



## wwPDB EM Validation Summary Report ⓘ

Jul 13, 2026 – 04:29 PM EDT

PDB ID : 11DG / pdb\_000011dg  
EMDB ID : EMD-75634  
Title : Chimeric Escherichia coli 70S ribosome containing an evolved Vibrio cholerae 16S rRNA (VC-S4.4)  
Authors : Raskar, T.; Badran, A.; Fraser, J.; Costello, A.  
Deposited on : 2026-02-18  
Resolution : 3.00 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](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev133  
Mogul : 2022.3.0, CSD as543be (2022)  
MolProbity : 4-5-2 with Phenix2.0  
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) : 2.50

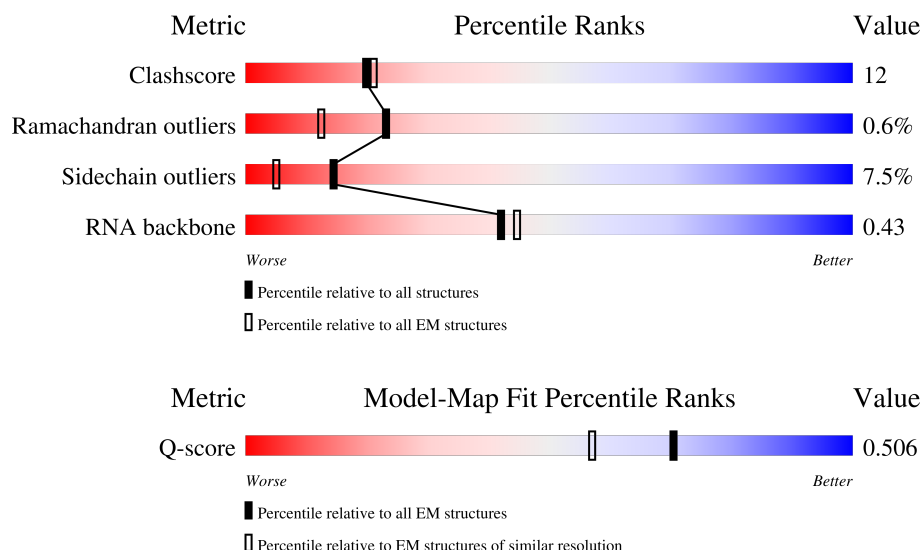
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.00 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<br>(#Entries) | EM structures<br>(#Entries) | Similar EM resolution<br>(#Entries, resolution range(Å)) |
|-----------------------|-----------------------------|-----------------------------|----------------------------------------------------------|
| Clashscore            | 229148                      | 23984                       | -                                                        |
| Ramachandran outliers | 224038                      | 23583                       | -                                                        |
| Sidechain outliers    | 223484                      | 23102                       | -                                                        |
| RNA backbone          | 8273                        | 3508                        | -                                                        |
| Q-score               | -                           | 25397                       | 14081 ( 2.50 - 3.50 )                                    |

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  $\geq 3$ , 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 EM map (all-atom inclusion  $< 40\%$ ). The numeric value is given above the bar.

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1   | 0     | 55     |                  |
| 2   | 1     | 46     |                  |
| 3   | 2     | 65     |                  |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 4   | 3     | 38     |                  |
| 5   | 4     | 70     |                  |
| 6   | 5     | 2      |                  |
| 7   | A     | 1517   |                  |
| 8   | B     | 241    |                  |
| 9   | C     | 233    |                  |
| 10  | D     | 206    |                  |
| 11  | E     | 167    |                  |
| 12  | F     | 135    |                  |
| 13  | G     | 179    |                  |
| 14  | H     | 130    |                  |
| 15  | I     | 130    |                  |
| 16  | J     | 103    |                  |
| 17  | K     | 129    |                  |
| 18  | L     | 124    |                  |
| 19  | M     | 118    |                  |
| 20  | N     | 101    |                  |
| 21  | O     | 89     |                  |
| 22  | P     | 82     |                  |
| 23  | Q     | 84     |                  |
| 24  | R     | 75     |                  |
| 25  | S     | 92     |                  |
| 26  | T     | 87     |                  |
| 27  | U     | 71     |                  |
| 28  | a     | 2904   |                  |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 29  | b     | 120    |                  |
| 30  | c     | 273    |                  |
| 31  | d     | 209    |                  |
| 32  | e     | 201    |                  |
| 33  | f     | 179    |                  |
| 34  | g     | 177    |                  |
| 35  | h     | 149    |                  |
| 36  | i     | 142    |                  |
| 37  | j     | 123    |                  |
| 38  | k     | 144    |                  |
| 39  | l     | 136    |                  |
| 40  | m     | 127    |                  |
| 41  | n     | 117    |                  |
| 42  | o     | 115    |                  |
| 43  | p     | 118    |                  |
| 44  | q     | 103    |                  |
| 45  | r     | 110    |                  |
| 46  | s     | 100    |                  |
| 47  | t     | 104    |                  |
| 48  | u     | 94     |                  |
| 49  | v     | 85     |                  |
| 50  | w     | 78     |                  |
| 51  | x     | 63     |                  |
| 52  | y     | 59     |                  |
| 53  | z     | 57     |                  |

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 |
|-----|------|-------|-----|-----------|----------|---------|------------------|
| 17  | BH2  | K     | 119 | X         | -        | -       | -                |

## 2 Entry composition

There are 57 unique types of molecules in this entry. The entry contains 138477 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     | 51       | Total | C   | N  | O  | 0       | 0     |
|     |       |          | 417   | 269 | 76 | 72 |         |       |

- Molecule 2 is a protein called Large ribosomal subunit protein bL34.

| Mol | Chain | Residues | Atoms |     |    |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 2   | 1     | 46       | Total | C   | N  | O  | S | 0       | 0     |
|     |       |          | 377   | 228 | 90 | 57 | 2 |         |       |

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

| Mol | Chain | Residues | Atoms |     |     |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---|---------|-------|
| 3   | 2     | 64       | Total | C   | N   | O  | S | 0       | 0     |
|     |       |          | 504   | 323 | 105 | 74 | 2 |         |       |

- Molecule 4 is a protein called Large ribosomal subunit protein bL36A.

| Mol | Chain | Residues | Atoms |     |    |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 4   | 3     | 38       | Total | C   | N  | O  | S | 0       | 0     |
|     |       |          | 302   | 185 | 65 | 48 | 4 |         |       |

- Molecule 5 is a protein called 50S ribosomal protein L31.

| Mol | Chain | Residues | Atoms |     |    |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 5   | 4     | 60       | Total | C   | N  | O  | S | 0       | 0     |
|     |       |          | 480   | 299 | 90 | 85 | 6 |         |       |

- Molecule 6 is a RNA chain called E-site tRNA.

| Mol | Chain | Residues | Atoms |    |   |    |   | AltConf | Trace |
|-----|-------|----------|-------|----|---|----|---|---------|-------|
| 6   | 5     | 2        | Total | C  | N | O  | P | 0       | 0     |
|     |       |          | 42    | 19 | 8 | 13 | 2 |         |       |

- Molecule 7 is a RNA chain called 16S ribosomal RNA.

| Mol | Chain | Residues | Atoms |       |      |       |      | AltConf | Trace |
|-----|-------|----------|-------|-------|------|-------|------|---------|-------|
| 7   | A     | 1517     | Total | C     | N    | O     | P    | 0       | 0     |
|     |       |          | 32564 | 14534 | 5976 | 10537 | 1517 |         |       |

- Molecule 8 is a protein called Small ribosomal subunit protein uS2.

| Mol | Chain | Residues | Atoms |      |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 8   | B     | 224      | Total | C    | N   | O   | S | 0       | 0     |
|     |       |          | 1753  | 1109 | 315 | 321 | 8 |         |       |

- Molecule 9 is a protein called 30S ribosomal protein S3.

| Mol | Chain | Residues | Atoms |      |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 9   | C     | 206      | Total | C    | N   | O   | S | 0       | 0     |
|     |       |          | 1624  | 1028 | 305 | 288 | 3 |         |       |

- Molecule 10 is a protein called Small ribosomal subunit protein uS4.

| Mol | Chain | Residues | Atoms |      |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 10  | D     | 205      | Total | C    | N   | O   | S | 0       | 0     |
|     |       |          | 1642  | 1026 | 315 | 297 | 4 |         |       |

- Molecule 11 is a protein called Small ribosomal subunit protein uS5.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 11  | E     | 156      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1152  | 717 | 217 | 212 | 6 |         |       |

- Molecule 12 is a protein called Small ribosomal subunit protein bS6.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 12  | F     | 103      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 839   | 530 | 151 | 151 | 7 |         |       |

- Molecule 13 is a protein called Small ribosomal subunit protein uS7.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 13  | G     | 153      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1203  | 750 | 231 | 218 | 4 |         |       |

- Molecule 14 is a protein called Small ribosomal subunit protein uS8.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 14  | H     | 129      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 978   | 616 | 173 | 183 | 6 |         |       |

- Molecule 15 is a protein called Small ribosomal subunit protein uS9.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 15  | I     | 127      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1022  | 634 | 206 | 179 | 3 |         |       |

- Molecule 16 is a protein called Small ribosomal subunit protein uS10.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 16  | J     | 98       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 786   | 493 | 150 | 142 | 1 |         |       |

- Molecule 17 is a protein called Small ribosomal subunit protein uS11.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 17  | K     | 117      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 876   | 540 | 173 | 160 | 3 |         |       |

There is a discrepancy between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment  | Reference  |
|-------|---------|----------|--------|----------|------------|
| K     | 119     | BH2      | ASN    | conflict | UNP A7ZSI6 |

- Molecule 18 is a protein called 30S ribosomal protein S12.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 18  | L     | 123      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 957   | 591 | 196 | 165 | 5 |         |       |

- Molecule 19 is a protein called Small ribosomal subunit protein uS13.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 19  | M     | 115      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 891   | 552 | 179 | 157 | 3 |         |       |

- Molecule 20 is a protein called Small ribosomal subunit protein uS14.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 20  | N     | 100      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 805   | 499 | 164 | 139 | 3 |         |       |

- Molecule 21 is a protein called Small ribosomal subunit protein uS15.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 21  | O     | 88       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 714   | 439 | 144 | 130 | 1 |         |       |

- Molecule 22 is a protein called Small ribosomal subunit protein bS16.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 22  | P     | 81       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 643   | 403 | 127 | 112 | 1 |         |       |

- Molecule 23 is a protein called Small ribosomal subunit protein uS17.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 23  | Q     | 79       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 641   | 406 | 120 | 112 | 3 |         |       |

- Molecule 24 is a protein called Small ribosomal subunit protein bS18.

| Mol | Chain | Residues | Atoms |     |     |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---|---------|-------|
| 24  | R     | 66       | Total | C   | N   | O  | S | 0       | 0     |
|     |       |          | 544   | 345 | 102 | 96 | 1 |         |       |

- Molecule 25 is a protein called Small ribosomal subunit protein uS19.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 25  | S     | 84       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 668   | 427 | 127 | 112 | 2 |         |       |

- Molecule 26 is a protein called Small ribosomal subunit protein bS20.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 26  | T     | 86       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 669   | 412 | 138 | 115 | 4 |         |       |

There is a discrepancy between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment  | Reference  |
|-------|---------|----------|--------|----------|------------|
| T     | 58      | CYS      | VAL    | conflict | UNP A7ZHB2 |

- Molecule 27 is a protein called Small ribosomal subunit protein bS21.

| Mol | Chain | Residues | Atoms |     |     |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---|---------|-------|
| 27  | U     | 70       | Total | C   | N   | O  | S | 0       | 0     |
|     |       |          | 589   | 366 | 125 | 97 | 1 |         |       |

- Molecule 28 is a RNA chain called 23S ribosomal RNA.

| Mol | Chain | Residues | Atoms |       |       |       |      | AltConf | Trace |
|-----|-------|----------|-------|-------|-------|-------|------|---------|-------|
| 28  | a     | 2753     | Total | C     | N     | O     | P    | 0       | 0     |
|     |       |          | 59130 | 26384 | 10897 | 19096 | 2753 |         |       |

- Molecule 29 is a RNA chain called 5S ribosomal RNA.

| Mol | Chain | Residues | Atoms |      |     |     |     | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|-----|---------|-------|
| 29  | b     | 119      | Total | C    | N   | O   | P   | 0       | 0     |
|     |       |          | 2549  | 1135 | 466 | 829 | 119 |         |       |

- Molecule 30 is a protein called 50S ribosomal protein L2.

| Mol | Chain | Residues | Atoms |      |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 30  | c     | 271      | Total | C    | N   | O   | S | 0       | 0     |
|     |       |          | 2082  | 1288 | 423 | 364 | 7 |         |       |

- Molecule 31 is a protein called 50S ribosomal protein L3.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 31  | d     | 209      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1566  | 980 | 288 | 294 | 4 |         |       |

- Molecule 32 is a protein called Large ribosomal subunit protein uL4.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 32  | e     | 201      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1552  | 974 | 283 | 290 | 5 |         |       |

- Molecule 33 is a protein called Large ribosomal subunit protein uL5.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 33  | f     | 177      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1410  | 899 | 249 | 256 | 6 |         |       |

- Molecule 34 is a protein called Large ribosomal subunit protein uL6.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 34  | g     | 176      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1323  | 832 | 243 | 246 | 2 |         |       |

- Molecule 35 is a protein called Large ribosomal subunit protein bL9.

| Mol | Chain | Residues | Atoms |     |    |    |   | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 35  | h     | 41       | Total | C   | N  | O  | S | 0       | 0     |
|     |       |          | 303   | 194 | 54 | 54 | 1 |         |       |

- Molecule 36 is a protein called Large ribosomal subunit protein uL13.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 36  | i     | 142      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1129  | 714 | 212 | 199 | 4 |         |       |

- Molecule 37 is a protein called Large ribosomal subunit protein uL14.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 37  | j     | 123      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 946   | 593 | 181 | 166 | 6 |         |       |

- Molecule 38 is a protein called Large ribosomal subunit protein uL15.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 38  | k     | 144      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1053  | 654 | 207 | 190 | 2 |         |       |

- Molecule 39 is a protein called Large ribosomal subunit protein uL16.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 39  | l     | 136      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 1075  | 686 | 205 | 177 | 7 |         |       |

- Molecule 40 is a protein called Large ribosomal subunit protein bL17.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 40  | m     | 118      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 945   | 585 | 194 | 161 | 5 |         |       |

- Molecule 41 is a protein called Large ribosomal subunit protein uL18.

| Mol | Chain | Residues | Atoms |     |     |     |  | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|--|---------|-------|
| 41  | n     | 116      | Total | C   | N   | O   |  | 0       | 0     |
|     |       |          | 892   | 552 | 178 | 162 |  |         |       |

- Molecule 42 is a protein called Large ribosomal subunit protein bL19.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 42  | o     | 114      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 917   | 574 | 179 | 163 | 1 |         |       |

- Molecule 43 is a protein called 50S ribosomal protein L20.

| Mol | Chain | Residues | Atoms |     |     |     |  | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|--|---------|-------|
| 43  | p     | 117      | Total | C   | N   | O   |  | 0       | 0     |
|     |       |          | 947   | 604 | 192 | 151 |  |         |       |

- Molecule 44 is a protein called Large ribosomal subunit protein bL21.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 44  | q     | 103      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 816   | 516 | 153 | 145 | 2 |         |       |

- Molecule 45 is a protein called Large ribosomal subunit protein uL22.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 45  | r     | 110      | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 857   | 532 | 166 | 156 | 3 |         |       |

- Molecule 46 is a protein called 50S ribosomal protein L23.

| Mol | Chain | Residues | Atoms |     |     |     |   | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 46  | s     | 93       | Total | C   | N   | O   | S | 0       | 0     |
|     |       |          | 738   | 466 | 139 | 131 | 2 |         |       |

- Molecule 47 is a protein called 50S ribosomal protein L24.

| Mol | Chain | Residues | Atoms |     |     |     | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 47  | t     | 102      | Total | C   | N   | O   | 0       | 0     |
|     |       |          | 779   | 492 | 146 | 141 |         |       |

- Molecule 48 is a protein called Large ribosomal subunit protein bL25.

| Mol | Chain | Residues | Atoms |     |     |     | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 48  | u     | 94       | Total | C   | N   | O   | S       | 0     |
|     |       |          | 753   | 479 | 137 | 134 | 3       | 0     |

- Molecule 49 is a protein called 50S ribosomal protein L27.

| Mol | Chain | Residues | Atoms |     |     |     | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 49  | v     | 78       | Total | C   | N   | O   | S       | 0     |
|     |       |          | 586   | 362 | 116 | 107 | 1       | 0     |

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

| Mol | Chain | Residues | Atoms |     |     |     | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 50  | w     | 77       | Total | C   | N   | O   | S       | 0     |
|     |       |          | 625   | 388 | 129 | 106 | 2       | 0     |

- Molecule 51 is a protein called Large ribosomal subunit protein uL29.

| Mol | Chain | Residues | Atoms |     |    |    | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---------|-------|
| 51  | x     | 62       | Total | C   | N  | O  | S       | 0     |
|     |       |          | 501   | 308 | 98 | 94 | 1       | 0     |

- Molecule 52 is a protein called 50S ribosomal protein L30.

| Mol | Chain | Residues | Atoms |     |    |    | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---------|-------|
| 52  | y     | 58       | Total | C   | N  | O  | S       | 0     |
|     |       |          | 449   | 281 | 87 | 79 | 2       | 0     |

- Molecule 53 is a protein called 50S ribosomal protein L32.

| Mol | Chain | Residues | Atoms |     |    |    | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---------|-------|
| 53  | z     | 56       | Total | C   | N  | O  | S       | 0     |
|     |       |          | 444   | 269 | 94 | 80 | 1       | 0     |

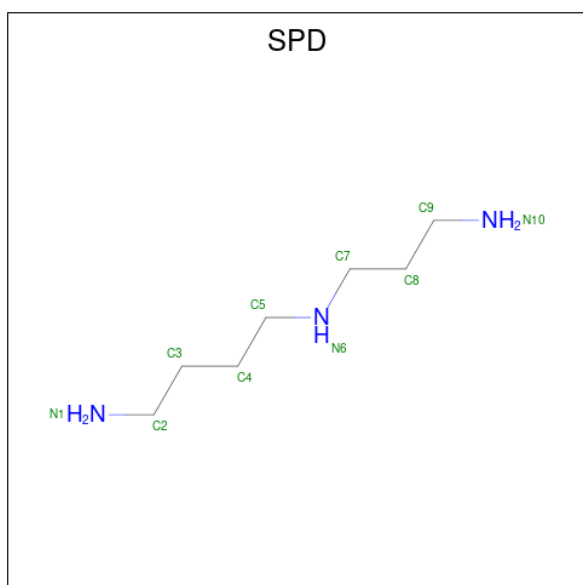
- Molecule 54 is ZINC ION (CCD ID: ZN) (formula: Zn).

| Mol | Chain | Residues | Atoms |    | AltConf |
|-----|-------|----------|-------|----|---------|
| 54  | 3     | 1        | Total | Zn | 0       |
|     |       |          | 1     | 1  |         |
| 54  | 4     | 1        | Total | Zn | 0       |
|     |       |          | 1     | 1  |         |

- Molecule 55 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

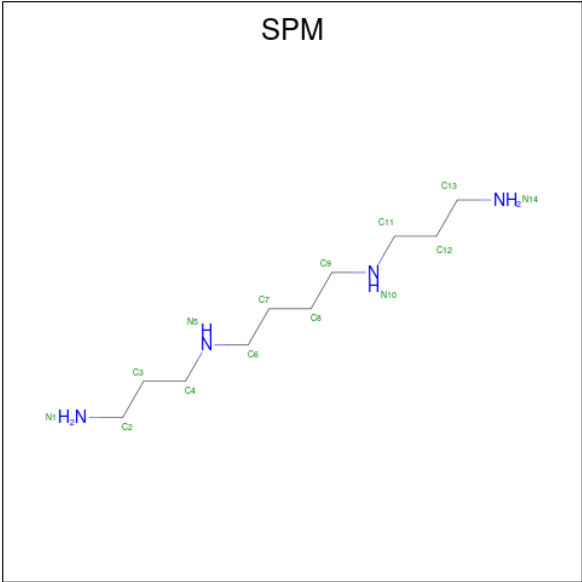
| Mol | Chain | Residues | Atoms |     | AltConf |
|-----|-------|----------|-------|-----|---------|
| 55  | A     | 55       | Total | Mg  | 0       |
|     |       |          | 55    | 55  |         |
| 55  | Q     | 1        | Total | Mg  | 0       |
|     |       |          | 1     | 1   |         |
| 55  | a     | 207      | Total | Mg  | 0       |
|     |       |          | 207   | 207 |         |
| 55  | b     | 5        | Total | Mg  | 0       |
|     |       |          | 5     | 5   |         |
| 55  | c     | 1        | Total | Mg  | 0       |
|     |       |          | 1     | 1   |         |
| 55  | m     | 1        | Total | Mg  | 0       |
|     |       |          | 1     | 1   |         |
| 55  | n     | 1        | Total | Mg  | 0       |
|     |       |          | 1     | 1   |         |
| 55  | z     | 1        | Total | Mg  | 0       |
|     |       |          | 1     | 1   |         |

- Molecule 56 is SPERMIDINE (CCD ID: SPD) (formula: C<sub>7</sub>H<sub>19</sub>N<sub>3</sub>).



| Mol | Chain | Residues | Atoms |   |   | AltConf |
|-----|-------|----------|-------|---|---|---------|
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |
| 56  | a     | 1        | Total | C | N | 0       |
|     |       |          | 10    | 7 | 3 |         |

- Molecule 57 is SPERMINE (CCD ID: SPM) (formula:  $C_{10}H_{26}N_4$ ).

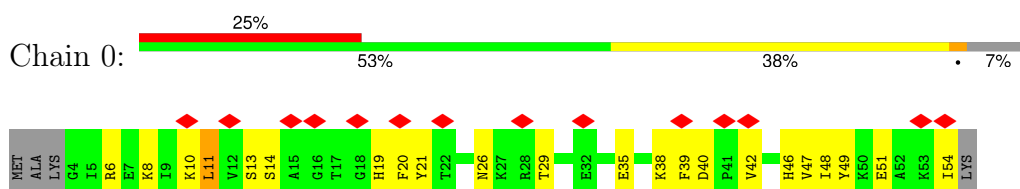


| Mol | Chain | Residues | Atoms |    |   | AltConf |
|-----|-------|----------|-------|----|---|---------|
| 57  | a     | 1        | Total | C  | N | 0       |
|     |       |          | 14    | 10 | 4 |         |

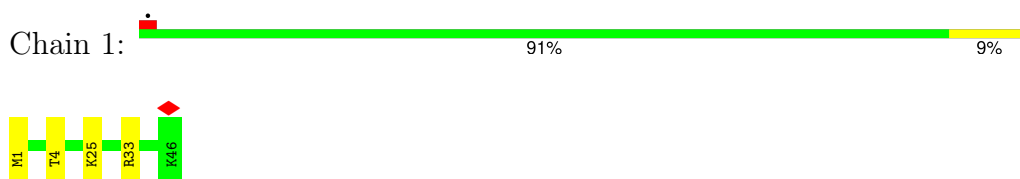
### 3 Residue-property plots [i](#)

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.

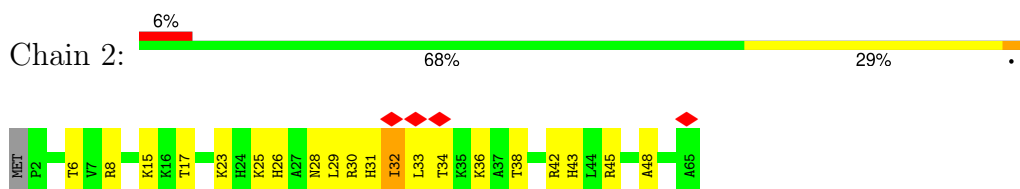
- Molecule 1: 50S ribosomal protein L33



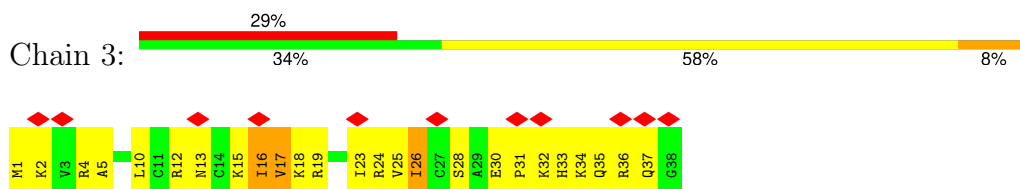
- Molecule 2: Large ribosomal subunit protein bL34



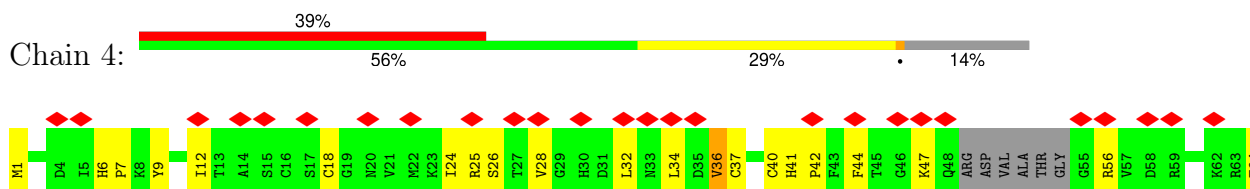
- Molecule 3: 50S ribosomal protein L35



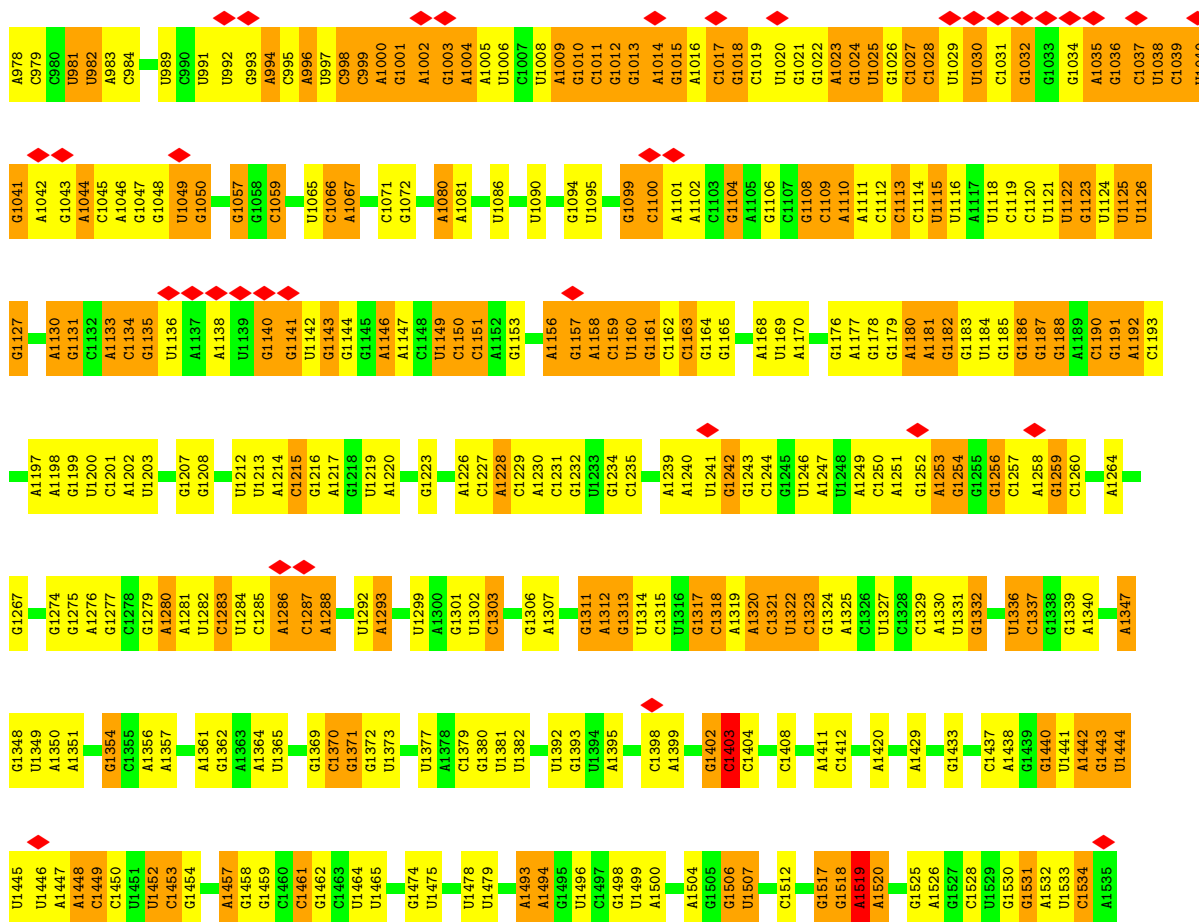
- Molecule 4: Large ribosomal subunit protein bL36A



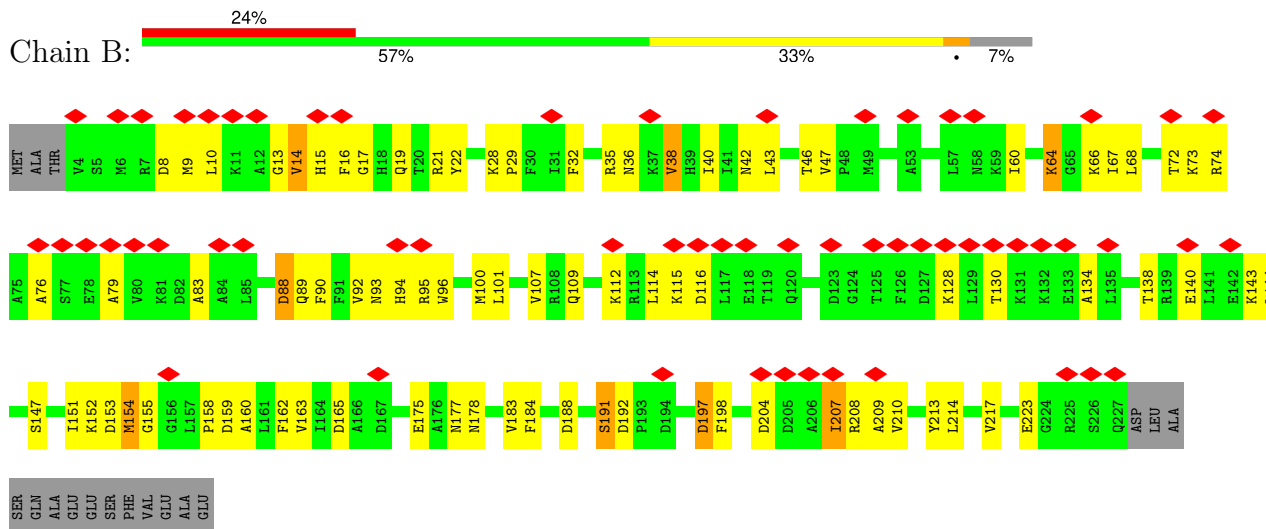
- Molecule 5: 50S ribosomal protein L31





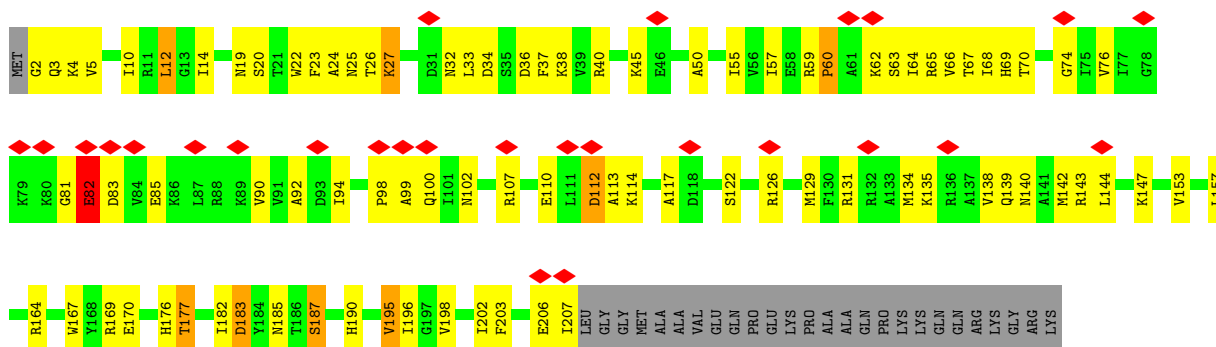


• Molecule 8: Small ribosomal subunit protein uS2

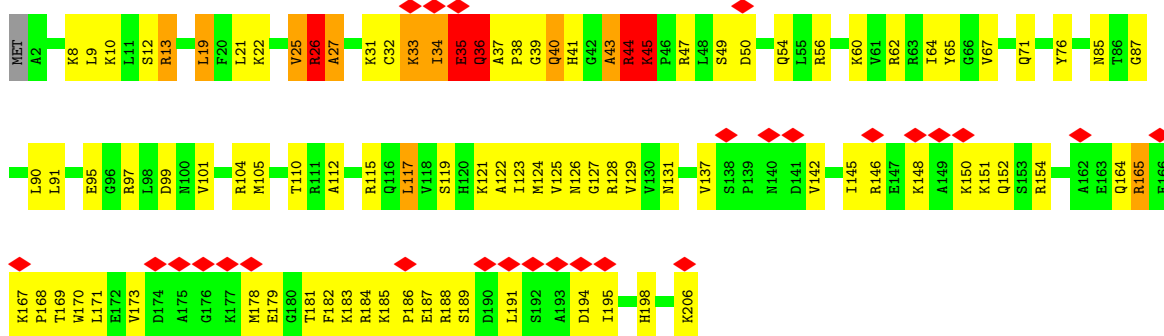


• Molecule 9: 30S ribosomal protein S3

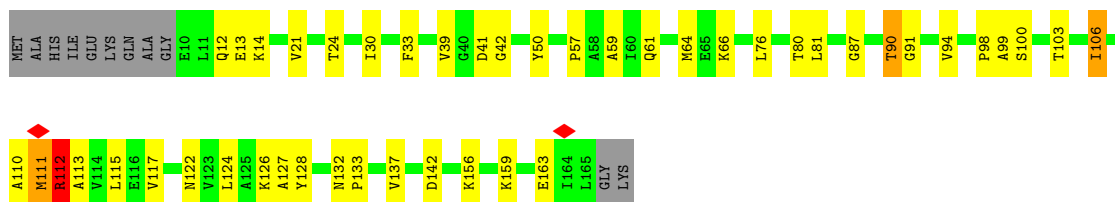




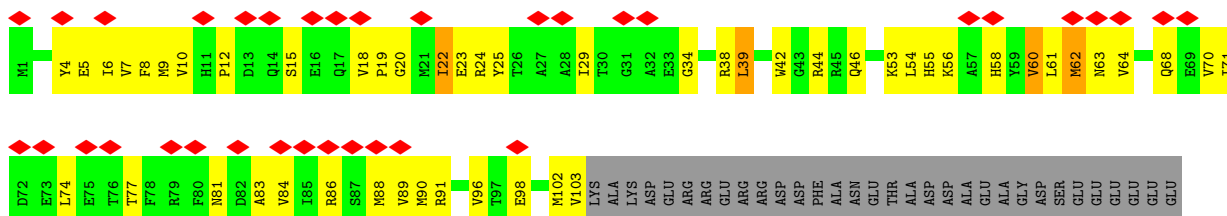
- Molecule 10: Small ribosomal subunit protein uS4



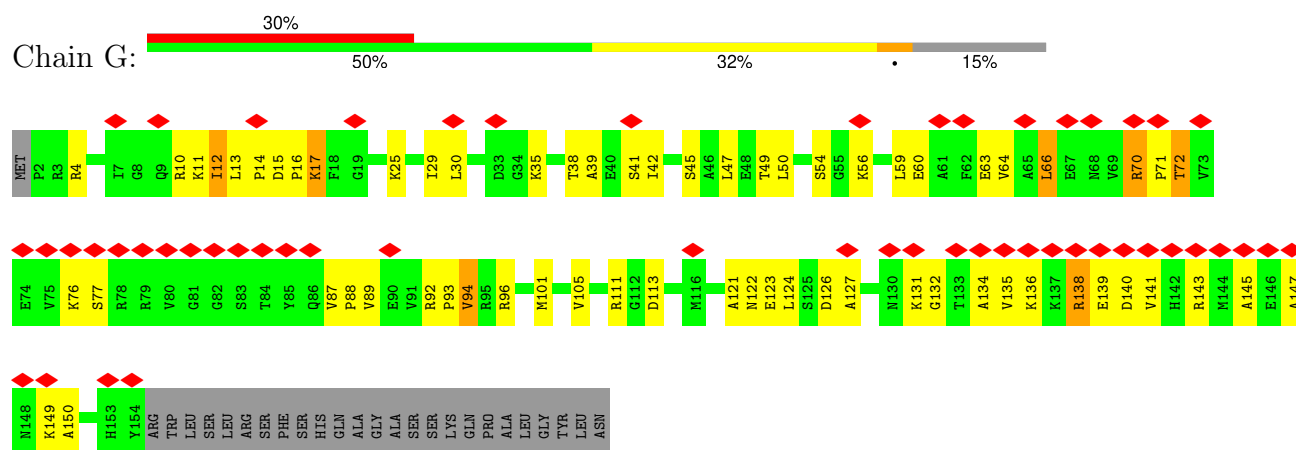
- Molecule 11: Small ribosomal subunit protein uS5



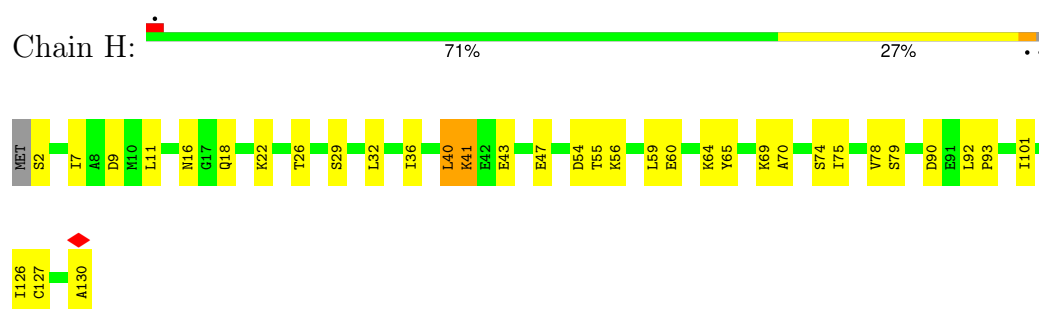
- Molecule 12: Small ribosomal subunit protein bS6



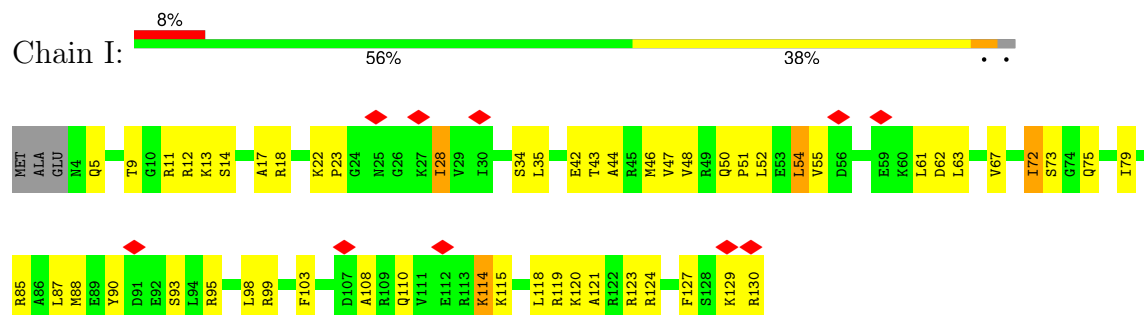
- Molecule 13: Small ribosomal subunit protein uS7



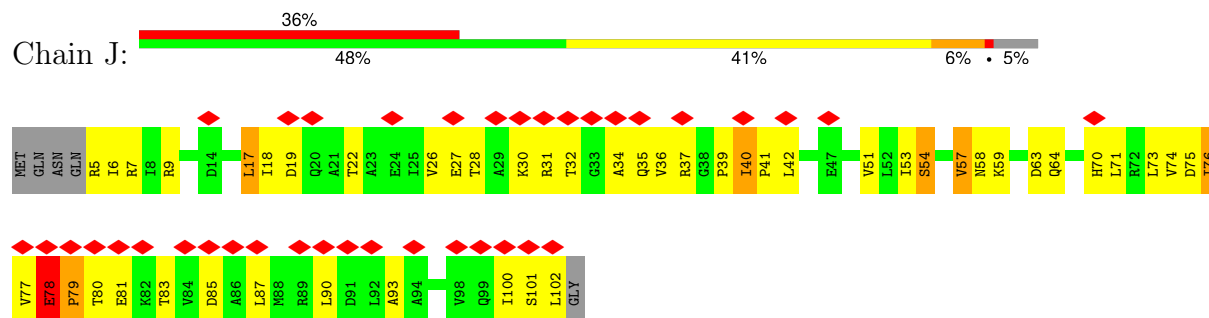
- Molecule 14: Small ribosomal subunit protein uS8



- Molecule 15: Small ribosomal subunit protein uS9

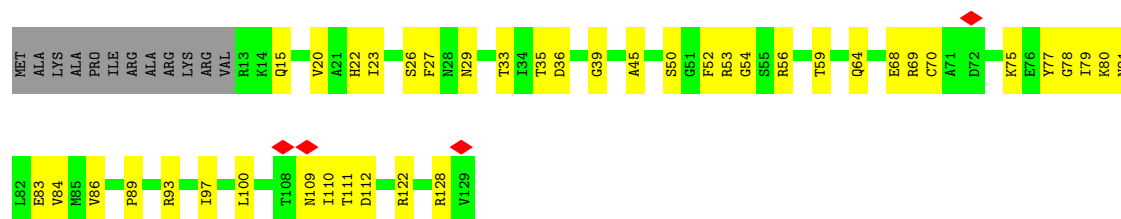


- Molecule 16: Small ribosomal subunit protein uS10



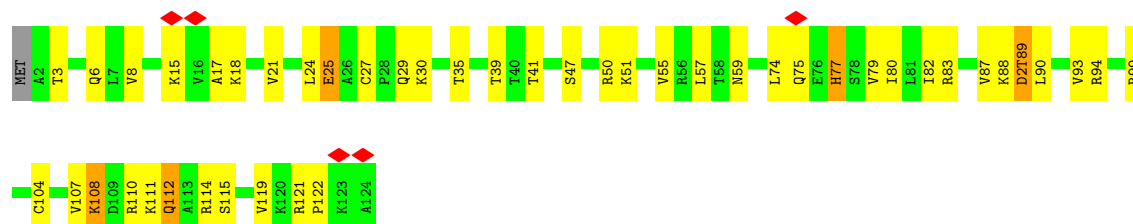
- Molecule 17: Small ribosomal subunit protein uS11





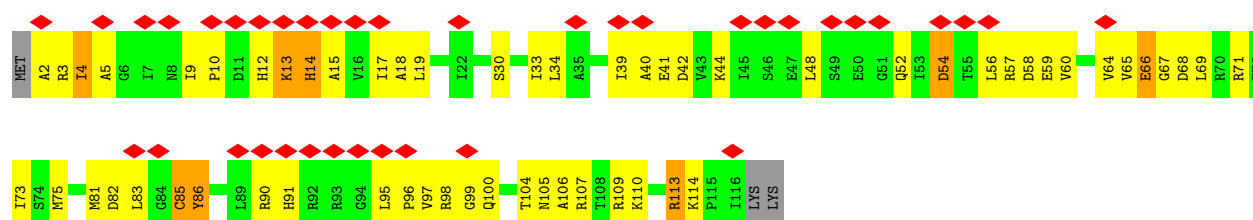
- Molecule 18: 30S ribosomal protein S12

Chain L: 62% 33%



- Molecule 19: Small ribosomal subunit protein uS13

Chain M: 32% 47% 43% 7%



- Molecule 20: Small ribosomal subunit protein uS14

Chain N: 5% 69% 27%



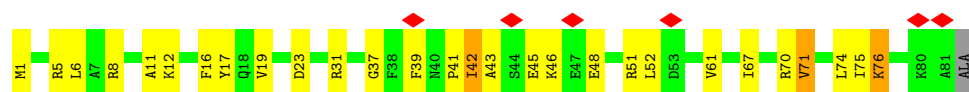
- Molecule 21: Small ribosomal subunit protein uS15

Chain O: 71% 25%

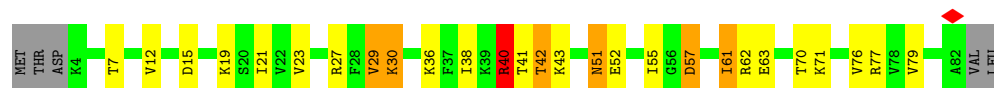


- Molecule 22: Small ribosomal subunit protein bS16

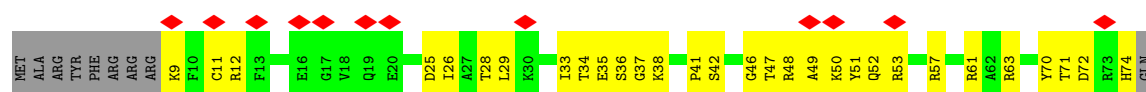
Chain P: 7% 65% 30%



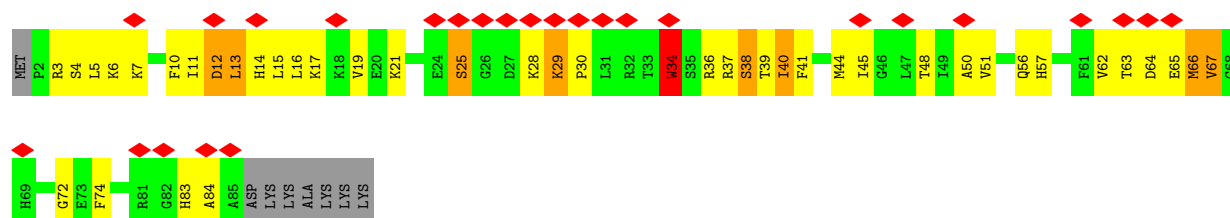
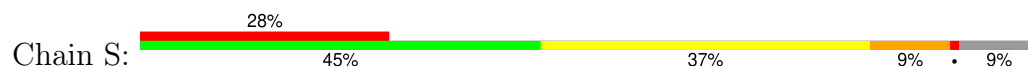
- Molecule 23: Small ribosomal subunit protein uS17



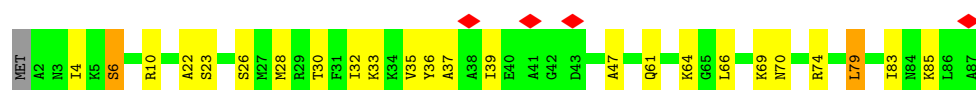
- Molecule 24: Small ribosomal subunit protein bS18



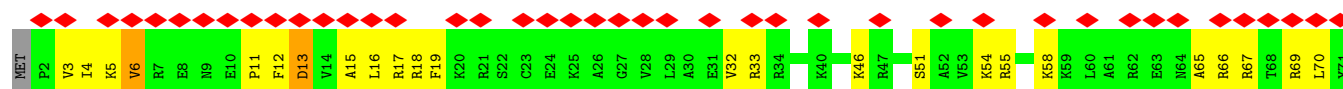
- Molecule 25: Small ribosomal subunit protein uS19



- Molecule 26: Small ribosomal subunit protein bS20



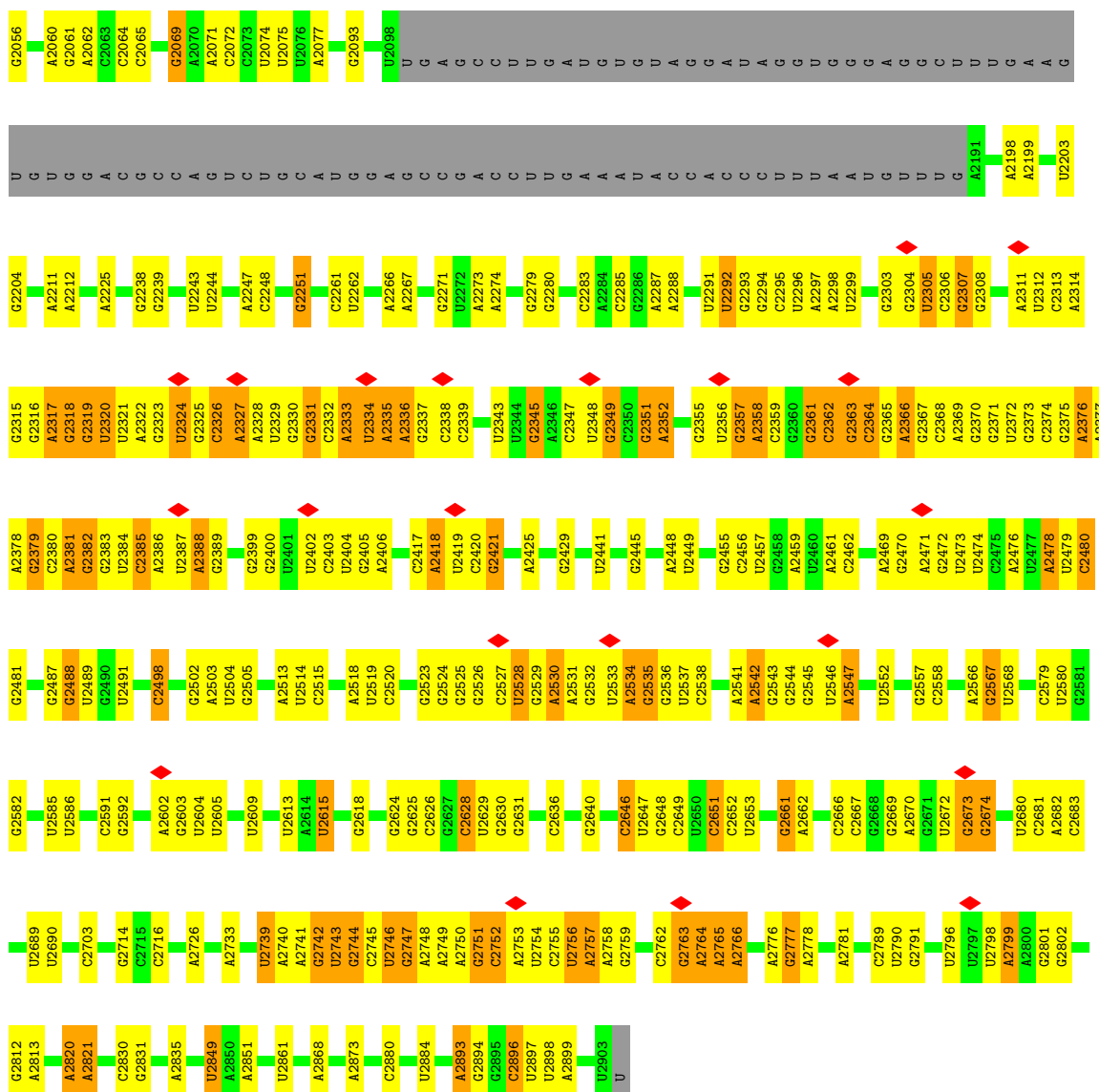
- Molecule 27: Small ribosomal subunit protein bS21



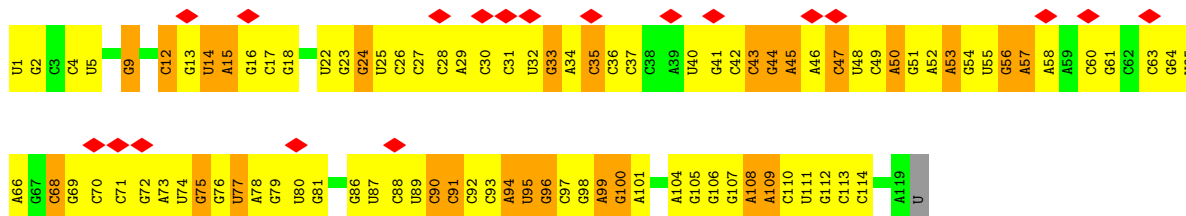
- Molecule 28: 23S ribosomal RNA



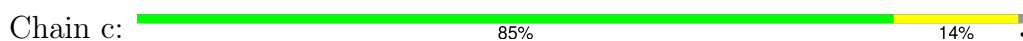


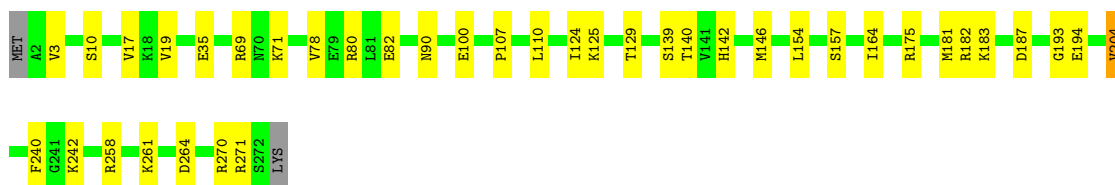


• Molecule 29: 5S ribosomal RNA

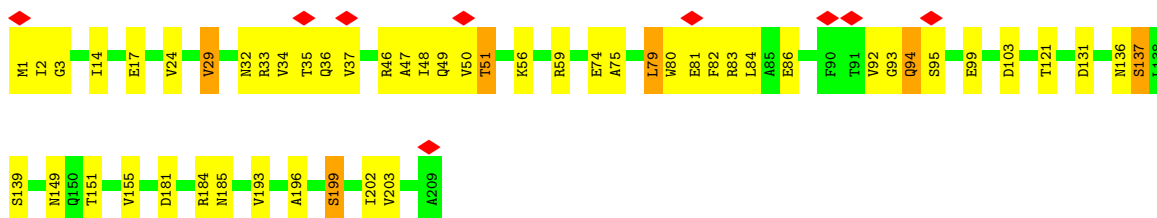


• Molecule 30: 50S ribosomal protein L2

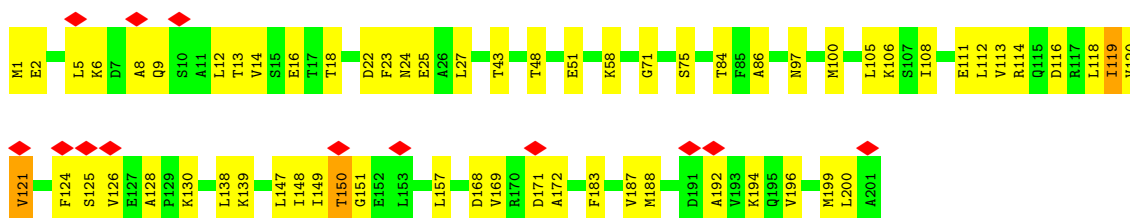




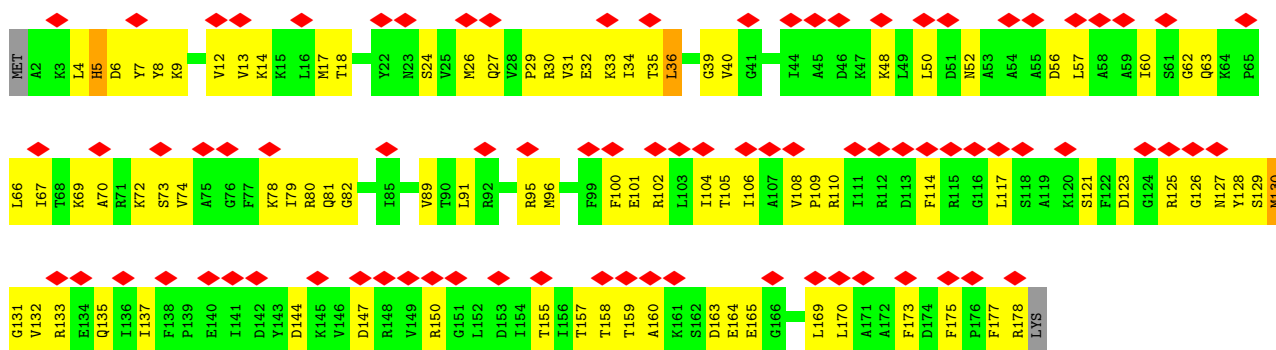
- Molecule 31: 50S ribosomal protein L3



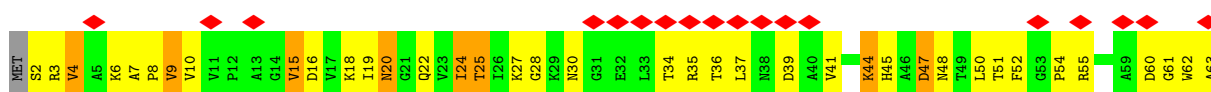
- Molecule 32: Large ribosomal subunit protein uL4



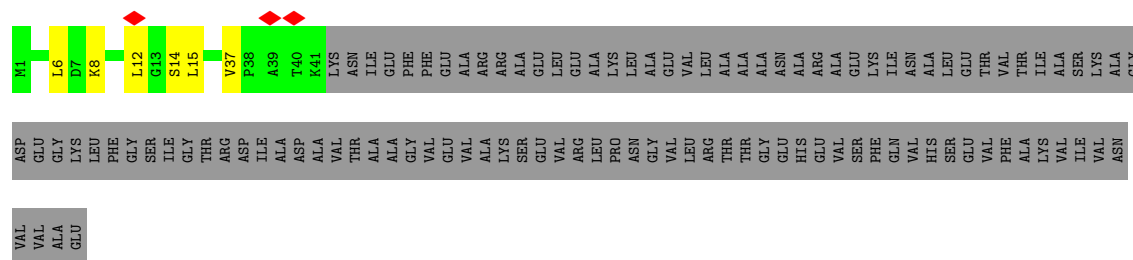
- Molecule 33: Large ribosomal subunit protein uL5



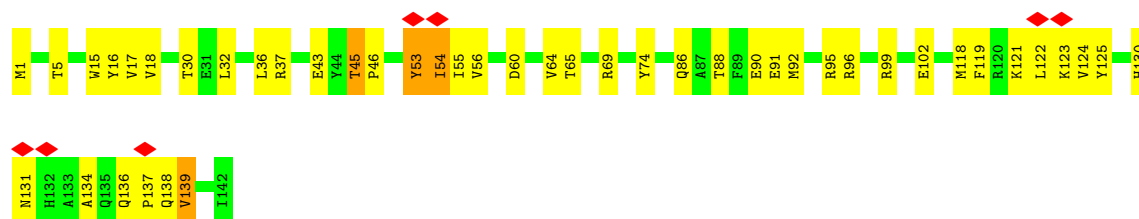
- Molecule 34: Large ribosomal subunit protein uL6



- Molecule 35: Large ribosomal subunit protein bL9



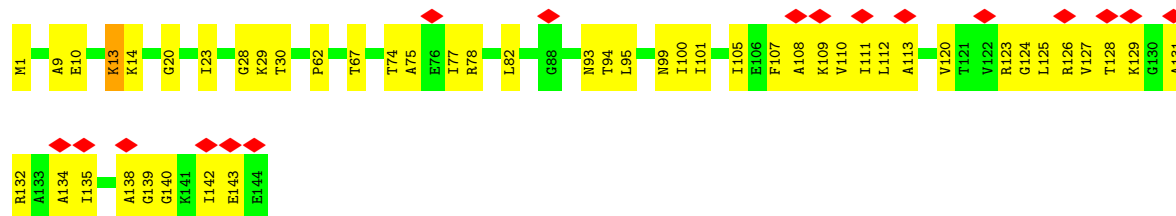
- Molecule 36: Large ribosomal subunit protein uL13



- Molecule 37: Large ribosomal subunit protein uL14

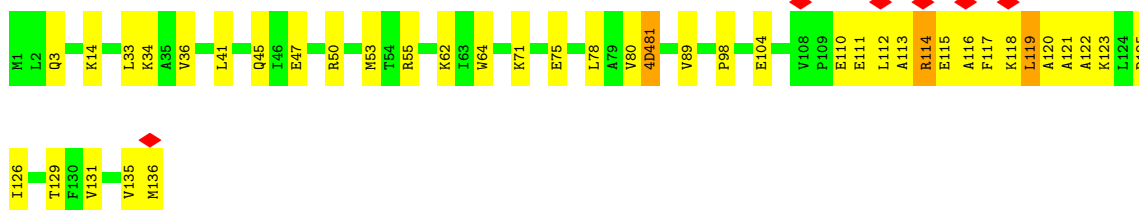


- Molecule 38: Large ribosomal subunit protein uL15



- Molecule 39: Large ribosomal subunit protein uL16

Chain l: 



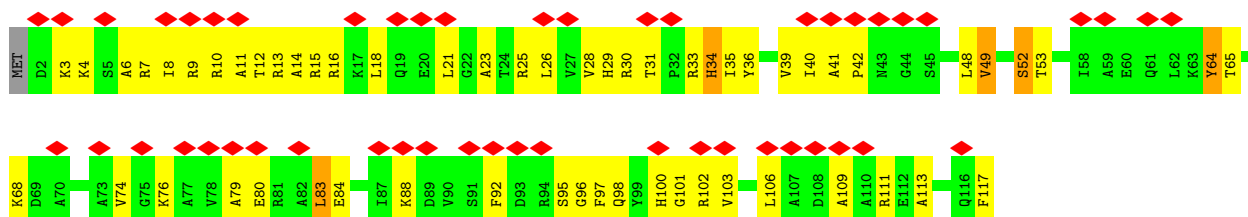
- Molecule 40: Large ribosomal subunit protein bL17

Chain m: 



- Molecule 41: Large ribosomal subunit protein uL18

Chain n: 



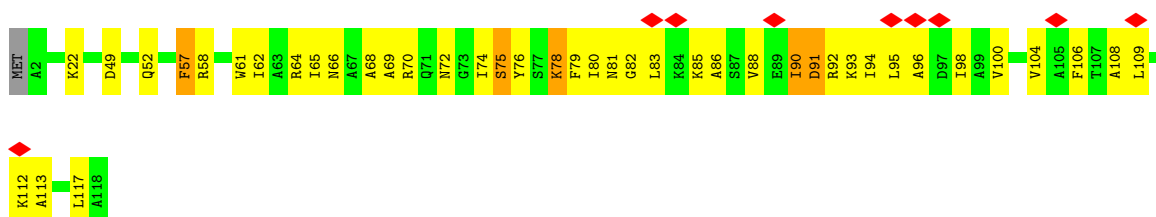
- Molecule 42: Large ribosomal subunit protein bL19

Chain o: 



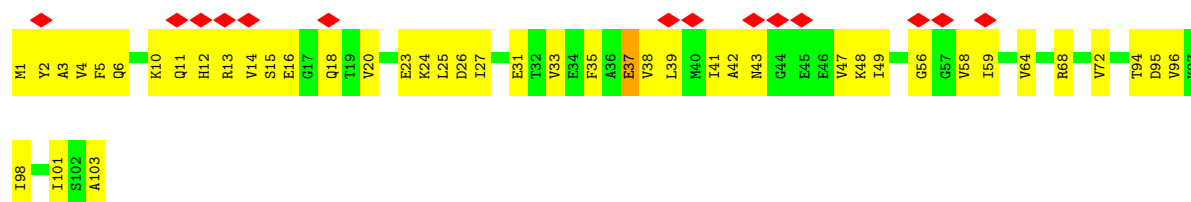
- Molecule 43: 50S ribosomal protein L20

Chain p: 

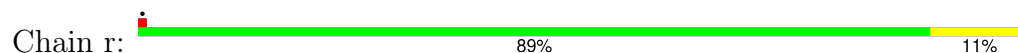


- Molecule 44: Large ribosomal subunit protein bL21

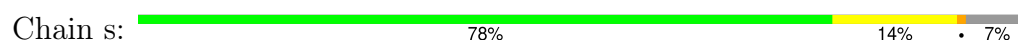
Chain q: 



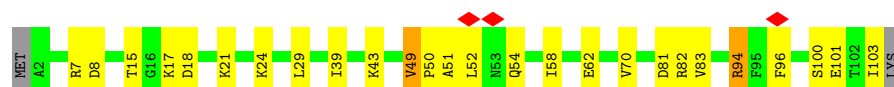
- Molecule 45: Large ribosomal subunit protein uL22



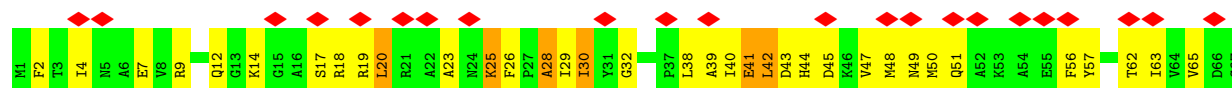
- Molecule 46: 50S ribosomal protein L23



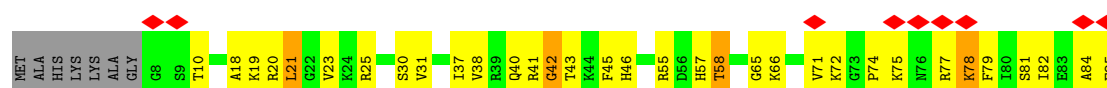
- Molecule 47: 50S ribosomal protein L24



- Molecule 48: Large ribosomal subunit protein bL25



- Molecule 49: 50S ribosomal protein L27



- Molecule 50: 50S ribosomal protein L28

Chain w:  78% 19% ..



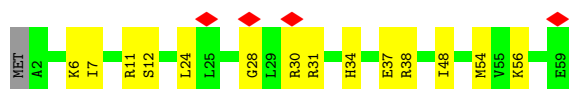
- Molecule 51: Large ribosomal subunit protein uL29

Chain x:  78% 21% .



- Molecule 52: 50S ribosomal protein L30

Chain y:  7% 75% 24% .



- Molecule 53: 50S ribosomal protein L32

Chain z:  79% 19% .



## 4 Experimental information

| Property                             | Value                                   | Source    |
|--------------------------------------|-----------------------------------------|-----------|
| EM reconstruction method             | SINGLE PARTICLE                         | Depositor |
| Imposed symmetry                     | POINT, Not provided                     |           |
| Number of particles used             | 450287                                  | Depositor |
| Resolution determination method      | FSC 0.143 CUT-OFF                       | Depositor |
| CTF correction method                | PHASE FLIPPING AND AMPLITUDE CORRECTION | Depositor |
| Microscope                           | TFS GLACIOS                             | Depositor |
| Voltage (kV)                         | 200                                     | Depositor |
| Electron dose ( $e^-/\text{\AA}^2$ ) | 60                                      | Depositor |
| Minimum defocus (nm)                 | 500                                     | Depositor |
| Maximum defocus (nm)                 | 2500                                    | Depositor |
| Magnification                        | Not provided                            |           |
| Image detector                       | FEI FALCON IV (4k x 4k)                 | Depositor |
| Maximum map value                    | 43.573                                  | Depositor |
| Minimum map value                    | -23.645                                 | Depositor |
| Average map value                    | 0.000                                   | Depositor |
| Map value standard deviation         | 1.000                                   | Depositor |
| Recommended contour level            | 2.8                                     | Depositor |
| Map size ( $\text{\AA}$ )            | 579.36, 579.36, 579.36                  | wwPDB     |
| Map dimensions                       | 480, 480, 480                           | wwPDB     |
| Map angles ( $^\circ$ )              | 90.0, 90.0, 90.0                        | wwPDB     |
| Pixel spacing ( $\text{\AA}$ )       | 1.207, 1.207, 1.207                     | Depositor |

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 5MC, UR3, MA6, 1MG, OMU, MG, BH2, H2U, SPM, OMC, 6MZ, 5MU, MS6, PSU, OMG, G7M, 2MA, SPD, 2MG, D2T, 3TD, ZN, 4OC, MEQ, 4D4

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     | 0.25         | 0/424         | 0.58        | 1/565 (0.2%)    |
| 2   | 1     | 0.24         | 0/380         | 0.38        | 0/498           |
| 3   | 2     | 0.27         | 0/513         | 0.45        | 0/676           |
| 4   | 3     | 0.33         | 0/303         | 0.76        | 0/397           |
| 5   | 4     | 0.23         | 0/488         | 0.49        | 0/649           |
| 6   | 5     | 0.10         | 0/46          | 0.09        | 0/69            |
| 7   | A     | 0.25         | 0/36183       | 0.45        | 10/56432 (0.0%) |
| 8   | B     | 0.26         | 0/1784        | 0.64        | 2/2403 (0.1%)   |
| 9   | C     | 0.24         | 0/1651        | 0.63        | 1/2225 (0.0%)   |
| 10  | D     | 0.41         | 0/1664        | 0.80        | 2/2226 (0.1%)   |
| 11  | E     | 0.25         | 0/1165        | 0.65        | 1/1568 (0.1%)   |
| 12  | F     | 0.25         | 0/858         | 0.60        | 0/1160          |
| 13  | G     | 0.55         | 2/1219 (0.2%) | 0.92        | 4/1635 (0.2%)   |
| 14  | H     | 0.22         | 0/988         | 0.59        | 1/1326 (0.1%)   |
| 15  | I     | 0.22         | 0/1034        | 0.61        | 0/1375          |
| 16  | J     | 0.55         | 1/796 (0.1%)  | 0.99        | 3/1077 (0.3%)   |
| 17  | K     | 0.20         | 0/884         | 0.54        | 0/1191          |
| 18  | L     | 0.21         | 0/960         | 0.52        | 1/1286 (0.1%)   |
| 19  | M     | 0.24         | 0/900         | 0.78        | 2/1204 (0.2%)   |
| 20  | N     | 0.21         | 0/817         | 0.59        | 0/1088          |
| 21  | O     | 0.37         | 0/722         | 0.55        | 1/964 (0.1%)    |
| 22  | P     | 0.24         | 0/653         | 0.65        | 0/877           |
| 23  | Q     | 0.30         | 0/650         | 0.71        | 0/871           |
| 24  | R     | 0.19         | 0/553         | 0.55        | 0/742           |
| 25  | S     | 0.26         | 0/685         | 0.75        | 1/922 (0.1%)    |
| 26  | T     | 0.20         | 0/675         | 0.58        | 1/893 (0.1%)    |
| 27  | U     | 0.24         | 0/597         | 0.55        | 0/792           |
| 28  | a     | 0.26         | 0/65651       | 0.36        | 0/102413        |
| 29  | b     | 0.26         | 0/2850        | 0.43        | 0/4444          |
| 30  | c     | 0.22         | 0/2121        | 0.36        | 0/2852          |
| 31  | d     | 0.26         | 0/1576        | 0.47        | 0/2119          |

| Mol | Chain | Bond lengths |                 | Bond angles |                  |
|-----|-------|--------------|-----------------|-------------|------------------|
|     |       | RMSZ         | # Z  >5         | RMSZ        | # Z  >5          |
| 32  | e     | 0.24         | 0/1571          | 0.48        | 0/2113           |
| 33  | f     | 0.21         | 0/1434          | 0.41        | 0/1926           |
| 34  | g     | 0.22         | 0/1343          | 0.53        | 0/1816           |
| 35  | h     | 0.20         | 0/306           | 0.54        | 0/413            |
| 36  | i     | 0.29         | 0/1152          | 0.54        | 0/1551           |
| 37  | j     | 0.22         | 0/955           | 0.37        | 0/1279           |
| 38  | k     | 0.23         | 0/1062          | 0.46        | 0/1413           |
| 39  | l     | 0.26         | 0/1073          | 0.54        | 0/1433           |
| 40  | m     | 0.23         | 0/958           | 0.39        | 0/1281           |
| 41  | n     | 0.53         | 1/902 (0.1%)    | 0.47        | 0/1209           |
| 42  | o     | 0.21         | 0/929           | 0.35        | 0/1242           |
| 43  | p     | 0.36         | 0/960           | 0.66        | 0/1278           |
| 44  | q     | 0.27         | 0/829           | 0.56        | 0/1107           |
| 45  | r     | 0.20         | 0/864           | 0.29        | 0/1156           |
| 46  | s     | 0.21         | 0/744           | 0.42        | 0/994            |
| 47  | t     | 0.21         | 0/787           | 0.49        | 0/1051           |
| 48  | u     | 0.33         | 1/766 (0.1%)    | 0.63        | 0/1025           |
| 49  | v     | 0.29         | 0/593           | 0.67        | 1/785 (0.1%)     |
| 50  | w     | 0.20         | 0/635           | 0.35        | 0/848            |
| 51  | x     | 0.22         | 0/502           | 0.52        | 0/667            |
| 52  | y     | 0.25         | 0/453           | 0.54        | 0/605            |
| 53  | z     | 0.21         | 0/450           | 0.37        | 0/599            |
| All | All   | 0.27         | 5/149058 (0.0%) | 0.45        | 32/222730 (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 |
|-----|-------|---------------------|---------------------|
| 10  | D     | 0                   | 3                   |
| 11  | E     | 0                   | 1                   |
| 16  | J     | 0                   | 1                   |
| 17  | K     | 1                   | 0                   |
| 20  | N     | 0                   | 1                   |
| 21  | O     | 0                   | 1                   |
| 23  | Q     | 0                   | 1                   |
| 34  | g     | 0                   | 1                   |
| 39  | l     | 0                   | 3                   |
| All | All   | 1                   | 12                  |

All (5) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z      | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|-------|--------|-------------|----------|
| 13  | G     | 14  | PRO  | CG-CD | -15.01 | 0.99        | 1.50     |
| 41  | n     | 34  | HIS  | C-N   | -13.82 | 1.25        | 1.33     |
| 16  | J     | 79  | PRO  | CG-CD | -12.36 | 1.08        | 1.50     |
| 13  | G     | 14  | PRO  | N-CD  | 6.37   | 1.56        | 1.47     |
| 48  | u     | 28  | ALA  | C-N   | -5.12  | 1.30        | 1.33     |

The worst 5 of 32 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms    | Z      | Observed(°) | Ideal(°) |
|-----|-------|-----|------|----------|--------|-------------|----------|
| 13  | G     | 14  | PRO  | N-CD-CG  | -18.57 | 75.34       | 103.20   |
| 16  | J     | 79  | PRO  | N-CD-CG  | -16.56 | 78.36       | 103.20   |
| 16  | J     | 79  | PRO  | CA-CB-CG | -12.96 | 79.88       | 104.50   |
| 13  | G     | 14  | PRO  | CA-CB-CG | -11.90 | 81.90       | 104.50   |
| 13  | G     | 14  | PRO  | CA-N-CD  | -9.30  | 98.98       | 112.00   |

All (1) chirality outliers are listed below:

| Mol | Chain | Res | Type | Atom |
|-----|-------|-----|------|------|
| 17  | K     | 119 | BH2  | CB   |

5 of 12 planarity outliers are listed below:

| Mol | Chain | Res | Type | Group     |
|-----|-------|-----|------|-----------|
| 10  | D     | 13  | ARG  | Sidechain |
| 10  | D     | 26  | ARG  | Sidechain |
| 10  | D     | 44  | ARG  | Sidechain |
| 11  | E     | 112 | ARG  | Sidechain |
| 16  | J     | 78  | GLU  | Peptide   |

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1   | 0     | 417   | 0        | 451      | 21      | 0            |
| 2   | 1     | 377   | 0        | 418      | 5       | 0            |
| 3   | 2     | 504   | 0        | 572      | 18      | 0            |
| 4   | 3     | 302   | 0        | 340      | 31      | 0            |
| 5   | 4     | 480   | 0        | 478      | 22      | 0            |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 6   | 5     | 42    | 0        | 23       | 0       | 0            |
| 7   | A     | 32564 | 0        | 16407    | 523     | 0            |
| 8   | B     | 1753  | 0        | 1780     | 58      | 0            |
| 9   | C     | 1624  | 0        | 1696     | 63      | 0            |
| 10  | D     | 1642  | 0        | 1704     | 73      | 0            |
| 11  | E     | 1152  | 0        | 1196     | 29      | 0            |
| 12  | F     | 839   | 0        | 833      | 42      | 0            |
| 13  | G     | 1203  | 0        | 1254     | 49      | 0            |
| 14  | H     | 978   | 0        | 1031     | 24      | 0            |
| 15  | I     | 1022  | 0        | 1070     | 46      | 0            |
| 16  | J     | 786   | 0        | 828      | 36      | 0            |
| 17  | K     | 876   | 0        | 882      | 35      | 0            |
| 18  | L     | 957   | 0        | 1016     | 27      | 0            |
| 19  | M     | 891   | 0        | 952      | 60      | 0            |
| 20  | N     | 805   | 0        | 844      | 28      | 0            |
| 21  | O     | 714   | 0        | 734      | 12      | 0            |
| 22  | P     | 643   | 0        | 661      | 23      | 0            |
| 23  | Q     | 641   | 0        | 682      | 16      | 0            |
| 24  | R     | 544   | 0        | 565      | 28      | 0            |
| 25  | S     | 668   | 0        | 693      | 43      | 0            |
| 26  | T     | 669   | 0        | 715      | 17      | 0            |
| 27  | U     | 589   | 0        | 629      | 27      | 0            |
| 28  | a     | 59130 | 0        | 29768    | 760     | 0            |
| 29  | b     | 2549  | 0        | 1291     | 121     | 0            |
| 30  | c     | 2082  | 0        | 2154     | 19      | 0            |
| 31  | d     | 1566  | 0        | 1617     | 46      | 0            |
| 32  | e     | 1552  | 0        | 1619     | 56      | 0            |
| 33  | f     | 1410  | 0        | 1444     | 79      | 0            |
| 34  | g     | 1323  | 0        | 1371     | 106     | 0            |
| 35  | h     | 303   | 0        | 327      | 4       | 0            |
| 36  | i     | 1129  | 0        | 1162     | 66      | 0            |
| 37  | j     | 946   | 0        | 1023     | 20      | 0            |
| 38  | k     | 1053  | 0        | 1129     | 49      | 0            |
| 39  | l     | 1075  | 0        | 1145     | 33      | 0            |
| 40  | m     | 945   | 0        | 989      | 11      | 0            |
| 41  | n     | 892   | 0        | 923      | 55      | 0            |
| 42  | o     | 917   | 0        | 962      | 9       | 0            |
| 43  | p     | 947   | 0        | 1019     | 67      | 0            |
| 44  | q     | 816   | 0        | 839      | 45      | 0            |
| 45  | r     | 857   | 0        | 922      | 6       | 0            |
| 46  | s     | 738   | 0        | 807      | 8       | 0            |
| 47  | t     | 779   | 0        | 831      | 22      | 0            |

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| Mol | Chain | Non-H  | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|--------|----------|----------|---------|--------------|
| 48  | u     | 753    | 0        | 780      | 77      | 0            |
| 49  | v     | 586    | 0        | 596      | 48      | 0            |
| 50  | w     | 625    | 0        | 652      | 5       | 0            |
| 51  | x     | 501    | 0        | 531      | 7       | 0            |
| 52  | y     | 449    | 0        | 488      | 11      | 0            |
| 53  | z     | 444    | 0        | 458      | 7       | 0            |
| 54  | 3     | 1      | 0        | 0        | 0       | 0            |
| 54  | 4     | 1      | 0        | 0        | 0       | 0            |
| 55  | A     | 55     | 0        | 0        | 0       | 0            |
| 55  | Q     | 1      | 0        | 0        | 0       | 0            |
| 55  | a     | 207    | 0        | 0        | 0       | 0            |
| 55  | b     | 5      | 0        | 0        | 0       | 0            |
| 55  | c     | 1      | 0        | 0        | 0       | 0            |
| 55  | m     | 1      | 0        | 0        | 0       | 0            |
| 55  | n     | 1      | 0        | 0        | 0       | 0            |
| 55  | z     | 1      | 0        | 0        | 0       | 0            |
| 56  | a     | 140    | 0        | 266      | 12      | 0            |
| 57  | a     | 14     | 0        | 26       | 1       | 0            |
| All | All   | 138477 | 0        | 93593    | 2821    | 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 12.

The worst 5 of 2821 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) |
|-----------------|----------------|--------------------------|-------------------|
| 29:b:78:A:H62   | 29:b:98:G:N2   | 1.20                     | 1.38              |
| 7:A:998:C:C6    | 7:A:1044:A:C5' | 2.07                     | 1.37              |
| 28:a:1418:G:H21 | 28:a:1580:A:N6 | 1.23                     | 1.37              |
| 28:a:1020:A:N6  | 28:a:1141:U:C2 | 2.00                     | 1.29              |
| 7:A:73:C:N4     | 7:A:96:U:H3    | 1.30                     | 1.29              |

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     | 49/55 (89%)   | 47 (96%)  | 2 (4%)   | 0        | 100         | 100 |
| 2   | 1     | 44/46 (96%)   | 43 (98%)  | 1 (2%)   | 0        | 100         | 100 |
| 3   | 2     | 62/65 (95%)   | 56 (90%)  | 6 (10%)  | 0        | 100         | 100 |
| 4   | 3     | 36/38 (95%)   | 30 (83%)  | 6 (17%)  | 0        | 100         | 100 |
| 5   | 4     | 56/70 (80%)   | 49 (88%)  | 7 (12%)  | 0        | 100         | 100 |
| 8   | B     | 222/241 (92%) | 196 (88%) | 24 (11%) | 2 (1%)   | 14          | 48  |
| 9   | C     | 204/233 (88%) | 159 (78%) | 40 (20%) | 5 (2%)   | 4           | 23  |
| 10  | D     | 203/206 (98%) | 150 (74%) | 41 (20%) | 12 (6%)  | 1           | 7   |
| 11  | E     | 154/167 (92%) | 129 (84%) | 24 (16%) | 1 (1%)   | 21          | 56  |
| 12  | F     | 101/135 (75%) | 91 (90%)  | 10 (10%) | 0        | 100         | 100 |
| 13  | G     | 151/179 (84%) | 130 (86%) | 21 (14%) | 0        | 100         | 100 |
| 14  | H     | 127/130 (98%) | 106 (84%) | 21 (16%) | 0        | 100         | 100 |
| 15  | I     | 125/130 (96%) | 103 (82%) | 22 (18%) | 0        | 100         | 100 |
| 16  | J     | 96/103 (93%)  | 81 (84%)  | 12 (12%) | 3 (3%)   | 3           | 19  |
| 17  | K     | 113/129 (88%) | 106 (94%) | 7 (6%)   | 0        | 100         | 100 |
| 18  | L     | 119/124 (96%) | 102 (86%) | 17 (14%) | 0        | 100         | 100 |
| 19  | M     | 113/118 (96%) | 81 (72%)  | 30 (26%) | 2 (2%)   | 6           | 31  |
| 20  | N     | 98/101 (97%)  | 88 (90%)  | 10 (10%) | 0        | 100         | 100 |
| 21  | O     | 86/89 (97%)   | 81 (94%)  | 3 (4%)   | 2 (2%)   | 5           | 25  |
| 22  | P     | 79/82 (96%)   | 59 (75%)  | 18 (23%) | 2 (2%)   | 4           | 23  |
| 23  | Q     | 77/84 (92%)   | 67 (87%)  | 10 (13%) | 0        | 100         | 100 |
| 24  | R     | 64/75 (85%)   | 55 (86%)  | 9 (14%)  | 0        | 100         | 100 |
| 25  | S     | 82/92 (89%)   | 57 (70%)  | 23 (28%) | 2 (2%)   | 4           | 24  |
| 26  | T     | 84/87 (97%)   | 83 (99%)  | 1 (1%)   | 0        | 100         | 100 |
| 27  | U     | 68/71 (96%)   | 63 (93%)  | 5 (7%)   | 0        | 100         | 100 |
| 30  | c     | 269/273 (98%) | 259 (96%) | 10 (4%)  | 0        | 100         | 100 |
| 31  | d     | 206/209 (99%) | 189 (92%) | 17 (8%)  | 0        | 100         | 100 |
| 32  | e     | 199/201 (99%) | 190 (96%) | 9 (4%)   | 0        | 100         | 100 |
| 33  | f     | 175/179 (98%) | 158 (90%) | 17 (10%) | 0        | 100         | 100 |
| 34  | g     | 174/177 (98%) | 150 (86%) | 23 (13%) | 1 (1%)   | 21          | 56  |

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| Mol | Chain | Analysed        | Favoured   | Allowed   | Outliers | Percentiles |     |
|-----|-------|-----------------|------------|-----------|----------|-------------|-----|
| 35  | h     | 39/149 (26%)    | 34 (87%)   | 5 (13%)   | 0        | 100         | 100 |
| 36  | i     | 140/142 (99%)   | 131 (94%)  | 9 (6%)    | 0        | 100         | 100 |
| 37  | j     | 121/123 (98%)   | 117 (97%)  | 4 (3%)    | 0        | 100         | 100 |
| 38  | k     | 142/144 (99%)   | 135 (95%)  | 7 (5%)    | 0        | 100         | 100 |
| 39  | l     | 132/136 (97%)   | 127 (96%)  | 5 (4%)    | 0        | 100         | 100 |
| 40  | m     | 116/127 (91%)   | 108 (93%)  | 8 (7%)    | 0        | 100         | 100 |
| 41  | n     | 114/117 (97%)   | 107 (94%)  | 7 (6%)    | 0        | 100         | 100 |
| 42  | o     | 112/115 (97%)   | 109 (97%)  | 3 (3%)    | 0        | 100         | 100 |
| 43  | p     | 115/118 (98%)   | 106 (92%)  | 9 (8%)    | 0        | 100         | 100 |
| 44  | q     | 101/103 (98%)   | 89 (88%)   | 12 (12%)  | 0        | 100         | 100 |
| 45  | r     | 108/110 (98%)   | 105 (97%)  | 3 (3%)    | 0        | 100         | 100 |
| 46  | s     | 91/100 (91%)    | 83 (91%)   | 8 (9%)    | 0        | 100         | 100 |
| 47  | t     | 100/104 (96%)   | 92 (92%)   | 8 (8%)    | 0        | 100         | 100 |
| 48  | u     | 92/94 (98%)     | 83 (90%)   | 9 (10%)   | 0        | 100         | 100 |
| 49  | v     | 76/85 (89%)     | 65 (86%)   | 11 (14%)  | 0        | 100         | 100 |
| 50  | w     | 75/78 (96%)     | 74 (99%)   | 1 (1%)    | 0        | 100         | 100 |
| 51  | x     | 60/63 (95%)     | 58 (97%)   | 2 (3%)    | 0        | 100         | 100 |
| 52  | y     | 56/59 (95%)     | 55 (98%)   | 1 (2%)    | 0        | 100         | 100 |
| 53  | z     | 54/57 (95%)     | 54 (100%)  | 0         | 0        | 100         | 100 |
| All | All   | 5480/5913 (93%) | 4890 (89%) | 558 (10%) | 32 (1%)  | 23          | 56  |

5 of 32 Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 8   | B     | 154 | MET  |
| 9   | C     | 26  | THR  |
| 10  | D     | 34  | ILE  |
| 10  | D     | 36  | GLN  |
| 10  | D     | 40  | GLN  |

### 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     | 46/49 (94%)    | 45 (98%)  | 1 (2%)   | 45          | 74  |
| 2   | 1     | 38/38 (100%)   | 38 (100%) | 0        | 100         | 100 |
| 3   | 2     | 51/52 (98%)    | 47 (92%)  | 4 (8%)   | 11          | 39  |
| 4   | 3     | 34/34 (100%)   | 30 (88%)  | 4 (12%)  | 5           | 22  |
| 5   | 4     | 55/62 (89%)    | 53 (96%)  | 2 (4%)   | 31          | 65  |
| 8   | B     | 186/199 (94%)  | 169 (91%) | 17 (9%)  | 9           | 33  |
| 9   | C     | 170/190 (90%)  | 155 (91%) | 15 (9%)  | 9           | 35  |
| 10  | D     | 171/173 (99%)  | 152 (89%) | 19 (11%) | 6           | 25  |
| 11  | E     | 119/126 (94%)  | 108 (91%) | 11 (9%)  | 8           | 33  |
| 12  | F     | 90/116 (78%)   | 83 (92%)  | 7 (8%)   | 11          | 39  |
| 13  | G     | 126/147 (86%)  | 112 (89%) | 14 (11%) | 6           | 25  |
| 14  | H     | 104/105 (99%)  | 99 (95%)  | 5 (5%)   | 23          | 57  |
| 15  | I     | 105/107 (98%)  | 98 (93%)  | 7 (7%)   | 15          | 46  |
| 16  | J     | 86/90 (96%)    | 73 (85%)  | 13 (15%) | 3           | 14  |
| 17  | K     | 89/98 (91%)    | 87 (98%)  | 2 (2%)   | 45          | 74  |
| 18  | L     | 102/103 (99%)  | 87 (85%)  | 15 (15%) | 3           | 15  |
| 19  | M     | 93/96 (97%)    | 86 (92%)  | 7 (8%)   | 12          | 41  |
| 20  | N     | 83/84 (99%)    | 77 (93%)  | 6 (7%)   | 13          | 43  |
| 21  | O     | 76/77 (99%)    | 68 (90%)  | 8 (10%)  | 6           | 27  |
| 22  | P     | 65/65 (100%)   | 60 (92%)  | 5 (8%)   | 12          | 40  |
| 23  | Q     | 73/78 (94%)    | 61 (84%)  | 12 (16%) | 2           | 12  |
| 24  | R     | 57/65 (88%)    | 56 (98%)  | 1 (2%)   | 51          | 77  |
| 25  | S     | 72/79 (91%)    | 59 (82%)  | 13 (18%) | 2           | 9   |
| 26  | T     | 65/66 (98%)    | 61 (94%)  | 4 (6%)   | 16          | 49  |
| 27  | U     | 60/61 (98%)    | 58 (97%)  | 2 (3%)   | 33          | 67  |
| 30  | c     | 216/218 (99%)  | 204 (94%) | 12 (6%)  | 19          | 52  |
| 31  | d     | 163/163 (100%) | 152 (93%) | 11 (7%)  | 15          | 46  |
| 32  | e     | 165/165 (100%) | 157 (95%) | 8 (5%)   | 23          | 57  |
| 33  | f     | 148/150 (99%)  | 142 (96%) | 6 (4%)   | 27          | 61  |
| 34  | g     | 137/138 (99%)  | 121 (88%) | 16 (12%) | 5           | 23  |

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| Mol | Chain | Analysed        | Rotameric  | Outliers | Percentiles |     |
|-----|-------|-----------------|------------|----------|-------------|-----|
| 35  | h     | 32/114 (28%)    | 32 (100%)  | 0        | 100         | 100 |
| 36  | i     | 116/116 (100%)  | 109 (94%)  | 7 (6%)   | 17          | 50  |
| 37  | j     | 104/104 (100%)  | 96 (92%)   | 8 (8%)   | 12          | 40  |
| 38  | k     | 103/103 (100%)  | 98 (95%)   | 5 (5%)   | 22          | 56  |
| 39  | l     | 107/107 (100%)  | 103 (96%)  | 4 (4%)   | 30          | 64  |
| 40  | m     | 98/103 (95%)    | 93 (95%)   | 5 (5%)   | 21          | 55  |
| 41  | n     | 86/87 (99%)     | 80 (93%)   | 6 (7%)   | 14          | 44  |
| 42  | o     | 99/100 (99%)    | 93 (94%)   | 6 (6%)   | 17          | 49  |
| 43  | p     | 89/90 (99%)     | 81 (91%)   | 8 (9%)   | 9           | 34  |
| 44  | q     | 84/84 (100%)    | 80 (95%)   | 4 (5%)   | 23          | 57  |
| 45  | r     | 93/93 (100%)    | 90 (97%)   | 3 (3%)   | 34          | 67  |
| 46  | s     | 80/84 (95%)     | 73 (91%)   | 7 (9%)   | 9           | 35  |
| 47  | t     | 83/85 (98%)     | 77 (93%)   | 6 (7%)   | 13          | 43  |
| 48  | u     | 78/78 (100%)    | 73 (94%)   | 5 (6%)   | 16          | 48  |
| 49  | v     | 58/63 (92%)     | 53 (91%)   | 5 (9%)   | 10          | 36  |
| 50  | w     | 67/68 (98%)     | 59 (88%)   | 8 (12%)  | 5           | 22  |
| 51  | x     | 54/55 (98%)     | 51 (94%)   | 3 (6%)   | 19          | 52  |
| 52  | y     | 48/49 (98%)     | 46 (96%)   | 2 (4%)   | 26          | 61  |
| 53  | z     | 47/48 (98%)     | 45 (96%)   | 2 (4%)   | 26          | 60  |
| All | All   | 4571/4825 (95%) | 4230 (92%) | 341 (8%) | 14          | 41  |

5 of 341 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 34  | g     | 4   | VAL  |
| 42  | o     | 47  | VAL  |
| 34  | g     | 60  | ASP  |
| 37  | j     | 91  | SER  |
| 45  | r     | 13  | SER  |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 44 such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 36  | i     | 40  | HIS  |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 41  | n     | 98  | GLN  |
| 36  | i     | 58  | ASN  |
| 37  | j     | 3   | GLN  |
| 44  | q     | 43  | ASN  |

### 5.3.3 RNA ⓘ

| Mol | Chain | Analysed        | Backbone Outliers | Pucker Outliers |
|-----|-------|-----------------|-------------------|-----------------|
| 28  | a     | 2747/2904 (94%) | 427 (15%)         | 0               |
| 29  | b     | 118/120 (98%)   | 36 (30%)          | 0               |
| 6   | 5     | 1/2 (50%)       | 1 (100%)          | 0               |
| 7   | A     | 1513/1517 (99%) | 564 (37%)         | 59 (3%)         |
| All | All   | 4379/4543 (96%) | 1028 (23%)        | 59 (1%)         |

5 of 1028 RNA backbone outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 6   | 5     | 76  | A    |
| 7   | A     | 4   | U    |
| 7   | A     | 6   | G    |
| 7   | A     | 7   | A    |
| 7   | A     | 8   | A    |

5 of 59 RNA pucker outliers are listed below:

| Mol | Chain | Res  | Type |
|-----|-------|------|------|
| 7   | A     | 996  | A    |
| 7   | A     | 1448 | A    |
| 7   | A     | 1040 | U    |
| 7   | A     | 1370 | C    |
| 7   | A     | 1306 | G    |

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

40 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$ |
| 28  | PSU  | a     | 1911 | 28    | 18,21,22     | 0.95 | 1 (5%)      | 21,30,33    | 1.98 | 4 (19%)     |
| 7   | 2MG  | A     | 1208 | 7     | 23,26,27     | 0.78 | 0           | 33,38,41    | 2.19 | 10 (30%)    |
| 28  | 6MZ  | a     | 2030 | 28    | 22,25,26     | 1.16 | 2 (9%)      | 29,36,39    | 2.21 | 9 (31%)     |
| 28  | PSU  | a     | 746  | 55,28 | 18,21,22     | 1.15 | 3 (16%)     | 21,30,33    | 1.94 | 4 (19%)     |
| 28  | OMU  | a     | 2552 | 28    | 19,22,23     | 1.11 | 3 (15%)     | 25,31,34    | 2.06 | 6 (24%)     |
| 7   | 5MC  | A     | 1408 | 7     | 19,22,23     | 1.37 | 2 (10%)     | 26,32,35    | 1.15 | 3 (11%)     |
| 7   | 4OC  | A     | 1403 | 7     | 20,23,24     | 0.36 | 0           | 25,32,35    | 0.65 | 1 (4%)      |
| 28  | 1MG  | a     | 745  | 28    | 23,26,27     | 1.05 | 1 (4%)      | 33,39,42    | 1.76 | 6 (18%)     |
| 28  | H2U  | a     | 2449 | 28    | 18,21,22     | 0.49 | 0           | 19,30,33    | 1.09 | 1 (5%)      |
| 28  | PSU  | a     | 2604 | 28    | 18,21,22     | 1.10 | 3 (16%)     | 21,30,33    | 2.02 | 4 (19%)     |
| 28  | G7M  | a     | 2069 | 28    | 23,26,27     | 0.58 | 0           | 34,39,42    | 0.93 | 2 (5%)      |
| 7   | MA6  | A     | 1520 | 7     | 23,26,27     | 2.51 | 5 (21%)     | 33,38,41    | 2.78 | 7 (21%)     |
| 28  | 3TD  | a     | 1915 | -     | 19,22,23     | 1.10 | 2 (10%)     | 23,32,35    | 1.95 | 3 (13%)     |
| 28  | 6MZ  | a     | 1618 | 28    | 22,25,26     | 1.11 | 2 (9%)      | 29,36,39    | 2.29 | 9 (31%)     |
| 28  | OMC  | a     | 2498 | 55,28 | 19,22,23     | 1.00 | 1 (5%)      | 25,31,34    | 1.28 | 2 (8%)      |
| 28  | 2MG  | a     | 1835 | 28    | 23,26,27     | 0.80 | 1 (4%)      | 33,38,41    | 2.30 | 11 (33%)    |
| 28  | PSU  | a     | 2580 | 28    | 18,21,22     | 1.22 | 3 (16%)     | 21,30,33    | 2.21 | 5 (23%)     |
| 7   | G7M  | A     | 527  | 7     | 23,26,27     | 0.51 | 0           | 34,39,42    | 0.94 | 1 (2%)      |
| 28  | PSU  | a     | 1917 | 28    | 18,21,22     | 1.00 | 1 (5%)      | 21,30,33    | 1.98 | 5 (23%)     |
| 28  | 5MU  | a     | 1939 | 28    | 19,22,23     | 1.21 | 4 (21%)     | 27,32,35    | 2.19 | 6 (22%)     |
| 28  | 5MU  | a     | 747  | 28    | 19,22,23     | 1.13 | 3 (15%)     | 27,32,35    | 2.16 | 7 (25%)     |
| 39  | MS6  | l     | 82   | 39    | 5,7,8        | 0.59 | 0           | 2,7,9       | 1.09 | 0           |
| 7   | 2MG  | A     | 1517 | 7     | 23,26,27     | 0.79 | 0           | 33,38,41    | 2.28 | 11 (33%)    |
| 28  | PSU  | a     | 2504 | 28    | 18,21,22     | 1.01 | 1 (5%)      | 21,30,33    | 1.92 | 4 (19%)     |
| 28  | PSU  | a     | 2457 | 28    | 18,21,22     | 1.14 | 4 (22%)     | 21,30,33    | 2.14 | 5 (23%)     |
| 7   | PSU  | A     | 516  | 7     | 18,21,22     | 0.99 | 1 (5%)      | 21,30,33    | 2.00 | 5 (23%)     |
| 18  | D2T  | L     | 89   | 18    | 8,9,10       | 1.78 | 2 (25%)     | 6,11,13     | 1.80 | 2 (33%)     |
| 39  | 4D4  | l     | 81   | 39    | 9,11,12      | 2.10 | 2 (22%)     | 7,13,15     | 2.19 | 3 (42%)     |
| 7   | 5MC  | A     | 967  | 7     | 19,22,23     | 1.28 | 2 (10%)     | 26,32,35    | 1.08 | 2 (7%)      |
| 31  | MEQ  | d     | 150  | 31    | 8,9,10       | 0.83 | 0           | 5,10,12     | 0.37 | 0           |
| 17  | BH2  | K     | 119  | 17    | 5,6,9        | 0.93 | 0           | 1,6,12      | 0.62 | 0           |
| 28  | OMG  | a     | 2251 | 28    | 23,26,27     | 0.77 | 1 (4%)      | 32,38,41    | 1.99 | 9 (28%)     |
| 28  | 2MG  | a     | 2445 | 28    | 23,26,27     | 0.91 | 2 (8%)      | 33,38,41    | 2.23 | 9 (27%)     |

| Mol | Type | Chain | Res  | Link  | Bond lengths |      |          | Bond angles |      |          |
|-----|------|-------|------|-------|--------------|------|----------|-------------|------|----------|
|     |      |       |      |       | Counts       | RMSZ | # Z  > 2 | Counts      | RMSZ | # Z  > 2 |
| 28  | 2MA  | a     | 2503 | 55,28 | 22,25,26     | 1.30 | 3 (13%)  | 32,37,40    | 2.07 | 7 (21%)  |
| 28  | 5MC  | a     | 1962 | 28    | 19,22,23     | 1.53 | 2 (10%)  | 26,32,35    | 1.20 | 2 (7%)   |
| 7   | 2MG  | A     | 966  | 7     | 23,26,27     | 0.74 | 1 (4%)   | 33,38,41    | 2.21 | 10 (30%) |
| 7   | UR3  | A     | 1499 | 7     | 19,22,23     | 3.89 | 6 (31%)  | 26,32,35    | 4.65 | 11 (42%) |
| 7   | MA6  | A     | 1519 | 7     | 23,26,27     | 2.53 | 5 (21%)  | 33,38,41    | 2.80 | 7 (21%)  |
| 28  | PSU  | a     | 2605 | 28    | 18,21,22     | 1.13 | 2 (11%)  | 21,30,33    | 1.97 | 3 (14%)  |
| 28  | PSU  | a     | 955  | 28    | 18,21,22     | 1.16 | 3 (16%)  | 21,30,33    | 2.08 | 4 (19%)  |

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   |
|-----|------|-------|------|-------|---------|------------|---------|
| 28  | PSU  | a     | 1911 | 28    | -       | 0/7/25/26  | 0/2/2/2 |
| 7   | 2MG  | A     | 1208 | 7     | -       | 0/9/27/28  | 0/3/3/3 |
| 28  | 6MZ  | a     | 2030 | 28    | -       | 2/9/27/28  | 0/3/3/3 |
| 28  | PSU  | a     | 746  | 55,28 | -       | 1/7/25/26  | 0/2/2/2 |
| 28  | OMU  | a     | 2552 | 28    | -       | 0/9/27/28  | 0/2/2/2 |
| 7   | 5MC  | A     | 1408 | 7     | -       | 0/7/25/26  | 0/2/2/2 |
| 7   | 4OC  | A     | 1403 | 7     | -       | 2/9/29/30  | 0/2/2/2 |
| 28  | 1MG  | a     | 745  | 28    | -       | 0/7/25/26  | 0/3/3/3 |
| 28  | H2U  | a     | 2449 | 28    | -       | 0/7/38/39  | 0/2/2/2 |
| 28  | PSU  | a     | 2604 | 28    | -       | 1/7/25/26  | 0/2/2/2 |
| 28  | G7M  | a     | 2069 | 28    | -       | 2/7/25/26  | 0/3/3/3 |
| 7   | MA6  | A     | 1520 | 7     | -       | 2/11/29/30 | 0/3/3/3 |
| 28  | 3TD  | a     | 1915 | -     | -       | 4/7/25/26  | 0/2/2/2 |
| 28  | 6MZ  | a     | 1618 | 28    | -       | 0/9/27/28  | 0/3/3/3 |
| 28  | OMC  | a     | 2498 | 55,28 | -       | 3/9/27/28  | 0/2/2/2 |
| 28  | 2MG  | a     | 1835 | 28    | -       | 0/9/27/28  | 0/3/3/3 |
| 28  | PSU  | a     | 2580 | 28    | -       | 1/7/25/26  | 0/2/2/2 |
| 7   | G7M  | A     | 527  | 7     | -       | 2/7/25/26  | 0/3/3/3 |
| 28  | PSU  | a     | 1917 | 28    | -       | 0/7/25/26  | 0/2/2/2 |
| 28  | 5MU  | a     | 1939 | 28    | -       | 0/7/25/26  | 0/2/2/2 |
| 28  | 5MU  | a     | 747  | 28    | -       | 0/7/25/26  | 0/2/2/2 |
| 39  | MS6  | l     | 82   | 39    | -       | 2/4/6/8    | -       |
| 7   | 2MG  | A     | 1517 | 7     | -       | 0/9/27/28  | 0/3/3/3 |
| 28  | PSU  | a     | 2504 | 28    | -       | 2/7/25/26  | 0/2/2/2 |
| 28  | PSU  | a     | 2457 | 28    | -       | 0/7/25/26  | 0/2/2/2 |

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| Mol | Type | Chain | Res  | Link  | Chirals | Torsions   | Rings   |
|-----|------|-------|------|-------|---------|------------|---------|
| 7   | PSU  | A     | 516  | 7     | -       | 1/7/25/26  | 0/2/2/2 |
| 18  | D2T  | L     | 89   | 18    | -       | 5/7/12/14  | -       |
| 39  | 4D4  | l     | 81   | 39    | -       | 2/11/12/14 | -       |
| 7   | 5MC  | A     | 967  | 7     | -       | 1/7/25/26  | 0/2/2/2 |
| 31  | MEQ  | d     | 150  | 31    | -       | 3/8/9/11   | -       |
| 17  | BH2  | K     | 119  | 17    | 1/1/1/4 | 1/4/5/12   | -       |
| 28  | OMG  | a     | 2251 | 28    | -       | 1/9/27/28  | 0/3/3/3 |
| 28  | 2MG  | a     | 2445 | 28    | -       | 0/9/27/28  | 0/3/3/3 |
| 28  | 2MA  | a     | 2503 | 55,28 | -       | 2/7/25/26  | 0/3/3/3 |
| 28  | 5MC  | a     | 1962 | 28    | -       | 0/7/25/26  | 0/2/2/2 |
| 7   | 2MG  | A     | 966  | 7     | -       | 0/9/27/28  | 0/3/3/3 |
| 7   | UR3  | A     | 1499 | 7     | -       | 2/7/25/26  | 0/2/2/2 |
| 7   | MA6  | A     | 1519 | 7     | -       | 3/11/29/30 | 0/3/3/3 |
| 28  | PSU  | a     | 2605 | 28    | -       | 0/7/25/26  | 0/2/2/2 |
| 28  | PSU  | a     | 955  | 28    | -       | 0/7/25/26  | 0/2/2/2 |

The worst 5 of 74 bond length outliers are listed below:

| Mol | Chain | Res  | Type | Atoms | Z     | Observed(Å) | Ideal(Å) |
|-----|-------|------|------|-------|-------|-------------|----------|
| 7   | A     | 1499 | UR3  | C2-N1 | 13.96 | 1.57        | 1.38     |
| 7   | A     | 1520 | MA6  | C5-N7 | 7.70  | 1.53        | 1.39     |
| 7   | A     | 1519 | MA6  | C5-N7 | 7.66  | 1.53        | 1.39     |
| 7   | A     | 1499 | UR3  | C6-N1 | 6.20  | 1.52        | 1.38     |
| 7   | A     | 1499 | UR3  | C2-N3 | 5.71  | 1.50        | 1.39     |

The worst 5 of 200 bond angle outliers are listed below:

| Mol | Chain | Res  | Type | Atoms    | Z      | Observed(°) | Ideal(°) |
|-----|-------|------|------|----------|--------|-------------|----------|
| 7   | A     | 1499 | UR3  | C6-N1-C2 | -14.80 | 109.71      | 121.80   |
| 7   | A     | 1499 | UR3  | C4-N3-C2 | -14.23 | 113.13      | 124.58   |
| 7   | A     | 1519 | MA6  | C4-N9-C8 | 13.53  | 119.95      | 105.74   |
| 7   | A     | 1520 | MA6  | C4-N9-C8 | 13.38  | 119.79      | 105.74   |
| 28  | a     | 1835 | 2MG  | C2-N3-C4 | 7.68   | 121.60      | 112.00   |

All (1) chirality outliers are listed below:

| Mol | Chain | Res | Type | Atom |
|-----|-------|-----|------|------|
| 17  | K     | 119 | BH2  | CB   |

5 of 45 torsion outliers are listed below:

| Mol | Chain | Res  | Type | Atoms           |
|-----|-------|------|------|-----------------|
| 18  | L     | 89   | D2T  | O-C-CA-CB       |
| 18  | L     | 89   | D2T  | CG-CB-SB-CB1    |
| 7   | A     | 527  | G7M  | O4'-C4'-C5'-O5' |
| 17  | K     | 119  | BH2  | CA-CB-CG-OD1    |
| 28  | a     | 1915 | 3TD  | O4'-C1'-C5-C4   |

There are no ring outliers.

8 monomers are involved in 10 short contacts:

| Mol | Chain | Res  | Type | Clashes | Symm-Clashes |
|-----|-------|------|------|---------|--------------|
| 7   | A     | 1403 | 4OC  | 3       | 0            |
| 7   | A     | 1520 | MA6  | 1       | 0            |
| 28  | a     | 1917 | PSU  | 1       | 0            |
| 7   | A     | 1517 | 2MG  | 1       | 0            |
| 18  | L     | 89   | D2T  | 1       | 0            |
| 7   | A     | 967  | 5MC  | 2       | 0            |
| 28  | a     | 2251 | OMG  | 1       | 0            |
| 7   | A     | 1519 | MA6  | 1       | 0            |

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 289 ligands modelled in this entry, 274 are monoatomic - leaving 15 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 | $\# Z  > 2$ |
| 57  | SPM  | a     | 6221 | -    | 13,13,13     | 0.34 | 0           | 12,12,12    | 0.90 | 0           |
| 56  | SPD  | a     | 6211 | -    | 9,9,9        | 0.32 | 0           | 8,8,8       | 0.76 | 0           |
| 56  | SPD  | a     | 6219 | -    | 9,9,9        | 0.36 | 0           | 8,8,8       | 0.82 | 0           |
| 56  | SPD  | a     | 6210 | -    | 9,9,9        | 0.31 | 0           | 8,8,8       | 0.85 | 0           |
| 56  | SPD  | a     | 6213 | -    | 9,9,9        | 0.32 | 0           | 8,8,8       | 0.83 | 0           |
| 56  | SPD  | a     | 6209 | -    | 9,9,9        | 0.33 | 0           | 8,8,8       | 0.70 | 0           |

| Mol | Type | Chain | Res  | Link | Bond lengths |      |          | Bond angles |      |          |
|-----|------|-------|------|------|--------------|------|----------|-------------|------|----------|
|     |      |       |      |      | Counts       | RMSZ | # Z  > 2 | Counts      | RMSZ | # Z  > 2 |
| 56  | SPD  | a     | 6214 | -    | 9,9,9        | 0.33 | 0        | 8,8,8       | 0.96 | 0        |
| 56  | SPD  | a     | 6215 | -    | 9,9,9        | 0.33 | 0        | 8,8,8       | 0.82 | 0        |
| 56  | SPD  | a     | 6220 | -    | 9,9,9        | 0.33 | 0        | 8,8,8       | 0.79 | 0        |
| 56  | SPD  | a     | 6208 | -    | 9,9,9        | 0.32 | 0        | 8,8,8       | 0.77 | 0        |
| 56  | SPD  | a     | 6212 | -    | 9,9,9        | 0.33 | 0        | 8,8,8       | 0.81 | 0        |
| 56  | SPD  | a     | 6218 | -    | 9,9,9        | 0.29 | 0        | 8,8,8       | 0.83 | 0        |
| 56  | SPD  | a     | 6216 | -    | 9,9,9        | 0.32 | 0        | 8,8,8       | 0.83 | 0        |
| 56  | SPD  | a     | 6217 | -    | 9,9,9        | 0.33 | 0        | 8,8,8       | 0.83 | 0        |
| 56  | SPD  | a     | 6207 | -    | 9,9,9        | 0.33 | 0        | 8,8,8       | 0.81 | 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 |
|-----|------|-------|------|------|---------|------------|-------|
| 57  | SPM  | a     | 6221 | -    | -       | 3/11/11/11 | -     |
| 56  | SPD  | a     | 6211 | -    | -       | 0/7/7/7    | -     |
| 56  | SPD  | a     | 6219 | -    | -       | 2/7/7/7    | -     |
| 56  | SPD  | a     | 6210 | -    | -       | 0/7/7/7    | -     |
| 56  | SPD  | a     | 6213 | -    | -       | 1/7/7/7    | -     |
| 56  | SPD  | a     | 6209 | -    | -       | 2/7/7/7    | -     |
| 56  | SPD  | a     | 6214 | -    | -       | 1/7/7/7    | -     |
| 56  | SPD  | a     | 6215 | -    | -       | 2/7/7/7    | -     |
| 56  | SPD  | a     | 6220 | -    | -       | 1/7/7/7    | -     |
| 56  | SPD  | a     | 6208 | -    | -       | 3/7/7/7    | -     |
| 56  | SPD  | a     | 6212 | -    | -       | 2/7/7/7    | -     |
| 56  | SPD  | a     | 6218 | -    | -       | 1/7/7/7    | -     |
| 56  | SPD  | a     | 6216 | -    | -       | 2/7/7/7    | -     |
| 56  | SPD  | a     | 6217 | -    | -       | 3/7/7/7    | -     |
| 56  | SPD  | a     | 6207 | -    | -       | 3/7/7/7    | -     |

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 26 torsion outliers are listed below:

| Mol | Chain | Res  | Type | Atoms           |
|-----|-------|------|------|-----------------|
| 57  | a     | 6221 | SPM  | N10-C11-C12-C13 |

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| Mol | Chain | Res  | Type | Atoms       |
|-----|-------|------|------|-------------|
| 56  | a     | 6209 | SPD  | C8-C7-N6-C5 |
| 56  | a     | 6215 | SPD  | C8-C7-N6-C5 |
| 57  | a     | 6221 | SPM  | N5-C6-C7-C8 |
| 56  | a     | 6208 | SPD  | C4-C5-N6-C7 |

There are no ring outliers.

10 monomers are involved in 13 short contacts:

| Mol | Chain | Res  | Type | Clashes | Symm-Clashes |
|-----|-------|------|------|---------|--------------|
| 57  | a     | 6221 | SPM  | 1       | 0            |
| 56  | a     | 6211 | SPD  | 2       | 0            |
| 56  | a     | 6219 | SPD  | 1       | 0            |
| 56  | a     | 6210 | SPD  | 2       | 0            |
| 56  | a     | 6214 | SPD  | 2       | 0            |
| 56  | a     | 6215 | SPD  | 1       | 0            |
| 56  | a     | 6220 | SPD  | 1       | 0            |
| 56  | a     | 6208 | SPD  | 1       | 0            |
| 56  | a     | 6218 | SPD  | 1       | 0            |
| 56  | a     | 6217 | SPD  | 1       | 0            |

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

The following chains have linkage breaks:

| Mol | Chain | Number of breaks |
|-----|-------|------------------|
| 7   | A     | 3                |
| 28  | a     | 2                |
| 18  | L     | 1                |

The worst 5 of 6 chain breaks are listed below:

| Model | Chain | Residue-1 | Atom-1 | Residue-2 | Atom-2 | Distance (Å) |
|-------|-------|-----------|--------|-----------|--------|--------------|
| 1     | A     | 1267:G    | O3'    | 1271:G    | P      | 15.18        |
| 1     | A     | 840:C     | O3'    | 846:G     | P      | 13.53        |
| 1     | A     | 204:G     | O3'    | 214:C     | P      | 10.50        |
| 1     | a     | 1914:C    | O3'    | 1915:3TD  | P      | 5.23         |

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| Model | Chain | Residue-1 | Atom-1 | Residue-2 | Atom-2 | Distance (Å) |
|-------|-------|-----------|--------|-----------|--------|--------------|
| 1     | a     | 1915:3TD  | O3'    | 1916:A    | P      | 4.29         |

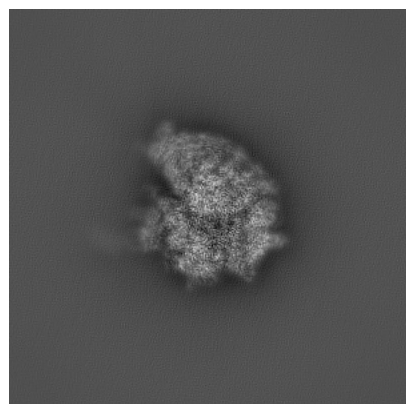
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-75634. These allow visual inspection of the internal detail of the map and identification of artifacts.

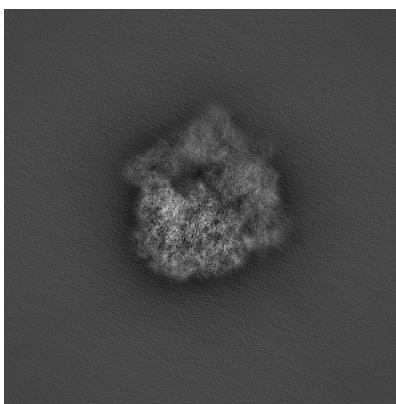
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

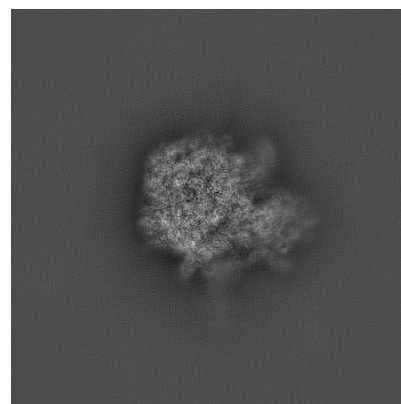
#### 6.1.1 Primary map



X

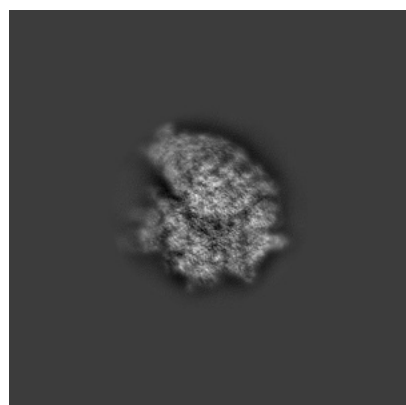


Y

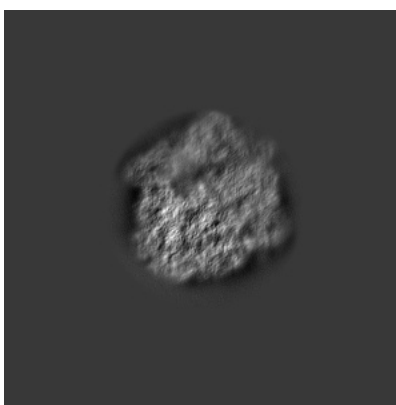


Z

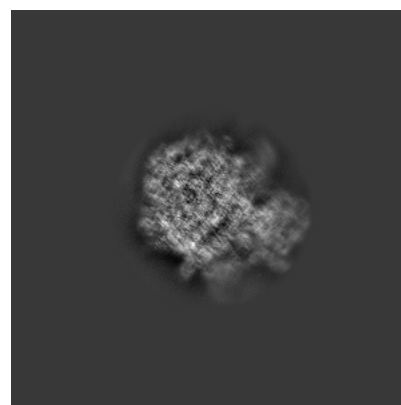
#### 6.1.2 Raw map



X



Y

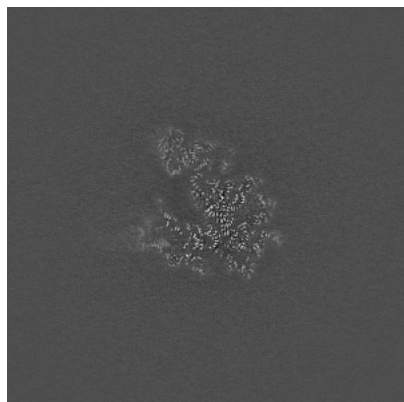


Z

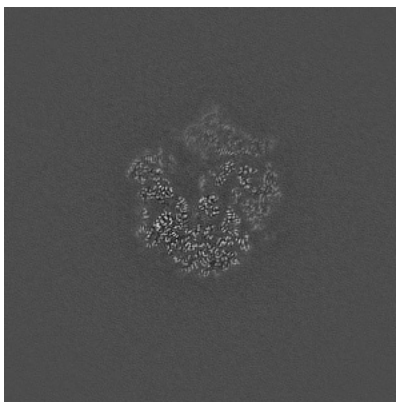
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

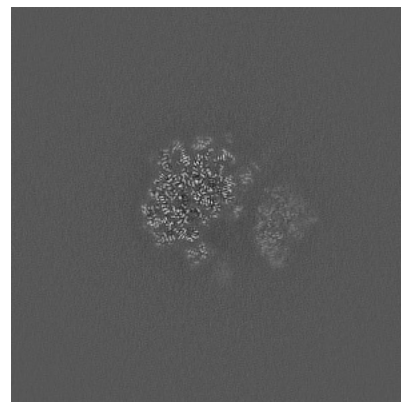
### 6.2.1 Primary map



X Index: 240

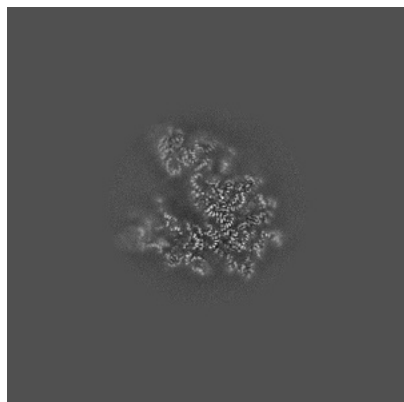


Y Index: 240

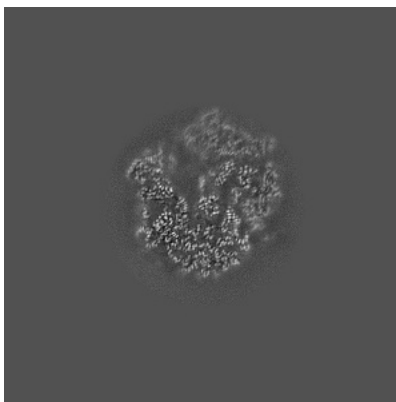


Z Index: 240

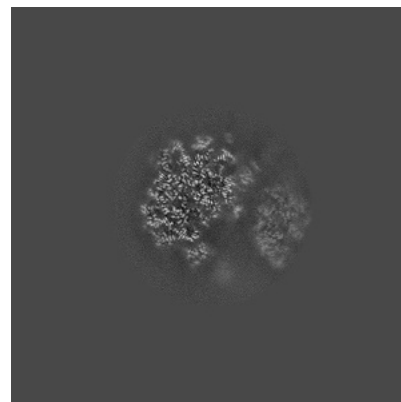
### 6.2.2 Raw map



X Index: 240



Y Index: 240

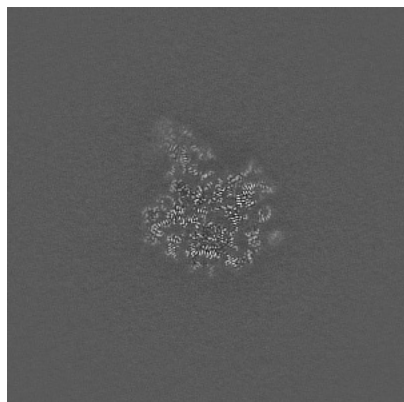


Z Index: 240

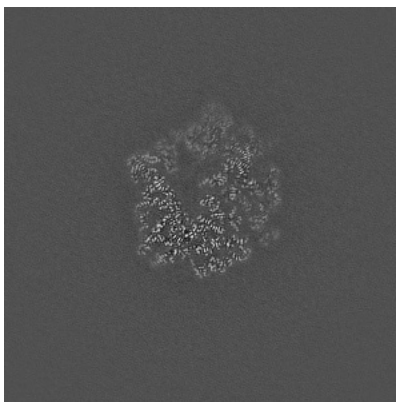
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

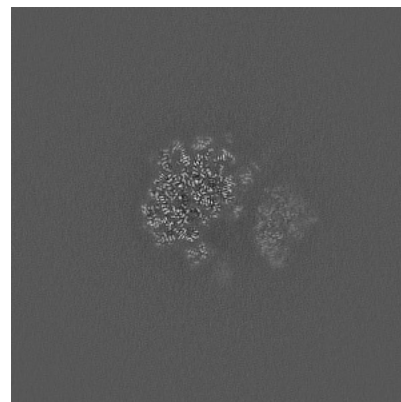
### 6.3.1 Primary map



X Index: 218

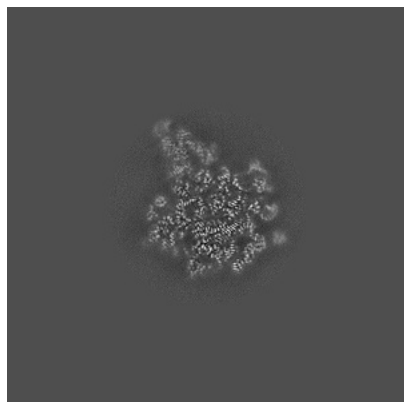


Y Index: 234

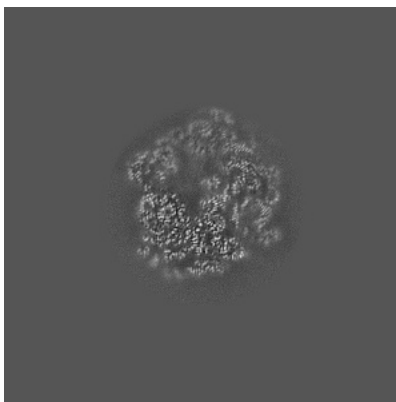


Z Index: 240

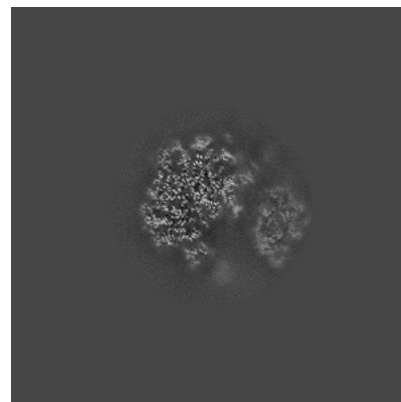
### 6.3.2 Raw map



X Index: 224



Y Index: 229

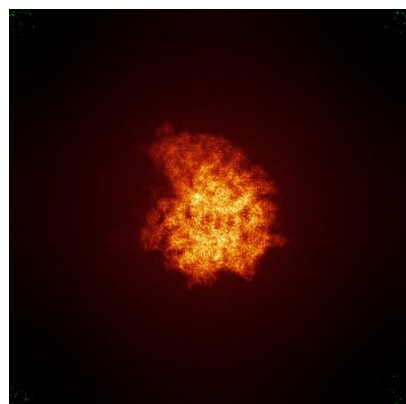


Z Index: 239

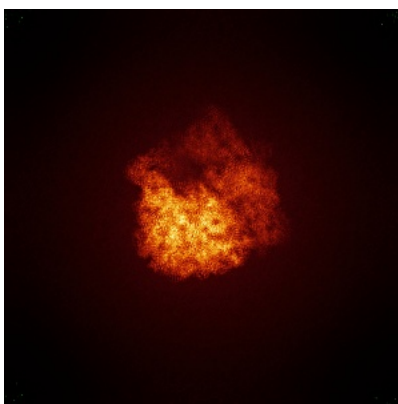
The images above show the largest variance slices of the map in three orthogonal directions.

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

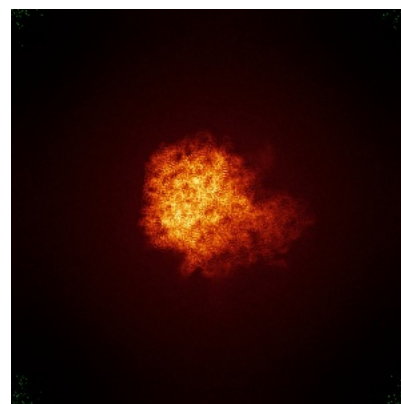
### 6.4.1 Primary map



X

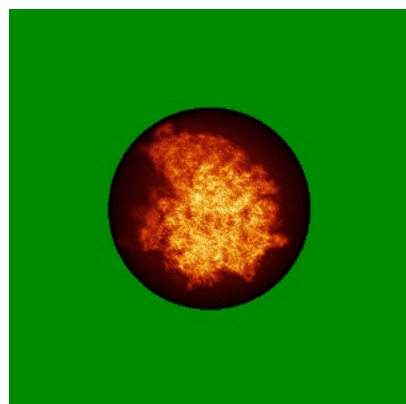


Y

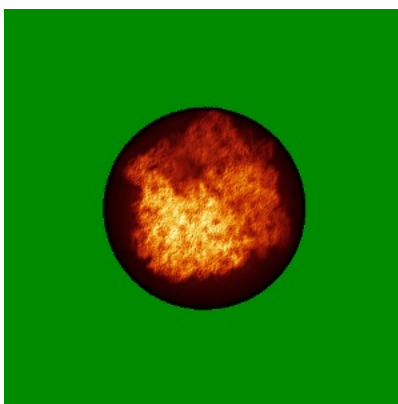


Z

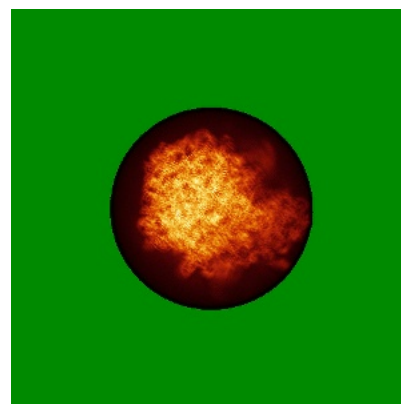
### 6.4.2 Raw map



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 2.8. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.5.2 Raw map



X



Y



Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

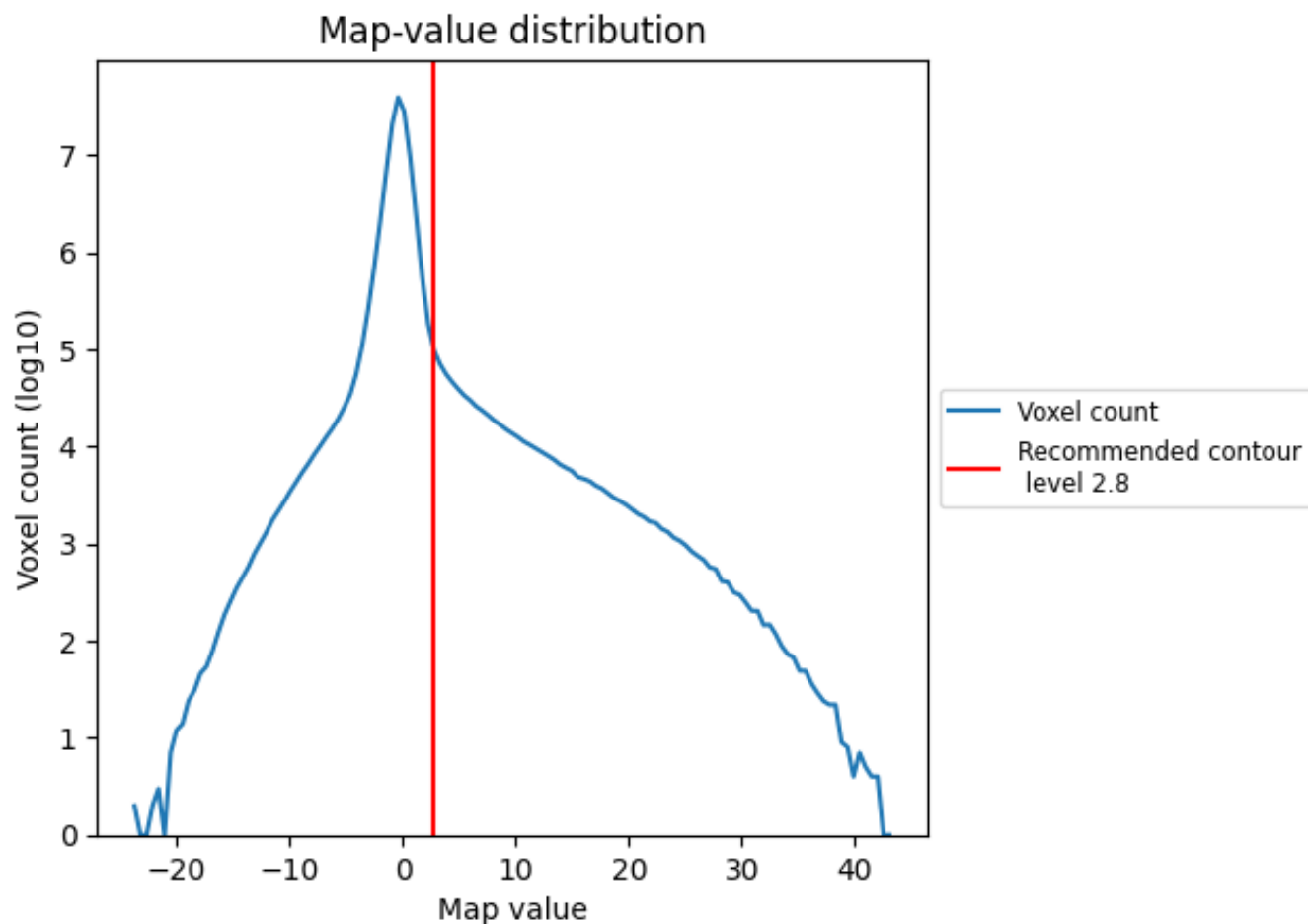
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

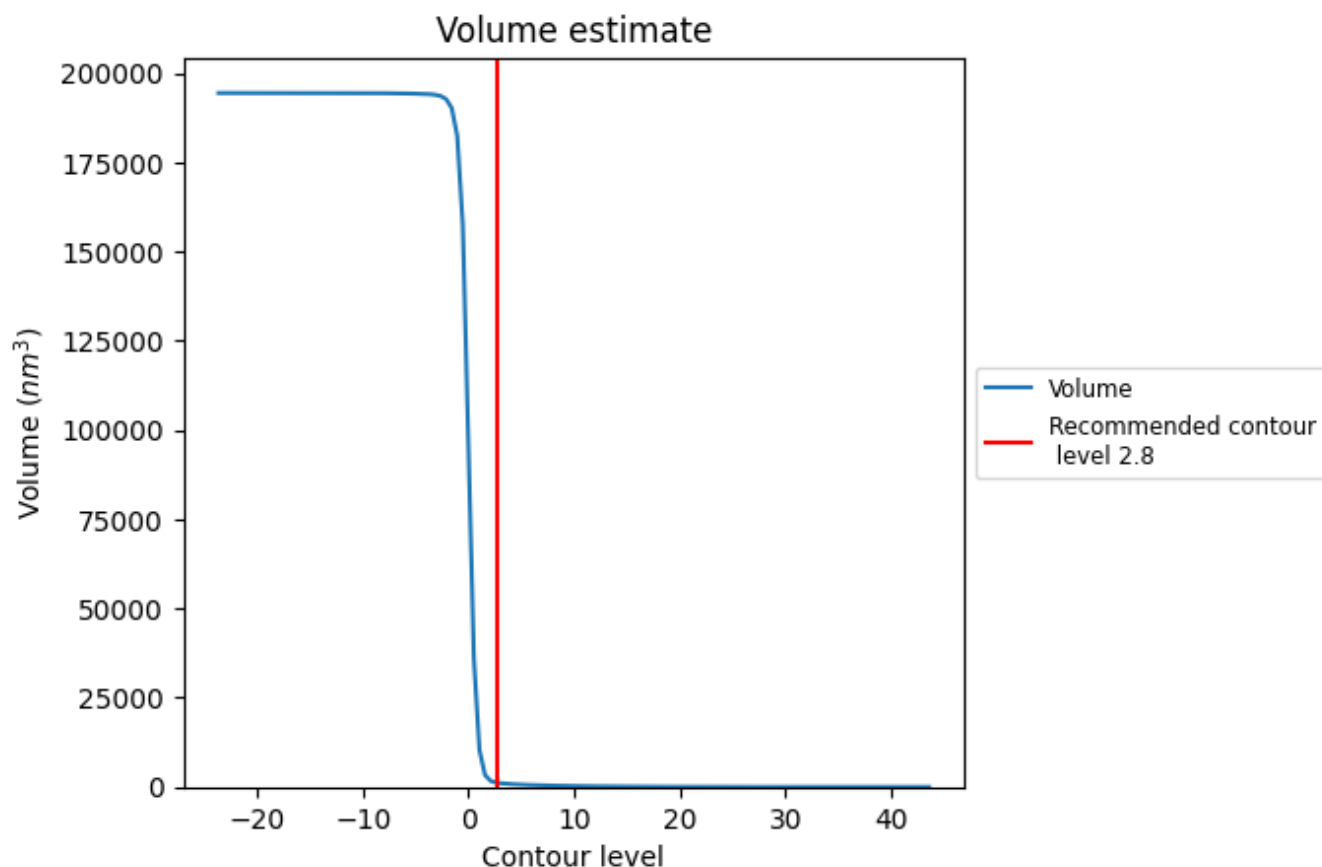
This section contains the results of statistical analysis of the map.

### 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.

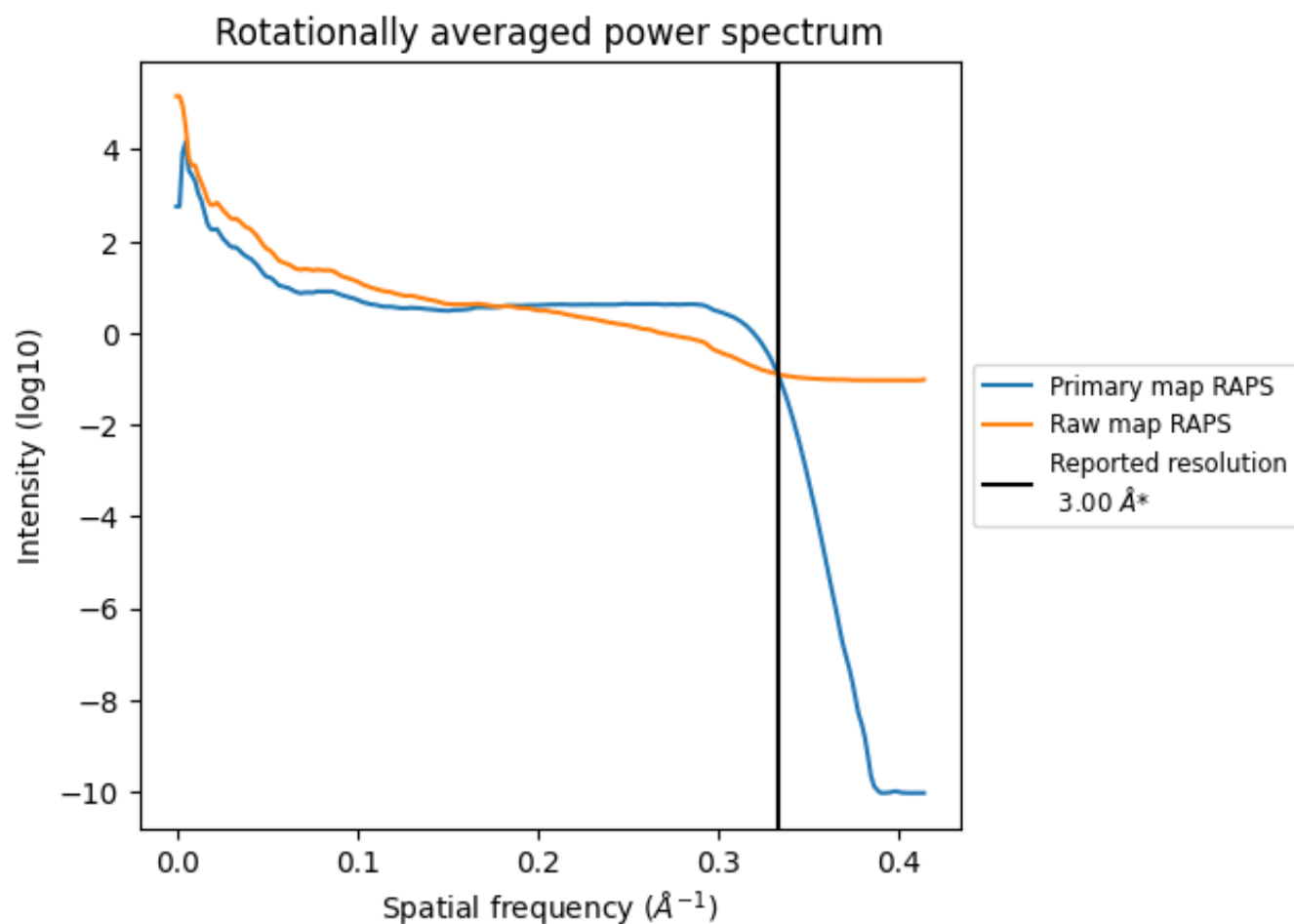
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1155  $\text{nm}^3$ ; this corresponds to an approximate mass of 1044 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 ⓘ

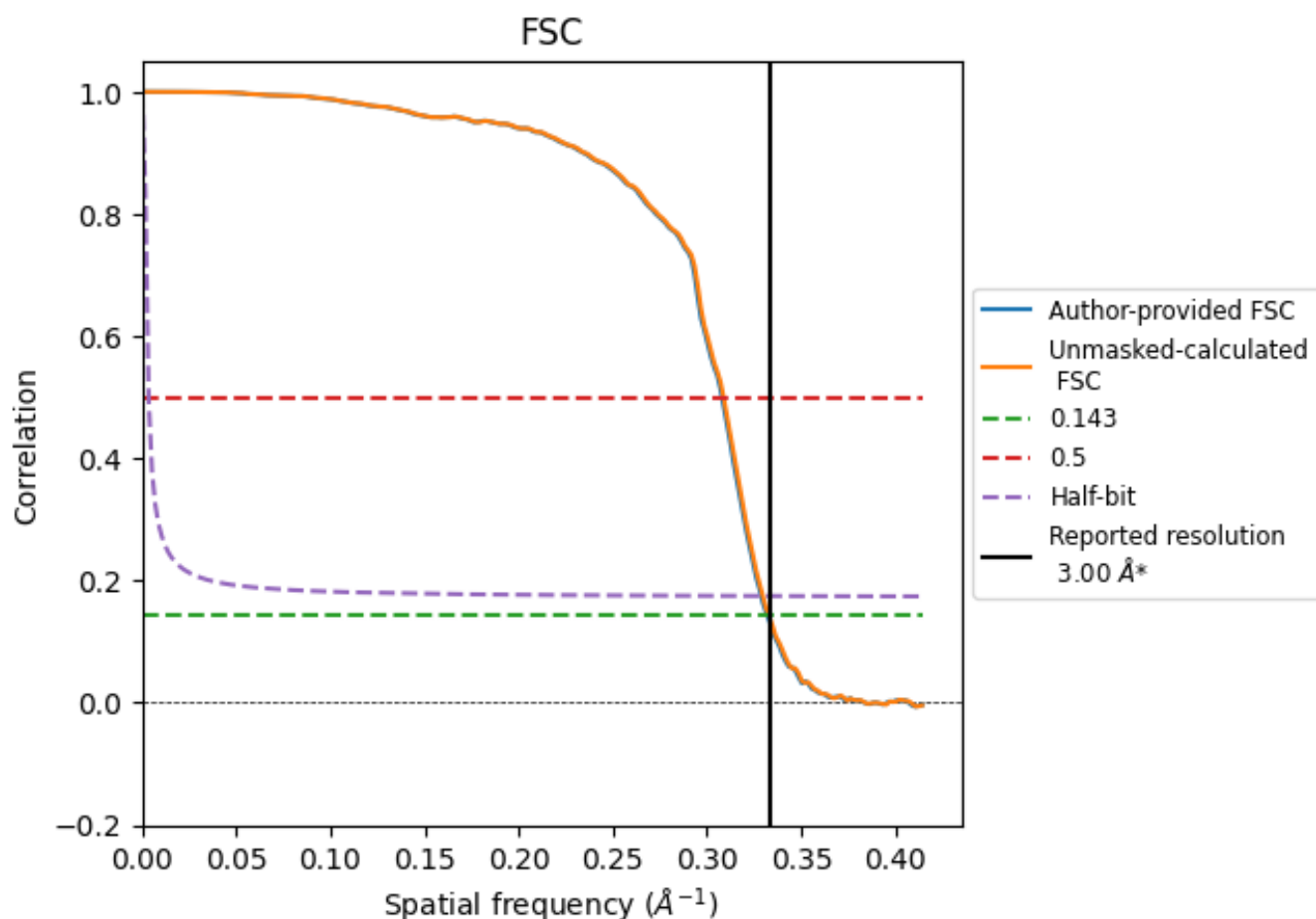


\*Reported resolution corresponds to spatial frequency of 0.333 Å<sup>-1</sup>

## 8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

### 8.1 FSC [i](#)



\*Reported resolution corresponds to spatial frequency of 0.333 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

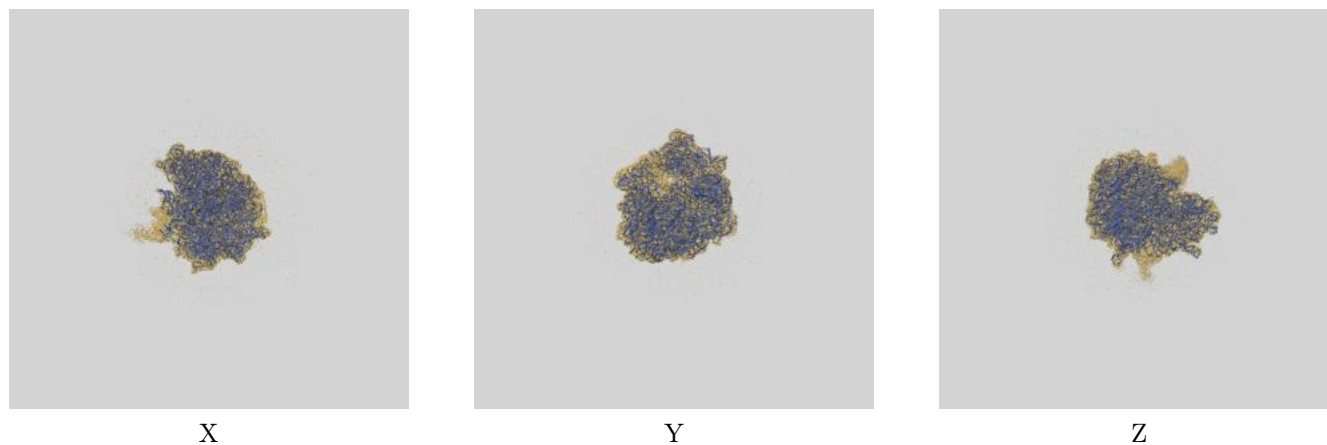
| Resolution estimate (Å)   | Estimation criterion (FSC cut-off) |      |          |
|---------------------------|------------------------------------|------|----------|
|                           | 0.143                              | 0.5  | Half-bit |
| Reported by author        | 3.00                               | -    | -        |
| Author-provided FSC curve | 3.01                               | 3.25 | 3.04     |
| Unmasked-calculated*      | 3.00                               | 3.24 | 3.03     |

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

## 9 Map-model fit [i](#)

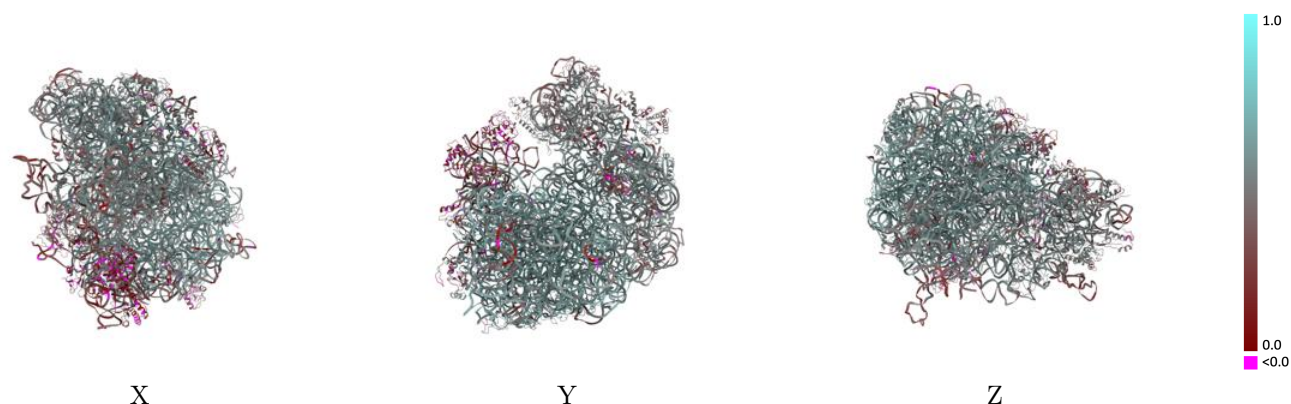
This section contains information regarding the fit between EMDB map EMD-75634 and PDB model 11DG. Per-residue inclusion information can be found in section [3](#) on page [17](#).

### 9.1 Map-model overlay [i](#)



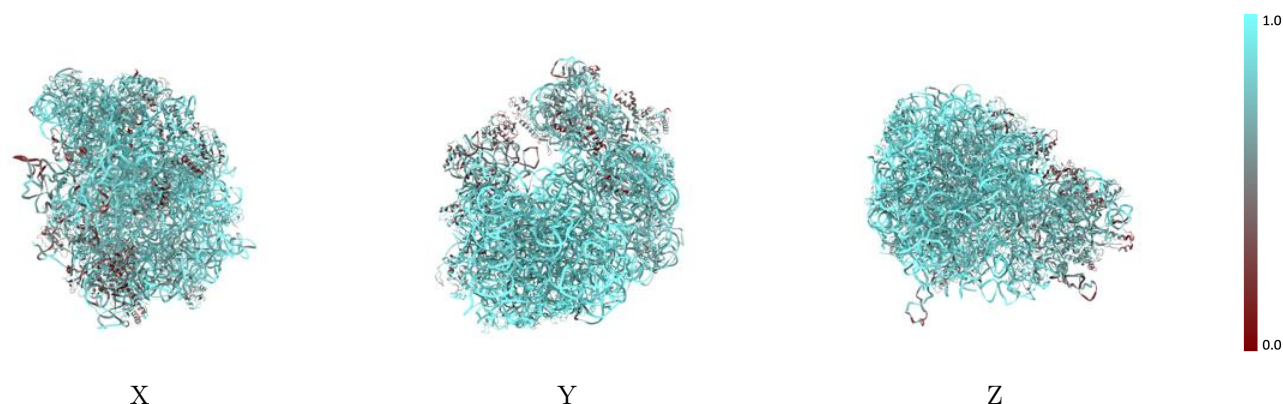
The images above show the 3D surface view of the map at the recommended contour level 2.8 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



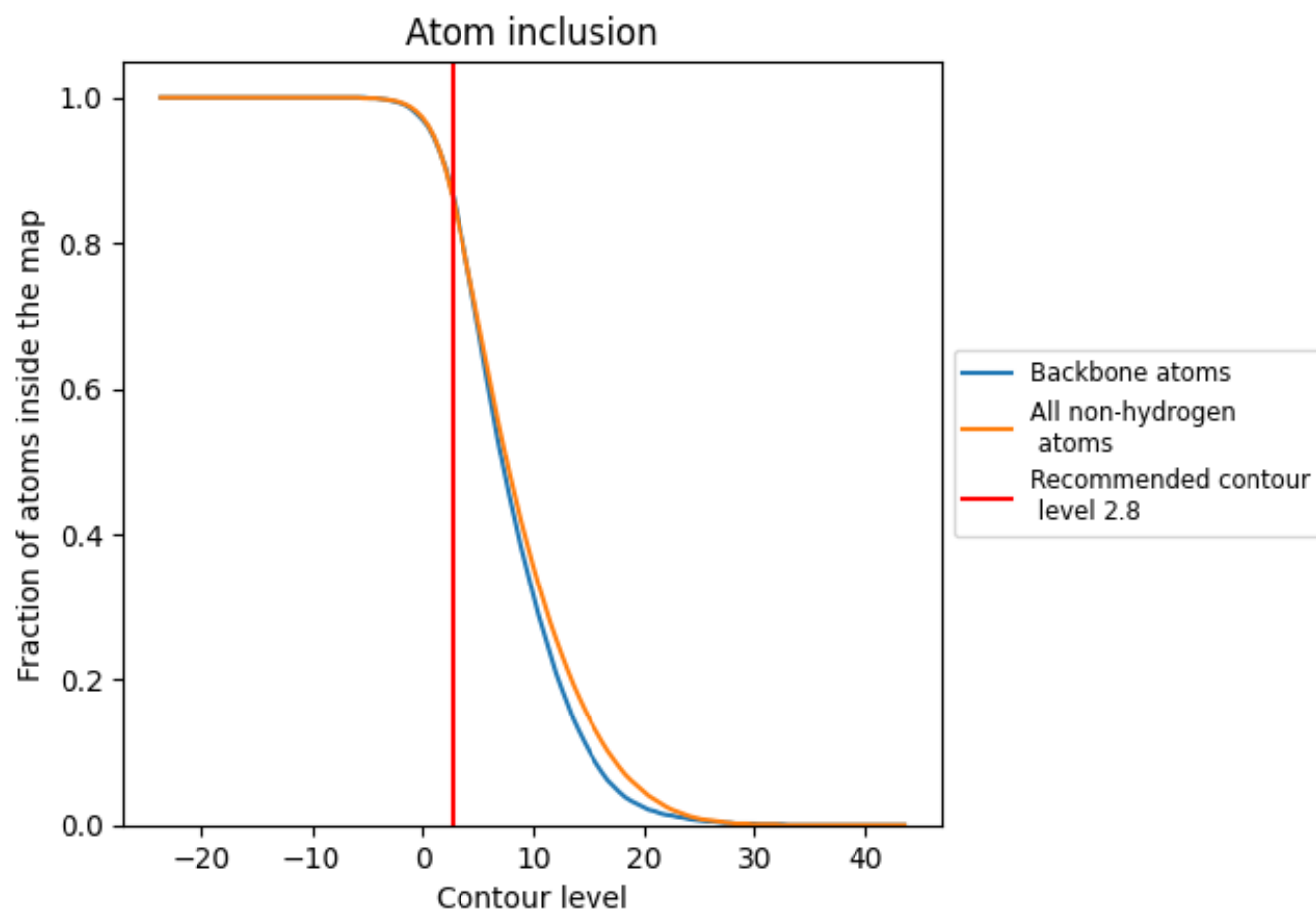
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (2.8).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 86% of all backbone atoms, 86% 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 (2.8) and Q-score for the entire model and for each chain.

| Chain | Atom inclusion                                                                             | Q-score                                                                                    |
|-------|--------------------------------------------------------------------------------------------|--------------------------------------------------------------------------------------------|
| All   |  0.8590   |  0.5060   |
| 0     |  0.5280   |  0.2890   |
| 1     |  0.9270   |  0.6130   |
| 2     |  0.8900   |  0.5760   |
| 3     |  0.6210   |  0.2870   |
| 4     |  0.4700   |  0.2220   |
| 5     |  0.5950   |  0.5370   |
| A     |  0.8880   |  0.5080   |
| B     |  0.5610   |  0.4390   |
| C     |  0.6610   |  0.4910   |
| D     |  0.6900   |  0.4990   |
| E     |  0.8300   |  0.5460   |
| F     |  0.5260   |  0.2720   |
| G     |  0.4750   |  0.4070   |
| H     |  0.8370  |  0.5570  |
| I     |  0.7250 |  0.4660 |
| J     |  0.5270 |  0.4340 |
| K     |  0.8120 |  0.5110 |
| L     |  0.8450 |  0.5510 |
| M     |  0.5470 |  0.3970 |
| N     |  0.7430 |  0.5010 |
| O     |  0.7670 |  0.4440 |
| P     |  0.7810 |  0.5160 |
| Q     |  0.7870 |  0.5260 |
| R     |  0.7250 |  0.3960 |
| S     |  0.5910 |  0.4310 |
| T     |  0.7940 |  0.5150 |
| U     |  0.3560 |  0.3510 |
| a     |  0.9390 |  0.5480 |
| b     |  0.6440 |  0.2780 |
| c     |  0.9470 |  0.6030 |
| d     |  0.8730 |  0.5300 |
| e     |  0.8300 |  0.4840 |
| f     |  0.4530 |  0.1560 |
| g     |  0.5260 |  0.2010 |



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| Chain | Atom inclusion                                                                           | Q-score                                                                                  |
|-------|------------------------------------------------------------------------------------------|------------------------------------------------------------------------------------------|
| h     |  0.7770 |  0.4560 |
| i     |  0.8360 |  0.4950 |
| j     |  0.8980 |  0.5870 |
| k     |  0.8070 |  0.4540 |
| l     |  0.8740 |  0.5310 |
| m     |  0.9590 |  0.5990 |
| n     |  0.4790 |  0.1840 |
| o     |  0.9060 |  0.5830 |
| p     |  0.8140 |  0.4450 |
| q     |  0.7590 |  0.4290 |
| r     |  0.8950 |  0.5820 |
| s     |  0.9040 |  0.5680 |
| t     |  0.8630 |  0.5180 |
| u     |  0.5150 |  0.1950 |
| v     |  0.7900 |  0.4710 |
| w     |  0.9270 |  0.5860 |
| x     |  0.8610 |  0.5300 |
| y     |  0.8720 |  0.5210 |
| z     |  0.9040 |  0.5920 |