



wwPDB EM Validation Summary Report ⓘ

Apr 30, 2024 – 06:22 pm BST

PDB ID : 4V8T
EMDB ID : EMD-2169
Title : Cryo-EM Structure of the 60S Ribosomal Subunit in Complex with Arx1 and Rei1
Authors : Greber, B.J.; Boehringer, D.; Montellese, C.; Ban, N.
Deposited on : 2012-08-07
Resolution : 8.10 Å (reported)
Based on initial models : 3U5I, 3U5H

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

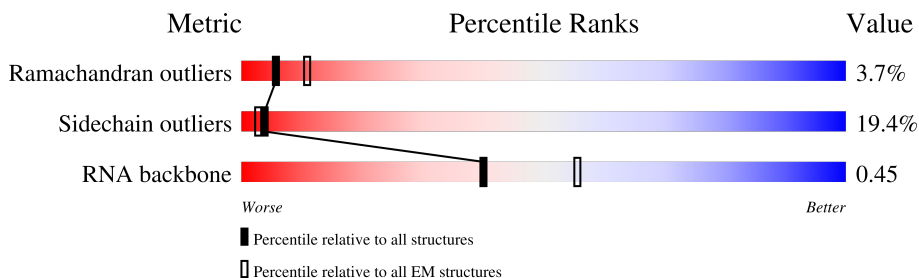
EMDB validation analysis : 0.0.1.dev92
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

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 8.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



| Metric | Whole archive (#Entries) | EM structures (#Entries) |
|-----------------------|--------------------------|--------------------------|
| Ramachandran outliers | 154571 | 4023 |
| Sidechain outliers | 154315 | 3826 |
| RNA backbone | 4643 | 859 |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1 | A | 254 | |
| 2 | B | 387 | |
| 3 | C | 362 | |
| 4 | D | 297 | |
| 5 | E | 176 | |
| 6 | F | 244 | |
| 7 | G | 256 | |
| 8 | H | 191 | |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------------------|
| 9 | I | 221 | 23% 72% 22% .. |
| 10 | J | 174 | 6% 72% 21% ... |
| 11 | K | 155 | 85% 97% . |
| 12 | L | 199 | 10% 76% 19% .. |
| 13 | M | 138 | . 80% 17% .. |
| 14 | N | 204 | 32% 80% 17% . |
| 15 | O | 219 | 5% 71% 20% 9% . |
| 16 | P | 184 | 17% 70% 12% . 16% |
| 17 | Q | 186 | 16% 78% 19% .. |
| 18 | R | 189 | 28% 79% 21% . |
| 19 | S | 172 | 7% 79% 19% . |
| 20 | T | 160 | 18% 80% 19% .. |
| 21 | U | 121 | 7% 60% 21% 19% |
| 22 | V | 137 | 26% 90% 8% .. |
| 23 | W | 155 | 50% 73% 13% . 13% |
| 24 | X | 142 | 11% 65% 18% . 15% |
| 25 | Y | 127 | 8% 76% 21% .. |
| 26 | Z | 136 | 10% 74% 21% 5% . |
| 27 | a | 149 | 19% 75% 22% ... |
| 28 | b | 59 | 24% 71% 24% .. |
| 29 | c | 105 | 5% 79% 15% . 5% |
| 30 | d | 113 | 13% 73% 21% .. |
| 31 | e | 130 | 28% 75% 20% ... |
| 32 | f | 107 | 8% 85% 12% .. |
| 33 | g | 121 | 26% 70% 21% . 7% |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 34 | h | 120 | |
| 35 | i | 100 | |
| 36 | j | 88 | |
| 37 | k | 78 | |
| 38 | l | 51 | |
| 39 | m | 128 | |
| 40 | n | 25 | |
| 41 | o | 106 | |
| 42 | p | 92 | |
| 43 | q | 312 | |
| 44 | r | 153 | |
| 45 | s | 46 | |
| 46 | t | 614 | |
| 47 | 1 | 114 | |
| 48 | 5 | 3396 | |
| 49 | 7 | 121 | |
| 50 | 8 | 158 | |

2 Entry composition [i](#)

There are 51 unique types of molecules in this entry. The entry contains 130050 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 60S ribosomal protein L2-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 1 | A | 252 | 1912 | 1190 | 388 | 333 | 1 | 0 | 0 |

- Molecule 2 is a protein called 60S ribosomal protein L3.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 2 | B | 386 | 3075 | 1950 | 584 | 533 | 8 | 0 | 0 |

- Molecule 3 is a protein called 60S ribosomal protein L4-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 3 | C | 361 | 2748 | 1729 | 522 | 494 | 3 | 0 | 0 |

- Molecule 4 is a protein called 60S ribosomal protein L5.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 4 | D | 294 | 2359 | 1489 | 412 | 456 | 2 | 0 | 0 |

- Molecule 5 is a protein called 60S ribosomal protein L6-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 5 | E | 157 | 1248 | 806 | 224 | 217 | 1 | 0 | 0 |

- Molecule 6 is a protein called 60S ribosomal protein L7-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 6 | F | 223 | 1791 | 1155 | 325 | 310 | 1 | 0 | 0 |

- Molecule 7 is a protein called 60S ribosomal protein L8-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 7 | G | 231 | 1763 | 1130 | 316 | 314 | 3 | 0 | 0 |

- Molecule 8 is a protein called 60S ribosomal protein L9-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 8 | H | 191 | 1518 | 963 | 274 | 277 | 4 | 0 | 0 |

- Molecule 9 is a protein called 60S ribosomal protein L10.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 9 | I | 213 | 1722 | 1094 | 325 | 297 | 6 | 0 | 0 |

- Molecule 10 is a protein called 60S ribosomal protein L11-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 10 | J | 169 | 1353 | 847 | 253 | 249 | 4 | 0 | 0 |

- Molecule 11 is a protein called 60S RIBOSOMAL PROTEIN L12.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| | | | Total | C | N | O | | |
| 11 | K | 150 | 750 | 450 | 150 | 150 | 0 | 0 |

- Molecule 12 is a protein called 60S ribosomal protein L13-A.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| | | | Total | C | N | O | | |
| 12 | L | 194 | 1548 | 965 | 316 | 267 | 0 | 0 |

- Molecule 13 is a protein called 60S ribosomal protein L14-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 13 | M | 137 | 1059 | 678 | 200 | 179 | 2 | 0 | 0 |

- Molecule 14 is a protein called 60S ribosomal protein L15-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 14 | N | 203 | 1720 | 1077 | 361 | 281 | 1 | 0 | 0 |

- Molecule 15 is a protein called Large ribosomal subunit protein uL13A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 15 | O | 197 | 3119 | 2008 | 581 | 528 | 2 | 197 | 0 |

- Molecule 16 is a protein called 60S ribosomal protein L17-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 16 | P | 155 | 1227 | 764 | 238 | 225 | | 0 | 0 |

- Molecule 17 is a protein called 60S ribosomal protein L18-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 17 | Q | 185 | 1441 | 908 | 290 | 241 | 2 | 0 | 0 |

- Molecule 18 is a protein called 60S ribosomal protein L19-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 18 | R | 188 | 1521 | 935 | 326 | 260 | | 0 | 0 |

- Molecule 19 is a protein called 60S ribosomal protein L20-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 19 | S | 172 | 1445 | 930 | 267 | 244 | 4 | 0 | 0 |

- Molecule 20 is a protein called 60S ribosomal protein L21-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 20 | T | 159 | 1276 | 805 | 246 | 221 | 4 | 0 | 0 |

- Molecule 21 is a protein called 60S ribosomal protein L22-A.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 21 | U | 98 | Total | C | N | O | 0 | 0 |
| | | | 778 | 505 | 127 | 146 | | |

- Molecule 22 is a protein called 60S ribosomal protein L23-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 22 | V | 136 | Total | C | N | O | S | 0 | 0 |
| | | | 1003 | 628 | 189 | 179 | 7 | | |

- Molecule 23 is a protein called 60S ribosomal protein L24-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 23 | W | 135 | Total | C | N | O | S | 0 | 0 |
| | | | 1038 | 651 | 206 | 180 | 1 | | |

- Molecule 24 is a protein called 60S ribosomal protein L25.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 24 | X | 120 | Total | C | N | O | S | 0 | 0 |
| | | | 959 | 617 | 168 | 172 | 2 | | |

- Molecule 25 is a protein called 60S ribosomal protein L26-A.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 25 | Y | 126 | Total | C | N | O | 0 | 0 |
| | | | 993 | 625 | 192 | 176 | | |

- Molecule 26 is a protein called 60S ribosomal protein L27-A.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 26 | Z | 135 | Total | C | N | O | 0 | 0 |
| | | | 1092 | 710 | 202 | 180 | | |

- Molecule 27 is a protein called 60S ribosomal protein L28.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 27 | a | 148 | Total | C | N | O | S | 0 | 0 |
| | | | 1173 | 749 | 231 | 190 | 3 | | |

- Molecule 28 is a protein called 60S ribosomal protein L29.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---------|-------|
| 28 | b | 58 | Total | C | N | O | 0 | 0 |
| | | | 462 | 289 | 100 | 73 | | |

- Molecule 29 is a protein called 60S ribosomal protein L30.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 29 | c | 100 | Total | C | N | O | S | 0 | 0 |
| | | | 767 | 492 | 128 | 146 | 1 | | |

- Molecule 30 is a protein called 60S ribosomal protein L31-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 30 | d | 109 | Total | C | N | O | S | 0 | 0 |
| | | | 883 | 559 | 167 | 156 | 1 | | |

- Molecule 31 is a protein called 60S ribosomal protein L32.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 31 | e | 127 | Total | C | N | O | S | 0 | 0 |
| | | | 1020 | 647 | 205 | 167 | 1 | | |

- Molecule 32 is a protein called 60S ribosomal protein L33-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 32 | f | 106 | Total | C | N | O | S | 0 | 0 |
| | | | 850 | 540 | 165 | 144 | 1 | | |

- Molecule 33 is a protein called 60S ribosomal protein L34-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 33 | g | 112 | Total | C | N | O | S | 0 | 0 |
| | | | 880 | 545 | 179 | 152 | 4 | | |

- Molecule 34 is a protein called 60S ribosomal protein L35-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 34 | h | 119 | Total | C | N | O | S | 0 | 0 |
| | | | 965 | 612 | 185 | 167 | 1 | | |

- Molecule 35 is a protein called 60S ribosomal protein L36-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 35 | i | 99 | Total | C | N | O | S | 0 | 0 |
| | | | 770 | 481 | 156 | 131 | 2 | | |

- Molecule 36 is a protein called 60S ribosomal protein L37-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 36 | j | 87 | Total | C | N | O | S | 0 | 0 |
| | | | 681 | 414 | 148 | 114 | 5 | | |

- Molecule 37 is a protein called 60S ribosomal protein L38.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---------|-------|
| 37 | k | 77 | Total | C | N | O | 0 | 0 |
| | | | 608 | 388 | 114 | 106 | | |

- Molecule 38 is a protein called 60S ribosomal protein L39.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 38 | l | 50 | Total | C | N | O | S | 0 | 0 |
| | | | 436 | 272 | 97 | 65 | 2 | | |

- Molecule 39 is a protein called Ubiquitin-60S ribosomal protein L40.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 39 | m | 52 | Total | C | N | O | S | 0 | 0 |
| | | | 417 | 259 | 86 | 67 | 5 | | |

- Molecule 40 is a protein called 60S ribosomal protein L41-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 40 | n | 25 | Total | C | N | O | S | 0 | 0 |
| | | | 233 | 142 | 63 | 27 | 1 | | |

- Molecule 41 is a protein called 60S ribosomal protein L42-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 41 | o | 105 | Total | C | N | O | S | 0 | 0 |
| | | | 847 | 534 | 170 | 138 | 5 | | |

- Molecule 42 is a protein called 60S ribosomal protein L43-A.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 42 | p | 91 | 694 | 429 | 138 | 121 | 6 | 0 | 0 |

- Molecule 43 is a protein called Large ribosomal subunit protein uL10.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| | | | Total | C | N | O | S | | |
| 43 | q | 143 | 1077 | 687 | 192 | 195 | 3 | 0 | 0 |

- Molecule 44 is a protein called RIBOSOMAL PROTEIN P1 ALPHA, Large ribosomal subunit protein P1A.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---------|-------|
| | | | Total | C | N | O | | |
| 44 | r | 47 | 235 | 141 | 47 | 47 | 0 | 0 |

- Molecule 45 is a protein called RIBOSOMAL PROTEIN P2 BETA.

| Mol | Chain | Residues | Atoms | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---------|-------|
| | | | Total | C | N | O | | |
| 45 | s | 46 | 230 | 138 | 46 | 46 | 0 | 0 |

- Molecule 46 is a protein called Probable metalloprotease ARX1.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|-------|
| | | | Total | C | N | O | S | | |
| 46 | t | 380 | 2938 | 1853 | 511 | 563 | 11 | 0 | 0 |

There are 21 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------------|------------|
| t | -20 | HIS | - | expression tag | UNP Q03862 |
| t | -19 | HIS | - | expression tag | UNP Q03862 |
| t | -18 | HIS | - | expression tag | UNP Q03862 |
| t | -17 | HIS | - | expression tag | UNP Q03862 |
| t | -16 | HIS | - | expression tag | UNP Q03862 |
| t | -15 | HIS | - | expression tag | UNP Q03862 |
| t | -14 | ASP | - | expression tag | UNP Q03862 |
| t | -13 | TYR | - | expression tag | UNP Q03862 |
| t | -12 | ASP | - | expression tag | UNP Q03862 |
| t | -11 | ILE | - | expression tag | UNP Q03862 |
| t | -10 | PRO | - | expression tag | UNP Q03862 |

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| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------------|------------|
| t | -9 | THR | - | expression tag | UNP Q03862 |
| t | -8 | THR | - | expression tag | UNP Q03862 |
| t | -7 | GLU | - | expression tag | UNP Q03862 |
| t | -6 | ASN | - | expression tag | UNP Q03862 |
| t | -5 | LEU | - | expression tag | UNP Q03862 |
| t | -4 | TYR | - | expression tag | UNP Q03862 |
| t | -3 | PHE | - | expression tag | UNP Q03862 |
| t | -2 | GLN | - | expression tag | UNP Q03862 |
| t | -1 | GLY | - | expression tag | UNP Q03862 |
| t | 0 | ALA | - | expression tag | UNP Q03862 |

- Molecule 47 is a RNA chain called ES27 OF THE 25S RRNA.

| Mol | Chain | Residues | Atoms | | AltConf | Trace |
|-----|-------|----------|-------|-----|---------|-------|
| 47 | 1 | 114 | Total | P | 0 | 114 |
| | | | 114 | 114 | | |

- Molecule 48 is a RNA chain called 25S RIBOSOMAL RNA.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-------|-------|-------|------|---------|-------|
| 48 | 5 | 3150 | Total | C | N | O | P | 0 | 0 |
| | | | 67376 | 30095 | 12145 | 21987 | 3149 | | |

- Molecule 49 is a RNA chain called 5S RIBOSOMAL RNA.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|-----|---------|-------|
| 49 | 7 | 121 | Total | C | N | O | P | 0 | 0 |
| | | | 2579 | 1152 | 461 | 845 | 121 | | |

- Molecule 50 is a RNA chain called 5.8S RIBOSOMAL RNA.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|------|-----|---------|-------|
| 50 | 8 | 158 | Total | C | N | O | P | 0 | 0 |
| | | | 3353 | 1500 | 586 | 1109 | 158 | | |

- Molecule 51 is ZINC ION (three-letter code: ZN) (formula: Zn).

| Mol | Chain | Residues | Atoms | | AltConf |
|-----|-------|----------|-------|----|---------|
| 51 | j | 1 | Total | Zn | 0 |
| | | | 1 | 1 | |
| 51 | m | 1 | Total | Zn | 0 |
| | | | 1 | 1 | |

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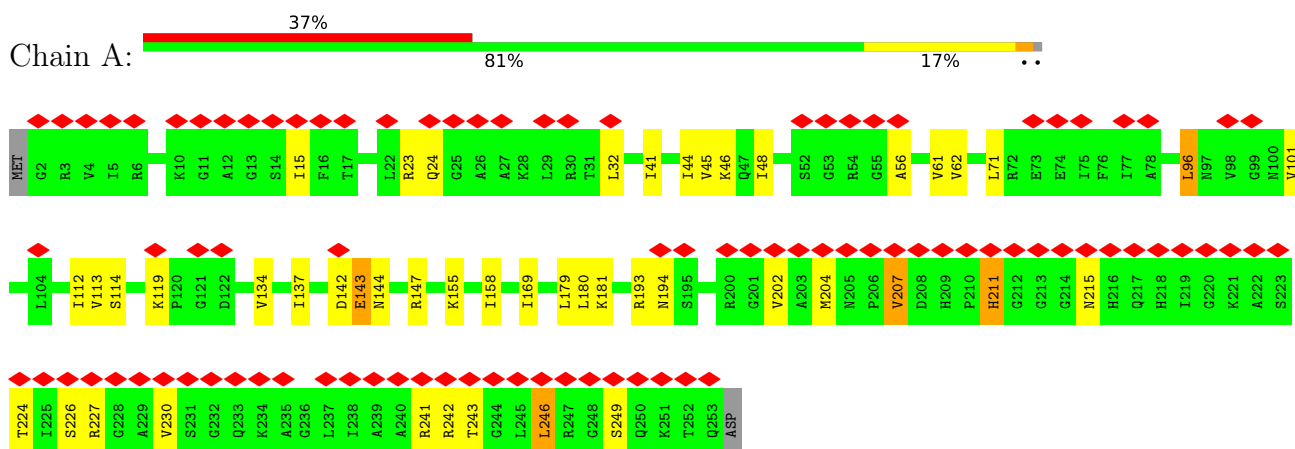
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| Mol | Chain | Residues | Atoms | | AltConf |
|------------|--------------|-----------------|--------------|---------|----------------|
| 51 | o | 1 | Total 1 | Zn 1 | 0 |
| 51 | p | 1 | Total 1 | Zn 1 | 0 |

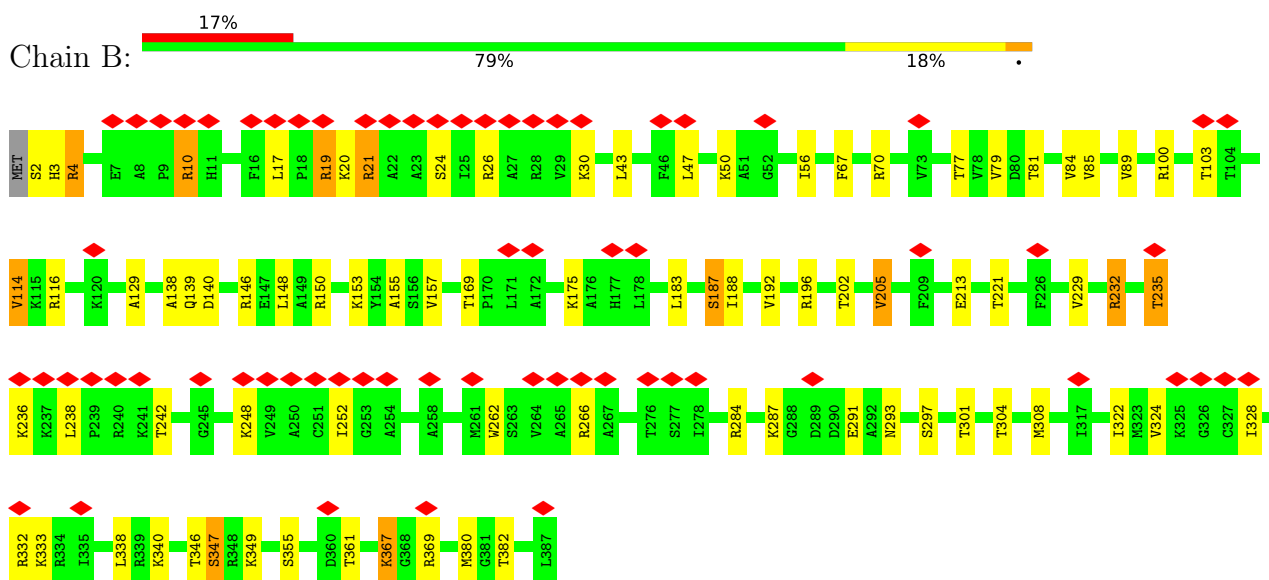
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and 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.

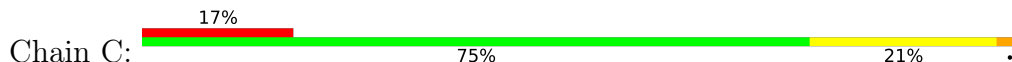
- Molecule 1: 60S ribosomal protein L2-A

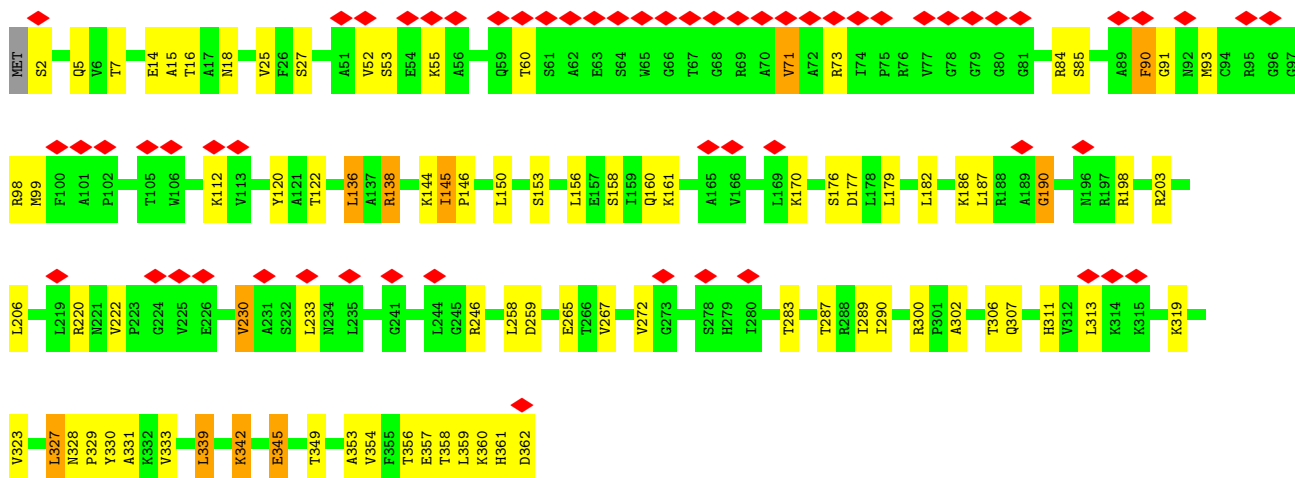


- Molecule 2: 60S ribosomal protein L3

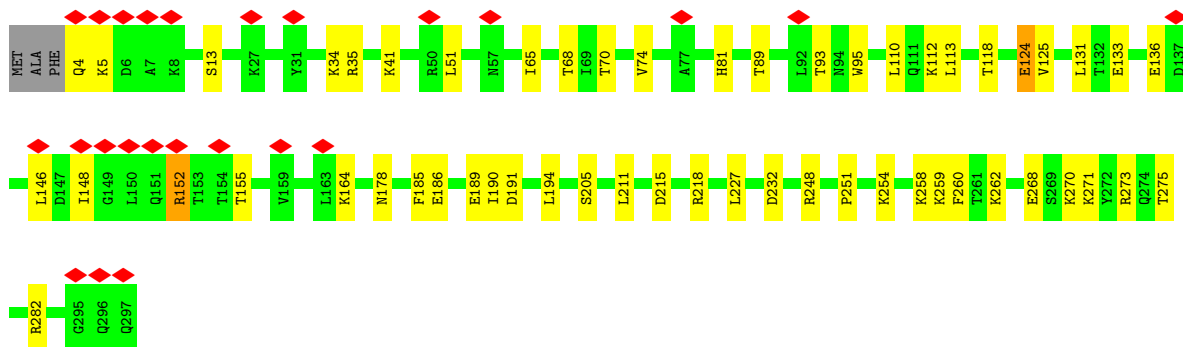
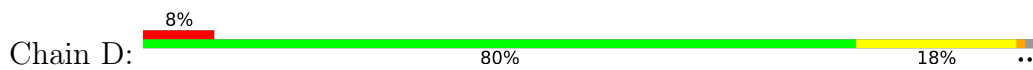


- Molecule 3: 60S ribosomal protein L4-A

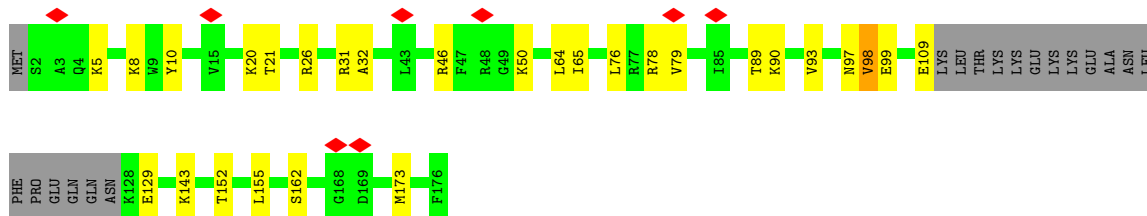
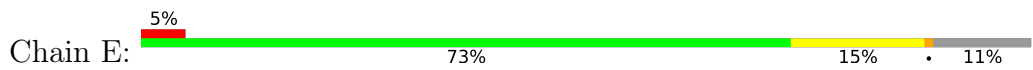




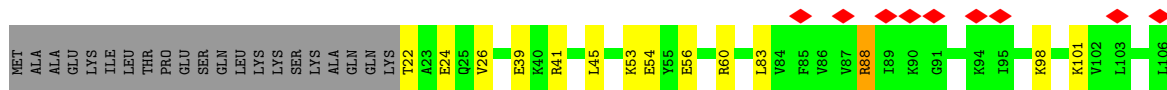
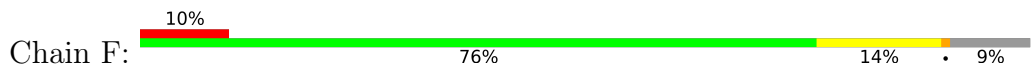
• Molecule 4: 60S ribosomal protein L5

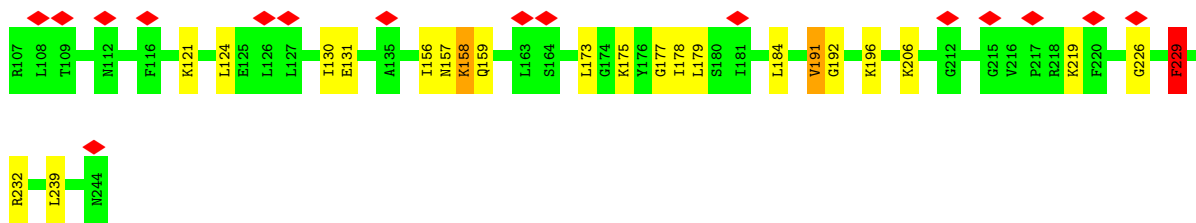


• Molecule 5: 60S ribosomal protein L6-A

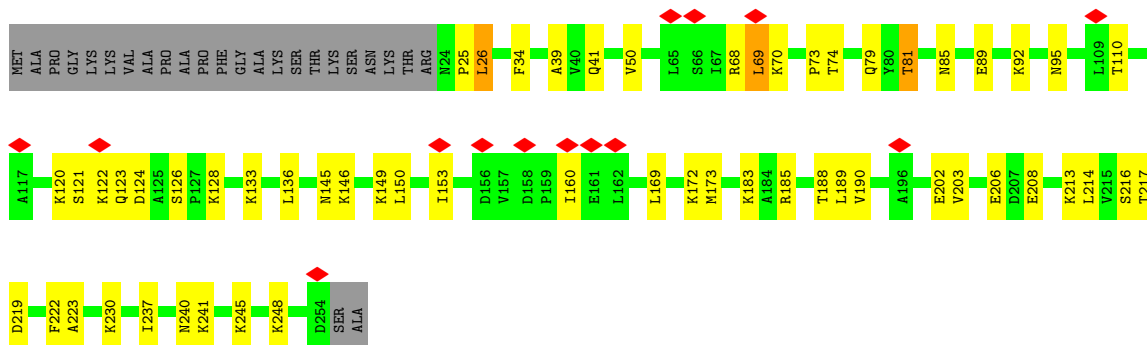


• Molecule 6: 60S ribosomal protein L7-A

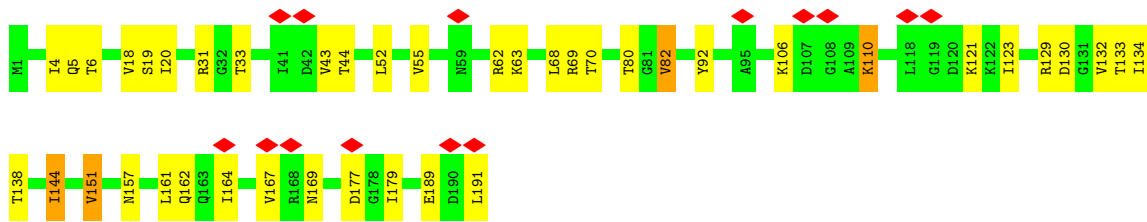
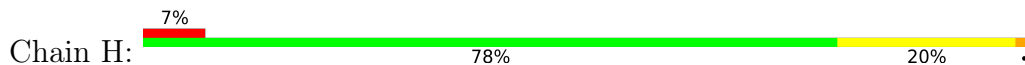




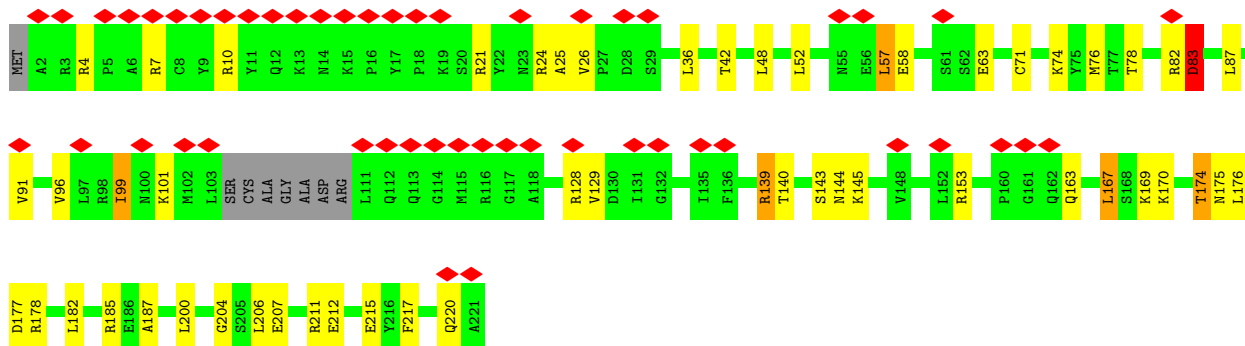
• Molecule 7: 60S ribosomal protein L8-A



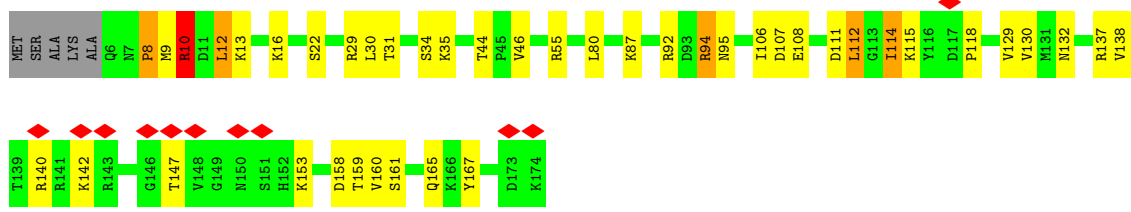
• Molecule 8: 60S ribosomal protein L9-A



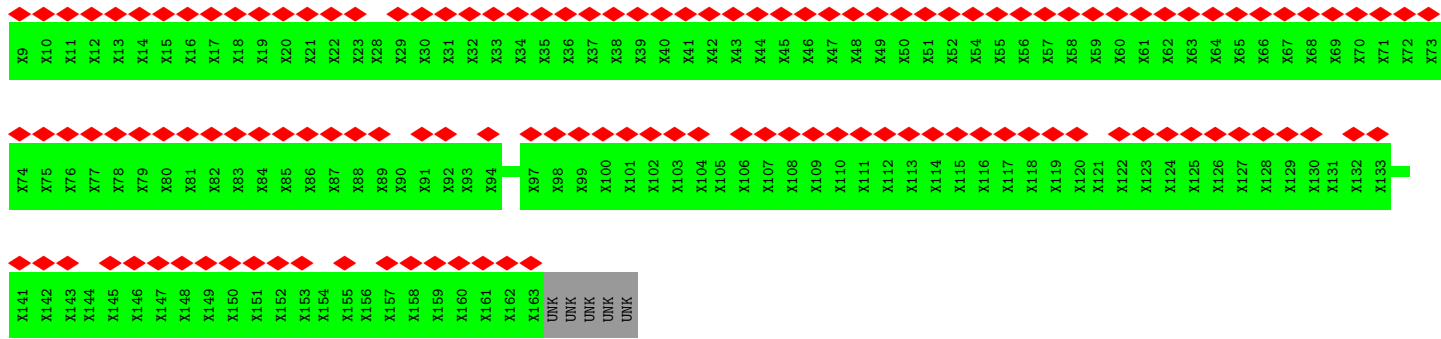
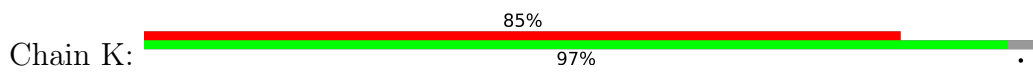
• Molecule 9: 60S ribosomal protein L10



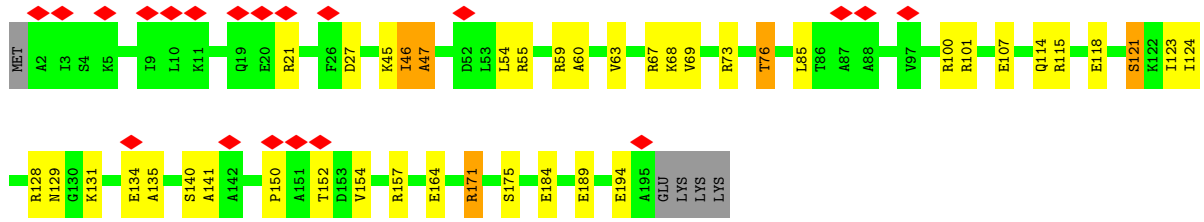
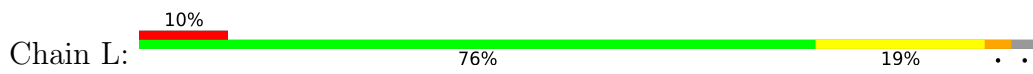
• Molecule 10: 60S ribosomal protein L11-A



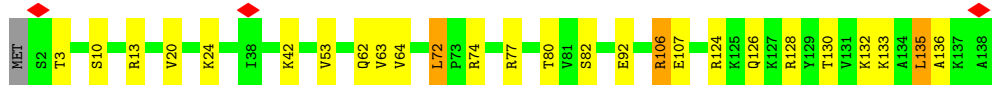
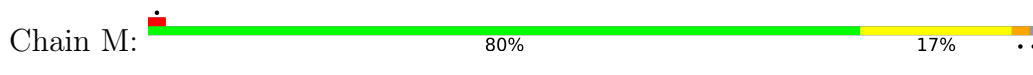
• Molecule 11: 60S RIBOSOMAL PROTEIN L12



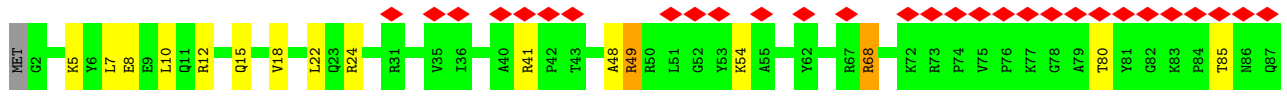
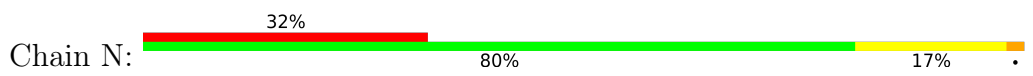
• Molecule 12: 60S ribosomal protein L13-A

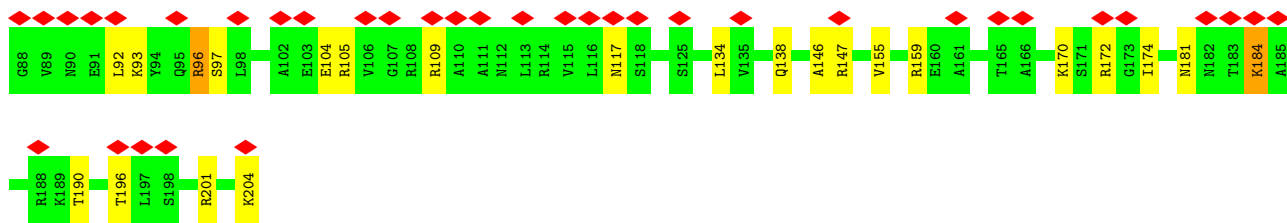


• Molecule 13: 60S ribosomal protein L14-A

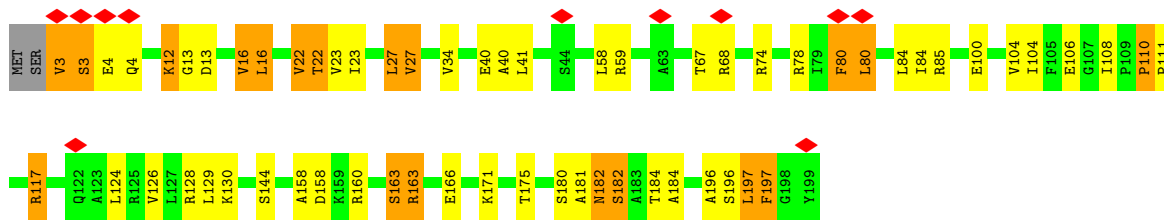


• Molecule 14: 60S ribosomal protein L15-A

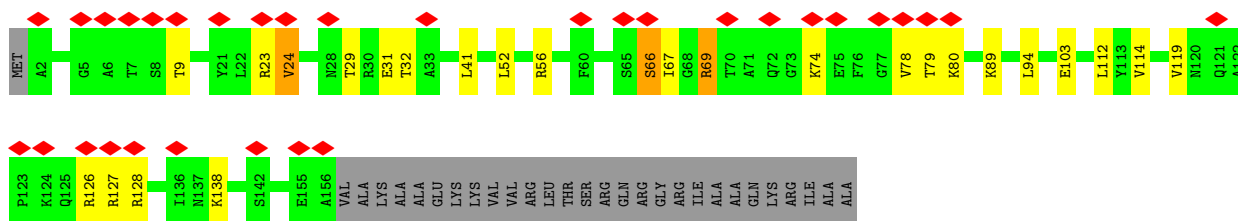




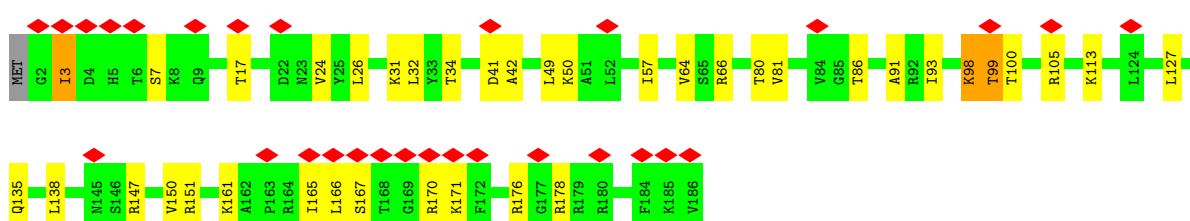
• Molecule 15: Large ribosomal subunit protein uL13A



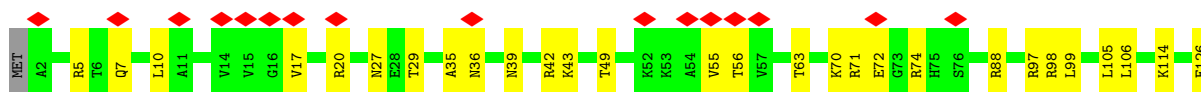
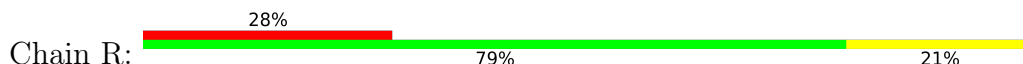
• Molecule 16: 60S ribosomal protein L17-A

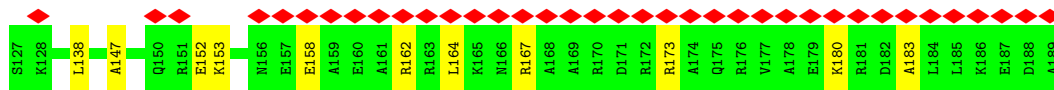


• Molecule 17: 60S ribosomal protein L18-A

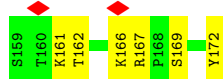
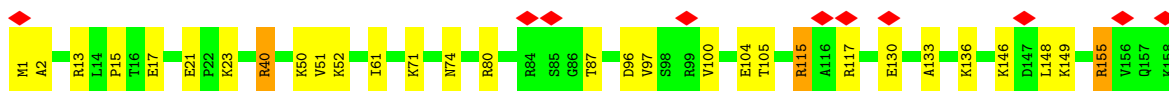
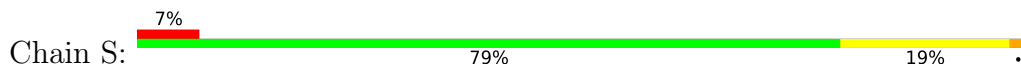


• Molecule 18: 60S ribosomal protein L19-A

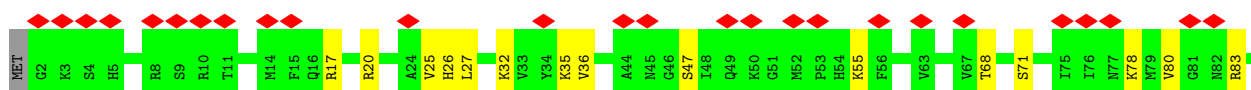
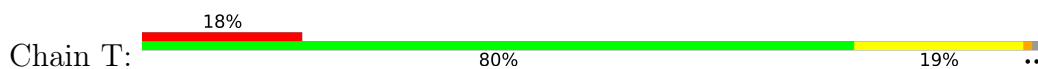




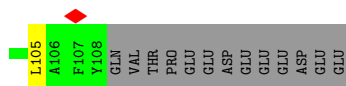
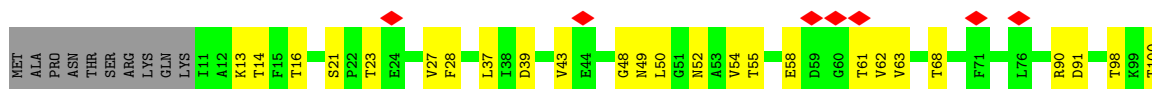
• Molecule 19: 60S ribosomal protein L20-A



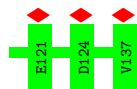
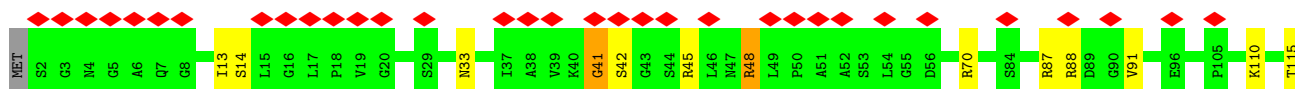
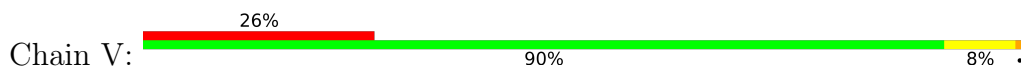
• Molecule 20: 60S ribosomal protein L21-A



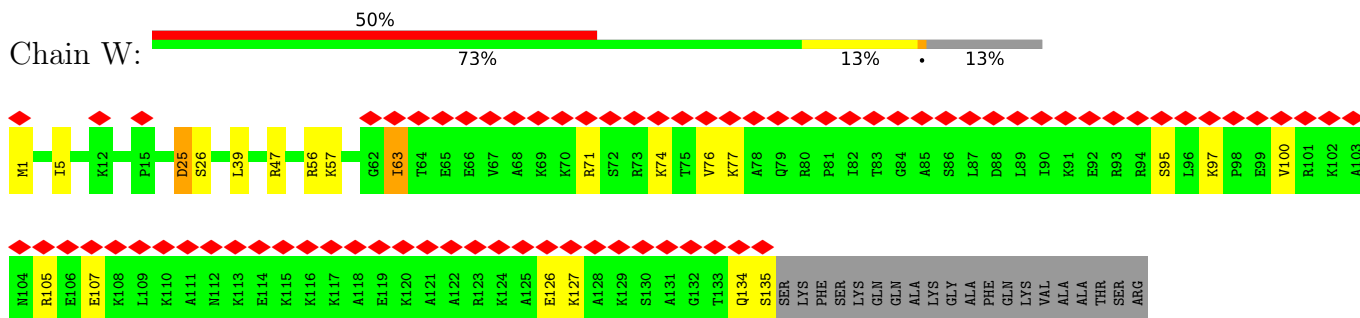
• Molecule 21: 60S ribosomal protein L22-A



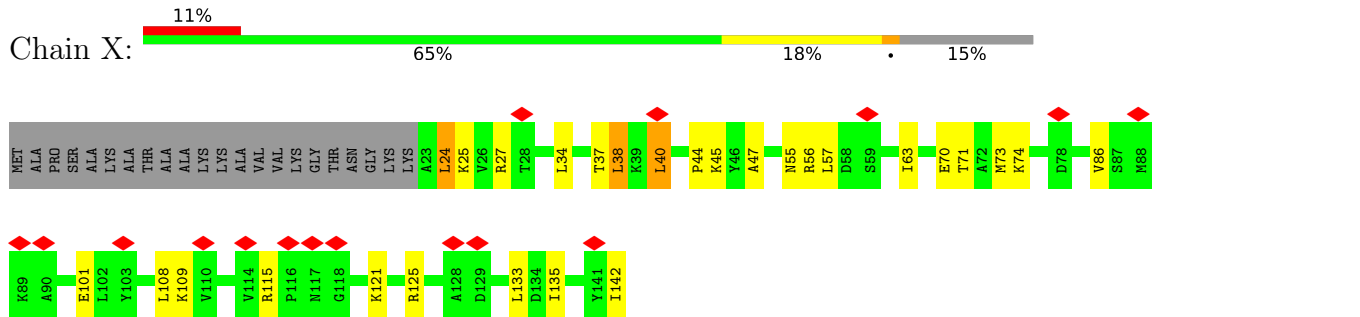
• Molecule 22: 60S ribosomal protein L23-A



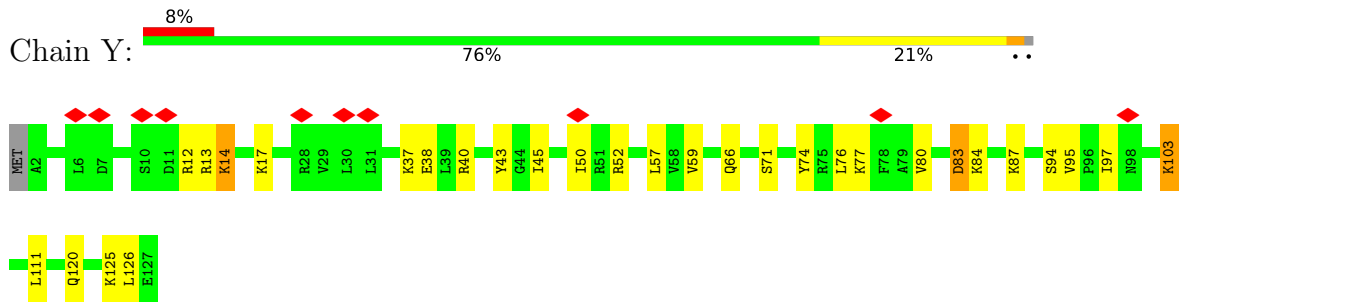
• Molecule 23: 60S ribosomal protein L24-A



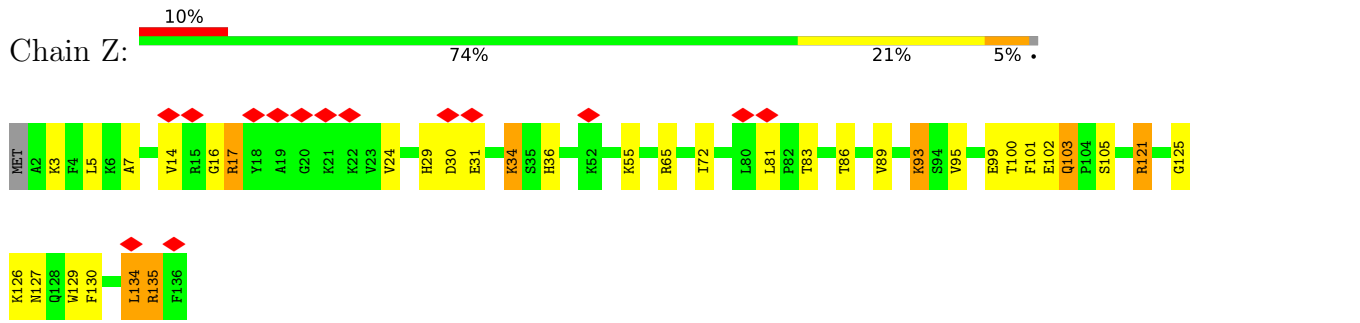
• Molecule 24: 60S ribosomal protein L25



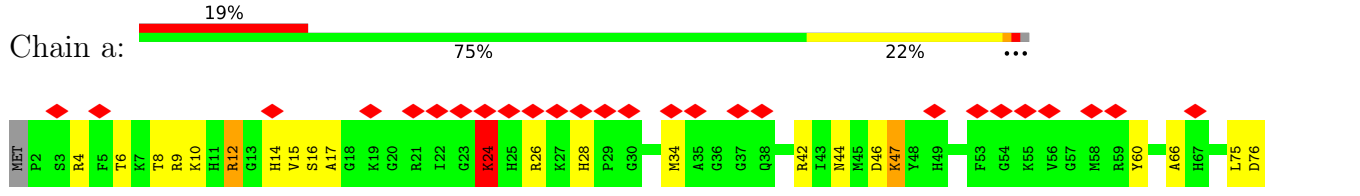
• Molecule 25: 60S ribosomal protein L26-A

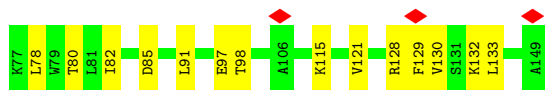


• Molecule 26: 60S ribosomal protein L27-A

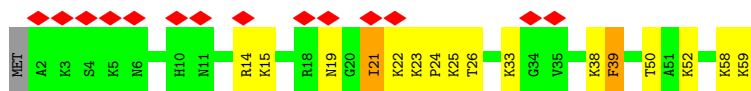


• Molecule 27: 60S ribosomal protein L28

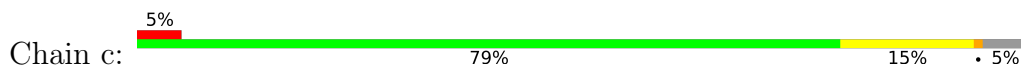




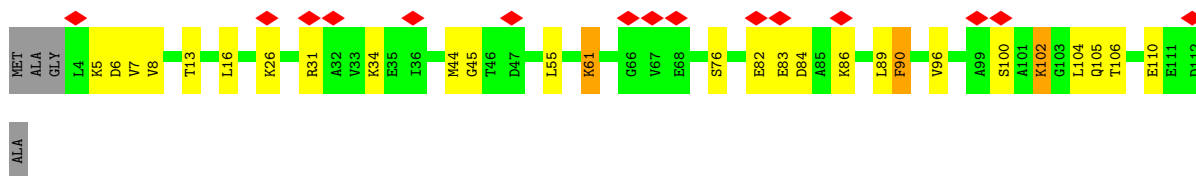
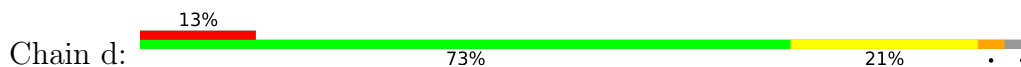
• Molecule 28: 60S ribosomal protein L29



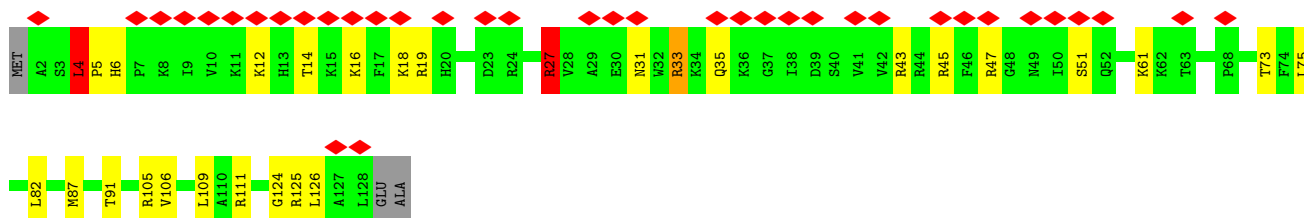
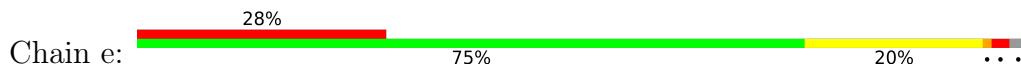
• Molecule 29: 60S ribosomal protein L30



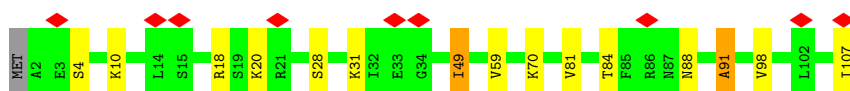
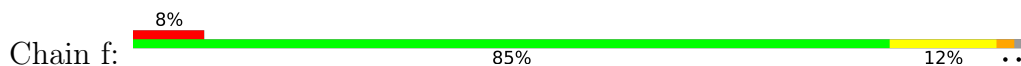
• Molecule 30: 60S ribosomal protein L31-A



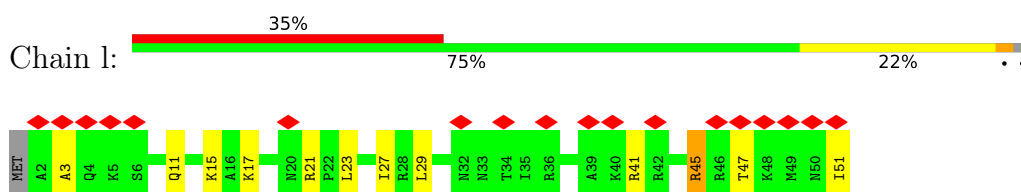
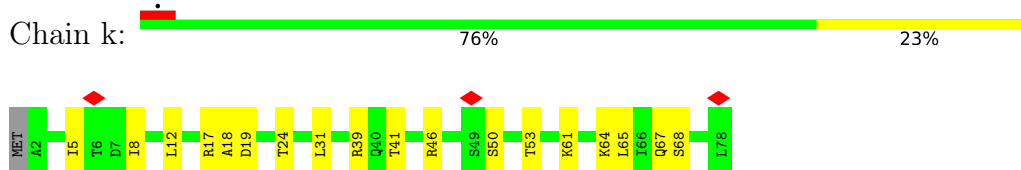
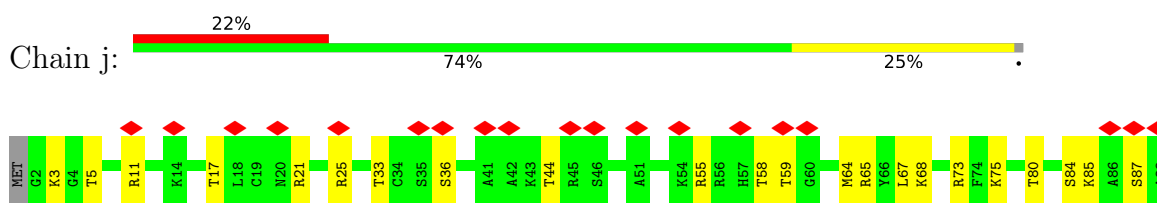
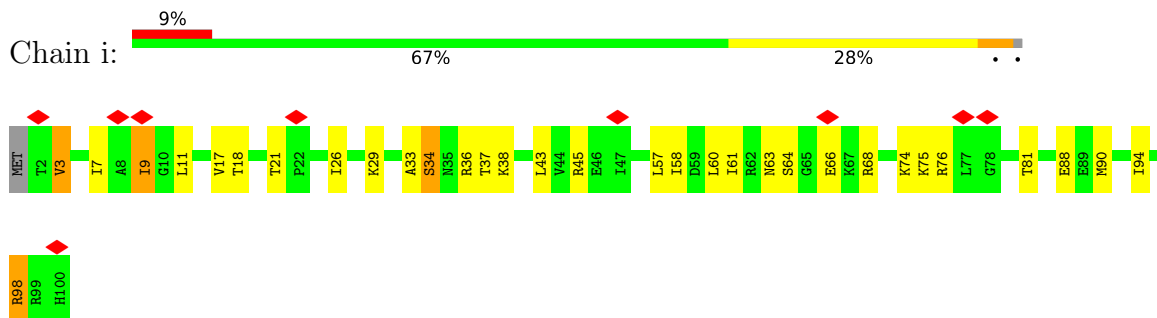
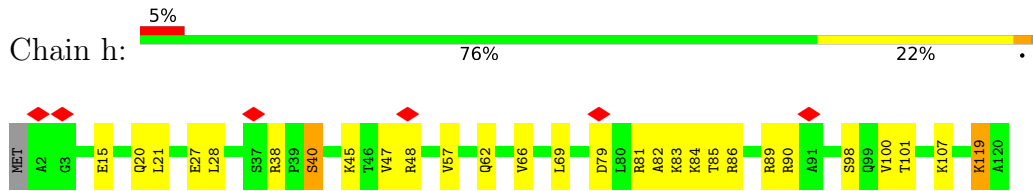
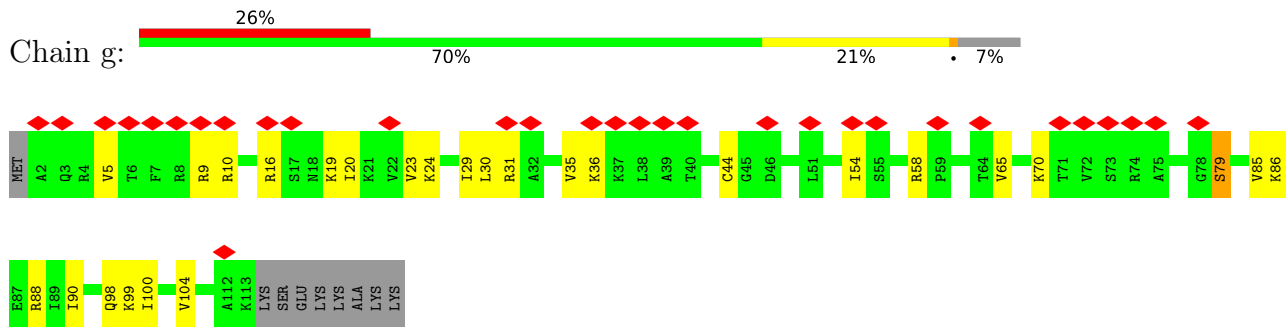
• Molecule 31: 60S ribosomal protein L32

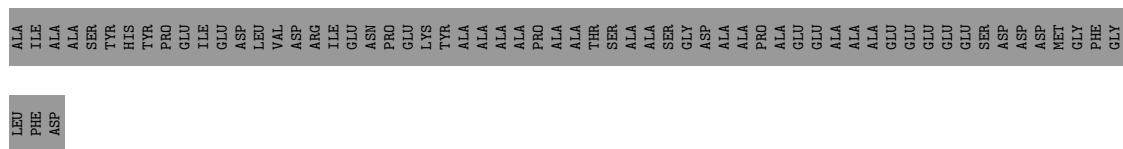


• Molecule 32: 60S ribosomal protein L33-A

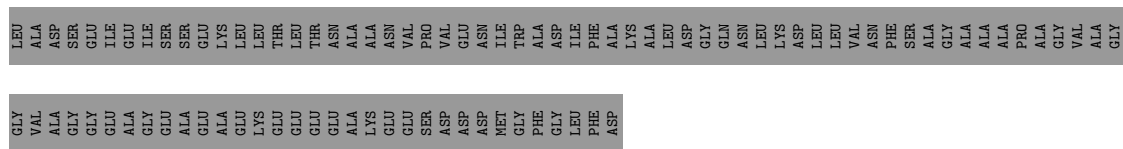
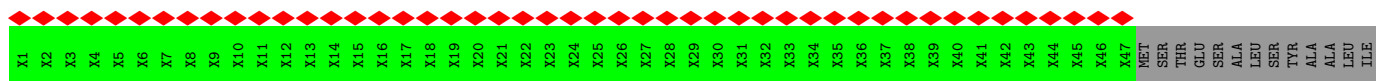


• Molecule 33: 60S ribosomal protein L34-A

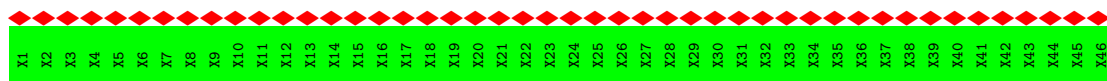




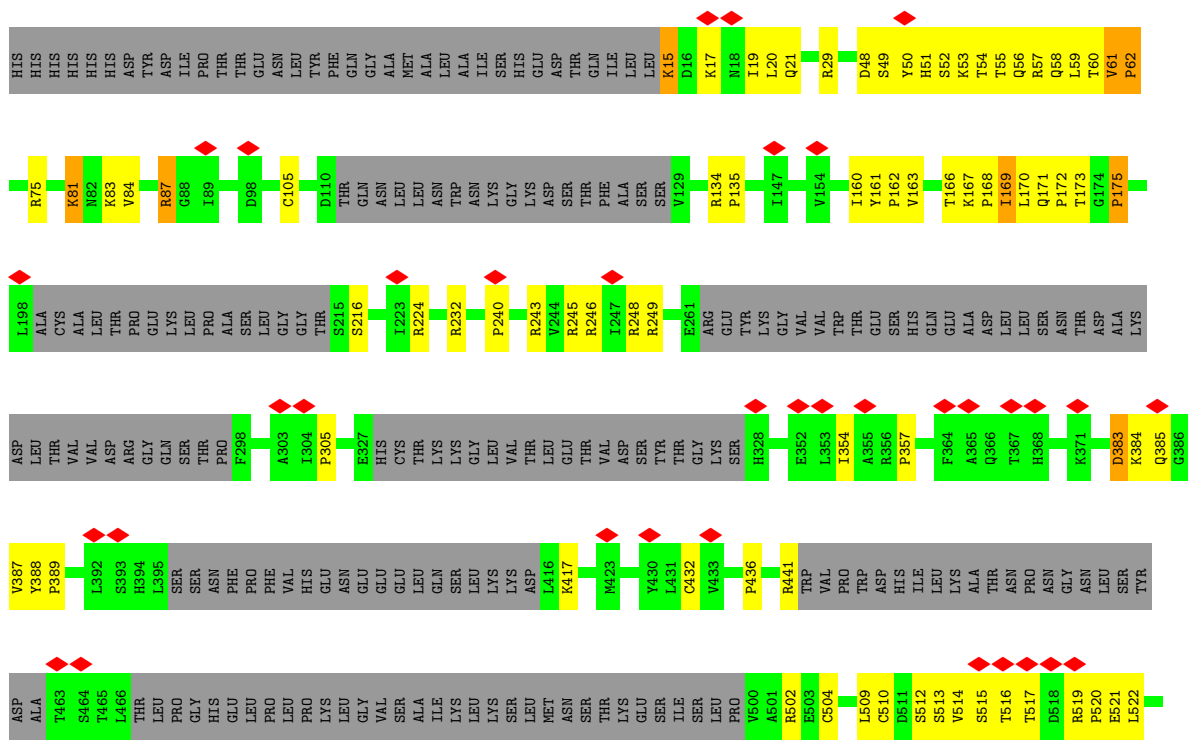
● Molecule 44: RIBOSOMAL PROTEIN P1 ALPHA, Large ribosomal subunit protein P1A

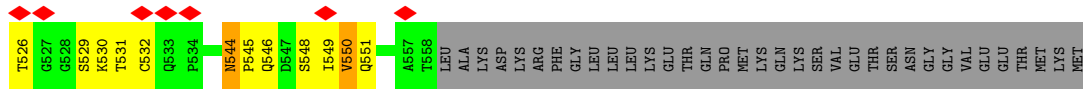


● Molecule 45: RIBOSOMAL PROTEIN P2 BETA

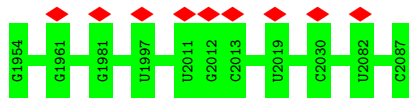


● Molecule 46: Probable metalloprotease ARX1

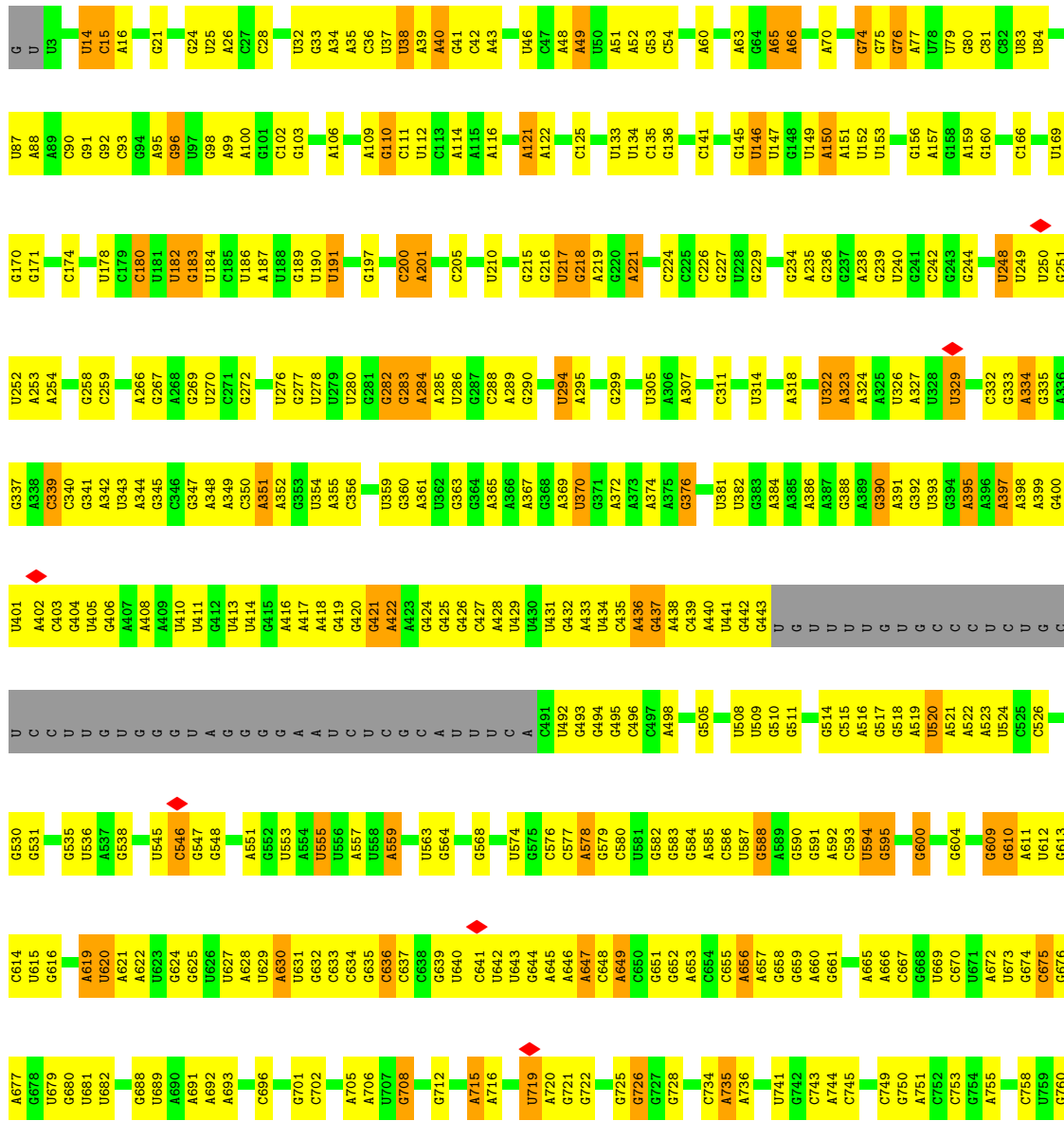


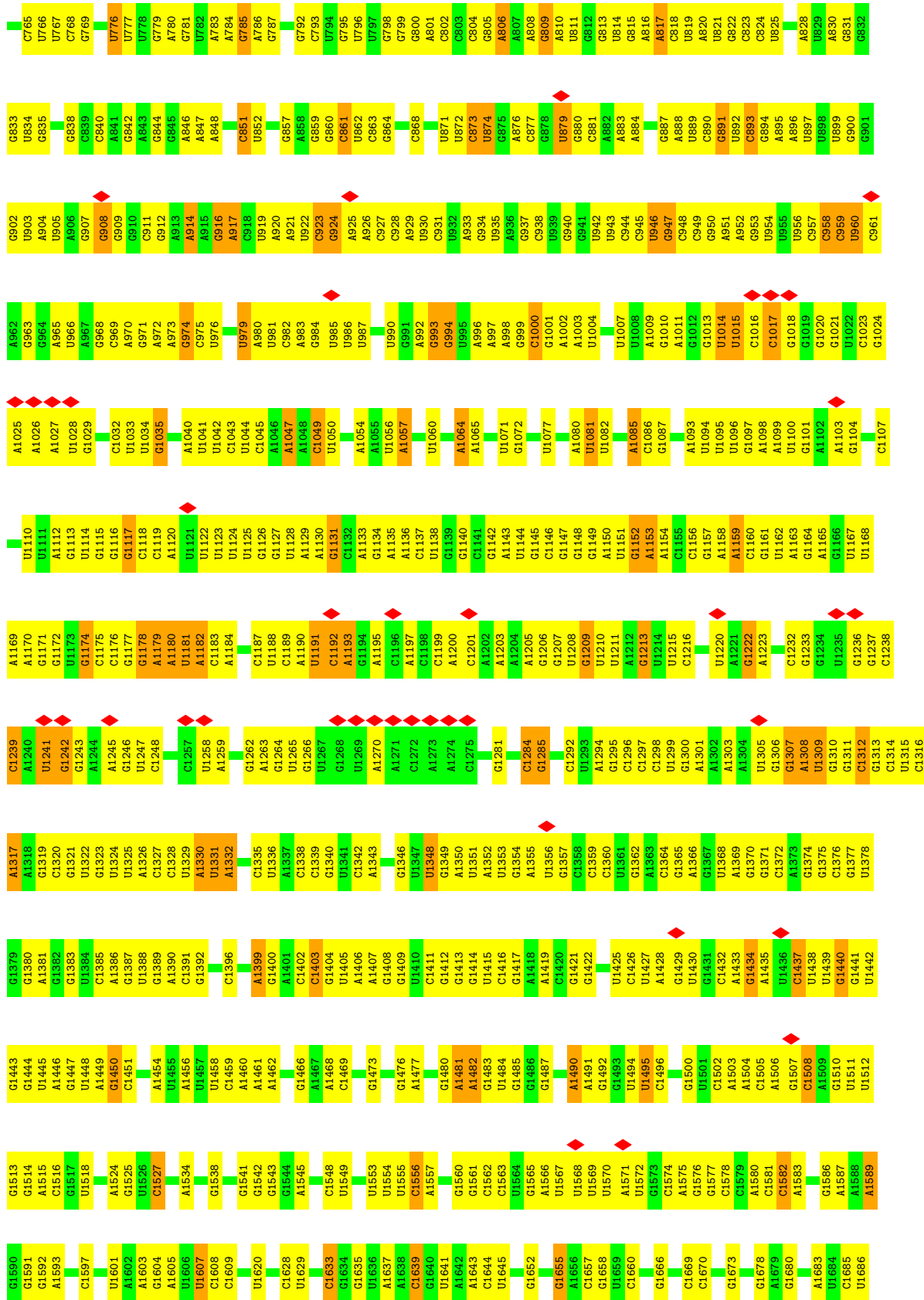


• Molecule 47: ES27 OF THE 25S RRNA

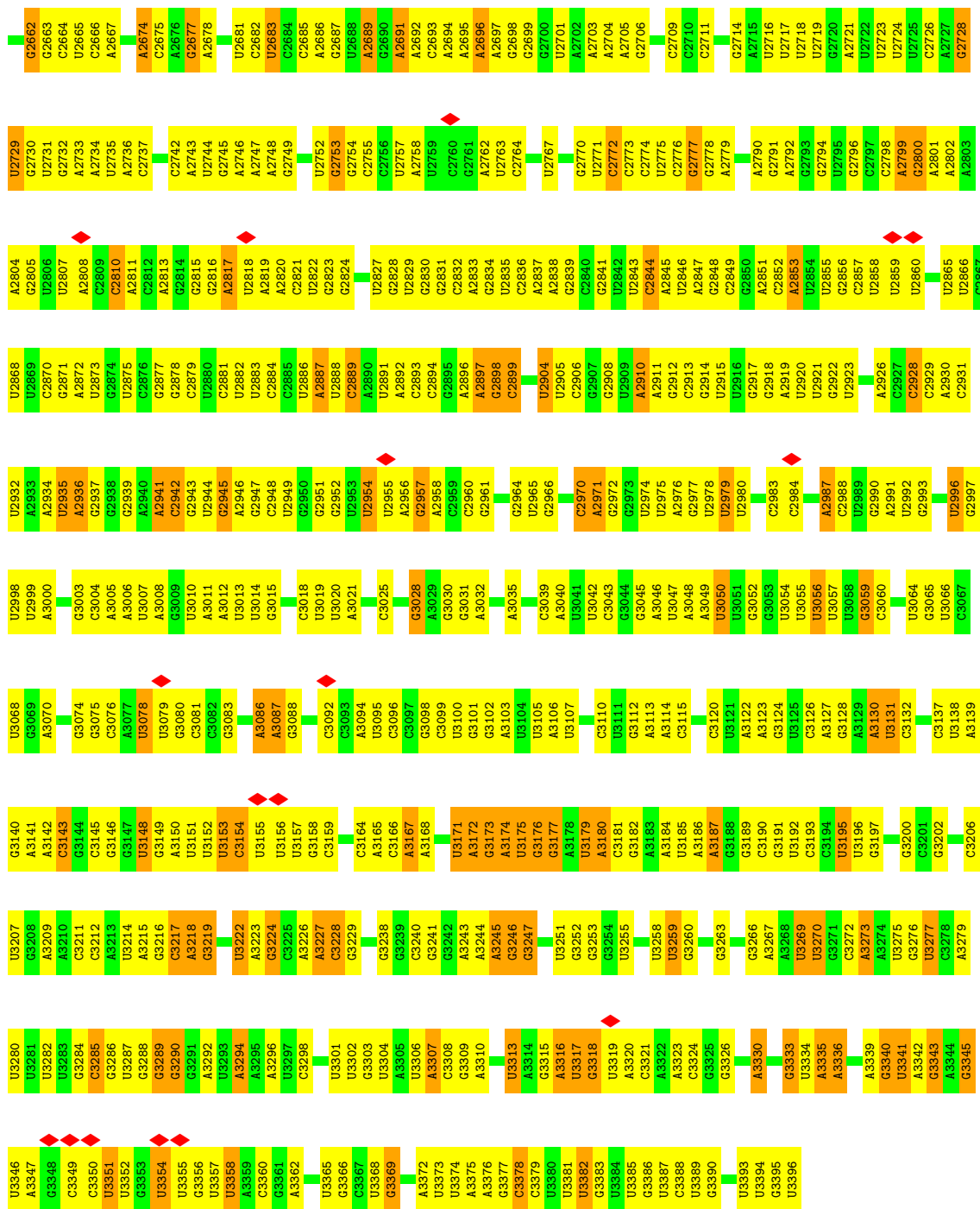


• Molecule 48: 25S RIBOSOMAL RNA

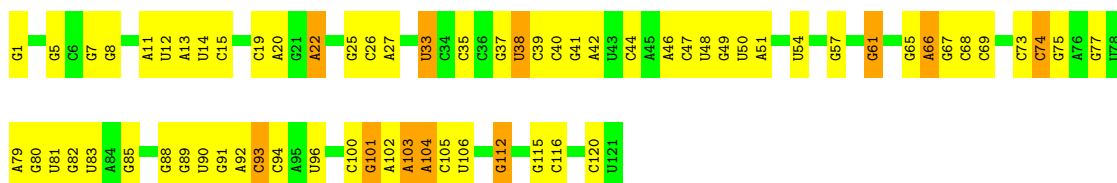




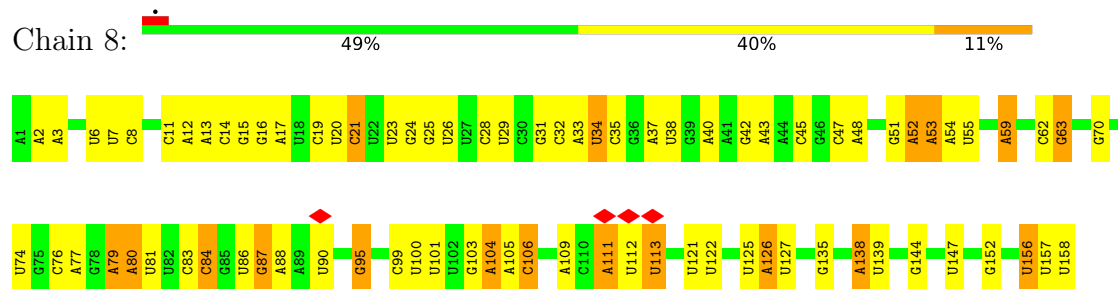
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|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|-------|
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| C2600 | U2517 | G2396 | U2336 | A2271 | G2206 | U2133 | A1950 | C1866 | G1780 | U1688 |
| G2603 | C2518 | A2397 | C2337 | G2272 | A2207 | U2134 | A1931 | U1867 | U1783 | C1690 |
| U2604 | A2398 | C2338 | C2339 | G2273 | A2208 | U2135 | A1932 | G1868 | U1784 | C1691 |
| G2605 | U2399 | C2339 | U2340 | G2274 | U2209 | U2136 | U1937 | G1869 | G1786 | C1692 |
| G2606 | A2401 | A2341 | A2342 | G2275 | G2210 | A2138 | U1938 | C1870 | A1787 | C1693 |
| G2607 | A2402 | U2343 | U2344 | C2276 | U2211 | A2139 | G1939 | U1871 | C1788 | U1694 |
| A2609 | G2403 | U2345 | U2346 | C2277 | C2212 | U2140 | G1940 | C1872 | U1789 | U1695 |
| G2610 | C2404 | U2346 | U2347 | A2280 | A2213 | A2141 | C1941 | U1791 | C1791 | A1699 |
| U2611 | C2405 | A2347 | A2348 | A2281 | A2214 | A2142 | U1942 | U1792 | C1792 | A1704 |
| U2612 | C2406 | U2347 | A2349 | A2282 | A2215 | A2143 | U1963 | C1793 | C1793 | A1704 |
| G2613 | A2348 | A2348 | U2349 | G2283 | G2216 | A2144 | G | U1797 | A1797 | A1744 |
| G2614 | U2349 | U2349 | C2350 | C2284 | U2217 | A2145 | G | U1877 | A1877 | A1744 |
| G2615 | C2350 | U2351 | A2222 | G2285 | A2222 | A2146 | U | G1878 | A1878 | A1715 |
| C2616 | U2352 | U2352 | A2223 | C2286 | A2223 | C2147 | U | U1880 | A1880 | U1716 |
| U2617 | G2412 | A2353 | A2224 | U2287 | A2224 | U2148 | A | U1881 | A1881 | U1717 |
| C2618 | A2413 | C2354 | U2225 | G2288 | U2225 | U2149 | U | G1882 | A1882 | G1718 |
| U2619 | G2414 | G2355 | U2226 | G2289 | U2226 | A2158 | G | A1883 | G1808 | G1719 |
| U2620 | C2415 | A2356 | C2228 | C2290 | C2227 | C2163 | G | U1884 | A1809 | U1722 |
| U2621 | U2416 | A2357 | A2229 | A2291 | A2228 | A2164 | G | U1885 | A1810 | U1723 |
| C2622 | G2417 | U2358 | U2229 | U2292 | A2229 | G2165 | G | A1886 | G1811 | U1724 |
| G2623 | U2418 | C2359 | C2230 | C2293 | C2230 | A2166 | G | A1887 | G1812 | U1724 |
| C2624 | A2419 | C2360 | C2231 | U2294 | C2231 | A2167 | C | U1888 | A1813 | C1725 |
| C2625 | C2420 | A2361 | G2294 | U2298 | G2239 | A2168 | C | G1889 | A1814 | C1726 |
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| C2627 | A2422 | G2363 | G2236 | G2300 | G2241 | U2170 | C | G1892 | A1816 | G1736 |
| U2628 | U2423 | A2364 | U2237 | U2301 | A2171 | A2172 | C | U1893 | A1817 | G1736 |
| U2629 | A2424 | C2365 | C2238 | G2302 | G2238 | U2173 | C | U1894 | A1818 | G1736 |
| C2630 | G2425 | C2366 | A2093 | A2303 | A2174 | G2174 | C | A1895 | A1819 | U1739 |
| A2631 | U2426 | A2367 | A2100 | U2298 | U2175 | U2176 | C | G1896 | A1820 | U1740 |
| U2632 | G2429 | G2368 | C2101 | G2304 | G2240 | U2177 | C | U1897 | U1821 | U1740 |
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| U2634 | C2431 | C2370 | C2108 | U2306 | U2177 | A2178 | C | U1901 | A1828 | G1744 |
| A2635 | A2432 | G2371 | U2109 | A2242 | A2243 | C2179 | C | G1902 | A1832 | A1749 |
| C2636 | A2433 | A2372 | C2109 | A2244 | A2244 | G2180 | C | U1903 | C1832 | A1750 |
| U2637 | U2434 | C2373 | C2110 | A2245 | A2245 | C2181 | C | G1904 | U1833 | A1750 |
| U2638 | G2435 | G2374 | U2111 | C2246 | C2246 | C2182 | C | G1905 | U1834 | G1751 |
| U2639 | U2437 | C2375 | U2112 | G2247 | G2247 | A2183 | C | G1906 | A1835 | A1752 |
| A2640 | A2438 | U2376 | U2113 | C2248 | C2248 | U2184 | C | U1907 | C1836 | G1754 |
| U2641 | U2439 | U2377 | U2114 | C2249 | C2249 | U2185 | C | A1908 | U1837 | G1754 |
| A2642 | G2440 | A2378 | G2115 | G2250 | G2250 | A2186 | C | A1909 | G1838 | G1758 |
| U2643 | A2441 | C2379 | G2116 | G2251 | G2251 | U2187 | C | U1910 | A1839 | C1759 |
| G2644 | G2442 | U2380 | G2117 | A2252 | A2252 | C2192 | C | A1911 | U1840 | A1760 |
| C2645 | A2443 | G2381 | U2118 | G2253 | G2253 | U2193 | C | U1912 | A1841 | C1761 |
| C2646 | U | A2382 | U2119 | U2254 | U2254 | A2194 | C | A1913 | A1842 | C1762 |
| A2647 | A | A2383 | U2120 | A2255 | A2255 | G2195 | C | G1914 | C1843 | U1763 |
| G2648 | G | U2384 | G2121 | U2256 | U2256 | G2196 | C | U1917 | U1765 | U1765 |
| U2649 | A | A2385 | G2122 | A2257 | A2257 | A2197 | C | C1918 | C1846 | G1766 |
| U2650 | G | U2386 | G2123 | C2258 | C2258 | U2198 | C | U1919 | A1847 | C1767 |
| G2651 | C | A2387 | G2124 | U2259 | U2259 | A2199 | C | U1920 | A1848 | C1767 |
| C2652 | U | U2388 | U2125 | U2260 | U2260 | G2201 | C | A1921 | C1849 | G1770 |
| U2653 | A | G2389 | G2126 | G2261 | G2261 | C2202 | C | U1925 | A1850 | G1771 |
| U2654 | G | A2390 | U2127 | U2262 | U2262 | U2203 | C | U1926 | U1851 | C1772 |
| U2655 | C | G2391 | C2128 | U2263 | U2263 | C2204 | C | G1927 | U1855 | C1773 |
| U2656 | U | U2392 | U2129 | U2264 | U2264 | C2263 | C | U1856 | C1857 | G1778 |
| U2657 | C | C2393 | G2130 | U2265 | U2265 | A2131 | C | C1857 | C1857 | G1778 |
| A2658 | C | G2394 | A2131 | U2266 | U2266 | C2264 | C | C1857 | C1857 | G1778 |
| G2659 | U | C2395 | U2267 | U2267 | U2267 | C2265 | C | C1857 | C1857 | G1778 |



● Molecule 49: 5S RIBOSOMAL RNA



● Molecule 50: 5.8S RIBOSOMAL RNA



4 Experimental information

| Property | Value | Source |
|--------------------------------------|--------------------------------|-----------|
| EM reconstruction method | SINGLE PARTICLE | Depositor |
| Imposed symmetry | POINT, C1 | Depositor |
| Number of particles used | 84113 | Depositor |
| Resolution determination method | Not provided | |
| CTF correction method | PER FRAME | Depositor |
| Microscope | FEI TECNAI 20 | Depositor |
| Voltage (kV) | 200 | Depositor |
| Electron dose ($e^-/\text{\AA}^2$) | 20 | Depositor |
| Minimum defocus (nm) | 1500 | Depositor |
| Maximum defocus (nm) | 4500 | Depositor |
| Magnification | 83000 | Depositor |
| Image detector | GATAN ULTRASCAN 4000 (4k x 4k) | Depositor |
| Maximum map value | 425.313 | Depositor |
| Minimum map value | -221.732 | Depositor |
| Average map value | -13.190 | Depositor |
| Map value standard deviation | 29.801 | Depositor |
| Recommended contour level | 35.0 | Depositor |
| Map size (\AA) | 405.44, 405.44, 405.44 | wwPDB |
| Map dimensions | 224, 224, 224 | wwPDB |
| Map angles ($^\circ$) | 90.0, 90.0, 90.0 | wwPDB |
| Pixel spacing (\AA) | 1.81, 1.81, 1.81 | Depositor |

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Chain | Bond lengths | | Bond angles | |
|-----|-------|--------------|----------------|-------------|----------------|
| | | RMSZ | # Z >5 | RMSZ | # Z >5 |
| 1 | A | 0.87 | 1/1946 (0.1%) | 1.05 | 4/2614 (0.2%) |
| 2 | B | 1.02 | 4/3146 (0.1%) | 1.11 | 14/4228 (0.3%) |
| 3 | C | 0.87 | 0/2800 | 1.07 | 11/3790 (0.3%) |
| 4 | D | 0.89 | 1/2408 (0.0%) | 0.96 | 3/3248 (0.1%) |
| 5 | E | 0.90 | 1/1269 (0.1%) | 1.00 | 3/1705 (0.2%) |
| 6 | F | 0.99 | 1/1828 (0.1%) | 1.04 | 6/2461 (0.2%) |
| 7 | G | 0.64 | 0/1795 | 0.81 | 1/2429 (0.0%) |
| 8 | H | 0.97 | 2/1539 (0.1%) | 1.01 | 1/2073 (0.0%) |
| 9 | I | 0.92 | 1/1758 (0.1%) | 1.08 | 11/2358 (0.5%) |
| 10 | J | 0.81 | 1/1374 (0.1%) | 0.99 | 3/1842 (0.2%) |
| 12 | L | 0.82 | 0/1573 | 1.04 | 6/2113 (0.3%) |
| 13 | M | 0.95 | 0/1074 | 1.01 | 4/1446 (0.3%) |
| 14 | N | 0.83 | 0/1757 | 1.00 | 6/2354 (0.3%) |
| 15 | O | 0.98 | 11/3159 (0.3%) | 1.02 | 25/4205 (0.6%) |
| 16 | P | 1.05 | 1/1250 (0.1%) | 1.09 | 5/1683 (0.3%) |
| 17 | Q | 0.89 | 1/1465 (0.1%) | 1.12 | 8/1965 (0.4%) |
| 18 | R | 0.78 | 1/1538 (0.1%) | 0.87 | 3/2050 (0.1%) |
| 19 | S | 1.02 | 0/1481 | 1.09 | 7/1990 (0.4%) |
| 20 | T | 1.01 | 2/1300 (0.2%) | 1.01 | 1/1743 (0.1%) |
| 21 | U | 0.56 | 0/794 | 0.77 | 0/1076 |
| 22 | V | 0.98 | 0/1018 | 1.09 | 4/1369 (0.3%) |
| 23 | W | 0.80 | 0/1052 | 0.90 | 1/1398 (0.1%) |
| 24 | X | 0.72 | 0/974 | 0.86 | 0/1314 |
| 25 | Y | 0.79 | 1/1004 (0.1%) | 0.98 | 2/1341 (0.1%) |
| 26 | Z | 0.55 | 0/1118 | 0.83 | 2/1497 (0.1%) |
| 27 | a | 0.95 | 2/1204 (0.2%) | 1.14 | 9/1612 (0.6%) |
| 28 | b | 0.91 | 0/473 | 1.14 | 1/629 (0.2%) |
| 29 | c | 0.61 | 0/775 | 0.77 | 0/1040 |
| 30 | d | 0.94 | 2/897 (0.2%) | 0.95 | 1/1205 (0.1%) |
| 31 | e | 1.04 | 0/1041 | 1.27 | 12/1394 (0.9%) |
| 32 | f | 1.12 | 1/868 (0.1%) | 1.09 | 3/1168 (0.3%) |
| 33 | g | 0.72 | 0/890 | 0.92 | 0/1189 |

| Mol | Chain | Bond lengths | | Bond angles | |
|-----|-------|--------------|-------------------|-------------|--------------------|
| | | RMSZ | # Z >5 | RMSZ | # Z >5 |
| 34 | h | 0.67 | 0/974 | 0.79 | 0/1297 |
| 35 | i | 0.67 | 0/777 | 0.85 | 0/1033 |
| 36 | j | 0.87 | 0/696 | 1.04 | 3/923 (0.3%) |
| 37 | k | 0.50 | 0/614 | 0.70 | 0/822 |
| 38 | l | 0.90 | 0/443 | 1.02 | 1/588 (0.2%) |
| 39 | m | 1.08 | 2/423 (0.5%) | 1.13 | 1/562 (0.2%) |
| 40 | n | 0.90 | 0/234 | 1.15 | 1/300 (0.3%) |
| 41 | o | 0.83 | 0/860 | 0.88 | 1/1136 (0.1%) |
| 42 | p | 0.86 | 0/701 | 0.98 | 1/934 (0.1%) |
| 43 | q | 0.53 | 0/1092 | 0.73 | 1/1474 (0.1%) |
| 46 | t | 5.64 | 18/2985 (0.6%) | 4.15 | 194/4053 (4.8%) |
| 48 | 5 | 1.46 | 609/75414 (0.8%) | 1.88 | 3517/117575 (3.0%) |
| 49 | 7 | 1.38 | 13/2883 (0.5%) | 1.80 | 118/4491 (2.6%) |
| 50 | 8 | 1.16 | 5/3746 (0.1%) | 1.70 | 132/5832 (2.3%) |
| All | All | 1.49 | 681/138410 (0.5%) | 1.70 | 4127/203549 (2.0%) |

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 |
|-----|-------|---------------------|---------------------|
| 1 | A | 0 | 2 |
| 3 | C | 0 | 1 |
| 4 | D | 0 | 1 |
| 5 | E | 0 | 1 |
| 6 | F | 0 | 2 |
| 15 | O | 0 | 2 |
| 19 | S | 0 | 1 |
| 22 | V | 0 | 1 |
| 25 | Y | 0 | 1 |
| 26 | Z | 0 | 1 |
| 27 | a | 0 | 3 |
| 28 | b | 0 | 1 |
| 46 | t | 0 | 6 |
| 48 | 5 | 0 | 1 |
| All | All | 0 | 24 |

The worst 5 of 681 bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|-------|--------|-------------|----------|
| 46 | t | 168 | PRO | N-CD | 120.75 | 3.16 | 1.47 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|-------|--------|-------------|----------|
| 46 | t | 545 | PRO | N-CD | 120.48 | 3.16 | 1.47 |
| 46 | t | 162 | PRO | N-CD | 120.12 | 3.16 | 1.47 |
| 46 | t | 172 | PRO | N-CD | 118.10 | 3.13 | 1.47 |
| 46 | t | 520 | PRO | N-CD | 117.23 | 3.12 | 1.47 |

The worst 5 of 4127 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|-----|------|---------|--------|-------------|----------|
| 46 | t | 81 | LYS | O-C-N | -73.85 | 4.54 | 122.70 |
| 46 | t | 15 | LYS | O-C-N | -50.53 | 41.85 | 122.70 |
| 46 | t | 544 | ASN | O-C-N | -46.67 | 32.42 | 121.10 |
| 46 | t | 162 | PRO | N-CA-CB | 37.84 | 148.71 | 103.30 |
| 46 | t | 168 | PRO | N-CA-CB | 37.79 | 148.65 | 103.30 |

There are no chirality outliers.

5 of 24 planarity outliers are listed below:

| Mol | Chain | Res | Type | Group |
|-----|-------|-----|------|---------|
| 1 | A | 143 | GLU | Peptide |
| 1 | A | 211 | HIS | Peptide |
| 3 | C | 91 | GLY | Peptide |
| 4 | D | 271 | LYS | Peptide |
| 5 | E | 129 | GLU | Peptide |

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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 | A | 250/254 (98%) | 213 (85%) | 30 (12%) | 7 (3%) | 5 | 30 |
| 2 | B | 384/387 (99%) | 341 (89%) | 34 (9%) | 9 (2%) | 6 | 34 |
| 3 | C | 359/362 (99%) | 306 (85%) | 32 (9%) | 21 (6%) | 1 | 18 |
| 4 | D | 292/297 (98%) | 267 (91%) | 19 (6%) | 6 (2%) | 7 | 36 |
| 5 | E | 153/176 (87%) | 134 (88%) | 15 (10%) | 4 (3%) | 5 | 31 |
| 6 | F | 221/244 (91%) | 201 (91%) | 15 (7%) | 5 (2%) | 6 | 34 |
| 7 | G | 229/256 (90%) | 181 (79%) | 27 (12%) | 21 (9%) | 1 | 11 |
| 8 | H | 189/191 (99%) | 172 (91%) | 13 (7%) | 4 (2%) | 7 | 36 |
| 9 | I | 209/221 (95%) | 175 (84%) | 22 (10%) | 12 (6%) | 1 | 18 |
| 10 | J | 167/174 (96%) | 135 (81%) | 19 (11%) | 13 (8%) | 1 | 13 |
| 12 | L | 192/199 (96%) | 161 (84%) | 20 (10%) | 11 (6%) | 1 | 18 |
| 13 | M | 135/138 (98%) | 124 (92%) | 10 (7%) | 1 (1%) | 22 | 63 |
| 14 | N | 201/204 (98%) | 182 (90%) | 13 (6%) | 6 (3%) | 4 | 28 |
| 15 | O | 352/219 (161%) | 324 (92%) | 18 (5%) | 10 (3%) | 5 | 30 |
| 16 | P | 153/184 (83%) | 142 (93%) | 9 (6%) | 2 (1%) | 12 | 48 |
| 17 | Q | 183/186 (98%) | 168 (92%) | 9 (5%) | 6 (3%) | 4 | 26 |
| 18 | R | 186/189 (98%) | 167 (90%) | 16 (9%) | 3 (2%) | 9 | 44 |
| 19 | S | 170/172 (99%) | 163 (96%) | 6 (4%) | 1 (1%) | 25 | 66 |
| 20 | T | 157/160 (98%) | 146 (93%) | 9 (6%) | 2 (1%) | 12 | 48 |
| 21 | U | 96/121 (79%) | 80 (83%) | 13 (14%) | 3 (3%) | 4 | 27 |
| 22 | V | 134/137 (98%) | 124 (92%) | 8 (6%) | 2 (2%) | 10 | 46 |
| 23 | W | 133/155 (86%) | 106 (80%) | 19 (14%) | 8 (6%) | 1 | 17 |
| 24 | X | 118/142 (83%) | 103 (87%) | 7 (6%) | 8 (7%) | 1 | 15 |
| 25 | Y | 124/127 (98%) | 107 (86%) | 12 (10%) | 5 (4%) | 3 | 23 |
| 26 | Z | 133/136 (98%) | 107 (80%) | 13 (10%) | 13 (10%) | 0 | 10 |
| 27 | a | 146/149 (98%) | 123 (84%) | 18 (12%) | 5 (3%) | 3 | 26 |
| 28 | b | 56/59 (95%) | 44 (79%) | 7 (12%) | 5 (9%) | 1 | 11 |
| 29 | c | 98/105 (93%) | 87 (89%) | 8 (8%) | 3 (3%) | 4 | 27 |
| 30 | d | 107/113 (95%) | 88 (82%) | 13 (12%) | 6 (6%) | 2 | 19 |
| 31 | e | 125/130 (96%) | 109 (87%) | 10 (8%) | 6 (5%) | 2 | 21 |
| 32 | f | 104/107 (97%) | 96 (92%) | 5 (5%) | 3 (3%) | 4 | 29 |
| 33 | g | 110/121 (91%) | 93 (84%) | 13 (12%) | 4 (4%) | 3 | 25 |

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| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Percentiles | |
|-----|-------|-----------------|------------|----------|----------|-------------|-----|
| 34 | h | 117/120 (98%) | 99 (85%) | 14 (12%) | 4 (3%) | 3 | 26 |
| 35 | i | 97/100 (97%) | 77 (79%) | 13 (13%) | 7 (7%) | 1 | 14 |
| 36 | j | 85/88 (97%) | 75 (88%) | 8 (9%) | 2 (2%) | 6 | 33 |
| 37 | k | 75/78 (96%) | 61 (81%) | 10 (13%) | 4 (5%) | 2 | 19 |
| 38 | l | 48/51 (94%) | 41 (85%) | 6 (12%) | 1 (2%) | 7 | 36 |
| 39 | m | 50/128 (39%) | 48 (96%) | 1 (2%) | 1 (2%) | 7 | 38 |
| 40 | n | 23/25 (92%) | 22 (96%) | 0 | 1 (4%) | 2 | 22 |
| 41 | o | 103/106 (97%) | 90 (87%) | 11 (11%) | 2 (2%) | 8 | 38 |
| 42 | p | 89/92 (97%) | 81 (91%) | 8 (9%) | 0 | 100 | 100 |
| 43 | q | 139/312 (45%) | 111 (80%) | 21 (15%) | 7 (5%) | 2 | 20 |
| 46 | t | 376/614 (61%) | 354 (94%) | 14 (4%) | 8 (2%) | 7 | 36 |
| All | All | 6868/7529 (91%) | 6028 (88%) | 588 (9%) | 252 (4%) | 6 | 24 |

5 of 252 Ramachandran outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | A | 96 | LEU |
| 2 | B | 129 | ALA |
| 2 | B | 140 | ASP |
| 2 | B | 347 | SER |
| 3 | C | 14 | GLU |

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all 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 | A | 192/196 (98%) | 154 (80%) | 38 (20%) | 1 | 8 |
| 2 | B | 321/323 (99%) | 251 (78%) | 70 (22%) | 1 | 6 |
| 3 | C | 288/289 (100%) | 222 (77%) | 66 (23%) | 1 | 4 |
| 4 | D | 243/245 (99%) | 196 (81%) | 47 (19%) | 1 | 8 |
| 5 | E | 135/153 (88%) | 115 (85%) | 20 (15%) | 3 | 15 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|----------------|-----------|----------|-------------|----|
| 6 | F | 187/205 (91%) | 158 (84%) | 29 (16%) | 2 | 14 |
| 7 | G | 177/208 (85%) | 138 (78%) | 39 (22%) | 1 | 5 |
| 8 | H | 171/171 (100%) | 132 (77%) | 39 (23%) | 1 | 5 |
| 9 | I | 179/187 (96%) | 142 (79%) | 37 (21%) | 1 | 6 |
| 10 | J | 147/150 (98%) | 114 (78%) | 33 (22%) | 1 | 5 |
| 12 | L | 154/159 (97%) | 124 (80%) | 30 (20%) | 1 | 8 |
| 13 | M | 108/109 (99%) | 84 (78%) | 24 (22%) | 1 | 6 |
| 14 | N | 175/176 (99%) | 143 (82%) | 32 (18%) | 1 | 10 |
| 15 | O | 323/179 (180%) | 267 (83%) | 56 (17%) | 2 | 11 |
| 16 | P | 125/146 (86%) | 103 (82%) | 22 (18%) | 2 | 11 |
| 17 | Q | 150/151 (99%) | 123 (82%) | 27 (18%) | 1 | 10 |
| 18 | R | 153/154 (99%) | 121 (79%) | 32 (21%) | 1 | 6 |
| 19 | S | 156/156 (100%) | 123 (79%) | 33 (21%) | 1 | 6 |
| 20 | T | 136/137 (99%) | 109 (80%) | 27 (20%) | 1 | 7 |
| 21 | U | 85/107 (79%) | 62 (73%) | 23 (27%) | 0 | 3 |
| 22 | V | 104/105 (99%) | 96 (92%) | 8 (8%) | 13 | 37 |
| 23 | W | 100/129 (78%) | 85 (85%) | 15 (15%) | 3 | 15 |
| 24 | X | 104/118 (88%) | 81 (78%) | 23 (22%) | 1 | 5 |
| 25 | Y | 109/110 (99%) | 85 (78%) | 24 (22%) | 1 | 5 |
| 26 | Z | 115/116 (99%) | 89 (77%) | 26 (23%) | 1 | 5 |
| 27 | a | 118/119 (99%) | 95 (80%) | 23 (20%) | 1 | 8 |
| 28 | b | 46/47 (98%) | 35 (76%) | 11 (24%) | 0 | 4 |
| 29 | c | 84/88 (96%) | 69 (82%) | 15 (18%) | 2 | 10 |
| 30 | d | 94/97 (97%) | 73 (78%) | 21 (22%) | 1 | 5 |
| 31 | e | 109/111 (98%) | 89 (82%) | 20 (18%) | 1 | 10 |
| 32 | f | 90/91 (99%) | 79 (88%) | 11 (12%) | 5 | 20 |
| 33 | g | 95/103 (92%) | 71 (75%) | 24 (25%) | 0 | 3 |
| 34 | h | 103/105 (98%) | 77 (75%) | 26 (25%) | 0 | 3 |
| 35 | i | 80/82 (98%) | 51 (64%) | 29 (36%) | 0 | 1 |
| 36 | j | 70/71 (99%) | 53 (76%) | 17 (24%) | 0 | 4 |
| 37 | k | 67/69 (97%) | 53 (79%) | 14 (21%) | 1 | 6 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|-----------------|------------|------------|-------------|----|
| 38 | l | 45/46 (98%) | 34 (76%) | 11 (24%) | 0 | 4 |
| 39 | m | 47/116 (40%) | 34 (72%) | 13 (28%) | 0 | 3 |
| 40 | n | 23/23 (100%) | 16 (70%) | 7 (30%) | 0 | 2 |
| 41 | o | 90/91 (99%) | 74 (82%) | 16 (18%) | 2 | 10 |
| 42 | p | 71/72 (99%) | 61 (86%) | 10 (14%) | 3 | 16 |
| 43 | q | 105/254 (41%) | 76 (72%) | 29 (28%) | 0 | 3 |
| 46 | t | 332/539 (62%) | 330 (99%) | 2 (1%) | 86 | 92 |
| All | All | 5806/6303 (92%) | 4687 (81%) | 1119 (19%) | 4 | 8 |

5 of 1119 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 33 | g | 23 | VAL |
| 34 | h | 47 | VAL |
| 33 | g | 20 | ILE |
| 38 | l | 47 | THR |
| 10 | J | 16 | LYS |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 31 such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 16 | P | 55 | GLN |
| 46 | t | 82 | ASN |
| 20 | T | 49 | GLN |
| 46 | t | 394 | HIS |
| 34 | h | 20 | GLN |

5.3.3 RNA [i](#)

| Mol | Chain | Analysed | Backbone Outliers | Pucker Outliers |
|-----|-------|-----------------|-------------------|-----------------|
| 47 | 1 | 0/114 | - | - |
| 48 | 5 | 3145/3396 (92%) | 731 (23%) | 129 (4%) |
| 49 | 7 | 120/121 (99%) | 18 (15%) | 0 |
| 50 | 8 | 157/158 (99%) | 32 (20%) | 3 (1%) |
| All | All | 3422/3789 (90%) | 781 (22%) | 132 (3%) |

5 of 781 RNA backbone outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 48 | 5 | 14 | U |
| 48 | 5 | 15 | C |
| 48 | 5 | 16 | A |
| 48 | 5 | 26 | A |
| 48 | 5 | 38 | U |

5 of 132 RNA pucker outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|------|------|
| 48 | 5 | 3155 | U |
| 48 | 5 | 3228 | C |
| 50 | 8 | 126 | A |
| 48 | 5 | 1284 | C |
| 48 | 5 | 1241 | U |

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

The following chains have linkage breaks:

| Mol | Chain | Number of breaks |
|-----|-------|------------------|
| 11 | K | 2 |

All chain breaks are listed below:

| Model | Chain | Residue-1 | Atom-1 | Residue-2 | Atom-2 | Distance (Å) |
|-------|-------|-----------|--------|-----------|--------|--------------|
| 1 | K | 52:UNK | C | 54:UNK | N | 3.86 |
| 1 | K | 23:UNK | C | 28:UNK | N | 3.48 |

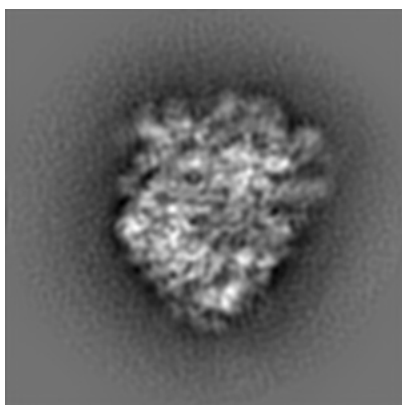
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-2169. These allow visual inspection of the internal detail of the map and identification of artifacts.

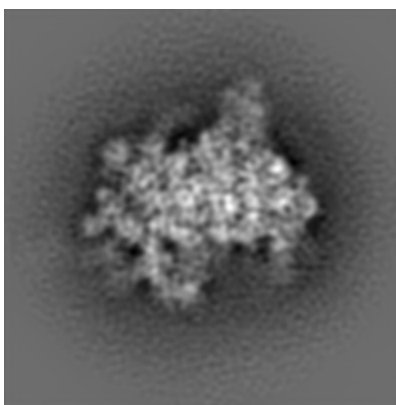
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

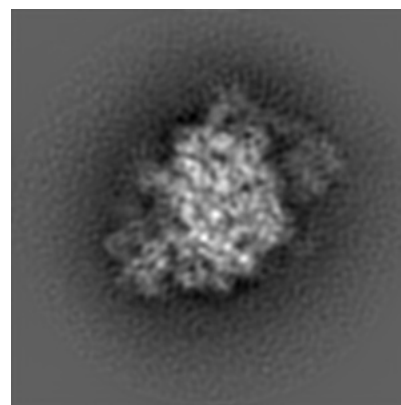
6.1.1 Primary map



X



Y

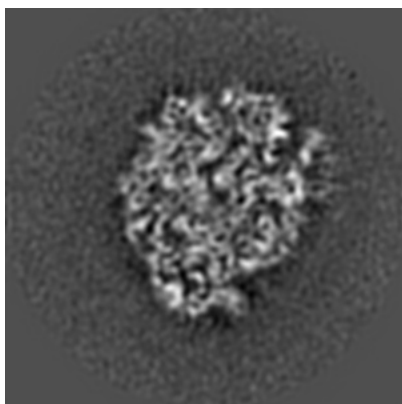


Z

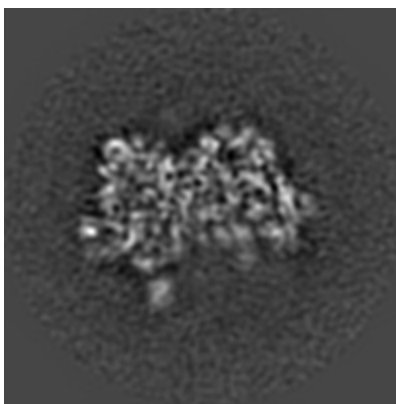
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

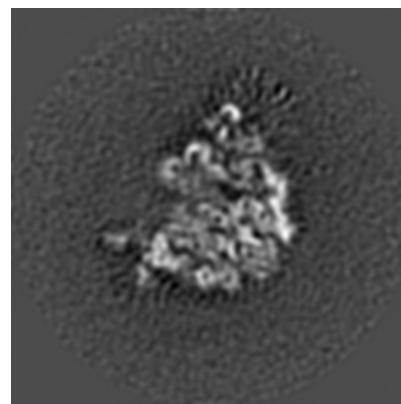
6.2.1 Primary map



X Index: 112



Y Index: 112

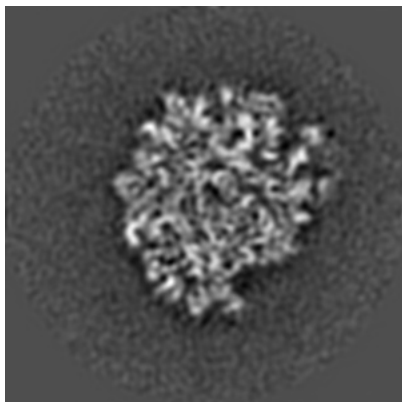


Z Index: 112

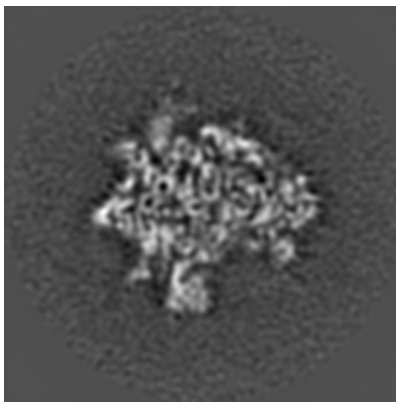
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

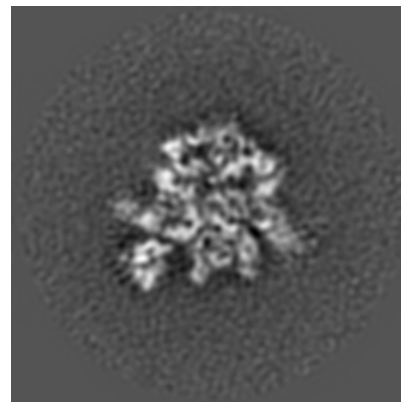
6.3.1 Primary map



X Index: 115



Y Index: 95

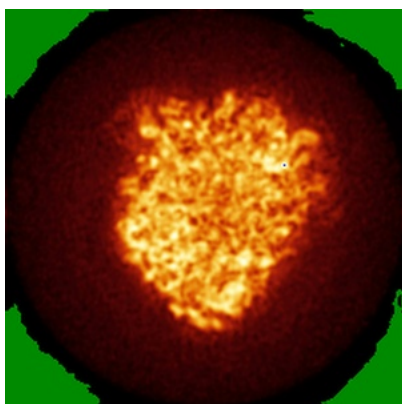


Z Index: 84

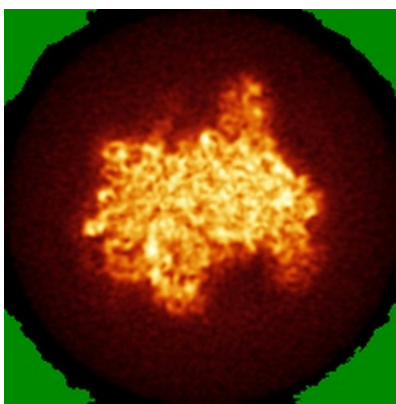
The images above show the largest variance slices of the map in three orthogonal directions.

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

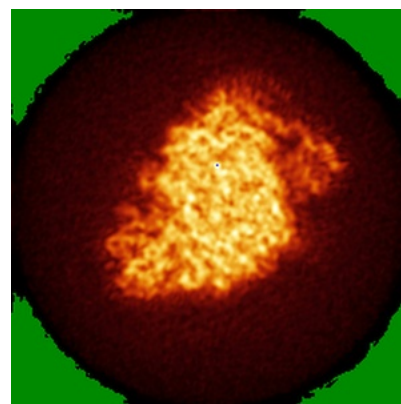
6.4.1 Primary map



X



Y

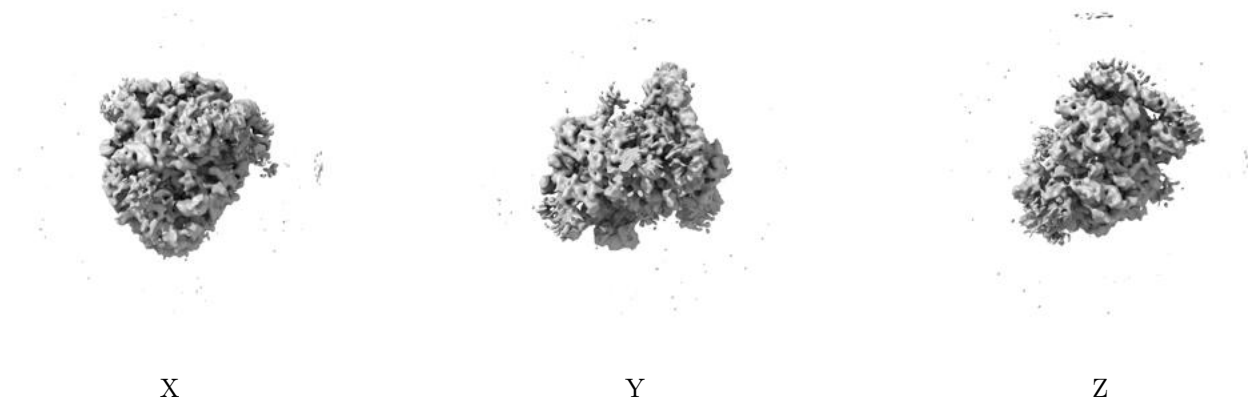


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 35.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

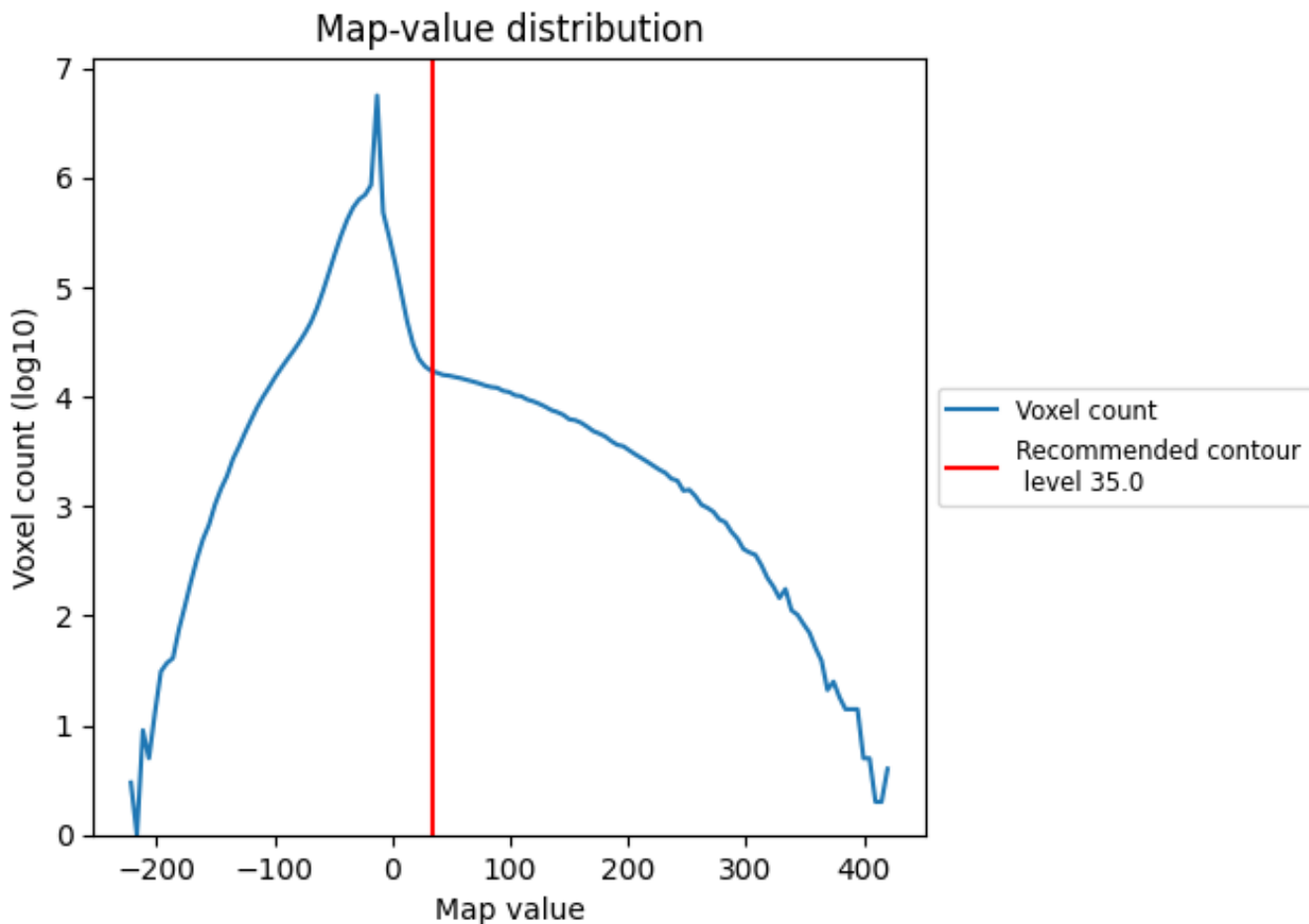
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

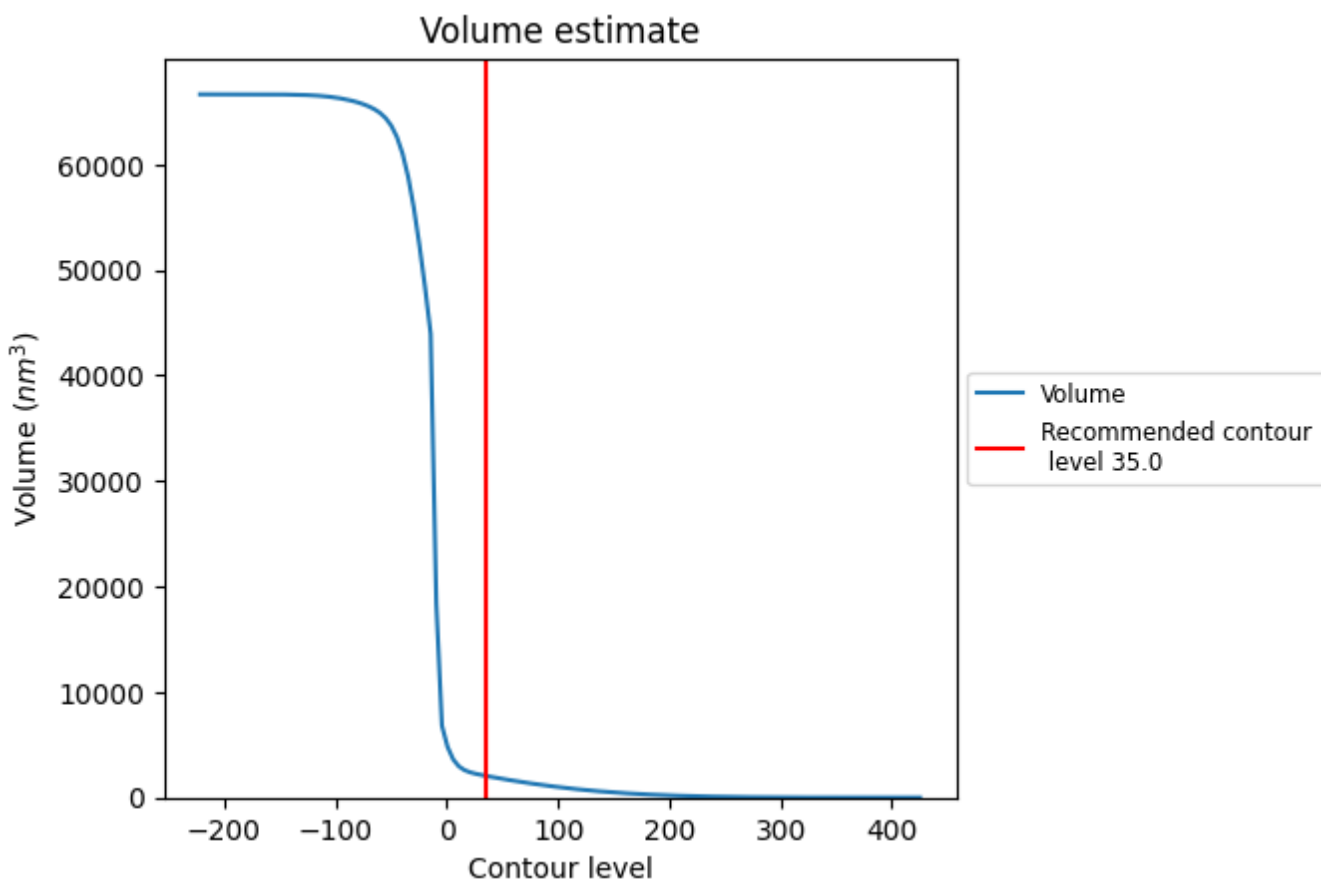
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

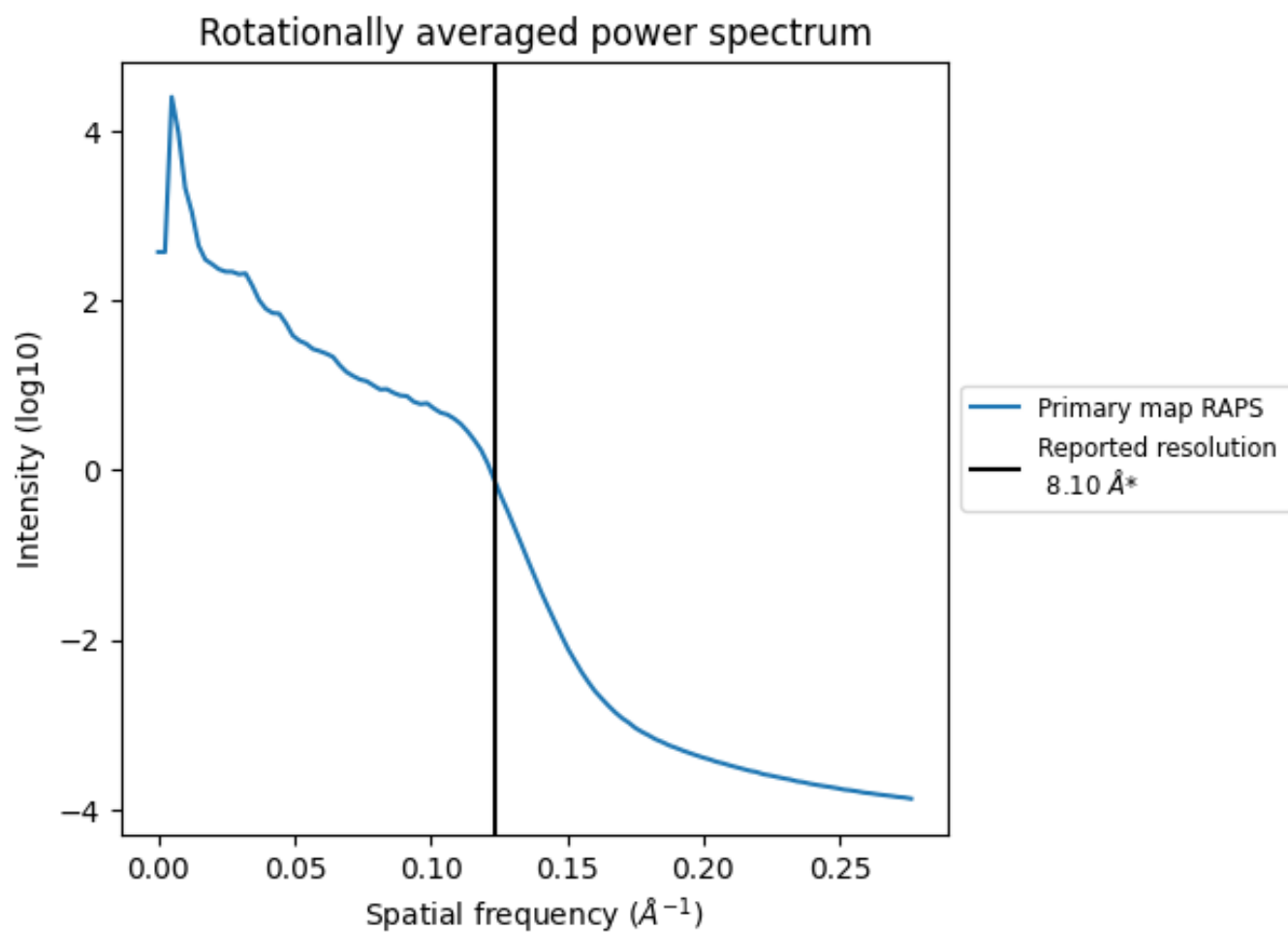
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2057 nm³; this corresponds to an approximate mass of 1858 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 [i](#)



*Reported resolution corresponds to spatial frequency of 0.123\AA^{-1}

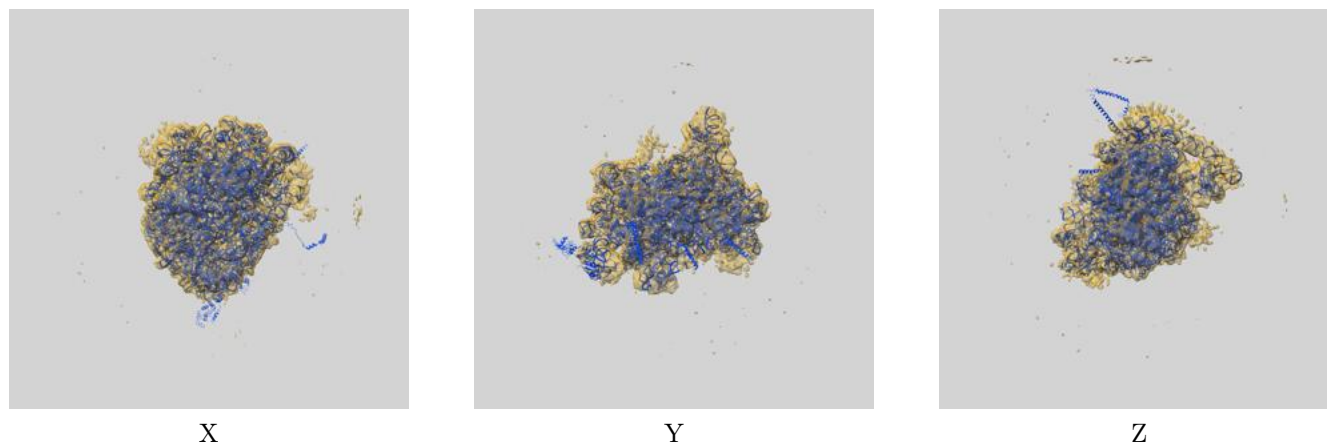
8 Fourier-Shell correlation

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

9 Map-model fit [i](#)

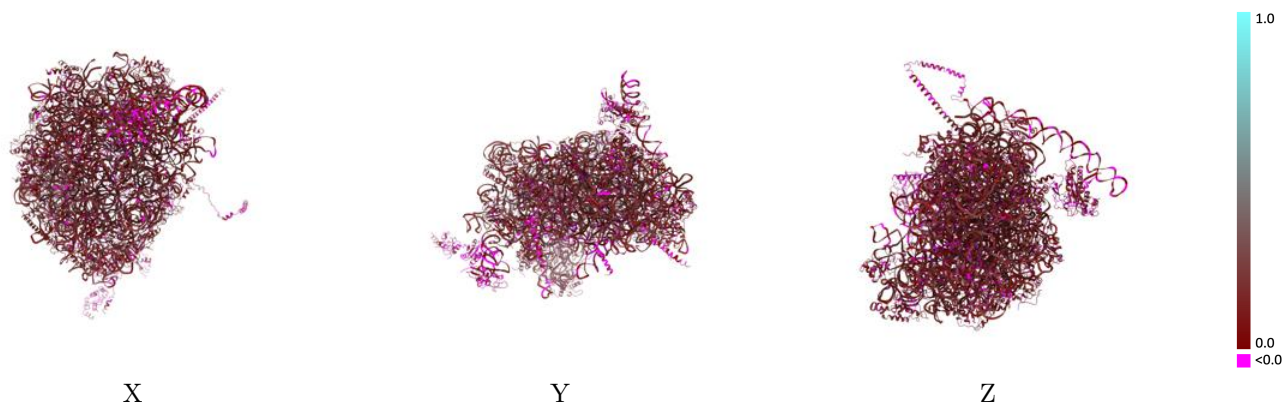
This section contains information regarding the fit between EMDB map EMD-2169 and PDB model 4V8T. Per-residue inclusion information can be found in section 3 on page 14.

9.1 Map-model overlay [i](#)



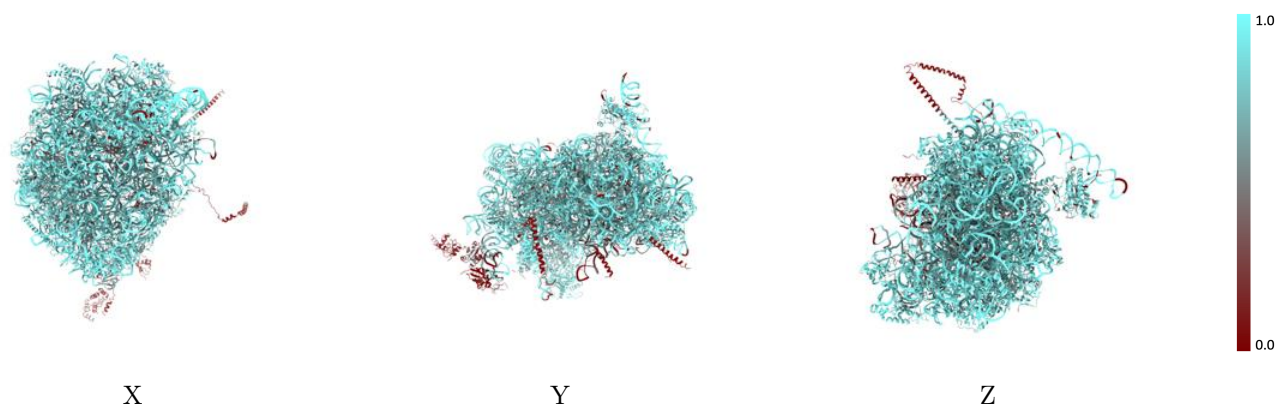
The images above show the 3D surface view of the map at the recommended contour level 35.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



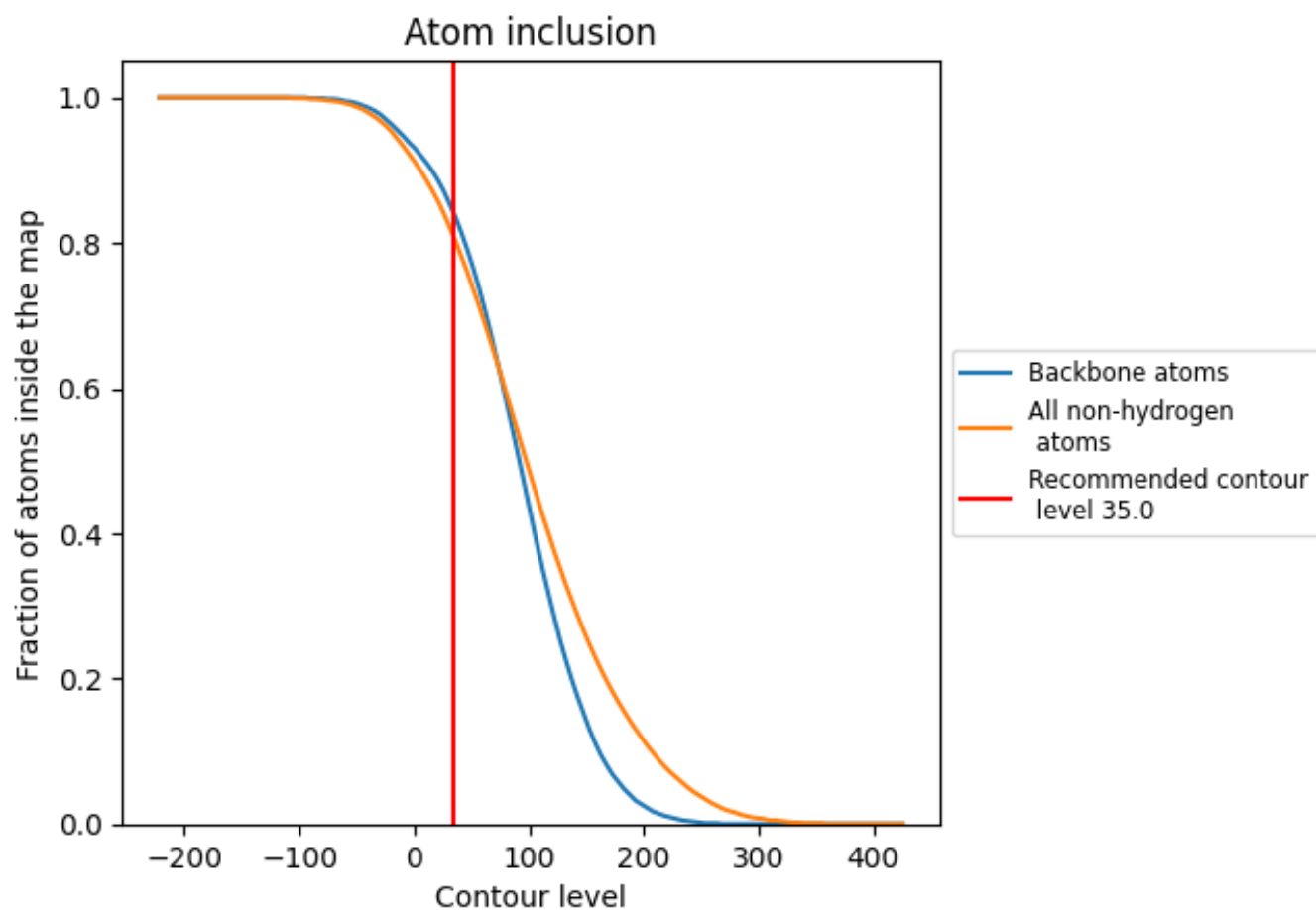
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (35.0).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 84% of all backbone atoms, 81% 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 (35.0) and Q-score for the entire model and for each chain.

| Chain | Atom inclusion | Q-score |
|-------|--|--|
| All |  0.8060 |  0.1350 |
| 1 |  0.9210 |  0.0730 |
| 5 |  0.8850 |  0.1540 |
| 7 |  0.9700 |  0.1670 |
| 8 |  0.9200 |  0.1690 |
| A |  0.5420 |  0.0710 |
| B |  0.7000 |  0.1060 |
| C |  0.7170 |  0.1120 |
| D |  0.8100 |  0.1160 |
| E |  0.8080 |  0.1310 |
| F |  0.7420 |  0.1380 |
| G |  0.8230 |  0.1300 |
| H |  0.8120 |  0.1290 |
| I |  0.6820 |  0.0940 |
| J |  0.8520 |  0.1170 |
| K |  0.1310 |  0.0320 |
| L |  0.7830 |  0.1380 |
| M |  0.8550 |  0.1500 |
| N |  0.5560 |  0.0810 |
| O |  0.7690 |  0.1420 |
| P |  0.6700 |  0.1020 |
| Q |  0.6930 |  0.1130 |
| R |  0.5720 |  0.1050 |
| S |  0.7620 |  0.1180 |
| T |  0.7100 |  0.1150 |
| U |  0.7660 |  0.1080 |
| V |  0.6610 |  0.1040 |
| W |  0.3990 |  0.0710 |
| X |  0.7100 |  0.1280 |
| Y |  0.7650 |  0.1130 |
| Z |  0.7740 |  0.1330 |
| a |  0.6720 |  0.1030 |
| b |  0.6130 |  0.1190 |
| c |  0.7960 |  0.1400 |
| d |  0.7210 |  0.1200 |



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| Chain | Atom inclusion | Q-score |
|-------|--|---|
| e |  0.6100 |  0.1120 |
| f |  0.7350 |  0.1050 |
| g |  0.5900 |  0.0950 |
| h |  0.7940 |  0.1410 |
| i |  0.7570 |  0.1360 |
| j |  0.6720 |  0.0960 |
| k |  0.8340 |  0.1360 |
| l |  0.5300 |  0.1020 |
| m |  0.7150 |  0.1130 |
| n |  0.0000 |  -0.0490 |
| o |  0.6250 |  0.0810 |
| p |  0.7700 |  0.1110 |
| q |  0.2200 |  0.0110 |
| r |  0.0000 |  0.0090 |
| s |  0.0000 |  0.0190 |
| t |  0.8060 |  0.0840 |