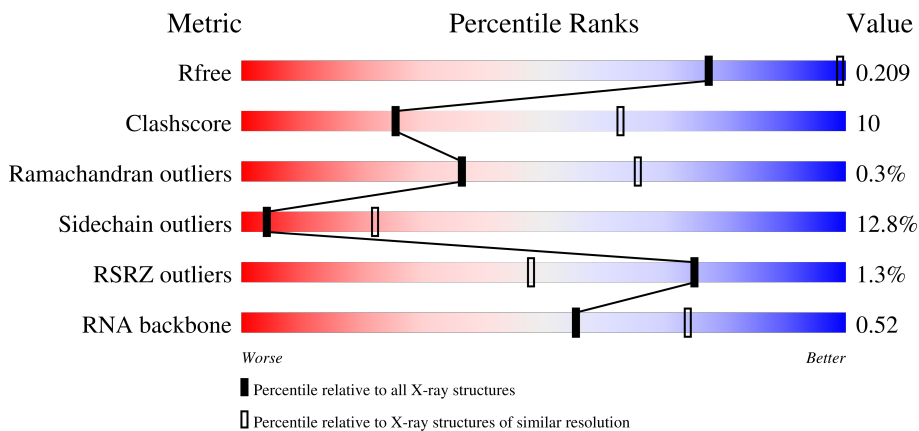
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.60 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.










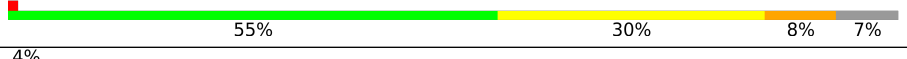



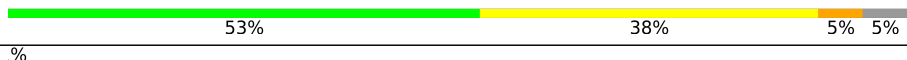
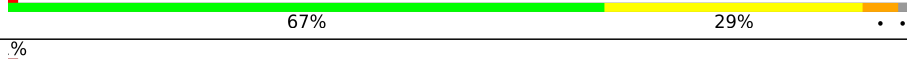


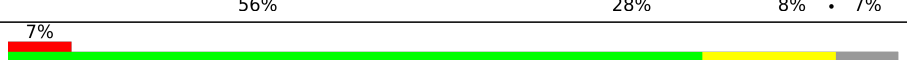
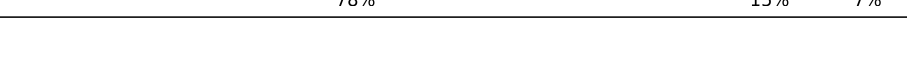
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)
RNA backbone	3983	1014 (4.14-3.06)

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 of 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  $\leq 5\%$ . 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	1522	 4% 60% 32% 7% ..
2	B	256	 1% 61% 28% • 8%
3	C	239	 61% 22% • 13%
4	D	209	 4% 66% 27% 7%

*Continued on next page...*

Continued from previous page...

Mol	Chain	Length	Quality of chain
5	E	162	
6	F	101	
7	G	156	
8	H	138	
9	I	128	
10	J	105	
11	K	129	
12	L	135	
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	U	27	

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
22	MG	A	1609	-	-	-	X
22	MG	A	1610	-	-	-	X
22	MG	A	1614	-	-	-	X

## 2 Entry composition [i](#)

There are 26 unique types of molecules in this entry. The entry contains 51888 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 rRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	1512	32504	14477	6011	10505	1511	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1534	C	A	conflict	GB M26923.1
A	1535	A	C	conflict	GB M26923.1

- Molecule 2 is a protein called ribosomal protein S2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	236	1874	1195	336	338	5	0	0	1

- Molecule 3 is a protein called ribosomal protein S3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	207	1613	1016	315	281	1	0	0	1

- Molecule 4 is a protein called ribosomal protein S4.

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

- Molecule 5 is a protein called ribosomal protein S5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	151	1147	724	218	201	4	0	0	1

- Molecule 6 is a protein called ribosomal protein S6.

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

- Molecule 7 is a protein called ribosomal protein S7.

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

- Molecule 8 is a protein called ribosomal protein S8.

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

- Molecule 9 is a protein called ribosomal protein S9.

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

- Molecule 10 is a protein called ribosomal protein S10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
10	J	99	793	498	157	137	1	0	0	1

- Molecule 11 is a protein called ribosomal protein S11.

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

- Molecule 12 is a protein called ribosomal protein S12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
12	L	125	973	612	196	163	2	0	0	1

- Molecule 13 is a protein called ribosomal protein S13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
13	M	118	937	579	193	163	2	0	0	0

- Molecule 14 is a protein called ribosomal protein S14.

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

- Molecule 15 is a protein called ribosomal protein S15.

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

- Molecule 16 is a protein called ribosomal protein S16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
16	P	84	701	443	140	117	1	0	0	1

- Molecule 17 is a protein called ribosomal protein S17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
17	Q	104	857	547	160	148	2	0	0	0

- Molecule 18 is a protein called ribosomal protein S18.

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

- Molecule 19 is a protein called ribosomal protein S19.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
19	S	81	648	414	120	112	2	0	0	1

- Molecule 20 is a protein called ribosomal protein S20.

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

- Molecule 21 is a protein called ribosomal protein THX.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
21	U	25	209	128	51	30	0	0	1

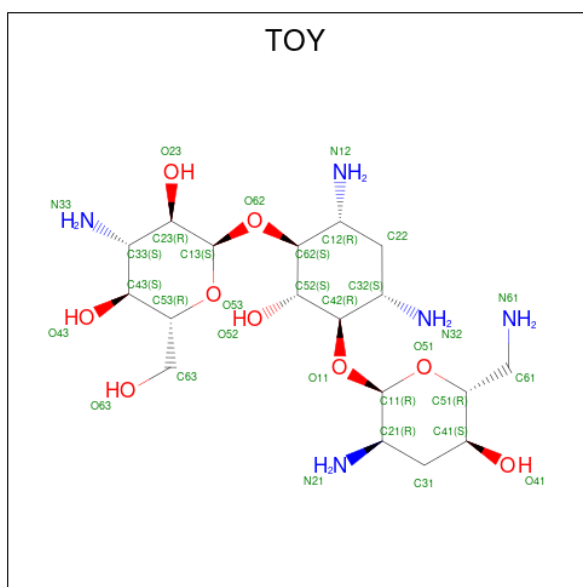
- Molecule 22 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
22	A	121	Total	Mg	0	0
			121	121		
22	B	1	Total	Mg	0	0
			1	1		
22	D	2	Total	Mg	0	0
			2	2		
22	E	1	Total	Mg	0	0
			1	1		
22	H	1	Total	Mg	0	0
			1	1		
22	J	1	Total	Mg	0	0
			1	1		
22	K	1	Total	Mg	0	0
			1	1		
22	M	1	Total	Mg	0	0
			1	1		
22	N	2	Total	Mg	0	0
			2	2		
22	Q	1	Total	Mg	0	0
			1	1		
22	S	1	Total	Mg	0	0
			1	1		

- Molecule 23 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
23	A	16	Total	K	0	0
			16	16		

- Molecule 24 is TOBRAMYCIN (CCD ID: TOY) (formula: C<sub>18</sub>H<sub>37</sub>N<sub>5</sub>O<sub>9</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
24	A	1	32	18	5	9	0	0

- Molecule 25 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

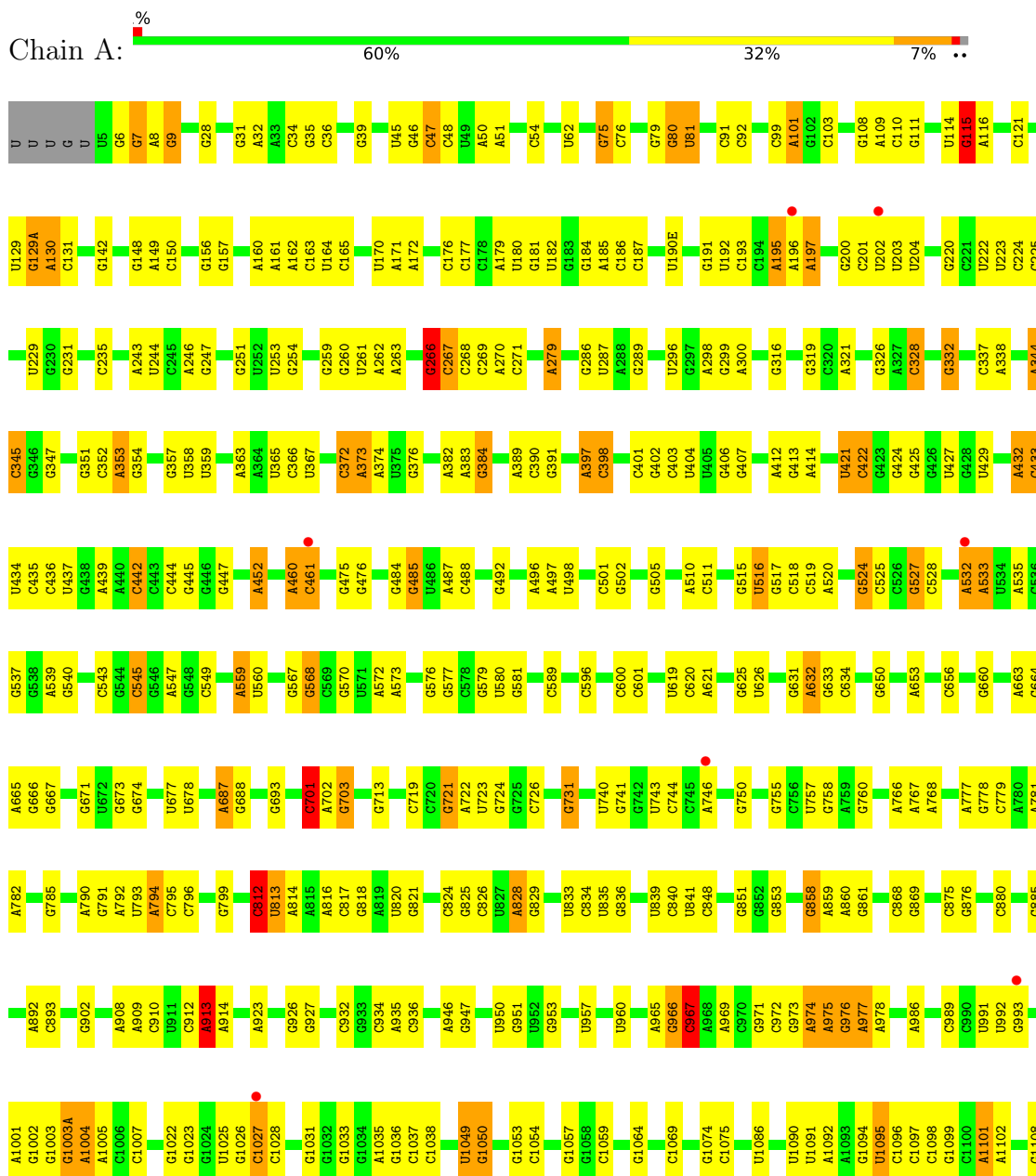
- Molecule 26 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
26	A	41	41	41	0	0
26	E	6	6	6	0	0
26	L	1	1	1	0	0

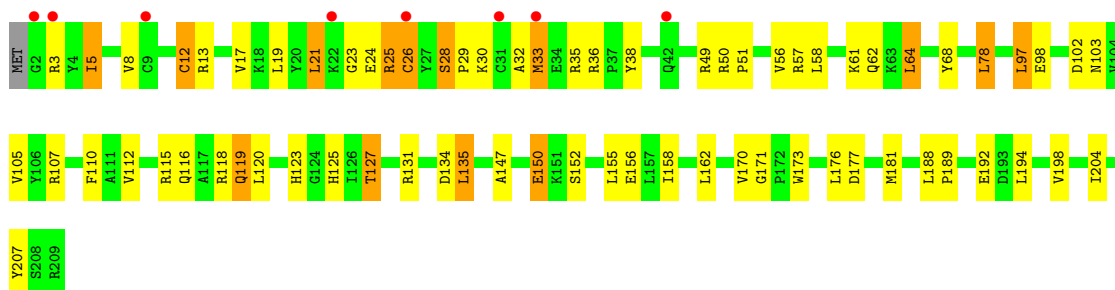
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

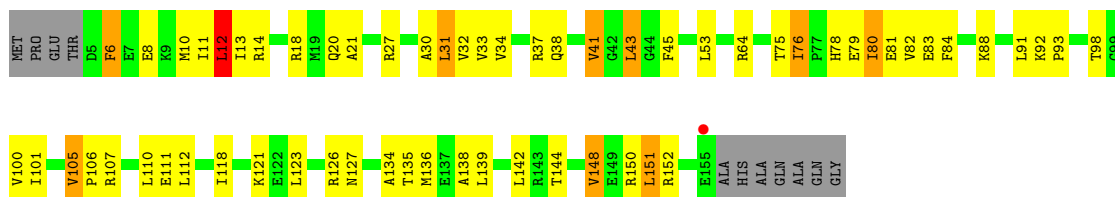
- Molecule 1: 16S rRNA



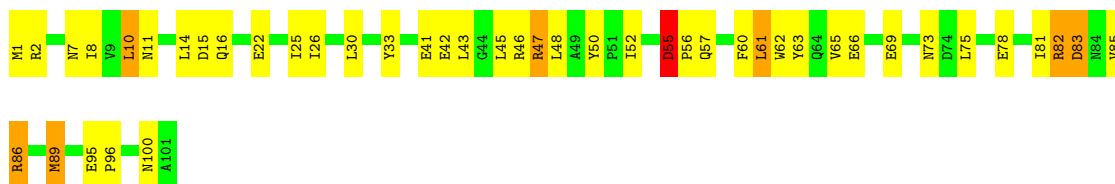




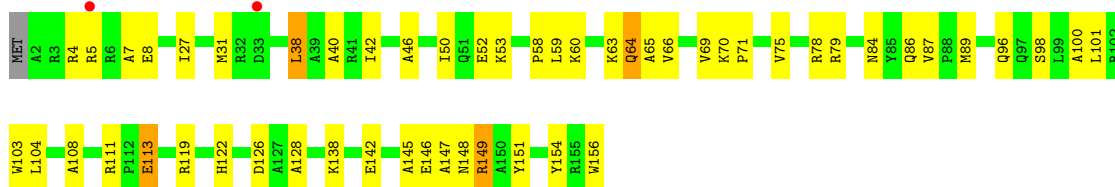
- Molecule 5: ribosomal protein S5



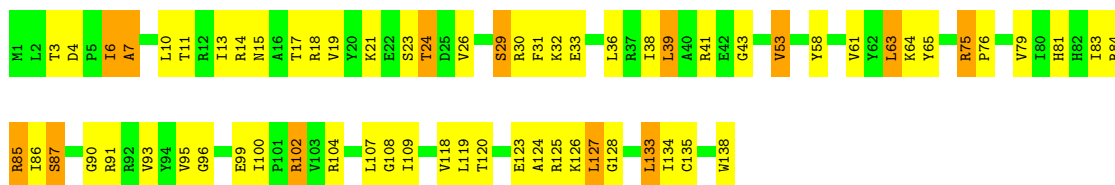
- Molecule 6: ribosomal protein S6



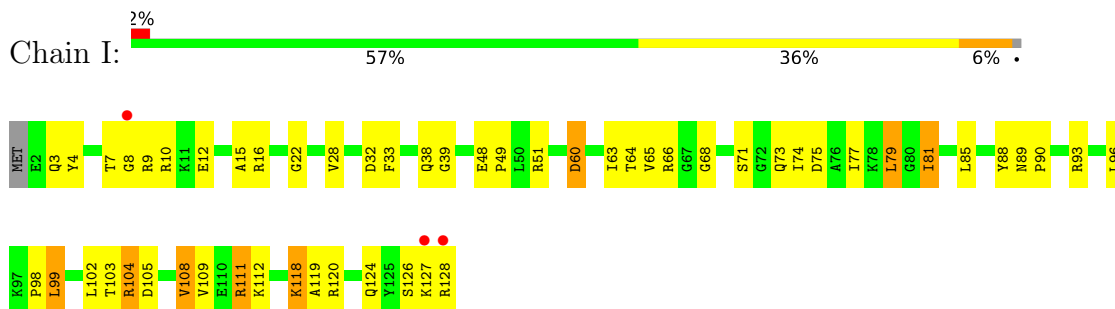
- Molecule 7: ribosomal protein S7



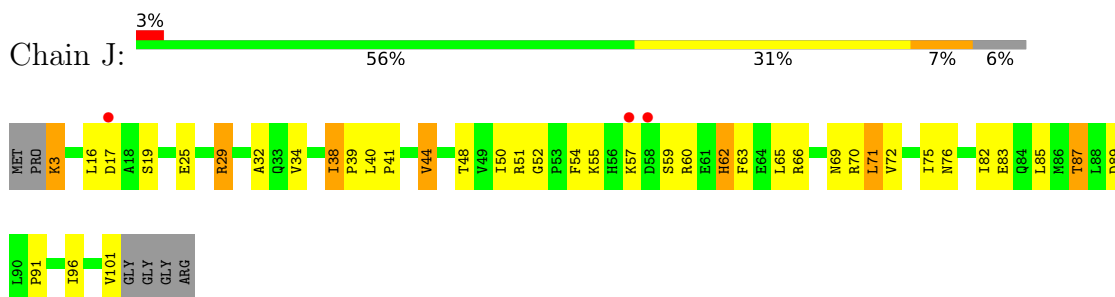
- Molecule 8: ribosomal protein S8



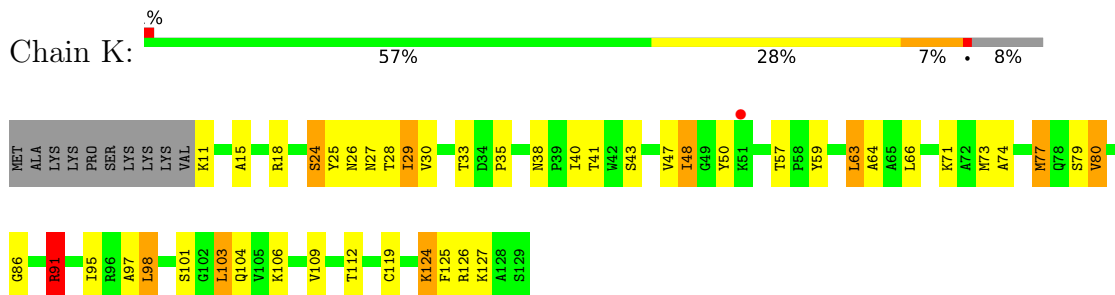
- Molecule 9: ribosomal protein S9



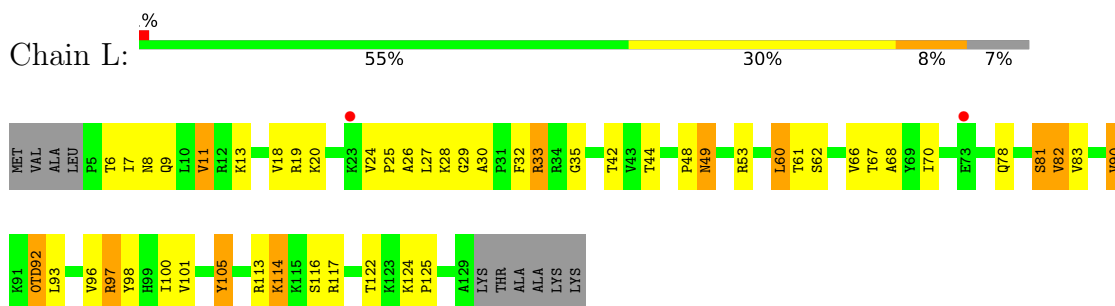
- Molecule 10: ribosomal protein S10



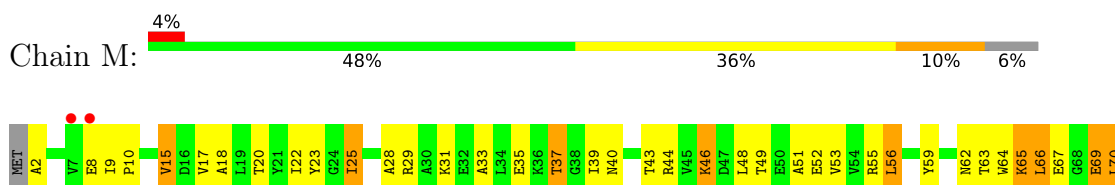
- Molecule 11: ribosomal protein S11



- Molecule 12: ribosomal protein S12

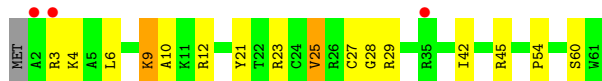


- Molecule 13: ribosomal protein S13





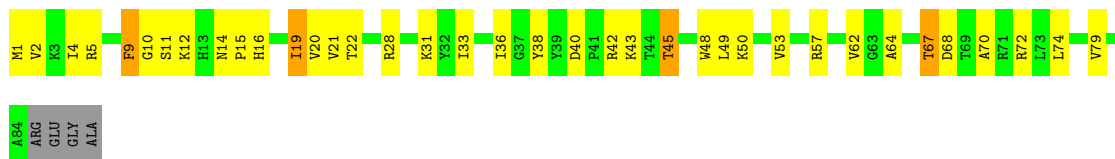
• Molecule 14: ribosomal protein S14



• Molecule 15: ribosomal protein S15



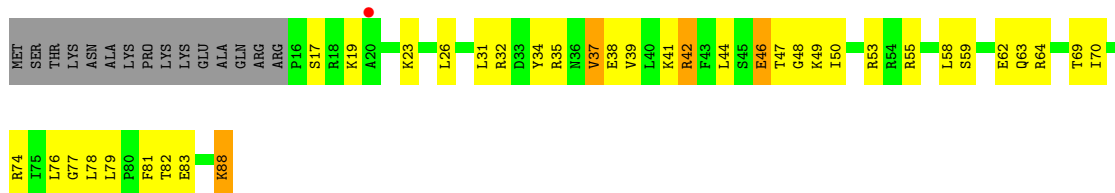
• Molecule 16: ribosomal protein S16



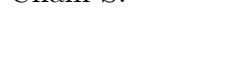
• Molecule 17: ribosomal protein S17

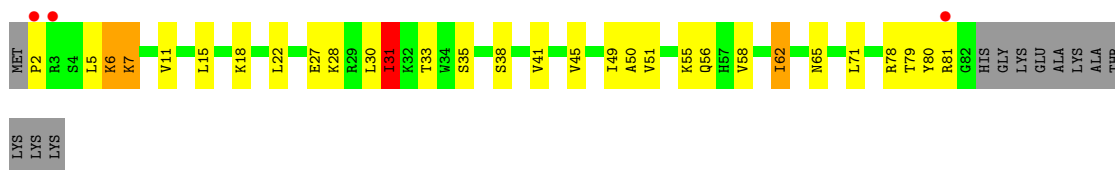


• Molecule 18: ribosomal protein S18

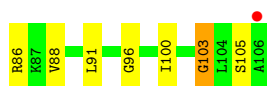
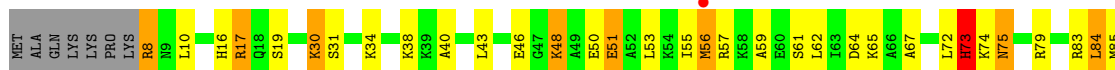


• Molecule 19: ribosomal protein S19

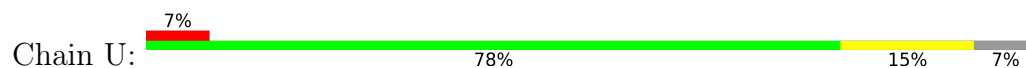




- Molecule 20: ribosomal protein S20



- Molecule 21: ribosomal protein THX



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	401.83Å 401.83Å 175.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.46 – 3.60 34.46 – 3.60	Depositor EDS
% Data completeness (in resolution range)	99.4 (34.46-3.60) 99.0 (34.46-3.60)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.86 (at 3.56Å)	Xtrriage
Refinement program	PHENIX dev_1119	Depositor
R, $R_{free}$	0.163 , 0.209 0.163 , 0.209	Depositor DCC
$R_{free}$ test set	8183 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	125.9	Xtrriage
Anisotropy	0.157	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 121.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	51888	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	133.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.57% 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 i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 4OC, MG, 2MG, MA6, UR3, K, 7MG, M2G, PSU, 0TD, TOY, 5MC

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	0.45	0/36037	0.63	10/56239 (0.0%)
2	B	0.73	2/1909 (0.1%)	1.04	9/2579 (0.3%)
3	C	0.58	0/1637	0.88	0/2207
4	D	0.67	0/1733	1.08	9/2318 (0.4%)
5	E	0.93	1/1163 (0.1%)	1.22	7/1566 (0.4%)
6	F	0.57	0/856	0.98	2/1154 (0.2%)
7	G	0.62	0/1276	0.95	2/1709 (0.1%)
8	H	0.96	0/1136	1.26	9/1527 (0.6%)
9	I	0.55	0/1029	1.02	3/1379 (0.2%)
10	J	0.62	0/806	1.04	5/1084 (0.5%)
11	K	0.82	0/900	1.19	6/1213 (0.5%)
12	L	0.78	1/978 (0.1%)	1.26	8/1308 (0.6%)
13	M	0.64	0/947	1.02	6/1270 (0.5%)
14	N	0.60	0/501	0.95	1/664 (0.2%)
15	O	0.77	0/745	0.98	0/992
16	P	0.72	0/717	1.10	6/965 (0.6%)
17	Q	0.93	1/870 (0.1%)	1.20	2/1159 (0.2%)
18	R	0.69	0/604	1.06	1/801 (0.1%)
19	S	0.56	0/662	0.97	2/892 (0.2%)
20	T	0.72	0/765	1.14	3/1007 (0.3%)
21	U	0.49	0/213	0.90	0/279
All	All	0.56	5/55484 (0.0%)	0.80	91/82312 (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
2	B	0	1
8	H	0	1

*Continued on next page...*

Continued from previous page...

Mol	Chain	#Chirality outliers	#Planarity outliers
9	I	0	1
10	J	0	1
All	All	0	4

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	13	ILE	CA-CB	-5.79	1.47	1.54
2	B	8	LYS	CA-C	5.71	1.58	1.52
2	B	9	GLU	CA-C	5.56	1.60	1.52
17	Q	9	VAL	CA-CB	-5.40	1.46	1.55
12	L	11	VAL	CA-CB	-5.06	1.48	1.54

The worst 5 of 91 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	I	39	GLY	N-CA-C	-13.05	96.21	115.00
4	D	26	CYS	N-CA-C	-10.15	100.92	113.20
12	L	30	ALA	CA-C-N	8.78	128.50	119.19
12	L	30	ALA	C-N-CA	8.78	128.50	119.19
8	H	75	ARG	CA-C-N	8.22	128.71	119.92

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	89	GLY	Peptide
8	H	90	GLY	Peptide
9	I	38	GLN	Peptide
10	J	87	THR	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	32504	0	16434	358	1

Continued on next page...

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1874	0	1887	42	0
3	C	1613	0	1677	37	0
4	D	1703	0	1763	45	0
5	E	1147	0	1207	40	0
6	F	843	0	857	35	0
7	G	1257	0	1296	32	0
8	H	1116	0	1177	39	0
9	I	1010	0	1037	39	1
10	J	793	0	835	24	0
11	K	885	0	904	24	0
12	L	973	0	1058	30	0
13	M	937	0	995	39	0
14	N	492	0	529	14	0
15	O	734	0	771	23	0
16	P	701	0	720	23	0
17	Q	857	0	928	22	0
18	R	598	0	670	24	0
19	S	648	0	673	19	0
20	T	763	0	861	37	0
21	U	209	0	221	2	0
22	A	121	0	0	0	0
22	B	1	0	0	0	0
22	D	2	0	0	0	0
22	E	1	0	0	0	0
22	H	1	0	0	0	0
22	J	1	0	0	0	0
22	K	1	0	0	0	0
22	M	1	0	0	0	0
22	N	2	0	0	0	0
22	Q	1	0	0	0	0
22	S	1	0	0	0	0
23	A	16	0	0	0	0
24	A	32	0	37	0	0
25	D	1	0	0	0	0
25	N	1	0	0	0	0
26	A	41	0	0	0	0
26	E	6	0	0	0	0
26	L	1	0	0	0	0
All	All	51888	0	36537	841	1

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 10.

The worst 5 of 841 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)
15:O:39:LEU:HD13	15:O:56:LEU:HB2	1.56	0.88
1:A:103:C:OP1	20:T:17:ARG:NH1	2.08	0.87
11:K:57:THR:HG22	11:K:59:TYR:H	1.42	0.84
16:P:57:ARG:NH1	16:P:79:VAL:O	2.11	0.83
7:G:27:ILE:HD12	7:G:40:ALA:HA	1.62	0.80

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	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:U:OP2	9:I:128:ARG:NH2[3_545]	2.19	0.01

## 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	
2	B	234/256 (91%)	214 (92%)	19 (8%)	1 (0%)	30	61
3	C	205/239 (86%)	190 (93%)	15 (7%)	0	100	100
4	D	206/209 (99%)	199 (97%)	6 (3%)	1 (0%)	24	57
5	E	149/162 (92%)	143 (96%)	6 (4%)	0	100	100
6	F	99/101 (98%)	96 (97%)	3 (3%)	0	100	100
7	G	153/156 (98%)	145 (95%)	8 (5%)	0	100	100
8	H	136/138 (99%)	132 (97%)	4 (3%)	0	100	100
9	I	125/128 (98%)	113 (90%)	11 (9%)	1 (1%)	16	49
10	J	97/105 (92%)	77 (79%)	18 (19%)	2 (2%)	5	31
11	K	117/129 (91%)	101 (86%)	16 (14%)	0	100	100
12	L	122/135 (90%)	113 (93%)	8 (7%)	1 (1%)	16	49

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
13	M	116/126 (92%)	107 (92%)	9 (8%)	0	100	100
14	N	58/61 (95%)	55 (95%)	3 (5%)	0	100	100
15	O	86/89 (97%)	82 (95%)	4 (5%)	0	100	100
16	P	82/88 (93%)	79 (96%)	3 (4%)	0	100	100
17	Q	102/105 (97%)	95 (93%)	7 (7%)	0	100	100
18	R	71/88 (81%)	65 (92%)	6 (8%)	0	100	100
19	S	79/93 (85%)	67 (85%)	11 (14%)	1 (1%)	9	39
20	T	97/106 (92%)	89 (92%)	7 (7%)	1 (1%)	12	45
21	U	23/27 (85%)	22 (96%)	1 (4%)	0	100	100
All	All	2357/2541 (93%)	2184 (93%)	165 (7%)	8 (0%)	36	65

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
12	L	28	LYS
19	S	31	ILE
10	J	72	VAL
9	I	119	ALA
20	T	73	HIS

### 5.3.2 Protein sidechains [i](#)

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	194/220 (88%)	173 (89%)	21 (11%)	6	28
3	C	160/188 (85%)	141 (88%)	19 (12%)	5	24
4	D	180/181 (99%)	162 (90%)	18 (10%)	7	30
5	E	115/123 (94%)	97 (84%)	18 (16%)	2	15
6	F	90/90 (100%)	78 (87%)	12 (13%)	4	21
7	G	126/127 (99%)	115 (91%)	11 (9%)	9	34
8	H	119/119 (100%)	101 (85%)	18 (15%)	3	17

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	I	98/99 (99%)	84 (86%)	14 (14%)	3	19
10	J	87/92 (95%)	77 (88%)	10 (12%)	5	25
11	K	90/99 (91%)	78 (87%)	12 (13%)	4	21
12	L	103/110 (94%)	86 (84%)	17 (16%)	2	14
13	M	94/101 (93%)	80 (85%)	14 (15%)	3	17
14	N	49/50 (98%)	45 (92%)	4 (8%)	10	36
15	O	79/80 (99%)	63 (80%)	16 (20%)	1	8
16	P	72/74 (97%)	64 (89%)	8 (11%)	6	27
17	Q	96/97 (99%)	85 (88%)	11 (12%)	5	25
18	R	64/77 (83%)	57 (89%)	7 (11%)	6	27
19	S	71/80 (89%)	61 (86%)	10 (14%)	3	19
20	T	76/82 (93%)	63 (83%)	13 (17%)	2	13
21	U	19/22 (86%)	19 (100%)	0	100	100
All	All	1982/2111 (94%)	1729 (87%)	253 (13%)	4	22

5 of 253 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
8	H	127	LEU
17	Q	92	ARG
11	K	48	ILE
17	Q	87	LYS
19	S	41	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
12	L	75	HIS
13	M	101	GLN
15	O	28	GLN
13	M	106	ASN
13	M	40	ASN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1510/1522 (99%)	241 (15%)	27 (1%)

5 of 241 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	7	G
1	A	9	G
1	A	31	G
1	A	32	A
1	A	39	G

5 of 27 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	965	A
1	A	1179	A
1	A	1347	G
1	A	1129	C
1	A	1182	G

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

15 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
1	5MC	A	1407	1	19,22,23	1.23	2 (10%)	26,32,35	1.15	2 (7%)
1	UR3	A	1498	1	19,22,23	0.93	0	26,32,35	1.10	1 (3%)
1	M2G	A	966	1	24,27,28	0.76	0	33,40,43	0.83	0
1	MA6	A	1518	1	23,26,27	1.03	1 (4%)	33,38,41	1.03	2 (6%)
1	PSU	A	1541	1	18,21,22	1.03	1 (5%)	21,30,33	1.79	5 (23%)
12	0TD	L	92	12	8,9,10	1.00	0	6,11,13	2.67	2 (33%)
1	5MC	A	1404	1	19,22,23	0.99	1 (5%)	26,32,35	1.01	2 (7%)
1	7MG	A	527	1	23,26,27	4.49	5 (21%)	27,39,42	2.69	10 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	MA6	A	1519	1	23,26,27	1.00	2 (8%)	33,38,41	0.97	0
1	PSU	A	516	22,1	18,21,22	1.33	1 (5%)	21,30,33	1.62	5 (23%)
1	4OC	A	1402	1	20,23,24	1.45	3 (15%)	25,32,35	0.80	1 (4%)
1	5MC	A	1400	1	19,22,23	1.16	2 (10%)	26,32,35	1.07	2 (7%)
1	5MC	A	967	1	19,22,23	1.16	2 (10%)	26,32,35	1.07	2 (7%)
1	PSU	A	1540	1	18,21,22	1.15	1 (5%)	21,30,33	1.65	2 (9%)
1	2MG	A	1207	1	23,26,27	1.44	6 (26%)	33,38,41	1.19	5 (15%)

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
1	5MC	A	1407	1	-	0/7/25/26	0/2/2/2
1	UR3	A	1498	1	-	0/7/25/26	0/2/2/2
1	M2G	A	966	1	-	3/11/29/30	0/3/3/3
1	MA6	A	1518	1	-	2/11/29/30	0/3/3/3
1	PSU	A	1541	1	-	1/7/25/26	0/2/2/2
12	0TD	L	92	12	-	3/7/12/14	-
1	5MC	A	1404	1	-	0/7/25/26	0/2/2/2
1	7MG	A	527	1	-	2/7/37/38	0/3/3/3
1	MA6	A	1519	1	-	2/11/29/30	0/3/3/3
1	PSU	A	516	22,1	-	0/7/25/26	0/2/2/2
1	4OC	A	1402	1	-	3/9/29/30	0/2/2/2
1	5MC	A	1400	1	-	2/7/25/26	0/2/2/2
1	5MC	A	967	1	-	2/7/25/26	0/2/2/2
1	PSU	A	1540	1	-	0/7/25/26	0/2/2/2
1	2MG	A	1207	1	-	0/9/27/28	0/3/3/3

The worst 5 of 27 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	527	7MG	C8-N9	-19.54	1.33	1.45
1	A	527	7MG	C5-N7	5.34	1.42	1.35
1	A	527	7MG	C2-N2	5.07	1.46	1.34
1	A	516	PSU	C6-C5	4.53	1.40	1.35
1	A	1402	4OC	C4-N4	4.11	1.44	1.36

The worst 5 of 41 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	527	7MG	C5-C6-N1	6.34	122.11	110.94
1	A	527	7MG	C2-N3-C4	5.82	122.33	112.30
12	L	92	0TD	CSB-SB-CB	-5.42	92.62	102.36
1	A	1540	PSU	C4-N3-C2	-4.69	119.91	126.37
1	A	1541	PSU	C4-N3-C2	-4.50	120.17	126.37

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	527	7MG	O4'-C4'-C5'-O5'
1	A	967	5MC	O4'-C4'-C5'-O5'
1	A	967	5MC	C3'-C4'-C5'-O5'
1	A	1402	4OC	O4'-C4'-C5'-O5'
1	A	527	7MG	C3'-C4'-C5'-O5'

There are no ring outliers.

10 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	1498	UR3	2	0
1	A	966	M2G	1	0
1	A	1518	MA6	2	0
12	L	92	0TD	3	0
1	A	1404	5MC	1	0
1	A	1519	MA6	2	0
1	A	516	PSU	1	0
1	A	1402	4OC	1	0
1	A	1400	5MC	1	0
1	A	967	5MC	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 152 ligands modelled in this entry, 151 are 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	$\# Z  > 2$
24	TOY	A	1738	-	33,34,34	1.14	4 (12%)	43,50,50	1.75	12 (27%)

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
24	TOY	A	1738	-	-	5/12/68/68	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
24	A	1738	TOY	C11-C21	2.71	1.57	1.52
24	A	1738	TOY	C23-C33	2.70	1.56	1.53
24	A	1738	TOY	C43-C53	2.13	1.57	1.53
24	A	1738	TOY	C41-C51	2.09	1.55	1.52

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	A	1738	TOY	C43-C33-C23	-3.75	102.92	111.06
24	A	1738	TOY	C62-C52-C42	3.62	116.44	109.11
24	A	1738	TOY	C13-C23-C33	-3.34	105.91	110.40
24	A	1738	TOY	O51-C51-C61	3.26	112.34	106.07
24	A	1738	TOY	O52-C52-C42	-2.98	102.29	109.94

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
24	A	1738	TOY	C41-C51-C61-N61
24	A	1738	TOY	O51-C51-C61-N61
24	A	1738	TOY	C43-C53-C63-O63

*Continued on next page...*

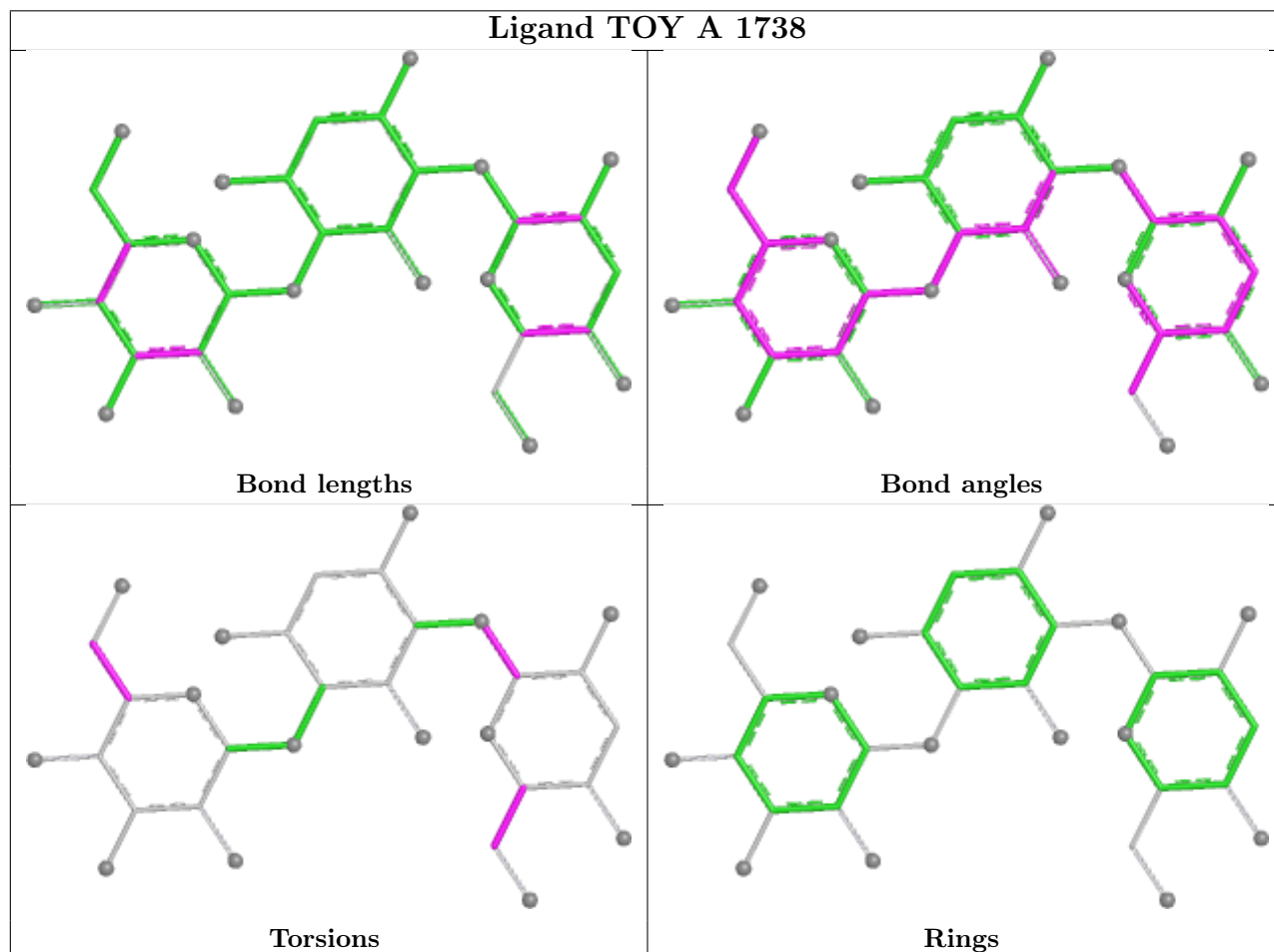
*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
24	A	1738	TOY	O51-C11-O11-C42
24	A	1738	TOY	O53-C53-C63-O63

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 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	1498/1522 (98%)	-0.37	10 (0%) 84 61	67, 118, 217, 321	0
2	B	236/256 (92%)	-0.32	2 (0%) 82 58	93, 136, 229, 255	0
3	C	207/239 (86%)	-0.14	0 100 100	76, 164, 205, 222	0
4	D	208/209 (99%)	-0.10	8 (3%) 44 25	86, 126, 175, 199	0
5	E	151/162 (93%)	-0.45	1 (0%) 84 61	71, 101, 141, 191	0
6	F	101/101 (100%)	-0.48	0 100 100	94, 142, 179, 222	0
7	G	155/156 (99%)	-0.18	2 (1%) 75 48	106, 149, 219, 255	0
8	H	138/138 (100%)	-0.64	0 100 100	62, 91, 129, 155	0
9	I	127/128 (99%)	0.12	3 (2%) 59 34	115, 170, 210, 233	0
10	J	99/105 (94%)	0.20	3 (3%) 52 30	104, 194, 292, 344	0
11	K	119/129 (92%)	-0.31	1 (0%) 82 58	82, 120, 159, 188	0
12	L	124/135 (91%)	-0.25	2 (1%) 70 43	61, 120, 159, 224	0
13	M	118/126 (93%)	-0.08	5 (4%) 40 22	124, 154, 198, 261	0
14	N	60/61 (98%)	0.07	3 (5%) 34 19	123, 150, 188, 221	0
15	O	88/89 (98%)	-0.29	0 100 100	74, 113, 157, 200	0
16	P	84/88 (95%)	-0.17	0 100 100	89, 122, 153, 219	0
17	Q	104/105 (99%)	-0.21	1 (0%) 79 54	72, 104, 164, 222	0
18	R	73/88 (82%)	-0.17	1 (1%) 73 46	85, 122, 187, 248	0
19	S	81/93 (87%)	0.22	3 (3%) 45 25	70, 182, 244, 270	0
20	T	99/106 (93%)	-0.05	2 (2%) 65 38	94, 123, 176, 217	0
21	U	25/27 (92%)	0.45	2 (8%) 18 13	72, 154, 194, 233	0
All	All	3895/4063 (95%)	-0.26	49 (1%) 75 48	61, 130, 211, 344	0

The worst 5 of 49 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
19	S	3	ARG	8.0
1	A	1129	C	6.3
4	D	31	CYS	5.9
4	D	2	GLY	5.0
9	I	127	LYS	4.8

## 6.2 Non-standard residues in protein, DNA, RNA chains [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. 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	B-factors(Å <sup>2</sup> )	Q<0.9
1	PSU	A	1540	20/21	0.89	0.14	204,217,231,234	0
1	PSU	A	1541	20/21	0.92	0.15	203,212,220,220	0
1	M2G	A	966	25/26	0.94	0.08	104,126,133,143	0
1	PSU	A	516	20/21	0.96	0.07	101,130,144,148	0
1	5MC	A	1407	21/22	0.97	0.08	93,112,121,124	0
1	UR3	A	1498	21/22	0.97	0.11	89,95,102,107	0
1	5MC	A	967	21/22	0.97	0.06	92,104,118,121	0
1	5MC	A	1400	21/22	0.97	0.12	86,107,119,121	0
12	0TD	L	92	10/11	0.97	0.13	92,125,143,316	0
1	7MG	A	527	24/25	0.98	0.06	104,111,119,126	0
1	MA6	A	1518	24/25	0.98	0.09	87,94,123,125	0
1	MA6	A	1519	24/25	0.98	0.11	83,89,94,107	0
1	4OC	A	1402	22/23	0.98	0.06	88,99,111,114	0
1	5MC	A	1404	21/22	0.98	0.06	86,89,97,102	0
1	2MG	A	1207	24/25	0.98	0.06	151,161,170,171	0

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides 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. 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	B-factors( $\text{\AA}^2$ )	Q<0.9
23	K	A	1707	1/1	0.53	0.24	148,148,148,148	0
22	MG	A	1725	1/1	0.54	0.34	129,129,129,129	0
22	MG	A	1614	1/1	0.57	0.47	107,107,107,107	0
22	MG	A	1718	1/1	0.61	0.40	86,86,86,86	0
22	MG	A	1608	1/1	0.63	0.32	93,93,93,93	0
23	K	A	1706	1/1	0.66	0.31	127,127,127,127	0
22	MG	A	1726	1/1	0.66	0.24	114,114,114,114	0
22	MG	N	102	1/1	0.70	0.37	96,96,96,96	0
22	MG	A	1724	1/1	0.72	0.24	89,89,89,89	0
23	K	A	1701	1/1	0.74	0.21	144,144,144,144	0
22	MG	A	1722	1/1	0.75	0.28	107,107,107,107	0
22	MG	A	1643	1/1	0.75	0.31	76,76,76,76	0
23	K	A	1712	1/1	0.75	0.23	130,130,130,130	0
22	MG	A	1656	1/1	0.76	0.24	102,102,102,102	0
22	MG	A	1606	1/1	0.76	0.36	97,97,97,97	0
23	K	A	1704	1/1	0.76	0.24	114,114,114,114	0
23	K	A	1714	1/1	0.77	0.18	147,147,147,147	0
22	MG	A	1730	1/1	0.78	0.22	126,126,126,126	0
22	MG	B	301	1/1	0.78	0.17	118,118,118,118	0
22	MG	A	1632	1/1	0.78	0.32	74,74,74,74	0
23	K	A	1700	1/1	0.78	0.20	131,131,131,131	0
23	K	A	1713	1/1	0.78	0.14	162,162,162,162	0
22	MG	A	1609	1/1	0.78	0.80	118,118,118,118	0
23	K	A	1711	1/1	0.79	0.14	163,163,163,163	0
22	MG	A	1610	1/1	0.79	0.45	67,67,67,67	0
22	MG	A	1717	1/1	0.79	0.34	88,88,88,88	0
22	MG	A	1616	1/1	0.79	0.25	69,69,69,69	0
22	MG	A	1650	1/1	0.80	0.32	61,61,61,61	0
22	MG	A	1720	1/1	0.80	0.30	90,90,90,90	0
22	MG	A	1645	1/1	0.81	0.18	88,88,88,88	0
23	K	A	1709	1/1	0.81	0.31	142,142,142,142	0
22	MG	A	1688	1/1	0.81	0.27	69,69,69,69	0
22	MG	A	1639	1/1	0.82	0.28	72,72,72,72	0
22	MG	A	1613	1/1	0.82	0.40	63,63,63,63	0
23	K	A	1708	1/1	0.83	0.27	164,164,164,164	0
22	MG	A	1652	1/1	0.83	0.33	86,86,86,86	0
22	MG	A	1661	1/1	0.83	0.48	81,81,81,81	0
22	MG	A	1731	1/1	0.84	0.06	361,361,361,361	0
23	K	A	1710	1/1	0.84	0.10	190,190,190,190	0
22	MG	A	1734	1/1	0.84	0.09	288,288,288,288	0
22	MG	A	1615	1/1	0.84	0.25	69,69,69,69	0
22	MG	A	1647	1/1	0.84	0.25	68,68,68,68	0
22	MG	A	1637	1/1	0.84	0.40	128,128,128,128	0

Continued on next page...

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
22	MG	D	303	1/1	0.85	0.17	99,99,99,99	0
22	MG	A	1658	1/1	0.85	0.24	99,99,99,99	0
23	K	A	1705	1/1	0.85	0.32	160,160,160,160	0
22	MG	A	1732	1/1	0.85	0.06	258,258,258,258	0
23	K	A	1715	1/1	0.85	0.17	155,155,155,155	0
22	MG	A	1636	1/1	0.86	0.45	94,94,94,94	0
22	MG	A	1697	1/1	0.86	0.18	79,79,79,79	0
22	MG	A	1651	1/1	0.86	0.35	100,100,100,100	0
22	MG	A	1633	1/1	0.86	0.27	81,81,81,81	0
22	MG	A	1681	1/1	0.86	0.17	105,105,105,105	0
22	MG	A	1675	1/1	0.87	0.54	90,90,90,90	0
22	MG	A	1670	1/1	0.88	0.37	100,100,100,100	0
22	MG	A	1655	1/1	0.88	0.34	76,76,76,76	0
22	MG	A	1602	1/1	0.88	0.63	94,94,94,94	0
22	MG	H	201	1/1	0.89	0.14	76,76,76,76	0
22	MG	M	201	1/1	0.89	0.14	123,123,123,123	0
22	MG	A	1716	1/1	0.89	0.19	84,84,84,84	0
22	MG	S	101	1/1	0.89	0.22	99,99,99,99	0
22	MG	A	1723	1/1	0.89	0.36	99,99,99,99	0
22	MG	A	1664	1/1	0.89	0.28	103,103,103,103	0
22	MG	A	1619	1/1	0.89	0.28	72,72,72,72	0
22	MG	A	1699	1/1	0.89	0.30	85,85,85,85	0
22	MG	A	1729	1/1	0.89	0.13	81,81,81,81	0
22	MG	A	1611	1/1	0.90	0.26	75,75,75,75	0
22	MG	A	1623	1/1	0.90	0.23	123,123,123,123	0
22	MG	D	302	1/1	0.90	0.12	74,74,74,74	0
23	K	A	1702	1/1	0.90	0.26	137,137,137,137	0
22	MG	A	1694	1/1	0.90	0.17	77,77,77,77	0
22	MG	A	1634	1/1	0.90	0.30	66,66,66,66	0
22	MG	A	1673	1/1	0.90	0.27	51,51,51,51	0
22	MG	A	1635	1/1	0.90	0.25	76,76,76,76	0
22	MG	A	1638	1/1	0.91	0.36	62,62,62,62	0
22	MG	A	1683	1/1	0.91	0.17	57,57,57,57	0
23	K	A	1703	1/1	0.91	0.21	138,138,138,138	0
22	MG	E	201	1/1	0.91	0.08	318,318,318,318	0
22	MG	A	1620	1/1	0.91	0.35	70,70,70,70	0
22	MG	A	1682	1/1	0.92	0.19	70,70,70,70	0
22	MG	A	1663	1/1	0.92	0.40	58,58,58,58	0
22	MG	A	1676	1/1	0.92	0.13	80,80,80,80	0
22	MG	N	103	1/1	0.92	0.15	100,100,100,100	0
22	MG	Q	201	1/1	0.92	0.12	91,91,91,91	0
22	MG	A	1692	1/1	0.92	0.20	118,118,118,118	0

*Continued on next page...*

*Continued from previous page...*

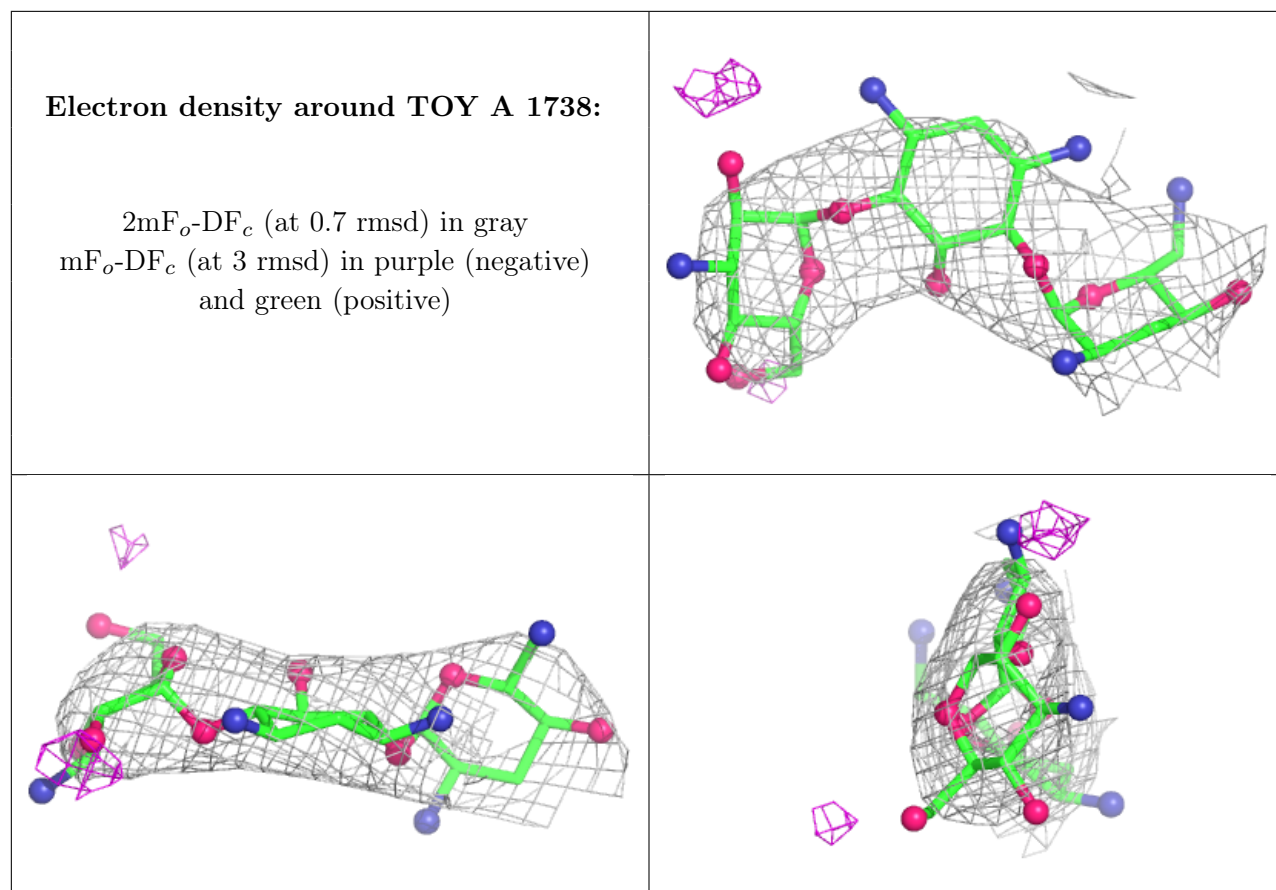
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
22	MG	A	1677	1/1	0.92	0.11	66,66,66,66	0
22	MG	A	1695	1/1	0.92	0.11	98,98,98,98	0
22	MG	A	1679	1/1	0.92	0.21	49,49,49,49	0
22	MG	A	1698	1/1	0.92	0.06	119,119,119,119	0
22	MG	A	1668	1/1	0.92	0.17	101,101,101,101	0
22	MG	A	1666	1/1	0.93	0.38	82,82,82,82	0
22	MG	A	1671	1/1	0.93	0.30	71,71,71,71	0
22	MG	A	1621	1/1	0.93	0.23	41,41,41,41	0
22	MG	A	1735	1/1	0.93	0.10	324,324,324,324	0
22	MG	A	1669	1/1	0.93	0.27	78,78,78,78	0
22	MG	A	1657	1/1	0.94	0.47	64,64,64,64	0
22	MG	A	1719	1/1	0.94	0.36	74,74,74,74	0
22	MG	A	1627	1/1	0.94	0.09	80,80,80,80	0
22	MG	A	1721	1/1	0.94	0.32	83,83,83,83	0
22	MG	A	1691	1/1	0.94	0.28	81,81,81,81	0
22	MG	A	1672	1/1	0.94	0.41	77,77,77,77	0
22	MG	A	1659	1/1	0.94	0.26	78,78,78,78	0
22	MG	A	1648	1/1	0.94	0.18	67,67,67,67	0
22	MG	A	1631	1/1	0.94	0.23	74,74,74,74	0
22	MG	A	1641	1/1	0.94	0.30	54,54,54,54	0
22	MG	A	1607	1/1	0.94	0.22	70,70,70,70	0
22	MG	A	1612	1/1	0.94	0.36	67,67,67,67	0
22	MG	A	1646	1/1	0.94	0.32	71,71,71,71	0
22	MG	A	1733	1/1	0.94	0.06	267,267,267,267	0
22	MG	A	1622	1/1	0.95	0.35	55,55,55,55	0
22	MG	A	1630	1/1	0.95	0.24	55,55,55,55	0
22	MG	A	1667	1/1	0.95	0.42	70,70,70,70	0
22	MG	A	1625	1/1	0.95	0.27	52,52,52,52	0
22	MG	K	201	1/1	0.96	0.09	122,122,122,122	0
22	MG	A	1674	1/1	0.96	0.18	72,72,72,72	0
22	MG	A	1626	1/1	0.96	0.38	72,72,72,72	0
22	MG	A	1653	1/1	0.96	0.66	90,90,90,90	0
22	MG	A	1629	1/1	0.96	0.09	55,55,55,55	0
22	MG	A	1736	1/1	0.96	0.05	238,238,238,238	0
22	MG	A	1727	1/1	0.96	0.26	99,99,99,99	0
22	MG	A	1728	1/1	0.96	0.22	116,116,116,116	0
22	MG	A	1684	1/1	0.96	0.20	59,59,59,59	0
22	MG	A	1686	1/1	0.96	0.17	69,69,69,69	0
22	MG	A	1678	1/1	0.96	0.27	45,45,45,45	0
22	MG	A	1687	1/1	0.97	0.31	72,72,72,72	0
22	MG	J	201	1/1	0.97	0.11	53,53,53,53	0
22	MG	A	1680	1/1	0.97	0.22	39,39,39,39	0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
22	MG	A	1603	1/1	0.97	0.33	37,37,37,37	0
22	MG	A	1662	1/1	0.97	0.07	60,60,60,60	0
22	MG	A	1693	1/1	0.97	0.05	80,80,80,80	0
24	TOY	A	1738	32/32	0.97	0.08	88,105,127,131	0
25	ZN	D	301	1/1	0.97	0.29	149,149,149,149	0
22	MG	A	1654	1/1	0.98	0.04	64,64,64,64	0
22	MG	A	1628	1/1	0.98	0.16	2,2,2,2	0
22	MG	A	1696	1/1	0.98	0.06	68,68,68,68	0
22	MG	A	1642	1/1	0.98	0.10	52,52,52,52	0
22	MG	A	1685	1/1	0.98	0.15	67,67,67,67	0
22	MG	A	1737	1/1	0.98	0.06	205,205,205,205	0
22	MG	A	1649	1/1	0.98	0.08	84,84,84,84	0
22	MG	A	1605	1/1	0.98	0.13	9,9,9,9	0
22	MG	A	1644	1/1	0.98	0.10	67,67,67,67	0
22	MG	A	1660	1/1	0.98	0.16	95,95,95,95	0
22	MG	A	1617	1/1	0.98	0.25	46,46,46,46	0
22	MG	A	1640	1/1	0.98	0.35	68,68,68,68	0
22	MG	A	1689	1/1	0.99	0.09	129,129,129,129	0
22	MG	A	1690	1/1	0.99	0.06	108,108,108,108	0
22	MG	A	1624	1/1	0.99	0.10	59,59,59,59	0
22	MG	A	1665	1/1	0.99	0.10	55,55,55,55	0
22	MG	A	1618	1/1	0.99	0.35	48,48,48,48	0
22	MG	A	1604	1/1	0.99	0.36	66,66,66,66	0
22	MG	A	1601	1/1	0.99	0.12	80,80,80,80	0
25	ZN	N	101	1/1	1.00	0.05	151,151,151,151	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.