



wwPDB EM Validation Summary Report ⓘ

Jun 24, 2026 – 05:17 PM EDT

PDB ID : 7UVV / pdb_00007uvv
EMDB ID : EMD-26817
Title : A. baumannii ribosome-Streptothricin-F complex: 70S with P-site tRNA
Authors : Morgan, C.E.; Yu, E.W.
Deposited on : 2022-05-02
Resolution : 2.50 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

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:

EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

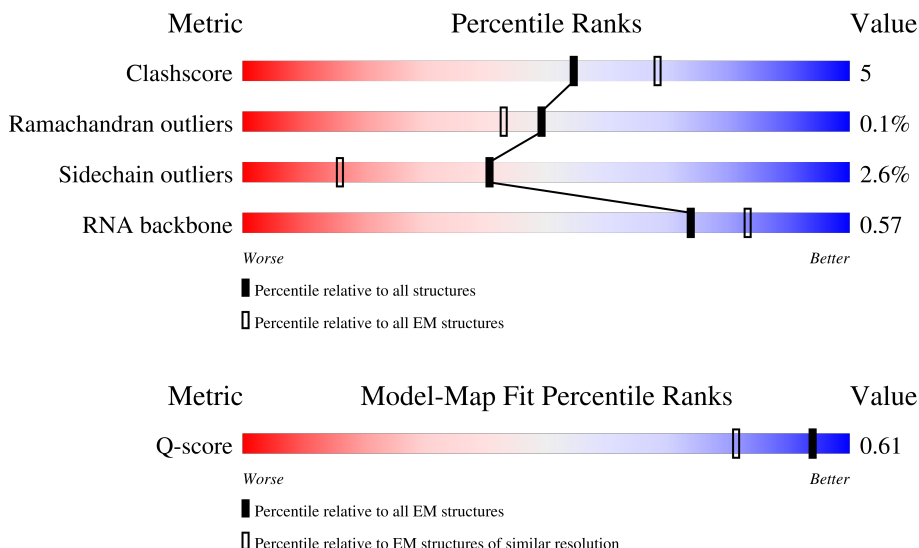
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.50 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	7115 (2.00 - 3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 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 EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	0	51	
2	1	44	
3	2	64	

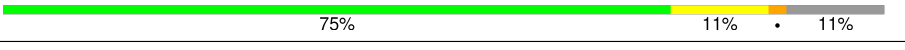
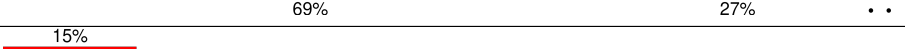
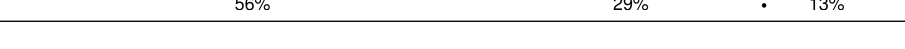
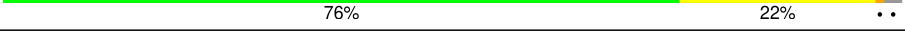
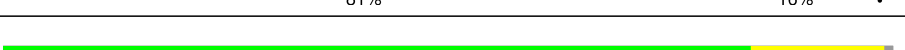
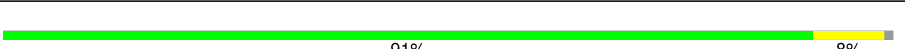
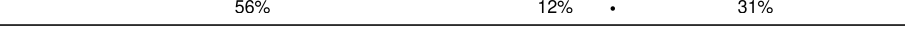
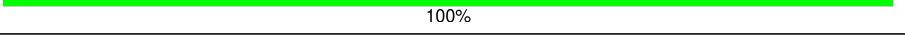
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	3	38	
5	A	2918	
6	B	115	
7	C	274	
8	D	212	
9	E	200	
10	F	178	
11	G	177	
12	H	148	
13	I	142	
14	J	122	
15	K	146	
16	L	137	
17	M	125	
18	N	116	
19	O	122	
20	P	119	
21	Q	103	
22	R	109	
23	S	106	
24	T	105	
25	U	98	
26	V	85	
27	W	78	
28	X	65	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
29	Y	58	
30	Z	61	
31	a	1544	
32	b	250	
33	c	250	
34	d	208	
35	e	165	
36	f	127	
37	g	156	
38	h	131	
39	i	128	
40	j	103	
41	k	128	
42	l	124	
43	m	118	
44	n	101	
45	o	89	
46	p	101	
47	q	85	
48	r	75	
49	s	91	
50	t	88	
51	u	71	
52	v	77	
53	w	3	

2 Entry composition

There are 55 unique types of molecules in this entry. The entry contains 138294 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 protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	0	49	Total	C	N	O	S	0	0
			409	263	74	70	2		

- Molecule 2 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	44	Total	C	N	O	S	0	0
			363	222	85	54	2		

- Molecule 3 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	63	Total	C	N	O	S	0	0
			509	319	110	76	4		

- Molecule 4 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	3	38	Total	C	N	O	S	0	0
			295	179	64	48	4		

- Molecule 5 is a RNA chain called 23s ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	A	2724	Total	C	N	O	P	0	0
			58438	26086	10700	18928	2724		

- Molecule 6 is a RNA chain called 5s ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	B	115	Total	C	N	O	P	0	0
			2450	1095	440	800	115		

- Molecule 7 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	C	272	Total	C	N	O	S	0	0
			2111	1302	436	365	8		

- Molecule 8 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	D	211	Total	C	N	O	S	0	0
			1572	972	297	300	3		

- Molecule 9 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	E	200	Total	C	N	O	S	0	0
			1516	952	281	278	5		

- Molecule 10 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	F	174	Total	C	N	O	S	0	0
			1370	871	243	248	8		

- Molecule 11 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	G	174	Total	C	N	O	S	0	0
			1318	832	236	249	1		

- Molecule 12 is a protein called 50S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	H	45	Total	C	N	O	S	0	0
			334	212	61	60	1		

- Molecule 13 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	I	142	Total	C	N	O	S	0	0
			1125	718	200	203	4		

- Molecule 14 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	J	122	Total	C	N	O	S	0	0
			946	592	180	169	5		

- Molecule 15 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	K	144	Total	C	N	O	S	0	0
			1071	663	213	195			

- Molecule 16 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	L	137	Total	C	N	O	S	0	0
			1087	687	210	185	5		

- Molecule 17 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	M	119	Total	C	N	O	S	0	0
			942	590	186	163	3		

- Molecule 18 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	N	114	Total	C	N	O	S	0	0
			857	528	173	155	1		

- Molecule 19 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	O	117	Total	C	N	O	S	0	0
			919	578	177	164			

- Molecule 20 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	P	117	Total	C	N	O	S	0	0
			934	589	197	146	2		

- Molecule 21 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	Q	103	Total	C	N	O	S	0	0
			807	506	155	143	3		

- Molecule 22 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	R	109	Total	C	N	O	S	0	0
			826	514	158	150	4		

- Molecule 23 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	S	91	Total	C	N	O	S	0	0
			710	452	128	129	1		

- Molecule 24 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	T	103	Total	C	N	O		0	0
			766	476	142	148			

- Molecule 25 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	U	97	Total	C	N	O	S	0	0
			760	477	143	139	1		

- Molecule 26 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	V	83	Total	C	N	O	S	0	0
			620	384	120	114	2		

- Molecule 27 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	W	77	Total	C	N	O	S	0	0
			632	395	130	105	2		

- Molecule 28 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	X	60	Total	C	N	O	S	0	0
			486	302	93	90	1		

- Molecule 29 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	Y	57	Total	C	N	O	S	0	0
			455	281	87	84	3		

- Molecule 30 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	Z	54	Total	C	N	O	S	0	0
			447	265	100	81	1		

- Molecule 31 is a RNA chain called 16s Ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	a	1528	Total	C	N	O	P	0	0
			32782	14631	5994	10630	1527		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
a	1007	U	C	conflict	GB 1211343212
a	1034	C	U	conflict	GB 1211343212

- Molecule 32 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	b	225	Total	C	N	O	S	0	0
			1769	1110	328	325	6		

- Molecule 33 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	c	215	Total	C	N	O	S	0	0
			1690	1065	318	299	8		

- Molecule 34 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	d	207	Total	C	N	O	S	0	0
			1631	1017	313	299	2		

- Molecule 35 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	e	155	Total	C	N	O	S	0	0
			1129	700	217	207	5		

- Molecule 36 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	f	104	Total	C	N	O	S	0	0
			867	546	158	159	4		

- Molecule 37 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	g	136	Total	C	N	O	S	0	0
			1073	671	201	194	7		

- Molecule 38 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	h	130	Total	C	N	O	S	0	0
			985	615	177	187	6		

- Molecule 39 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	i	126	Total	C	N	O	S	0	0
			991	618	197	175	1		

- Molecule 40 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	j	100	Total	C	N	O	S	0	0
			801	500	150	148	3		

- Molecule 41 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	k	117	Total	C	N	O	S	0	0
			862	535	167	159	1		

- Molecule 42 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	l	122	Total	C	N	O	S	0	0
			945	580	193	167	5		

- Molecule 43 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	m	115	Total	C	N	O	S	0	0
			903	558	184	158	3		

- Molecule 44 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	n	100	Total	C	N	O	S	0	0
			792	493	158	137	4		

- Molecule 45 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	o	88	Total	C	N	O	S	0	0
			705	434	144	126	1		

- Molecule 46 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	p	82	Total	C	N	O	S	0	0
			644	403	128	112	1		

- Molecule 47 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	q	79	Total	C	N	O	S	0	0
			621	390	116	114	1		

- Molecule 48 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				AltConf	Trace
48	r	52	Total	C	N	O	0	0
			426	273	74	79		

- Molecule 49 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	s	82	Total	C	N	O	S	0	0
			646	412	125	107	2		

- Molecule 50 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	t	86	Total	C	N	O	S	0	0
			663	409	139	113	2		

- Molecule 51 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	u	63	Total	C	N	O	S	0	0
			522	327	105	89	1		

- Molecule 52 is a RNA chain called tRNA-met.

Mol	Chain	Residues	Atoms						AltConf	Trace
52	v	77	Total	C	N	O	P	S	0	0
			1636	733	291	535	76	1		

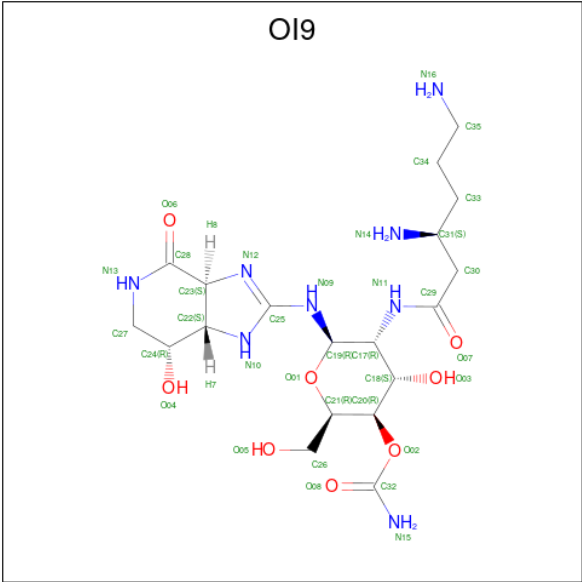
- Molecule 53 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	w	3	Total	C	N	O	P	0	0
			65	29	12	21	3		

- Molecule 54 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
54	3	1	Total	Zn	0
			1	1	

- Molecule 55 is Streptothricin F (CCD ID: OI9) (formula: C₁₉H₃₄N₈O₈) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
55	A	1	Total	C	N	O	0
			35	19	8	8	
55	a	1	Total	C	N	O	0
			35	19	8	8	

3 Residue-property plots [i](#)

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 atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: 50S ribosomal protein L33

Chain 0: 



- Molecule 2: 50S ribosomal protein L34

Chain 1: 

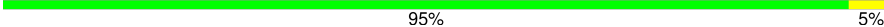


- Molecule 3: 50S ribosomal protein L35

Chain 2: 



- Molecule 4: 50S ribosomal protein L36

Chain 3: 

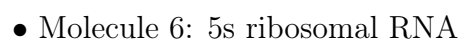


- Molecule 5: 23s ribosomal RNA

Chain A: 



A1915	U1792	A1596	U1502	U1415	G1245	U	G1014	G	A728	U596	U418	U283	
C1916	C1793	U1597	C1503	G1416	A1248	A	G1014	G	G740	A597	C419	U284	A147
G1917	C1796	C1598	U1504	A1422	A1505	U	G1023	U	U741	G598	G423	A285	A148
A1918	A1797	A1506	A1505	C1423	G1251	A	G1023	C	U741	A601	G424	G287	G149
C1920	A1798	A1607	U1506	C1423	G1261	A	U1030	U	U745	U605	C435	U288	U157
A1923	G1807	A1616	G1510	G1427	G1266	U	A1036	C	G773	A606	C435	G289	U158
A1924		U1632	A1511	A1429	A1267	U	A1037	C	G774	A607	C455	U290	U159
G1925	C1812	U1633	G1512	G1430	A1267	C	G1038	G	U774	U611	C455	G291	C160
G1926		A1634	U1513	G1431		C	G1039	C	A760	A612	C474	U292	A161
U1927	A1825	G1634	C1515	C1432	A1296	U	C1040	A	A760	A612	C474	U293	A162
A1928		A1635	A1518	A1435	A1311	U	A1043	U	G782	G618	A478	G297	A170
G1929	C1828	A1639	U1519	A1435	A1312	U	A1043	U	G783	G618	A478	G297	C171
A1932	C1829	G1640	U1520	A1437	U1312	A	G1044	A	A790	A625	A479	C305	U177
A1933		U1645	G1521	C1438	U1313	C	A1047	C	G803	A630	A481	U306	G177
A1934	G1835	U1646	G1522	C1439	C1314	C	A1047	C	G803	A630	A482	U307	U178
U1935	G1836	G1647	A1523	C1441	C1315	A	G1048	A	G803	A631	G493	U308	U179
				C1442	A1316	A	G1049	A	G803	A631	G493	U309	A182
U1940	U1840	G1665	C1526	G1445	A1317		U1050	A899	C810	A635	U498	A313	A183
U1951	G1841	G1665	U1527	G1446	U1324		A1051	C900	U811	A635	U498	A313	A183
U1959	G1842	G1672	G1528	G1447	G1333		G	G905	C812	A636	A499	A331	A188
A1962	A1844		U1529	U1447	G1333		G	G905	C812	U637	G499	A332	A189
C1963	G1845	G1680	A1530	U1448	C1344		G	A908	A817	C638	U503	A332	C190
A1966	U1681	U1681	C1531	U1455	C1344		G	A909	U825	U643	U503	A504	C190
U1967	A1682	A1682	U1532	C1456	C1345		U	C929	U826	U643	U507	A504	A198
G1968	C1683	C1683	U1533	G1462	U1347		U	C929	U826	U643	U507	A504	G199
				A1462	U1348		G	A938	U830	U650	C508	U336	A203
G1708	G1853	G1708	G1534	A1463	A1348		G	A938	A831	U651	A527	G345	A204
U1709	A1854	U1709	U1535	A1464	A1349		U	A942	G832	U652	A528	G346	C205
				A1465			C	G943	U837	A653	G529	A347	C206
G1983			G1536	U1471	C1352		U	A944	C838	U654	G529	A348	A206
U1987	G1865	U1718	C1537	A1472	G1353		A	A946	U837	U655	G531	G351	C215
G1988	U1866	U1719	A1538	A1472	C1357		C	U952	U837	U656	G532	G352	C216
A1989	C1726	G1727	C1540	A1472	C1357		G	U952	G843	U656	U533	G353	C216
U1989	G1727	G1727	G1541	U1477			A	U952	A844	A666	U544	A354	G222
G2008	C1870	C1731	A1542	G1478	A1360		C	G953	A845	G667	U544	A354	A223
A2009	A1871	C1732	A1543	U1479	G1363		C	U953	U846	G667	C545	A364	A228
A2010	U1733	U1733	G1544	A1480			A	C958	A847	A668	G546	U365	A229
A2011		G1734	G1558	U1481	A1373		C	U958	C848	A675	A554	A367	A229
	U1879	A1735		C1482	U1374		C	G966	G854	U684	A554	A367	U235
A2017	G1880	G1736	A1564	U1483	A1378		C	U967	G855	G857	A561	A370	U236
U2018	A1885			U1484	A1381		C	G971	G856	A667	A561	G371	U236
U2018	A1886	G1742	A1567	A1485	U1382		C	G971	G857	U688	G568	G371	A240
C2019		G1749	A1568	C1488			U	A980	G857	U688	U571	G385	U246
G2020	G1901		A1570	U1488	U1382		U	A980	G857	U688	U571	G385	U246
U2022	G1902	G1752		A1489	U1382		U	G990	G857	U688	U571	G385	U246
				A1489	U1382		U	G990	G857	U688	U571	G385	U246
A2026	G1906	U1576	U1576	A1490	U1400		U	A993	G857	U688	U571	G385	U246
A2027	U1907	A1577	A1577	A1491	U1401		U	A993	G857	U688	U571	G385	U246
G2028	A1908	U1578	U1492	U1492	C1402		U	A997	G857	U688	U571	G385	U246
A2029	A1909	A1769	G1579	C1493	G1403		A	A998	G857	U688	U571	G385	U246
	C1910		C1580		G1404		G	A998	G857	U688	U571	G385	U246
G2034	3TD1911	U1581	U1581	G1496	G1411		A	A1006	G857	U688	U571	G385	U246
U2035	U1912	U1582	C1583	G1497	C1412		A	A1006	G857	U688	U571	G385	U246
	U1913	C1583	U1499	G1498	G1413		C	A1009	G857	U688	U571	G385	U246
C2039	A1914	A1584	A1500	U1499	A1414		G	U1009	G857	U688	U571	G385	U246
		C1791		A1500			G	U1009	G857	U688	U571	G385	U246
				C1501			G	U1009	G857	U688	U571	G385	U246



G1	G22	G26	A27	G28	U38	C39	C40	G42	A43	A44	C45	U46	C47	A51	G52	U53	G54	A55	C59	U60	C61	A64	G65	C66	G67	A71	U86	A87	U104	C105	A106	C111	A112	C113	C114	U115
----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	------	------	------	------	------	------	------	------

- Molecule 7: 50S ribosomal protein L2

- Molecule 8: 50S ribosomal protein L3

MET
A2
I3
G4
L5
V6
C10
V30
R34
Q50
G54
S59
L79
V80
K81
E82
A88
E91
I99
D111
S142
Q151
Q176
S188
V189
I205
A212

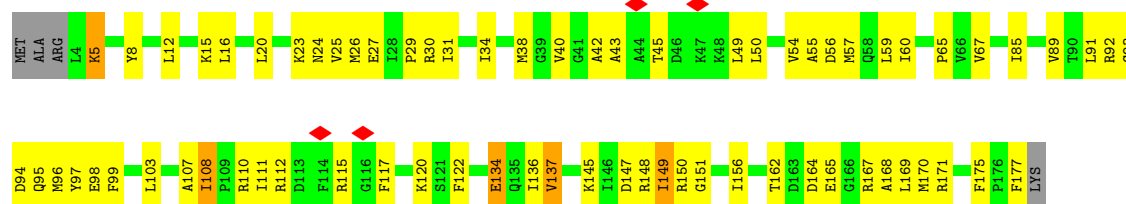
- Molecule 9: 50S ribosomal protein L4

Chain E:  91% 9%



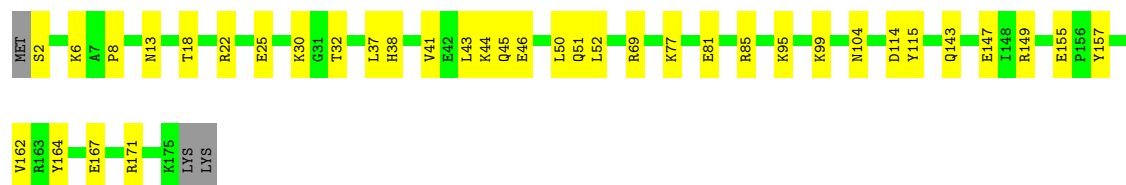
- Molecule 10: 50S ribosomal protein L5

Chain F:  58% 37%



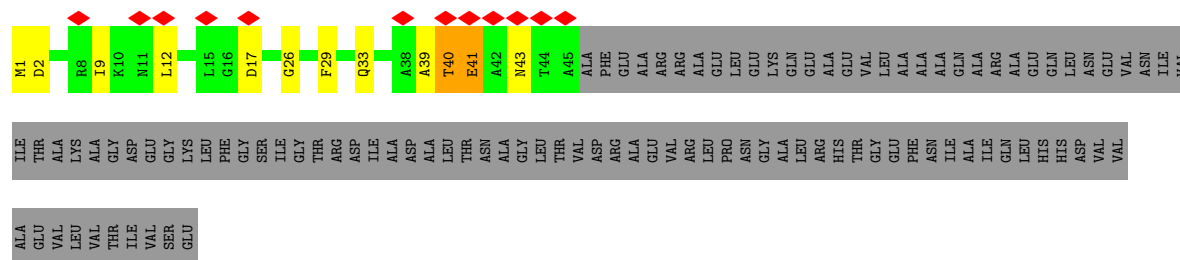
- Molecule 11: 50S ribosomal protein L6

Chain G:  77% 21%



- Molecule 12: 50S ribosomal protein L9

Chain H:  8% 22% 7% 70%



- Molecule 13: 50S ribosomal protein L13

Chain I:  89% 11%



- Molecule 14: 50S ribosomal protein L14

Chain J:  81% 17%



- Molecule 15: 50S ribosomal protein L15

Chain K: 92% 7%



- Molecule 16: 50S ribosomal protein L16

Chain L: 87% 13%



- Molecule 17: 50S ribosomal protein L17

Chain M: 90% 5% 5%



- Molecule 18: 50S ribosomal protein L18

Chain N: 79% 17% 4%



- Molecule 19: 50S ribosomal protein L19

Chain O: 87% 9% 4%



- Molecule 20: 50S ribosomal protein L20

Chain P: 90% 8% 2%



- Molecule 21: 50S ribosomal protein L21

Chain Q: 88% 12%



- Molecule 22: 50S ribosomal protein L22

Chain R: 94% 6%



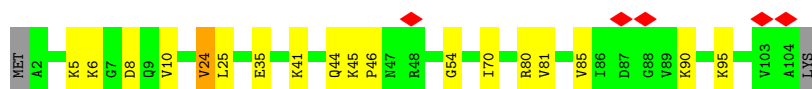
- Molecule 23: 50S ribosomal protein L23

Chain S: 77% 8% 14%



- Molecule 24: 50S ribosomal protein L24

Chain T: 5% 81% 16% ..



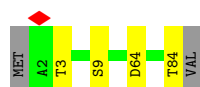
- Molecule 25: 50S ribosomal protein L25

Chain U: 88% 11% .



- Molecule 26: 50S ribosomal protein L27

Chain V: 93% 5% .



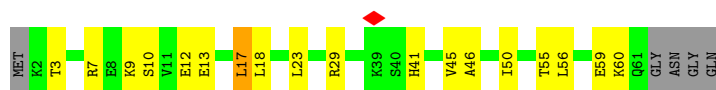
- Molecule 27: 50S ribosomal protein L28

Chain W: 90% 9% .



- Molecule 28: 50S ribosomal protein L29

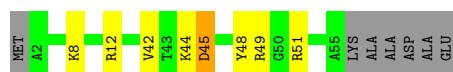
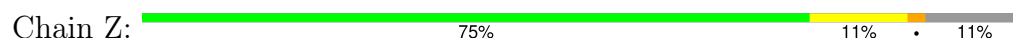
Chain X: 65% 26% 8%



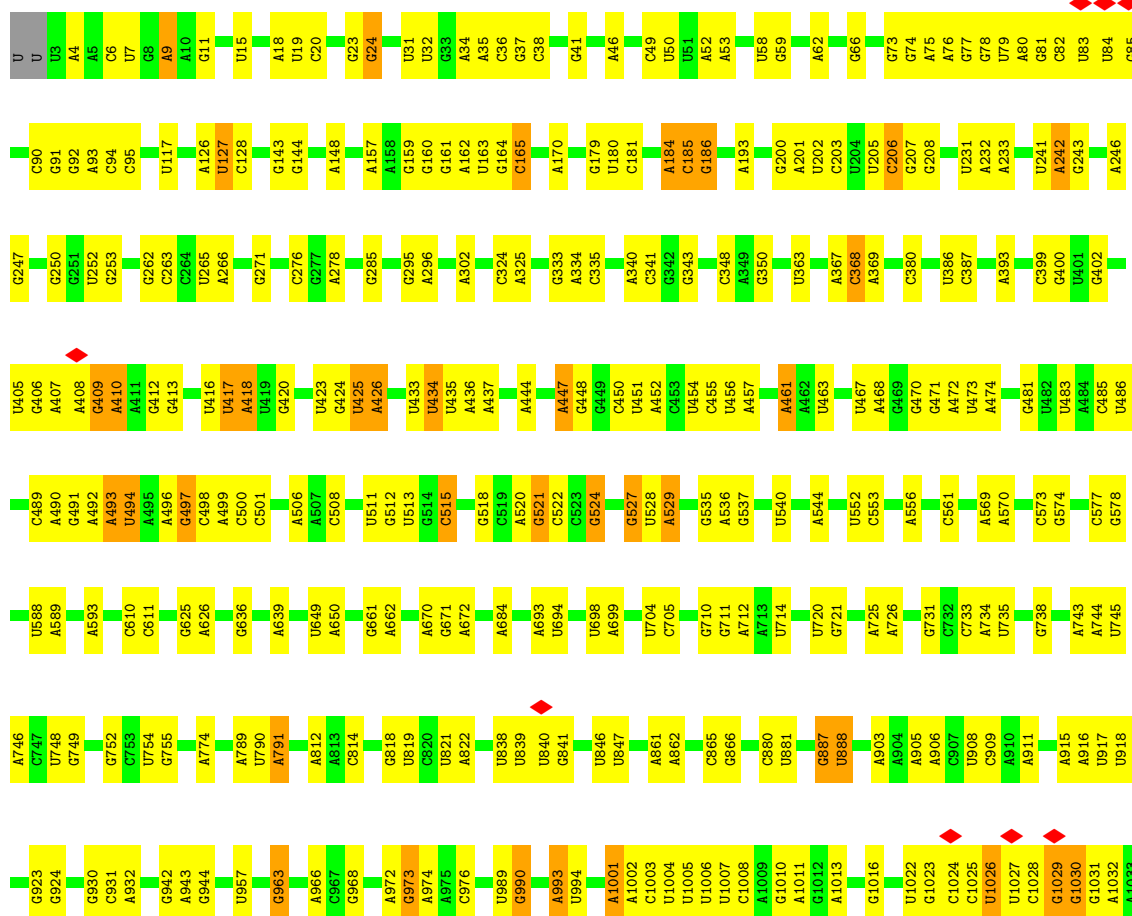
- Molecule 29: 50S ribosomal protein L30

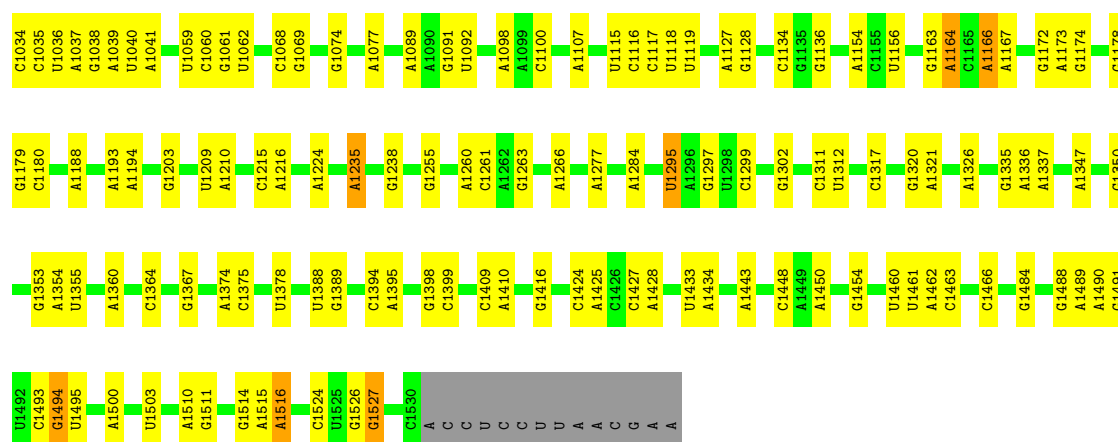


- Molecule 30: 50S ribosomal protein L32

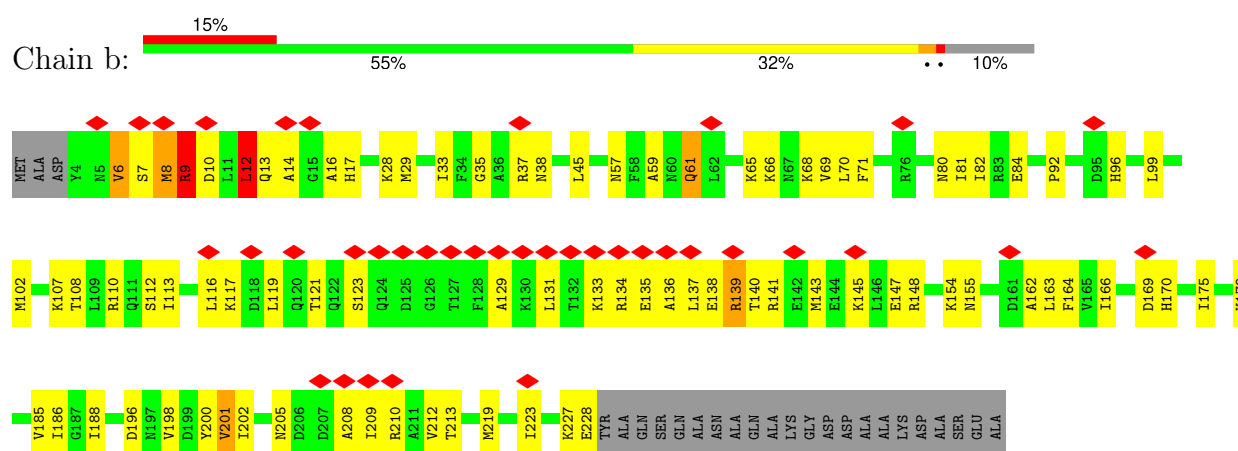


- Molecule 31: 16s Ribosomal RNA

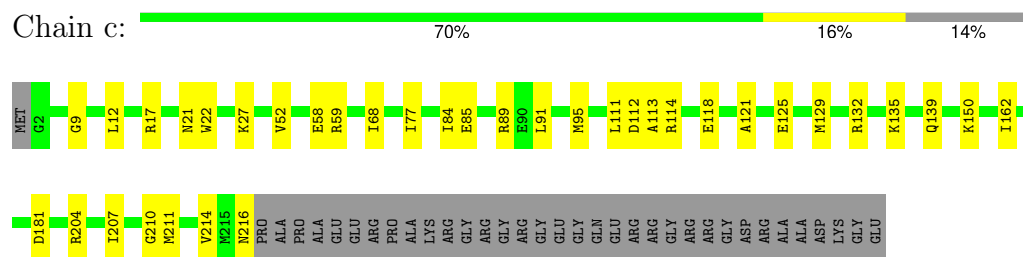




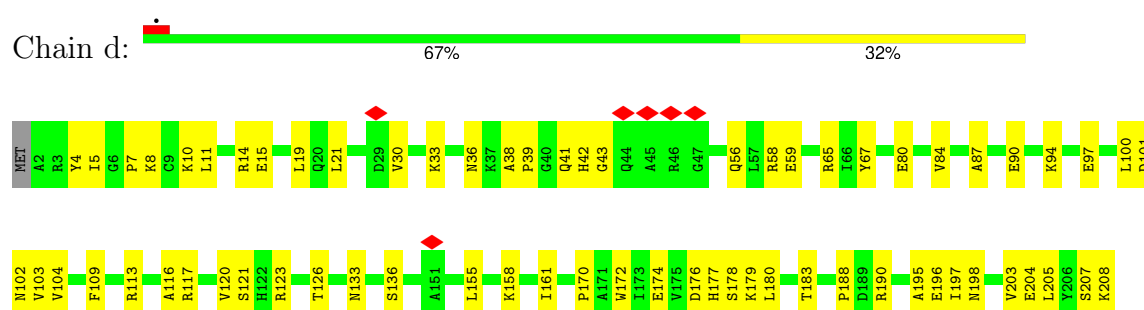
• Molecule 32: 30S ribosomal protein S2



• Molecule 33: 30S ribosomal protein S3



• Molecule 34: 30S ribosomal protein S4



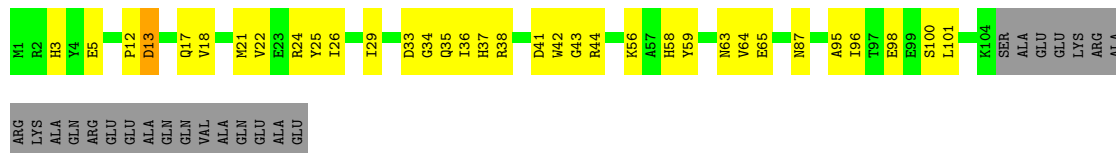
- Molecule 35: 30S ribosomal protein S5

Chain e: 



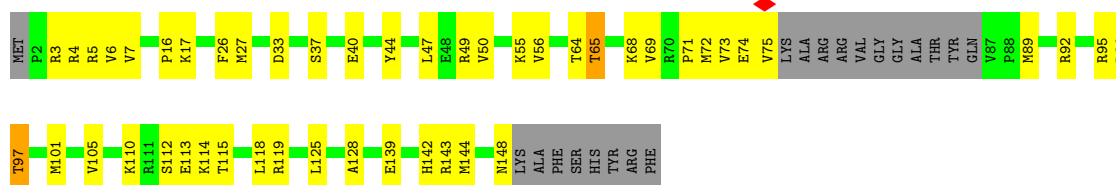
- Molecule 36: 30S ribosomal protein S6

Chain f: 



- Molecule 37: 30S ribosomal protein S7

Chain g: 



- Molecule 38: 30S ribosomal protein S8

Chain h: 



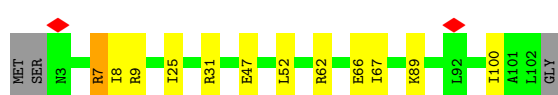
- Molecule 39: 30S ribosomal protein S9

Chain i: 



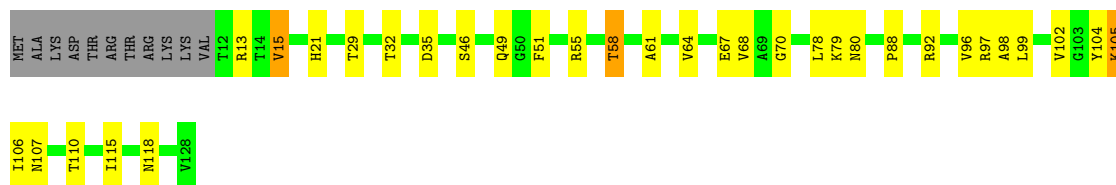
- Molecule 40: 30S ribosomal protein S10

Chain j: 



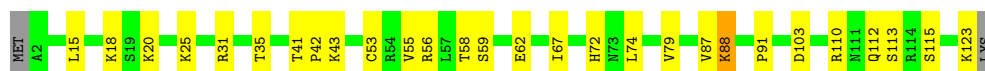
- Molecule 41: 30S ribosomal protein S11

Chain k:  66% 23% 9%



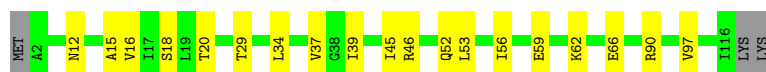
- Molecule 42: 30S ribosomal protein S12

Chain l:  76% 22% ..



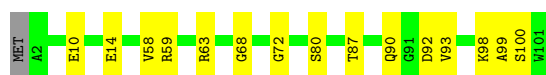
- Molecule 43: 30S ribosomal protein S13

Chain m:  81% 16% .



- Molecule 44: 30S ribosomal protein S14

Chain n:  84% 15% .



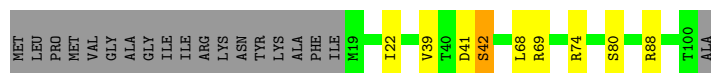
- Molecule 45: 30S ribosomal protein S15

Chain o:  91% 8% .



- Molecule 46: 30S ribosomal protein S16

Chain p:  72% 8% 19%



- Molecule 47: 30S ribosomal protein S17

Chain q:  76% 15% 7%



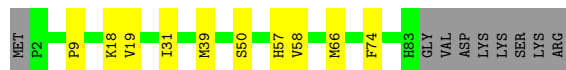
- Molecule 48: 30S ribosomal protein S18

Chain r: 



- Molecule 49: 30S ribosomal protein S19

Chain s: 



- Molecule 50: 30S ribosomal protein S20

Chain t: 



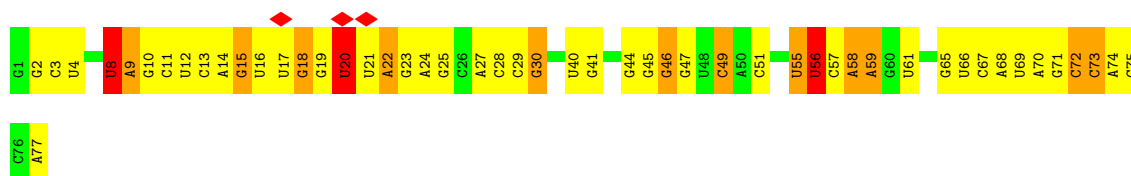
- Molecule 51: 30S ribosomal protein S21

Chain u: 



- Molecule 52: tRNA-met

Chain v: 



- Molecule 53: mRNA

Chain w: 

There are no outlier residues recorded for this chain.

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	53114	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.334	Depositor
Minimum map value	-0.044	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.035	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	434.176, 434.176, 434.176	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.848, 0.848, 0.848	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3TD, 6MZ, MA6, 5MC, OMU, 5MU, 2MG, 4OC, OMG, 7MG, 4SU, PSU, 2MA, UR3, OI9, H2U, 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	0	0.19	0/416	0.44	0/552
2	1	0.18	0/367	0.38	0/481
3	2	0.17	0/515	0.42	0/678
4	3	0.19	0/296	0.36	0/389
5	A	0.20	0/65086	0.32	0/101513
6	B	0.16	0/2739	0.27	0/4266
7	C	0.20	0/2152	0.44	0/2891
8	D	0.19	0/1590	0.38	0/2142
9	E	0.19	0/1537	0.39	0/2073
10	F	0.18	0/1390	0.48	0/1863
11	G	0.17	0/1337	0.41	0/1807
12	H	0.14	0/336	0.44	0/450
13	I	0.19	0/1151	0.42	0/1551
14	J	0.20	0/956	0.40	0/1286
15	K	0.19	0/1079	0.44	0/1439
16	L	0.19	0/1104	0.40	0/1475
17	M	0.21	0/956	0.41	0/1282
18	N	0.17	0/865	0.41	0/1156
19	O	0.18	0/931	0.41	0/1249
20	P	0.21	0/947	0.36	0/1262
21	Q	0.18	0/818	0.37	0/1094
22	R	0.18	0/831	0.36	0/1113
23	S	0.18	0/716	0.37	0/957
24	T	0.14	0/770	0.35	0/1034
25	U	0.18	0/770	0.45	0/1036
26	V	0.20	0/628	0.42	0/839
27	W	0.17	0/642	0.36	0/856
28	X	0.15	0/487	0.36	0/646
29	Y	0.19	0/460	0.34	0/614
30	Z	0.18	0/453	0.36	0/604
31	a	0.18	0/36476	0.30	0/56895

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	b	0.27	0/1799	0.64	5/2429 (0.2%)
33	c	0.18	0/1714	0.39	0/2304
34	d	0.17	0/1653	0.44	0/2213
35	e	0.19	0/1141	0.38	0/1537
36	f	0.17	0/882	0.45	0/1189
37	g	0.19	0/1087	0.47	0/1456
38	h	0.18	0/993	0.37	0/1331
39	i	0.20	0/1002	0.46	0/1339
40	j	0.19	0/811	0.46	0/1096
41	k	0.18	0/878	0.41	0/1189
42	l	0.18	0/958	0.46	0/1284
43	m	0.18	0/913	0.42	0/1226
44	n	0.19	0/803	0.36	0/1071
45	o	0.18	0/715	0.31	0/958
46	p	0.17	0/655	0.38	0/879
47	q	0.15	0/628	0.38	0/847
48	r	0.17	0/432	0.36	0/583
49	s	0.18	0/664	0.36	0/897
50	t	0.20	0/669	0.35	0/892
51	u	0.13	0/528	0.32	0/697
52	v	0.18	0/1739	0.33	0/2709
53	w	0.14	0/72	0.19	0/110
All	All	0.19	0/149537	0.34	5/223729 (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
21	Q	0	1
32	b	0	2
All	All	0	3

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
32	b	8	MET	CA-CB-CG	7.32	128.73	114.10
32	b	9	ARG	CA-CB-CG	6.90	127.89	114.10
32	b	12	LEU	CA-C-N	-5.91	112.78	122.54
32	b	12	LEU	C-N-CA	-5.91	112.78	122.54

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
32	b	8	MET	CB-CG-SD	5.55	129.36	112.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
21	Q	51	ALA	Peptide
32	b	139	ARG	Sidechain
32	b	9	ARG	Sidechain

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	0	409	0	437	6	0
2	1	363	0	401	2	0
3	2	509	0	566	4	0
4	3	295	0	326	1	0
5	A	58438	0	29389	334	0
6	B	2450	0	1241	16	0
7	C	2111	0	2176	27	0
8	D	1572	0	1610	11	0
9	E	1516	0	1569	11	0
10	F	1370	0	1420	54	0
11	G	1318	0	1373	22	0
12	H	334	0	363	8	0
13	I	1125	0	1148	10	0
14	J	946	0	1007	19	0
15	K	1071	0	1141	8	0
16	L	1087	0	1162	10	0
17	M	942	0	987	4	0
18	N	857	0	899	14	0
19	O	919	0	973	10	0
20	P	934	0	997	7	0
21	Q	807	0	842	6	0
22	R	826	0	894	4	0
23	S	710	0	768	7	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
24	T	766	0	816	11	0
25	U	760	0	783	9	0
26	V	620	0	629	2	0
27	W	632	0	667	3	0
28	X	486	0	528	14	0
29	Y	455	0	476	3	0
30	Z	447	0	435	7	0
31	a	32782	0	16508	244	0
32	b	1769	0	1787	62	0
33	c	1690	0	1774	23	0
34	d	1631	0	1691	42	0
35	e	1129	0	1174	16	0
36	f	867	0	868	19	0
37	g	1073	0	1118	34	0
38	h	985	0	1047	10	0
39	i	991	0	1048	10	0
40	j	801	0	832	7	0
41	k	862	0	877	21	0
42	l	945	0	998	17	0
43	m	903	0	962	11	0
44	n	792	0	833	8	0
45	o	705	0	712	5	0
46	p	644	0	655	6	0
47	q	621	0	665	7	0
48	r	426	0	447	9	0
49	s	646	0	663	4	0
50	t	663	0	715	4	0
51	u	522	0	565	10	0
52	v	1636	0	835	31	0
53	w	65	0	33	0	0
54	3	1	0	0	0	0
55	A	35	0	0	0	0
55	a	35	0	0	0	0
All	All	138294	0	92830	1168	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 5.

The worst 5 of 1168 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)
31:a:74:G:H1	31:a:93:A:N6	1.56	1.03
32:b:135:GLU:HG3	32:b:139:ARG:HH12	1.22	1.01
52:v:27:A:H61	52:v:45:G:H1	1.01	0.93
5:A:1036:A:H61	5:A:1113:G:H1	1.04	0.92
32:b:135:GLU:HG3	32:b:139:ARG:NH1	1.90	0.86

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	
1	0	47/51 (92%)	47 (100%)	0	0	100	100
2	1	42/44 (96%)	42 (100%)	0	0	100	100
3	2	61/64 (95%)	58 (95%)	3 (5%)	0	100	100
4	3	36/38 (95%)	35 (97%)	1 (3%)	0	100	100
7	C	270/274 (98%)	262 (97%)	8 (3%)	0	100	100
8	D	209/212 (99%)	200 (96%)	9 (4%)	0	100	100
9	E	198/200 (99%)	195 (98%)	3 (2%)	0	100	100
10	F	172/178 (97%)	160 (93%)	9 (5%)	3 (2%)	7	13
11	G	172/177 (97%)	166 (96%)	5 (3%)	1 (1%)	21	38
12	H	43/148 (29%)	43 (100%)	0	0	100	100
13	I	140/142 (99%)	138 (99%)	2 (1%)	0	100	100
14	J	120/122 (98%)	120 (100%)	0	0	100	100
15	K	142/146 (97%)	137 (96%)	5 (4%)	0	100	100
16	L	135/137 (98%)	132 (98%)	2 (2%)	1 (1%)	18	34
17	M	117/125 (94%)	114 (97%)	3 (3%)	0	100	100
18	N	112/116 (97%)	111 (99%)	1 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
19	O	115/122 (94%)	113 (98%)	2 (2%)	0	100	100
20	P	115/119 (97%)	115 (100%)	0	0	100	100
21	Q	101/103 (98%)	99 (98%)	1 (1%)	1 (1%)	12	24
22	R	107/109 (98%)	105 (98%)	2 (2%)	0	100	100
23	S	89/106 (84%)	87 (98%)	2 (2%)	0	100	100
24	T	101/105 (96%)	98 (97%)	3 (3%)	0	100	100
25	U	95/98 (97%)	92 (97%)	3 (3%)	0	100	100
26	V	81/85 (95%)	79 (98%)	2 (2%)	0	100	100
27	W	75/78 (96%)	75 (100%)	0	0	100	100
28	X	58/65 (89%)	58 (100%)	0	0	100	100
29	Y	55/58 (95%)	55 (100%)	0	0	100	100
30	Z	52/61 (85%)	50 (96%)	2 (4%)	0	100	100
32	b	223/250 (89%)	207 (93%)	16 (7%)	0	100	100
33	c	213/250 (85%)	207 (97%)	6 (3%)	0	100	100
34	d	205/208 (99%)	200 (98%)	5 (2%)	0	100	100
35	e	153/165 (93%)	151 (99%)	2 (1%)	0	100	100
36	f	102/127 (80%)	99 (97%)	3 (3%)	0	100	100
37	g	132/156 (85%)	126 (96%)	5 (4%)	1 (1%)	16	31
38	h	128/131 (98%)	125 (98%)	3 (2%)	0	100	100
39	i	124/128 (97%)	120 (97%)	4 (3%)	0	100	100
40	j	98/103 (95%)	95 (97%)	3 (3%)	0	100	100
41	k	115/128 (90%)	109 (95%)	6 (5%)	0	100	100
42	l	120/124 (97%)	115 (96%)	5 (4%)	0	100	100
43	m	113/118 (96%)	110 (97%)	2 (2%)	1 (1%)	14	27
44	n	98/101 (97%)	94 (96%)	4 (4%)	0	100	100
45	o	86/89 (97%)	84 (98%)	2 (2%)	0	100	100
46	p	80/101 (79%)	79 (99%)	1 (1%)	0	100	100
47	q	77/85 (91%)	76 (99%)	1 (1%)	0	100	100
48	r	50/75 (67%)	48 (96%)	2 (4%)	0	100	100
49	s	80/91 (88%)	78 (98%)	2 (2%)	0	100	100
50	t	84/88 (96%)	84 (100%)	0	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
51	u	61/71 (86%)	59 (97%)	2 (3%)	0	100	100
All	All	5402/5872 (92%)	5252 (97%)	142 (3%)	8 (0%)	49	68

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
16	L	79	LEU
10	F	137	VAL
11	G	13	ASN
43	m	66	GLU
10	F	5	LYS

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	
1	0	45/47 (96%)	43 (96%)	2 (4%)	25	50
2	1	36/36 (100%)	36 (100%)	0	100	100
3	2	52/53 (98%)	51 (98%)	1 (2%)	50	76
4	3	33/33 (100%)	33 (100%)	0	100	100
7	C	218/220 (99%)	213 (98%)	5 (2%)	44	72
8	D	166/167 (99%)	161 (97%)	5 (3%)	36	64
9	E	155/155 (100%)	151 (97%)	4 (3%)	40	68
10	F	144/147 (98%)	138 (96%)	6 (4%)	26	52
11	G	139/142 (98%)	138 (99%)	1 (1%)	76	89
12	H	34/112 (30%)	31 (91%)	3 (9%)	9	20
13	I	118/118 (100%)	117 (99%)	1 (1%)	73	88
14	J	103/103 (100%)	99 (96%)	4 (4%)	28	55
15	K	106/108 (98%)	105 (99%)	1 (1%)	70	87
16	L	113/113 (100%)	111 (98%)	2 (2%)	51	77
17	M	96/101 (95%)	95 (99%)	1 (1%)	68	86

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
18	N	83/85 (98%)	78 (94%)	5 (6%)	17	36
19	O	99/102 (97%)	98 (99%)	1 (1%)	68	86
20	P	85/86 (99%)	85 (100%)	0	100	100
21	Q	84/84 (100%)	82 (98%)	2 (2%)	43	70
22	R	88/88 (100%)	88 (100%)	0	100	100
23	S	77/87 (88%)	75 (97%)	2 (3%)	40	68
24	T	83/85 (98%)	80 (96%)	3 (4%)	31	58
25	U	79/80 (99%)	78 (99%)	1 (1%)	61	82
26	V	62/64 (97%)	61 (98%)	1 (2%)	55	79
27	W	69/70 (99%)	67 (97%)	2 (3%)	37	65
28	X	53/56 (95%)	51 (96%)	2 (4%)	29	56
29	Y	53/54 (98%)	53 (100%)	0	100	100
30	Z	46/50 (92%)	45 (98%)	1 (2%)	45	73
32	b	185/200 (92%)	177 (96%)	8 (4%)	26	51
33	c	175/198 (88%)	173 (99%)	2 (1%)	65	84
34	d	170/171 (99%)	163 (96%)	7 (4%)	27	53
35	e	113/120 (94%)	111 (98%)	2 (2%)	51	77
36	f	94/111 (85%)	91 (97%)	3 (3%)	34	62
37	g	113/128 (88%)	107 (95%)	6 (5%)	20	42
38	h	108/109 (99%)	106 (98%)	2 (2%)	50	76
39	i	99/100 (99%)	98 (99%)	1 (1%)	68	86
40	j	89/91 (98%)	87 (98%)	2 (2%)	45	73
41	k	88/98 (90%)	83 (94%)	5 (6%)	18	39
42	l	104/106 (98%)	100 (96%)	4 (4%)	29	56
43	m	95/98 (97%)	92 (97%)	3 (3%)	34	62
44	n	81/82 (99%)	78 (96%)	3 (4%)	30	57
45	o	71/72 (99%)	71 (100%)	0	100	100
46	p	63/77 (82%)	61 (97%)	2 (3%)	34	62
47	q	71/76 (93%)	67 (94%)	4 (6%)	19	39
48	r	46/66 (70%)	45 (98%)	1 (2%)	45	73
49	s	70/78 (90%)	68 (97%)	2 (3%)	37	65

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
50	t	65/67 (97%)	63 (97%)	2 (3%)	35 62
51	u	54/62 (87%)	53 (98%)	1 (2%)	50 76
All	All	4473/4756 (94%)	4357 (97%)	116 (3%)	41 68

5 of 116 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
32	b	12	LEU
48	r	24	LYS
34	d	204	GLU
47	q	69	SER
43	m	34	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
40	j	58	ASN
43	m	76	ASN
18	N	35	HIS
13	I	47	HIS
48	r	31	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
31	a	1527/1544 (98%)	184 (12%)	0
5	A	2720/2918 (93%)	369 (13%)	18 (0%)
52	v	76/77 (98%)	28 (36%)	0
53	w	2/3 (66%)	0	0
6	B	114/115 (99%)	11 (9%)	1 (0%)
All	All	4439/4657 (95%)	592 (13%)	19 (0%)

5 of 592 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
5	A	10	U
5	A	41	U
5	A	53	G
5	A	56	A

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
5	A	67	G

5 of 19 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
5	A	1519	U
5	A	2502	U
6	B	86	U
5	A	2443	G
5	A	782	G

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

28 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	$\# Z > 2$
5	PSU	A	2576	5	18,21,22	1.12	2 (11%)	21,30,33	2.02	6 (28%)
5	PSU	A	2453	5	18,21,22	1.07	1 (5%)	21,30,33	2.03	6 (28%)
31	PSU	a	513	31	18,21,22	1.08	1 (5%)	21,30,33	1.92	5 (23%)
5	OMG	A	2247	5,52	23,26,27	0.53	0	32,38,41	0.47	0
5	2MG	A	2441	5	23,26,27	0.52	0	33,38,41	0.46	0
5	5MU	A	1935	5	19,22,23	0.52	0	27,32,35	0.49	0
52	4SU	v	8	52	18,21,22	4.44	8 (44%)	25,30,33	2.30	4 (16%)
5	PSU	A	1913	5	18,21,22	1.09	1 (5%)	21,30,33	1.87	4 (19%)
5	3TD	A	1911	5	19,22,23	4.39	6 (31%)	23,32,35	1.75	3 (13%)
31	MA6	a	1516	31	23,26,27	1.43	4 (17%)	33,38,41	3.29	12 (36%)
5	PSU	A	2500	5	18,21,22	1.06	1 (5%)	21,30,33	1.88	4 (19%)
31	7MG	a	524	31	23,26,27	1.08	1 (4%)	27,39,42	0.89	2 (7%)
5	PSU	A	2601	5	18,21,22	1.08	1 (5%)	21,30,33	1.97	3 (14%)
31	4OC	a	1399	31	20,23,24	3.12	8 (40%)	25,32,35	0.97	1 (4%)
5	7MG	A	2065	5	23,26,27	1.05	1 (4%)	27,39,42	0.93	2 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
52	PSU	v	56	52	18,21,22	1.12	1 (5%)	21,30,33	1.83	5 (23%)
31	MA6	a	1515	31	23,26,27	1.42	4 (17%)	33,38,41	3.22	12 (36%)
52	H2U	v	20	52	18,21,22	0.55	0	19,30,33	0.99	1 (5%)
31	2MG	a	1204	31	23,26,27	0.48	0	33,38,41	0.49	0
31	2MG	a	963	31	23,26,27	0.49	0	33,38,41	0.50	0
5	6MZ	A	2026	5	22,25,26	2.48	4 (18%)	29,36,39	3.47	11 (37%)
5	PSU	A	952	5	18,21,22	1.06	1 (5%)	21,30,33	1.97	4 (19%)
52	5MU	v	55	52	19,22,23	0.47	0	27,32,35	0.71	0
5	PSU	A	1907	5	18,21,22	1.09	1 (5%)	21,30,33	1.93	5 (23%)
5	OMU	A	2548	5	19,22,23	3.02	8 (42%)	25,31,34	1.87	5 (20%)
31	UR3	a	1495	31	19,22,23	2.88	7 (36%)	26,32,35	1.63	3 (11%)
5	2MA	A	2499	5	22,25,26	3.87	9 (40%)	32,37,40	2.59	11 (34%)
31	5MC	a	964	31	19,22,23	0.58	0	26,32,35	0.63	0

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
5	PSU	A	2576	5	-	0/7/25/26	0/2/2/2
5	PSU	A	2453	5	-	0/7/25/26	0/2/2/2
31	PSU	a	513	31	-	0/7/25/26	0/2/2/2
5	OMG	A	2247	5,52	-	1/9/27/28	0/3/3/3
5	2MG	A	2441	5	-	2/9/27/28	0/3/3/3
5	5MU	A	1935	5	-	0/7/25/26	0/2/2/2
52	4SU	v	8	52	-	2/7/25/26	0/2/2/2
5	PSU	A	1913	5	-	0/7/25/26	0/2/2/2
5	3TD	A	1911	5	-	2/7/25/26	0/2/2/2
31	MA6	a	1516	31	-	0/11/29/30	0/3/3/3
5	PSU	A	2500	5	-	1/7/25/26	0/2/2/2
31	7MG	a	524	31	-	3/7/37/38	0/3/3/3
5	PSU	A	2601	5	-	0/7/25/26	0/2/2/2
31	4OC	a	1399	31	-	2/9/29/30	0/2/2/2
5	7MG	A	2065	5	-	1/7/37/38	0/3/3/3
52	PSU	v	56	52	-	4/7/25/26	0/2/2/2
31	MA6	a	1515	31	-	0/11/29/30	0/3/3/3
52	H2U	v	20	52	-	1/7/38/39	0/2/2/2
31	2MG	a	1204	31	-	0/9/27/28	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
31	2MG	a	963	31	-	2/9/27/28	0/3/3/3
5	6MZ	A	2026	5	-	1/9/27/28	0/3/3/3
5	PSU	A	952	5	-	0/7/25/26	0/2/2/2
52	5MU	v	55	52	-	3/7/25/26	0/2/2/2
5	PSU	A	1907	5	-	0/7/25/26	0/2/2/2
5	OMU	A	2548	5	-	0/9/27/28	0/2/2/2
31	UR3	a	1495	31	-	0/7/25/26	0/2/2/2
5	2MA	A	2499	5	-	2/7/25/26	0/3/3/3
31	5MC	a	964	31	-	0/7/25/26	0/2/2/2

The worst 5 of 70 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1911	3TD	C6-C5	13.21	1.49	1.35
5	A	2499	2MA	C4-N3	11.70	1.49	1.34
5	A	1911	3TD	C2-N1	10.48	1.50	1.37
5	A	2026	6MZ	C6-N6	9.74	1.45	1.34
52	v	8	4SU	C2-N1	9.61	1.53	1.38

The worst 5 of 109 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
31	a	1516	MA6	N1-C6-N6	-11.86	102.40	116.86
31	a	1515	MA6	N1-C6-N6	-11.53	102.81	116.86
5	A	2026	6MZ	C4-N9-C1'	-9.89	103.49	126.63
5	A	2026	6MZ	C1'-N9-C8	9.82	148.88	127.09
31	a	1516	MA6	C5-C6-N6	8.03	138.03	125.33

There are no chirality outliers.

5 of 27 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
52	v	8	4SU	O4'-C4'-C5'-O5'
52	v	55	5MU	O4'-C4'-C5'-O5'
52	v	56	PSU	O4'-C1'-C5-C4
52	v	56	PSU	O4'-C1'-C5-C6
5	A	2247	OMG	C1'-C2'-O2'-CM2

There are no ring outliers.

10 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2247	OMG	2	0
5	A	1935	5MU	1	0
52	v	8	4SU	1	0
5	A	1913	PSU	1	0
5	A	1911	3TD	1	0
52	v	56	PSU	1	0
52	v	20	H2U	2	0
31	a	963	2MG	1	0
5	A	2026	6MZ	1	0
52	v	55	5MU	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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
55	OI9	a	1601	-	36,37,37	1.40	6 (16%)	37,52,52	1.60	4 (10%)
55	OI9	A	3001	-	36,37,37	1.40	6 (16%)	37,52,52	1.56	5 (13%)

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
55	OI9	a	1601	-	-	5/22/67/67	0/3/3/3
55	OI9	A	3001	-	-	6/22/67/67	0/3/3/3

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
55	A	3001	OI9	C28-N13	3.96	1.40	1.33
55	a	1601	OI9	C28-N13	3.82	1.40	1.33
55	a	1601	OI9	C24-C22	-3.68	1.50	1.53
55	A	3001	OI9	C27-C24	3.64	1.55	1.52
55	a	1601	OI9	C27-C24	3.42	1.55	1.52

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
55	A	3001	OI9	C24-C27-N13	6.67	116.63	109.83
55	a	1601	OI9	C24-C27-N13	6.34	116.29	109.83
55	a	1601	OI9	O06-C28-N13	4.03	127.87	122.41
55	A	3001	OI9	C30-C29-N11	-3.28	111.78	116.25
55	a	1601	OI9	C30-C29-N11	-3.15	111.95	116.25

There are no chirality outliers.

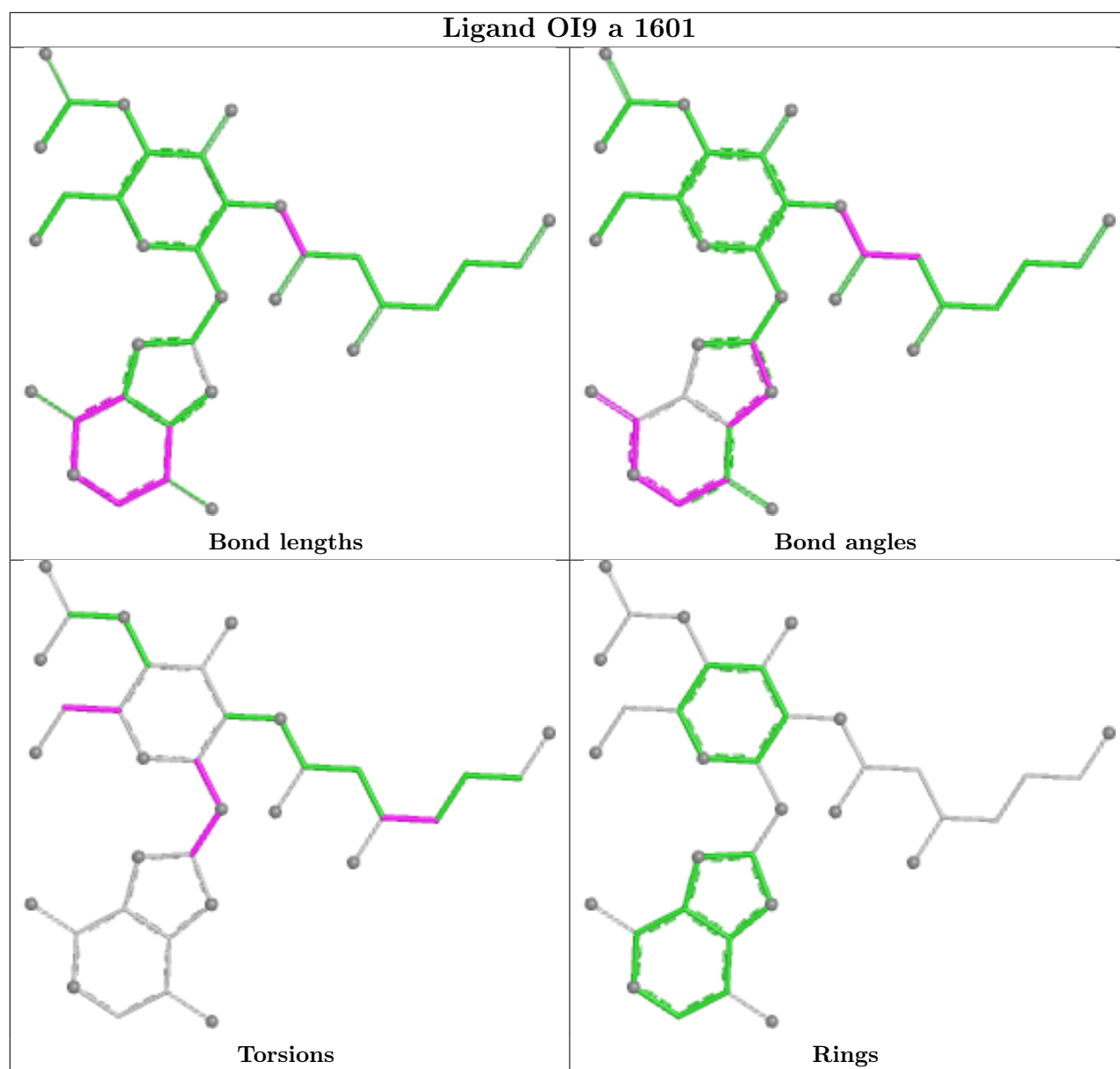
5 of 11 torsion outliers are listed below:

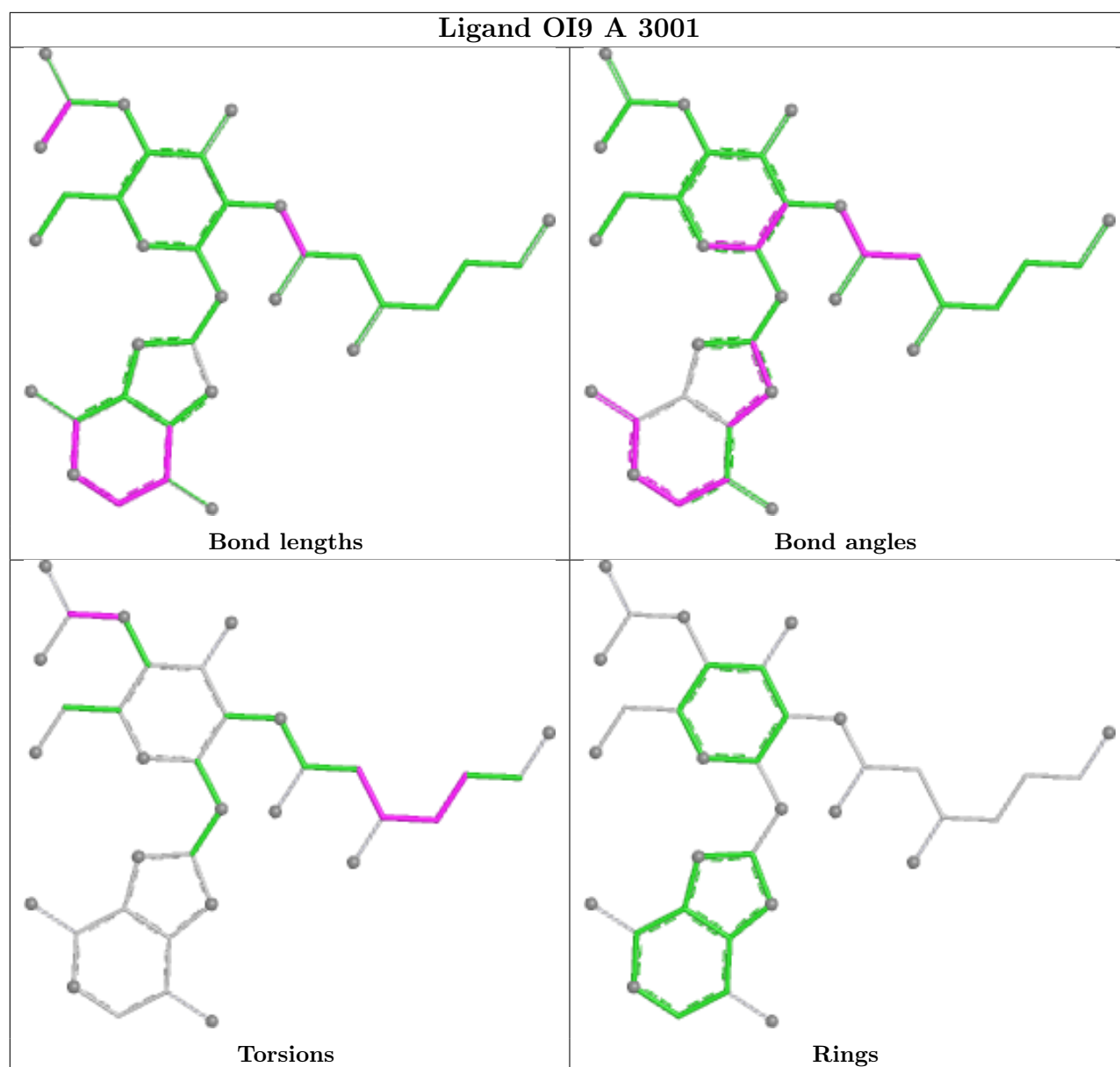
Mol	Chain	Res	Type	Atoms
55	A	3001	OI9	C30-C31-C33-C34
55	A	3001	OI9	N15-C32-O02-C20
55	A	3001	OI9	O08-C32-O02-C20
55	a	1601	OI9	O01-C21-C26-O05
55	a	1601	OI9	C20-C21-C26-O05

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.

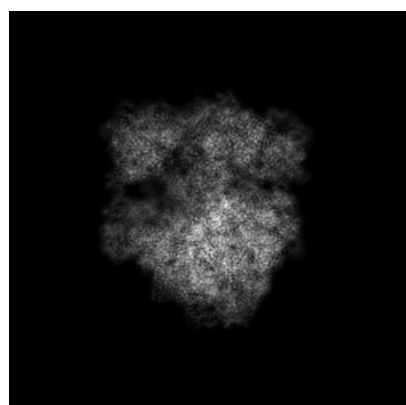
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26817. These allow visual inspection of the internal detail of the map and identification of artifacts.

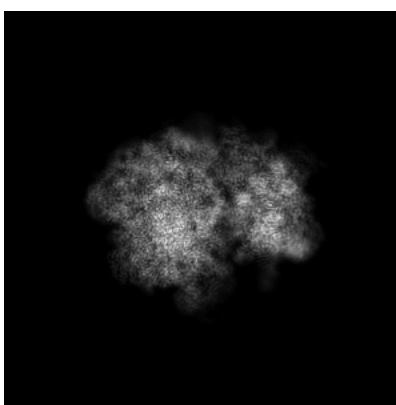
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

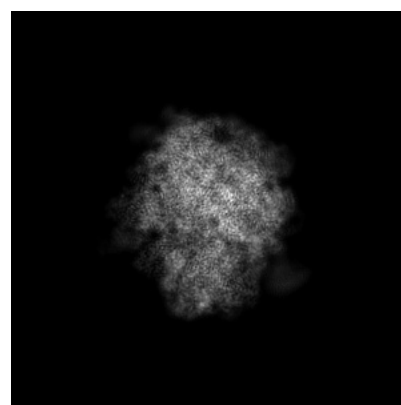
6.1.1 Primary map



X



Y



Z

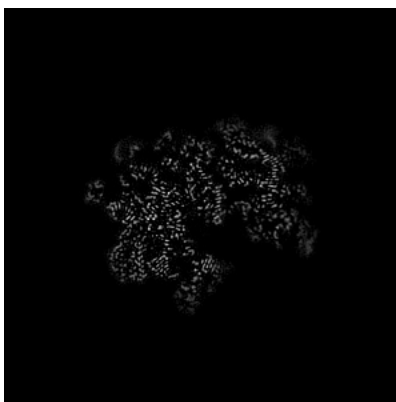
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

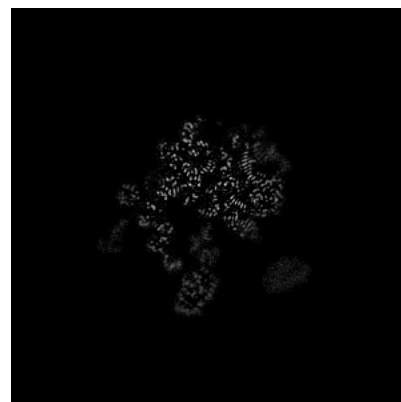
6.2.1 Primary map



X Index: 256



Y Index: 256

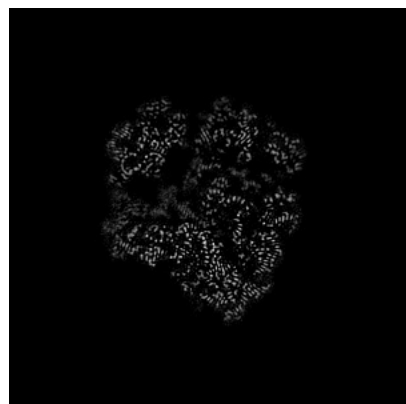


Z Index: 256

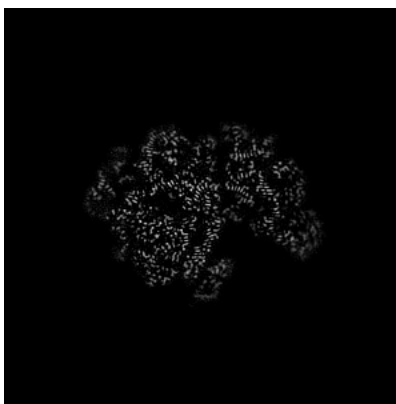
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

6.3.1 Primary map



X Index: 248



Y Index: 275

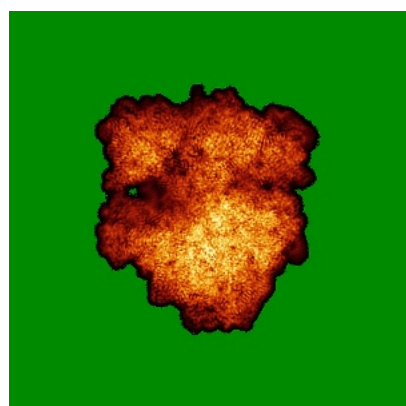


Z Index: 218

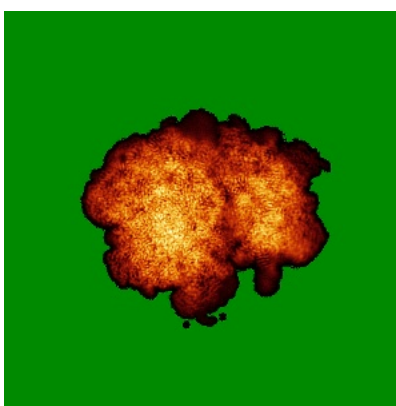
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

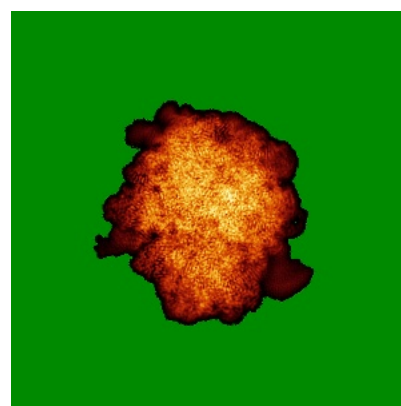
6.4.1 Primary map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

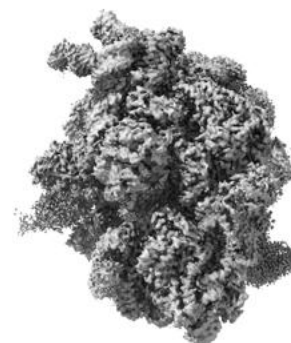
6.5.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

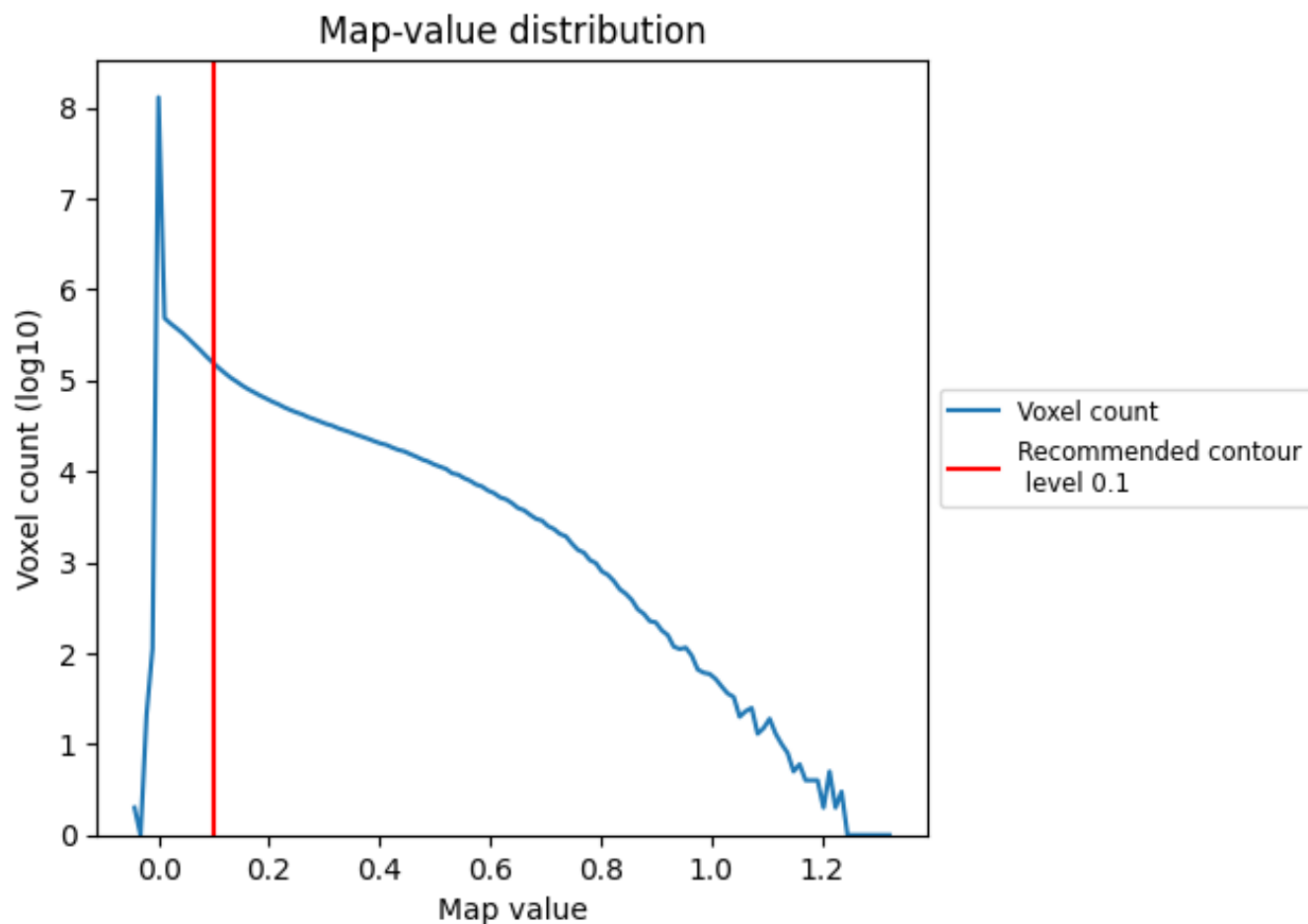
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

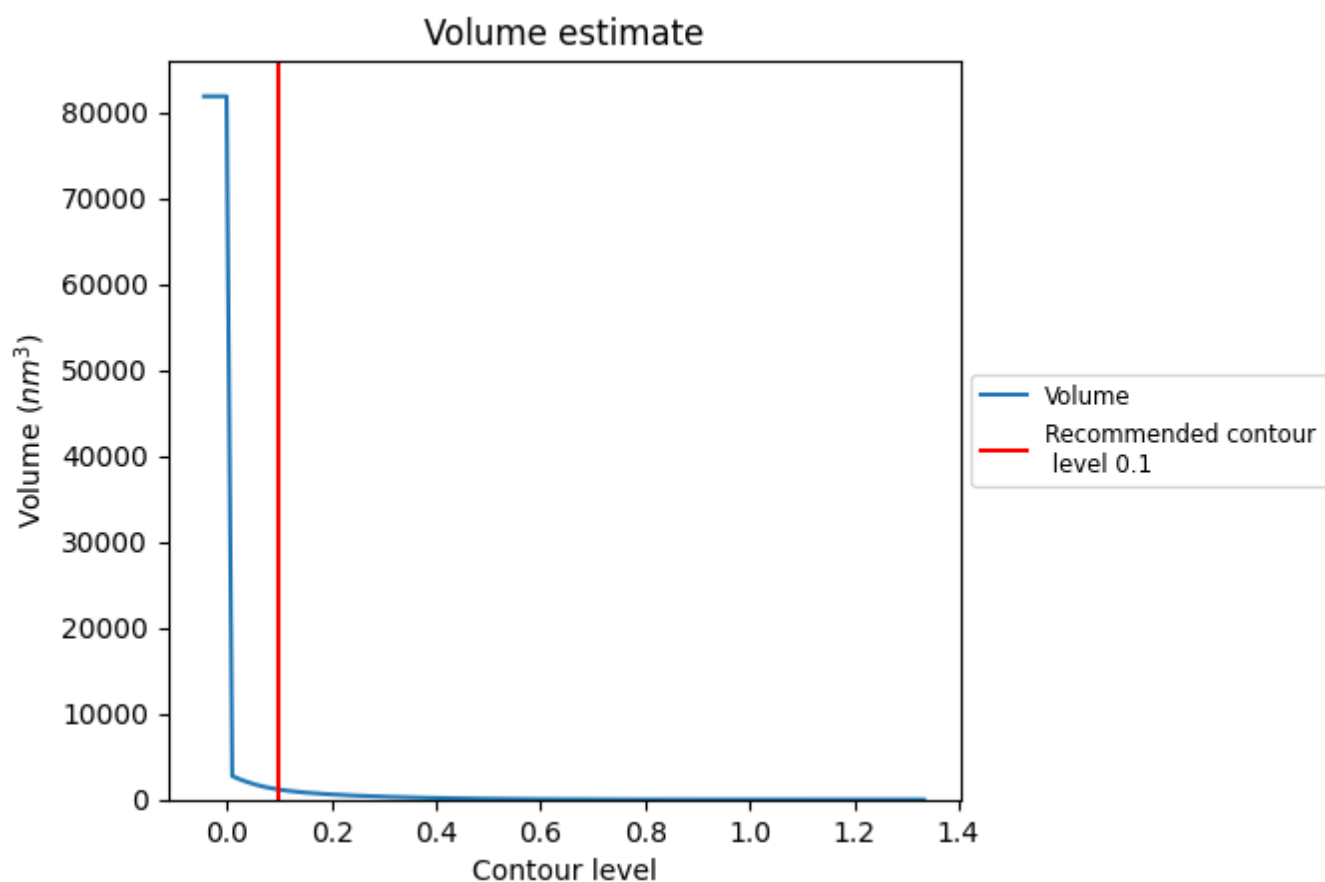
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

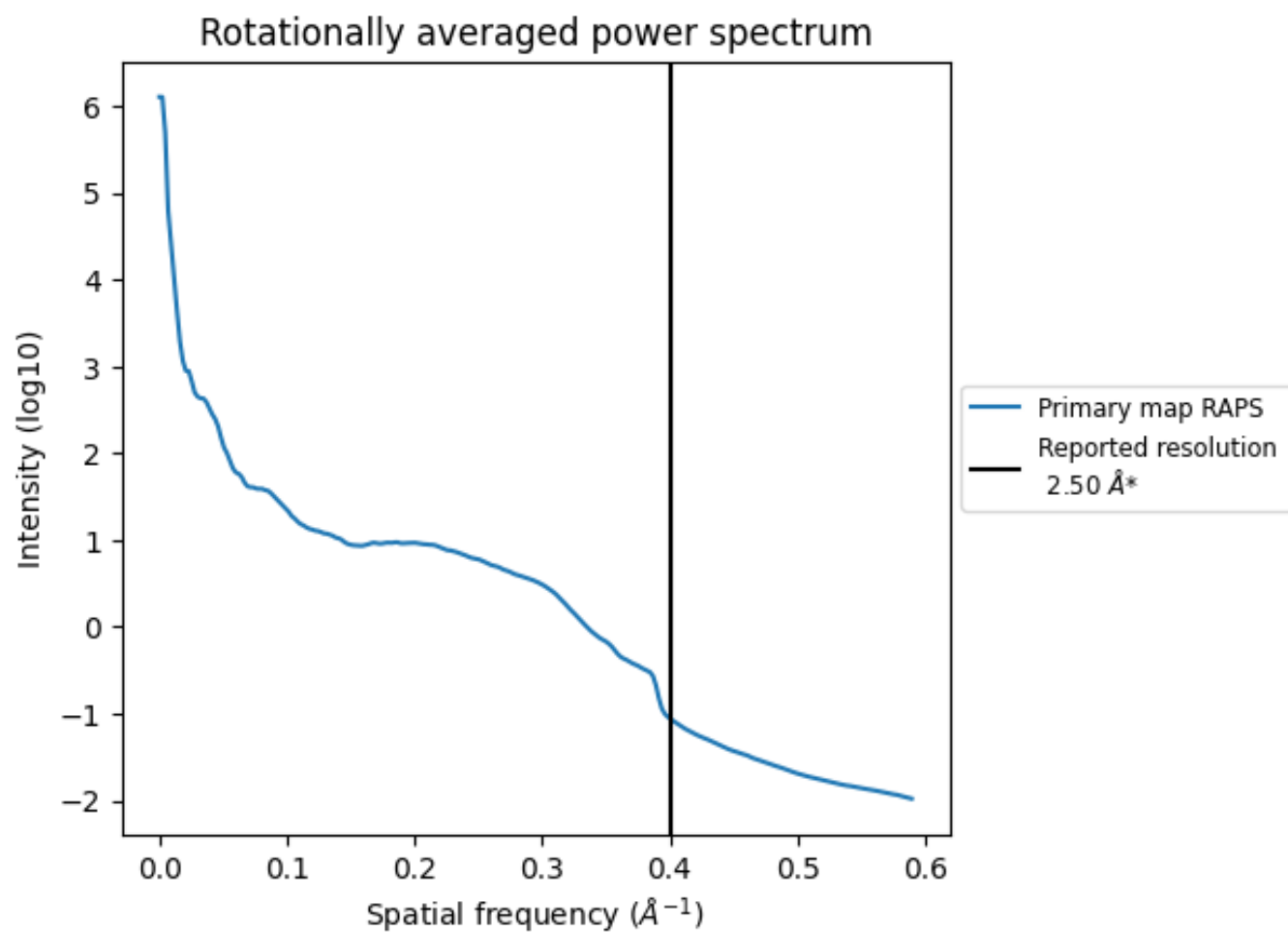
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1161 nm^3 ; this corresponds to an approximate mass of 1049 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.400 Å⁻¹

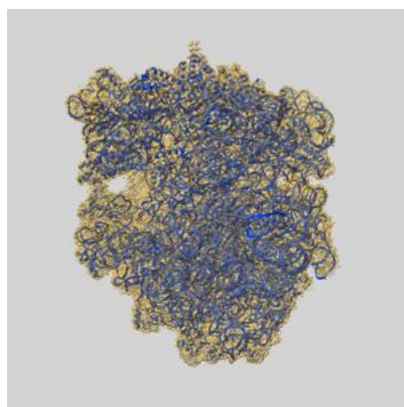
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

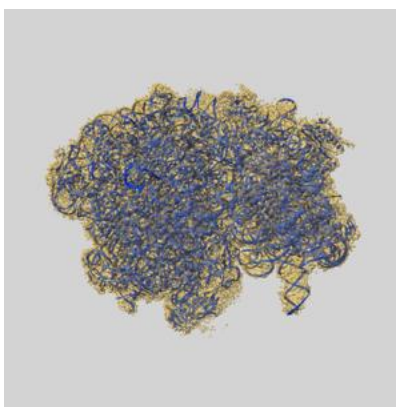
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26817 and PDB model 7UVV. Per-residue inclusion information can be found in section [3](#) on page [14](#).

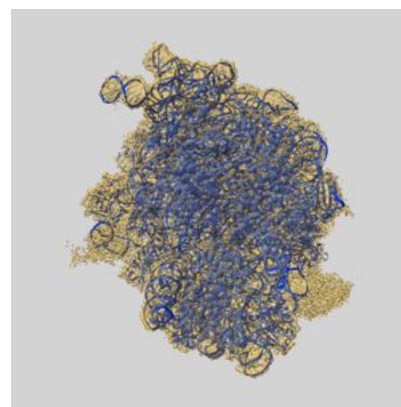
9.1 Map-model overlay [i](#)



X



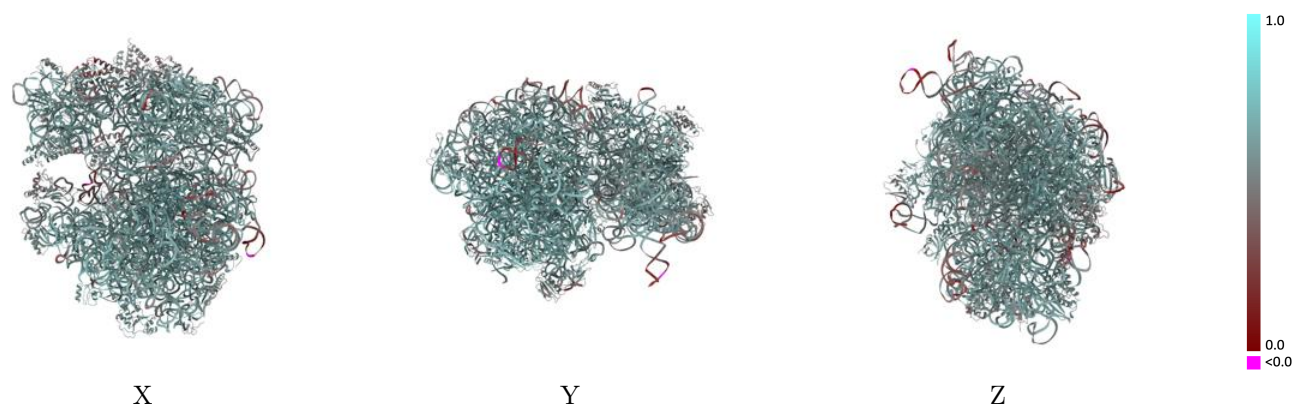
Y



Z

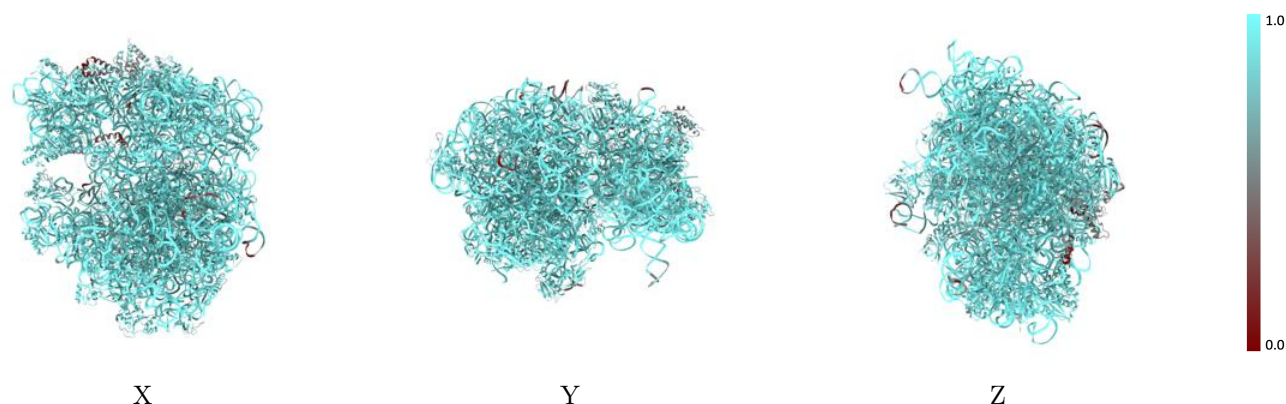
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



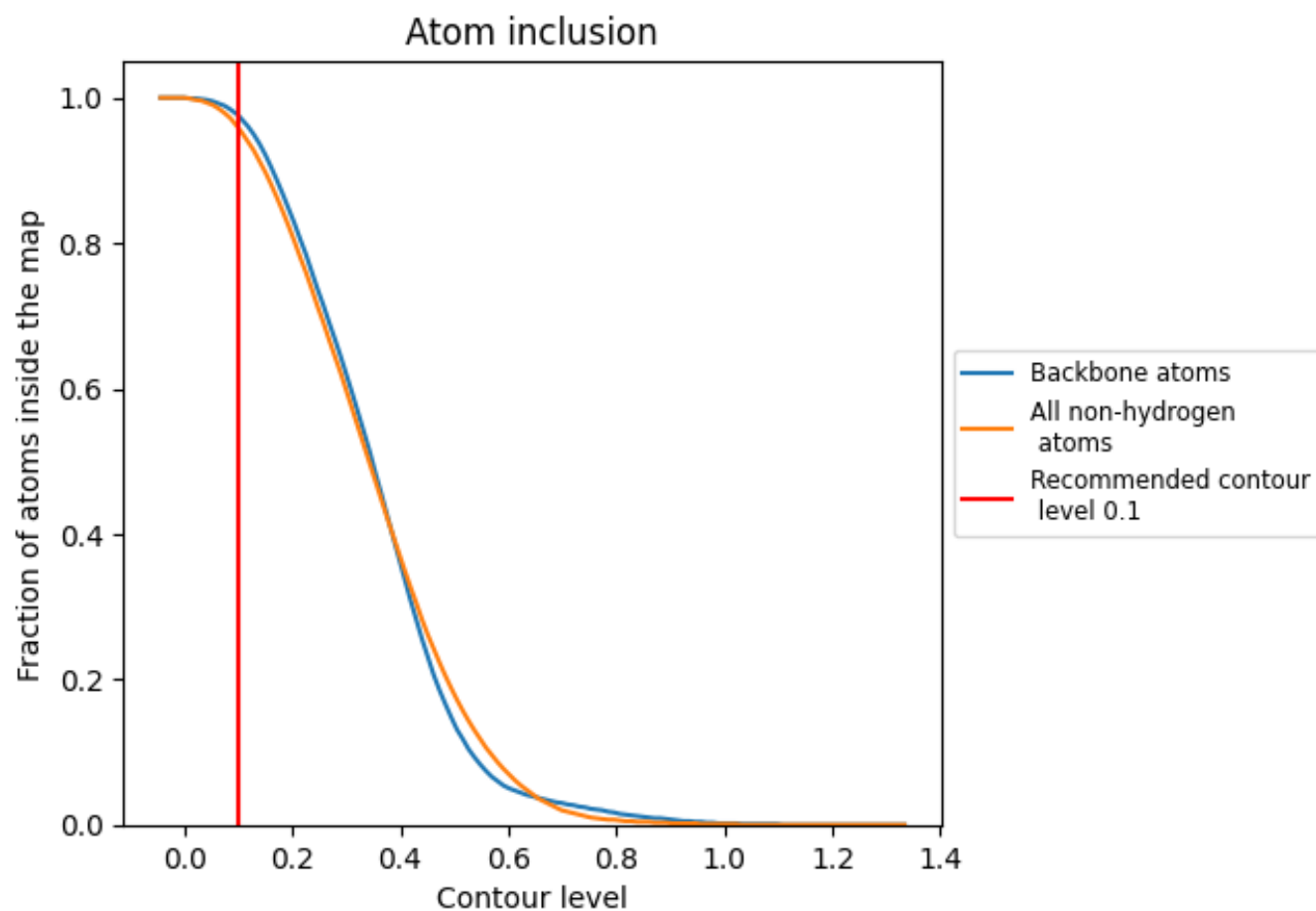
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.1).

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 96% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9590	 0.6100
0	 0.9420	 0.6300
1	 0.9800	 0.6820
2	 0.9880	 0.6850
3	 0.9830	 0.6510
A	 0.9800	 0.6310
B	 0.9940	 0.5840
C	 0.9660	 0.6530
D	 0.9690	 0.6670
E	 0.9430	 0.6340
F	 0.8420	 0.4770
G	 0.9080	 0.5420
H	 0.6560	 0.4650
I	 0.9730	 0.6590
J	 0.9270	 0.6370
K	 0.9780	 0.6550
L	 0.9580	 0.6500
M	 0.9790	 0.6680
N	 0.9630	 0.5890
O	 0.9370	 0.6360
P	 0.9960	 0.6770
Q	 0.9680	 0.6520
R	 0.9720	 0.6600
S	 0.9080	 0.6070
T	 0.8660	 0.5670
U	 0.9460	 0.6200
V	 0.9620	 0.6670
W	 0.9510	 0.6550
X	 0.8820	 0.5320
Y	 0.9690	 0.6600
Z	 0.9650	 0.6570
a	 0.9810	 0.5930
b	 0.6180	 0.4670
c	 0.9310	 0.5970
d	 0.8350	 0.4910



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
e	 0.9650	 0.6170
f	 0.8580	 0.5330
g	 0.8210	 0.5110
h	 0.9620	 0.6270
i	 0.9570	 0.6190
j	 0.9050	 0.5820
k	 0.9020	 0.5520
l	 0.9020	 0.5990
m	 0.9640	 0.6070
n	 0.9440	 0.6220
o	 0.9560	 0.6180
p	 0.9700	 0.6210
q	 0.9370	 0.5870
r	 0.9320	 0.5960
s	 0.9780	 0.6260
t	 0.9690	 0.6030
u	 0.4820	 0.4710
v	 0.8700	 0.4420
w	 1.0000	 0.6210