



Full wwPDB EM Validation Report ⓘ

Oct 13, 2024 – 03:30 PM EDT

PDB ID : 7SYR
EMDB ID : EMD-25538
Title : Structure of the wt IRES eIF2-containing 48S initiation complex, closed conformation. Structure 12(wt).
Authors : Brown, Z.P.; Abaeva, I.S.; De, S.; Hellen, C.U.T.; Pestova, T.V.; Frank, J.
Deposited on : 2021-11-25
Resolution : 3.60 Å(reported)
Based on initial models : 6D9J, 4KZZ, 5FLX, 6O85, 5K0Y

This is a Full wwPDB EM Validation 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.dev113
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

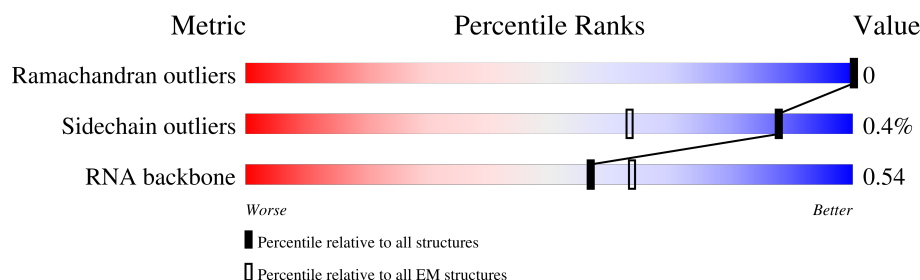
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.60 Å.

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



| Metric | Whole archive (#Entries) | EM structures (#Entries) |
|-----------------------|-----------------------------|-----------------------------|
| Ramachandran outliers | 207382 | 16835 |
| Sidechain outliers | 206894 | 16415 |
| RNA backbone | 6643 | 2191 |

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 | 2 | 1870 | |
| 2 | A | 144 | |
| 3 | B | 295 | |
| 4 | C | 264 | |
| 5 | D | 221 | |
| 6 | E | 281 | |
| 7 | F | 263 | |
| 8 | G | 204 | |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 9 | H | 249 | |
| 10 | I | 432 | |
| 11 | J | 208 | |
| 12 | K | 194 | |
| 13 | L | 149 | |
| 14 | M | 158 | |
| 15 | N | 132 | |
| 16 | O | 151 | |
| 17 | P | 168 | |
| 18 | Q | 145 | |
| 19 | R | 172 | |
| 20 | S | 135 | |
| 21 | T | 152 | |
| 22 | U | 145 | |
| 23 | V | 119 | |
| 24 | W | 83 | |
| 25 | X | 130 | |
| 26 | Y | 143 | |
| 27 | Z | 131 | |
| 28 | a | 124 | |
| 29 | b | 101 | |
| 30 | c | 84 | |
| 31 | d | 69 | |
| 32 | e | 56 | |
| 33 | f | 133 | |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|--|
| 34 | g | 188 | <div><div><div></div><div></div><div></div></div><div>31%36%64%</div></div> |
| 35 | h | 317 | <div><div><div></div><div></div><div></div></div><div>16%99%</div></div> |
| 36 | i | 75 | <div><div><div></div><div></div><div></div></div><div>17%80%19%</div></div> |
| 37 | j | 315 | <div><div><div></div><div></div><div></div></div><div>57%42%</div></div> |
| 38 | n | 25 | <div><div><div></div><div></div><div></div></div><div>20%100%</div></div> |
| 39 | z | 400 | <div><div><div></div><div></div><div></div></div><div>10%34%13%53%</div></div> |

2 Entry composition

There are 40 unique types of molecules in this entry. The entry contains 83005 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 RNA chain called 18S rRNA.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-------|------|-------|------|---------|-------|
| 1 | 2 | 1697 | Total | C | N | O | P | 0 | 0 |
| | | | 36227 | 16170 | 6504 | 11857 | 1696 | | |

- Molecule 2 is a protein called Eukaryotic translation initiation factor 1A, X-chromosomal.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 2 | A | 99 | Total | C | N | O | S | 0 | 0 |
| | | | 798 | 503 | 143 | 148 | 4 | | |

- Molecule 3 is a protein called uS2 (SA).

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 3 | B | 217 | Total | C | N | O | S | 0 | 0 |
| | | | 1710 | 1086 | 300 | 316 | 8 | | |

There is a discrepancy between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------|------------|
| B | 114 | THR | ALA | conflict | UNP G1TLT8 |

- Molecule 4 is a protein called eS1.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|-------|
| 4 | C | 213 | Total | C | N | O | S | 0 | 0 |
| | | | 1729 | 1098 | 309 | 308 | 14 | | |

- Molecule 5 is a protein called uS5.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 5 | D | 221 | Total | C | N | O | S | 0 | 0 |
| | | | 1716 | 1111 | 295 | 301 | 9 | | |

- Molecule 6 is a protein called uS3.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 6 | E | 228 | Total | C | N | O | S | 0 | 0 |
| | | | 1768 | 1126 | 318 | 316 | 8 | | |

- Molecule 7 is a protein called eS4 (S4 X isoform).

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 7 | F | 262 | Total | C | N | O | S | 0 | 0 |
| | | | 2076 | 1324 | 386 | 358 | 8 | | |

There are 4 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------|------------|
| F | 25 | GLY | SER | conflict | UNP G1TK17 |
| F | 51 | ARG | LYS | conflict | UNP G1TK17 |
| F | 78 | THR | ALA | conflict | UNP G1TK17 |
| F | 156 | VAL | MET | conflict | UNP G1TK17 |

- Molecule 8 is a protein called uS7.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 8 | G | 191 | Total | C | N | O | S | 0 | 0 |
| | | | 1509 | 943 | 286 | 273 | 7 | | |

- Molecule 9 is a protein called eS6.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 9 | H | 237 | Total | C | N | O | S | 0 | 0 |
| | | | 1923 | 1200 | 387 | 329 | 7 | | |

- Molecule 10 is a protein called eS7.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 10 | I | 185 | Total | C | N | O | S | 0 | 0 |
| | | | 1488 | 952 | 271 | 264 | 1 | | |

- Molecule 11 is a protein called eS8.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|---|---------|-------|
| 11 | J | 206 | Total | C | N | O | S | 1 | 0 |
| | | | 1691 | 1061 | 333 | 292 | 5 | | |

There is a discrepancy between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------|------------|
| J | 47 | ARG | GLY | conflict | UNP G1TJW1 |

- Molecule 12 is a protein called uS4.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 12 | K | 185 | Total | C | N | O | S | 0 | 0 |
| | | | 1525 | 969 | 306 | 248 | 2 | | |

- Molecule 13 is a protein called eS10.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 13 | L | 96 | Total | C | N | O | S | 0 | 0 |
| | | | 810 | 530 | 143 | 131 | 6 | | |

- Molecule 14 is a protein called uS17.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 14 | M | 151 | Total | C | N | O | S | 0 | 0 |
| | | | 1233 | 785 | 231 | 211 | 6 | | |

- Molecule 15 is a protein called eS12.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 15 | N | 117 | Total | C | N | O | S | 0 | 0 |
| | | | 908 | 570 | 161 | 169 | 8 | | |

- Molecule 16 is a protein called uS15.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 16 | O | 149 | Total | C | N | O | S | 0 | 0 |
| | | | 1202 | 770 | 228 | 203 | 1 | | |

- Molecule 17 is a protein called uS11.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 17 | P | 136 | Total | C | N | O | S | 0 | 0 |
| | | | 1016 | 621 | 199 | 190 | 6 | | |

- Molecule 18 is a protein called uS19.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 18 | Q | 120 | Total | C | N | O | S | 0 | 0 |
| | | | 997 | 635 | 187 | 168 | 7 | | |

- Molecule 19 is a protein called uS9.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 19 | R | 142 | Total | C | N | O | S | 0 | 0 |
| | | | 1128 | 717 | 213 | 195 | 3 | | |

- Molecule 20 is a protein called eS17.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 20 | S | 132 | Total | C | N | O | S | 0 | 0 |
| | | | 1068 | 670 | 199 | 195 | 4 | | |

- Molecule 21 is a protein called uS13.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 21 | T | 144 | Total | C | N | O | S | 0 | 0 |
| | | | 1190 | 746 | 241 | 202 | 1 | | |

- Molecule 22 is a protein called eS19.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 22 | U | 141 | Total | C | N | O | S | 0 | 0 |
| | | | 1097 | 688 | 211 | 195 | 3 | | |

- Molecule 23 is a protein called uS10.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 23 | V | 100 | Total | C | N | O | S | 0 | 0 |
| | | | 795 | 498 | 152 | 141 | 4 | | |

- Molecule 24 is a protein called eS21.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 24 | W | 83 | Total | C | N | O | S | 0 | 0 |
| | | | 636 | 393 | 117 | 121 | 5 | | |

There are 7 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------|------------|
| W | 3 | ASN | SER | conflict | UNP G1TM82 |
| W | 4 | ASP | ASN | conflict | UNP G1TM82 |
| W | 33 | GLN | PRO | conflict | UNP G1TM82 |
| W | 50 | PHE | SER | conflict | UNP G1TM82 |
| W | 75 | ALA | SER | conflict | UNP G1TM82 |
| W | 76 | ASP | HIS | conflict | UNP G1TM82 |
| W | 81 | LYS | GLN | conflict | UNP G1TM82 |

- Molecule 25 is a protein called uS8.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 25 | X | 129 | Total | C | N | O | S | 0 | 0 |
| | | | 1034 | 659 | 193 | 176 | 6 | | |

- Molecule 26 is a protein called uS12.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 26 | Y | 141 | Total | C | N | O | S | 0 | 0 |
| | | | 1098 | 693 | 219 | 183 | 3 | | |

- Molecule 27 is a protein called eS24.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 27 | Z | 124 | Total | C | N | O | S | 0 | 0 |
| | | | 1011 | 640 | 198 | 168 | 5 | | |

- Molecule 28 is a protein called eS25.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 28 | a | 77 | Total | C | N | O | S | 0 | 0 |
| | | | 614 | 393 | 114 | 106 | 1 | | |

- Molecule 29 is a protein called eS26.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 29 | b | 101 | Total | C | N | O | S | 0 | 0 |
| | | | 814 | 507 | 170 | 132 | 5 | | |

There are 2 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------|------------|
| b | 28 | ARG | CYS | conflict | UNP G1TFE8 |

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| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------|------------|
| b | 56 | ALA | VAL | conflict | UNP G1TFE8 |

- Molecule 30 is a protein called eS27.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 30 | c | 83 | Total | C | N | O | S | 0 | 0 |
| | | | 651 | 408 | 121 | 115 | 7 | | |

- Molecule 31 is a protein called eS28.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---|---------|-------|
| 31 | d | 67 | Total | C | N | O | S | 0 | 0 |
| | | | 530 | 321 | 108 | 99 | 2 | | |

- Molecule 32 is a protein called uS14.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 32 | e | 55 | Total | C | N | O | S | 0 | 0 |
| | | | 459 | 286 | 94 | 74 | 5 | | |

- Molecule 33 is a protein called eS30.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---|---------|-------|
| 33 | f | 57 | Total | C | N | O | S | 0 | 0 |
| | | | 457 | 282 | 101 | 73 | 1 | | |

- Molecule 34 is a protein called eS31.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|----|---|---------|-------|
| 34 | g | 68 | Total | C | N | O | S | 0 | 0 |
| | | | 555 | 351 | 103 | 94 | 7 | | |

- Molecule 35 is a protein called Receptor for Activated C Kinase 1 (RACK1).

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|-----|----|---------|-------|
| 35 | h | 313 | Total | C | N | O | S | 0 | 0 |
| | | | 2436 | 1535 | 424 | 465 | 12 | | |

- Molecule 36 is a RNA chain called Met-tRNA-i-Met.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|----|---------|-------|
| 36 | i | 75 | Total | C | N | O | P | 0 | 0 |
| | | | 1604 | 717 | 298 | 515 | 74 | | |

- Molecule 37 is a protein called Eukaryotic translation initiation factor 2 subunit 1.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|-------|
| 37 | j | 182 | Total | C | N | O | S | 0 | 0 |
| | | | 1245 | 795 | 234 | 213 | 3 | | |

- Molecule 38 is a protein called eL41.

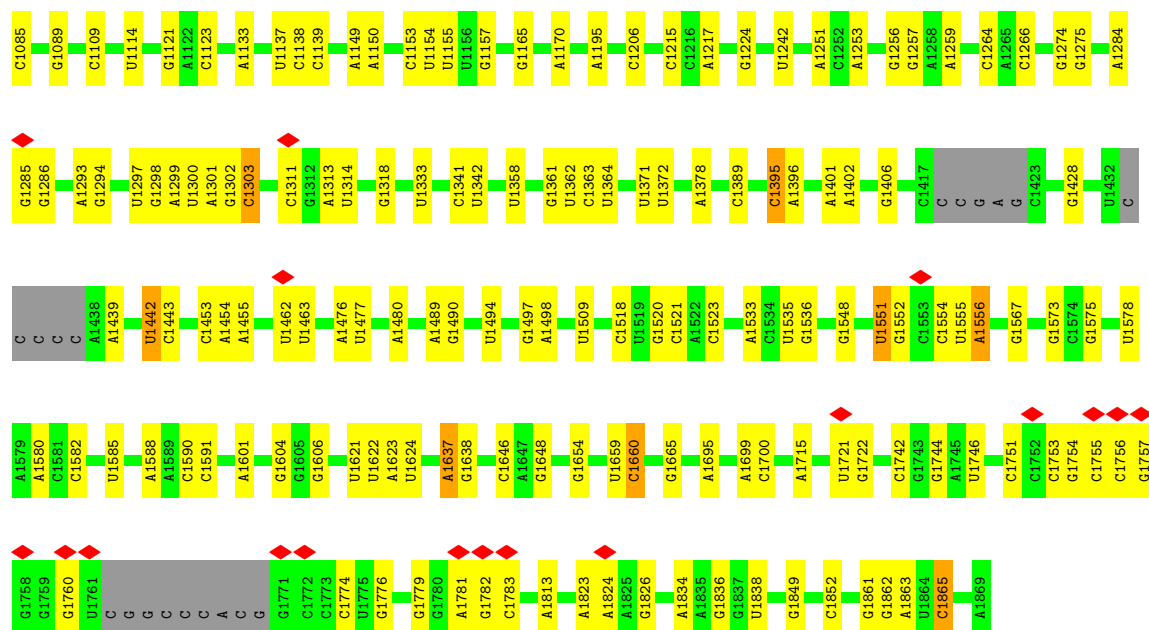
| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|-------|
| 38 | n | 25 | Total | C | N | O | S | 0 | 0 |
| | | | 239 | 145 | 64 | 27 | 3 | | |

- Molecule 39 is a RNA chain called HCV IRES.

| Mol | Chain | Residues | Atoms | | | | | AltConf | Trace |
|-----|-------|----------|-------|------|-----|------|-----|---------|-------|
| 39 | z | 188 | Total | C | N | O | P | 0 | 0 |
| | | | 4017 | 1789 | 721 | 1319 | 188 | | |

- Molecule 40 is ZINC ION (three-letter code: ZN) (formula: Zn).

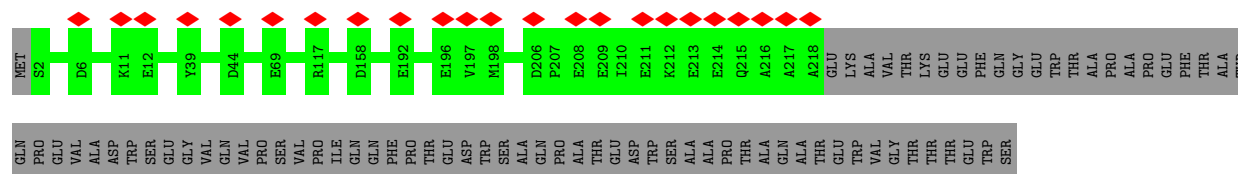
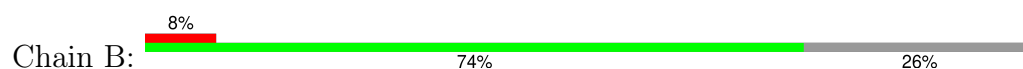
| Mol | Chain | Residues | Atoms | | AltConf |
|-----|-------|----------|-------|----|---------|
| 40 | b | 1 | Total | Zn | 0 |
| | | | 1 | 1 | |



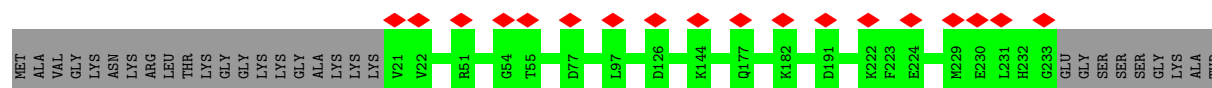
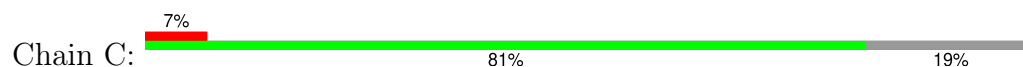
- Molecule 2: Eukaryotic translation initiation factor 1A, X-chromosomal



- Molecule 3: uS2 (SA)



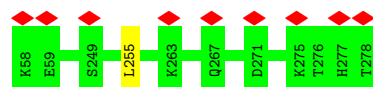
- Molecule 4: eS1




GLY ASP GLU THR GLY ALA LYS VAL GLU ARG ALA ASP GLY TYR GLU PRO PRO VAL GLN GLU SER VAL

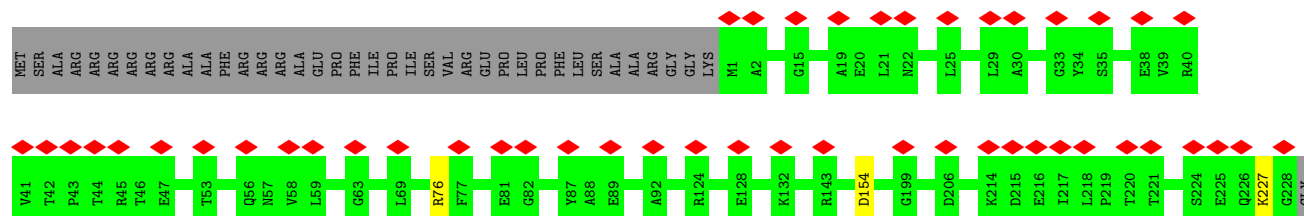
• Molecule 5: uS5

Chain D:  100%



• Molecule 6: uS3

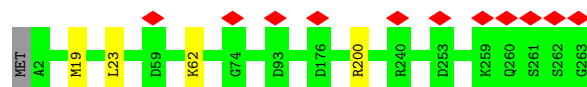
Chain E:  17% 80% 19%



LYS PRO GLU PRO PRO ALA MET PRO GLN PRO VAL PRO THR ALA

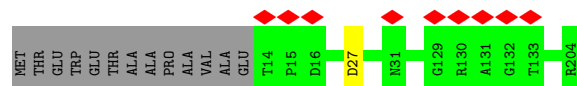
• Molecule 7: eS4 (S4 X isoform)

Chain F:  98%



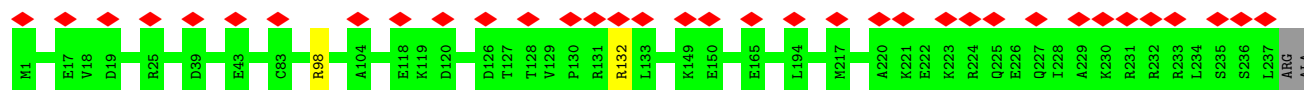
• Molecule 8: uS7

Chain G:  93% 6%



• Molecule 9: eS6

Chain H:  14% 94% 5%

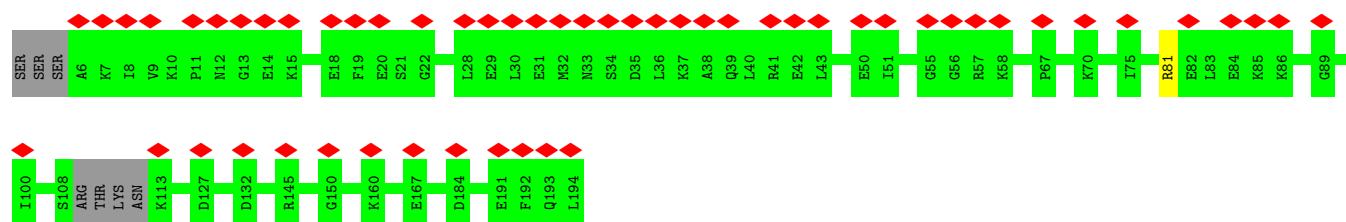


SER THR SER LYS SER SER SER GLN LYS

• Molecule 10: eS7



| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| GLU | PRO | ILE | THR | THR | ASN | ARG | VAL | PHE | GLN | ALA | LEU | THR | SER | SER | ASP | PHE | LYS | GLY | GLN | ASP | GLY | GLY | SER | PHE | SER | LEU | ARG | THR | ALA | HIS | LEU | SER | GLY | PHE | ARG | SER | ARG | PRO | ARG | ARG | THR | THR | THR | THR | ALA | ALA | THR | ARG | PRO | PRO | PRO | PRO | ILE | GLY | PRO | LEU | SER | ALA | ARG | GLY | GLY | THR | ALA | GLY | GLY | THR | TRP |
|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|



- Molecule 11: eS8



| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|
| MET | G2 | A34 | I76 | I79 | L96 | S115 | R123 | K124 | K125 | Q126 | A127 | K128 | L129 | T130 | P131 | E132 | E133 | E134 | E135 | I136 | L137 | K140 | K143 | K147 | K154 | N155 | G183 | R205 | K206 | G207 | LYS |
|-----|----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|

- Molecule 12: uS4



| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| MET | P2 | K22 | D26 | L29 | L60 | K66 | R79 | M92 | D95 | E107 | K139 | R150 | L151 | D152 | K180 | G181 | Q182 | G183 | G184 | A185 | G186 | ALA | GLY | ASP | ASP | ASP | GLU | GLU | GLU | ASP |
|-----|----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|-----|-----|-----|

- Molecule 13: eS10



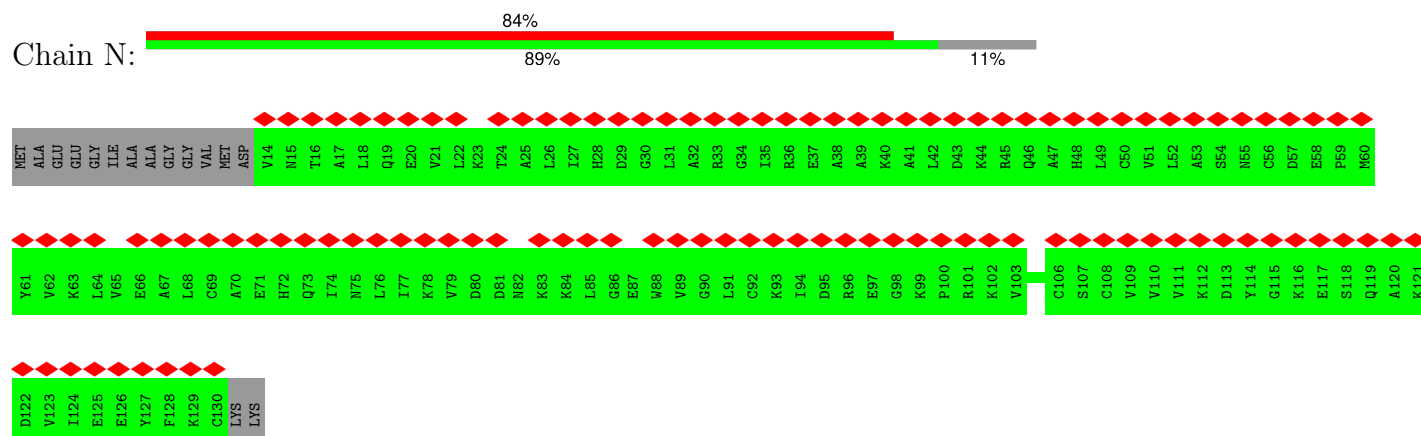
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| M1 | L2 | K17 | E18 | G19 | V20 | H23 | E34 | L35 | A36 | D37 | K38 | N39 | M46 | R55 | L71 | L79 | R80 | D81 | H82 | H84 | L85 | P86 | P87 | E88 | T89 | V90 | P91 | A92 | T93 | L94 | R95 | R96 | SER | ARG | ARG | PRO | PRO | GLU | THR | GLY | ARG | ARG | PRO | PRO | LYS | GLY | LEU | GLY | GLY | ARG | PRO | ALA |
|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|

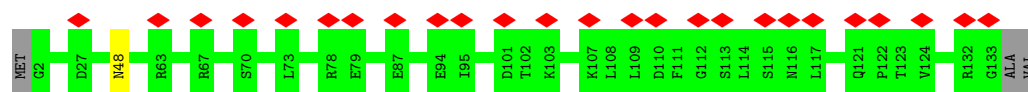
- Molecule 14: uS17



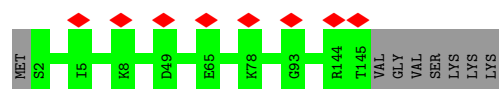
| | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|-----|-----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|
| MET | ALA | D3 | F17 | Q18 | N19 | R22 | V23 | L24 | L25 | G26 | E27 | T28 | G29 | K30 | E31 | K32 | R69 | E103 | R118 | A149 | G150 | T151 | K152 | K153 | GLN | PHE | GLN | LYS | PHE |
|-----|-----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|

- Molecule 15: eS12

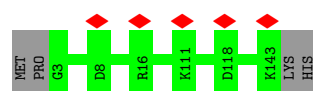




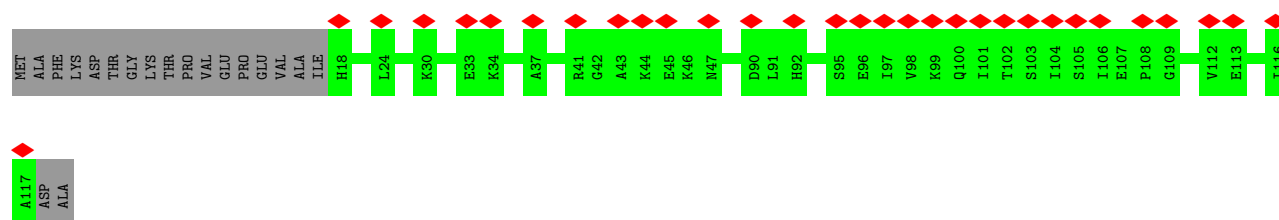
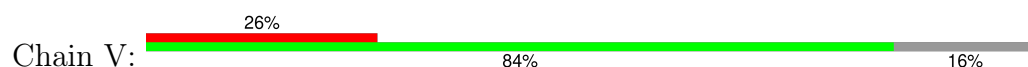
- Molecule 21: uS13



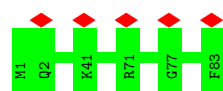
- Molecule 22: eS19



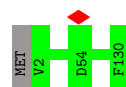
- Molecule 23: uS10



- Molecule 24: eS21

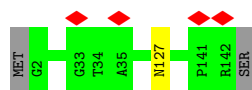


- Molecule 25: uS8

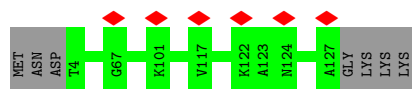


- Molecule 26: uS12

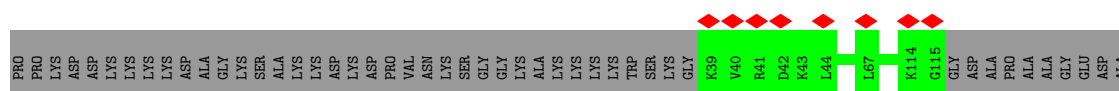




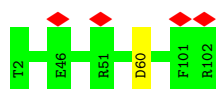
- Molecule 27: eS24



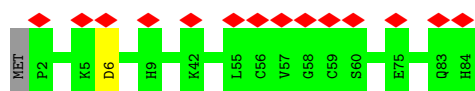
- Molecule 28: eS25



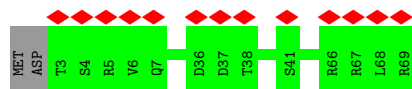
- Molecule 29: eS26



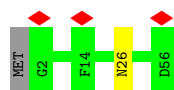
- Molecule 30: eS27



- Molecule 31: eS28



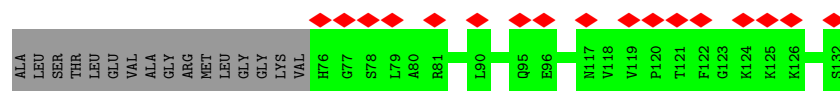
- Molecule 32: uS14



- Molecule 33: eS30



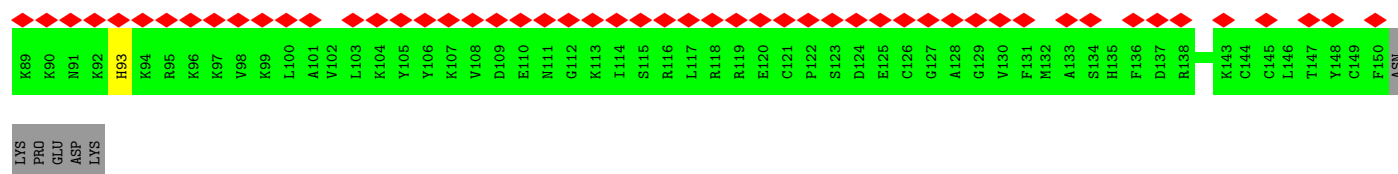
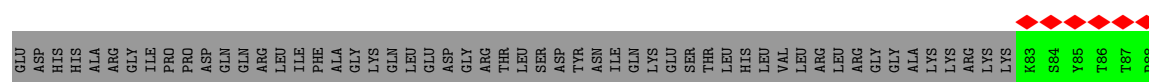
MET GLN LEU PHE THR VAL ARG ALA GLN LEU HIS THR LEU VAL THR ARG GLN THR VAL ALA GLN ILE LYS ALA HIS VAL SER LEU GLU ILE ALA PRO GLU ASP GLN VAL LEU LEU ALA GLY THR PRO LEU GLU ASP GLU THR LEU GLN CYS GLY VAL GLU



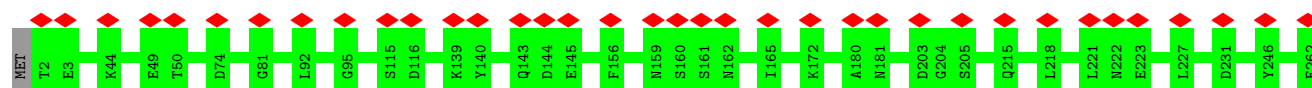
• Molecule 34: eS31



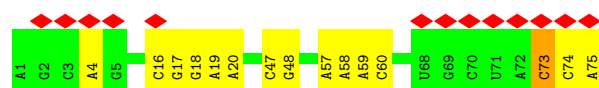
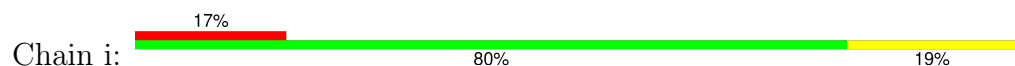
MET ALA ARG HIS LEU TVR GLN LEU PHE THR VAL ARG GLN LEU HIS THR LEU VAL THR ARG GLN CYS GLY VAL THR ARG GLN ILE LYS ALA HIS VAL SER LEU GLU ILE ALA PRO GLU ASP GLN VAL LEU LEU ALA GLY THR PRO LEU GLU ASP GLU THR LEU GLN CYS GLY VAL GLU



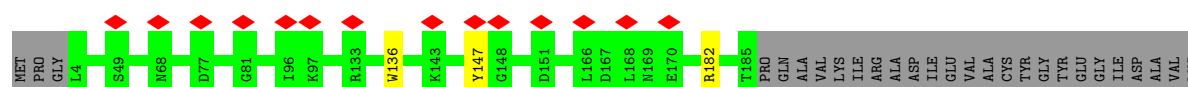
• Molecule 35: Receptor for Activated C Kinase 1 (RACK1)



• Molecule 36: Met-tRNA-i-Met



• Molecule 37: Eukaryotic translation initiation factor 2 subunit 1



4 Experimental information

| Property | Value | Source |
|--------------------------------------|---|-----------|
| EM reconstruction method | SINGLE PARTICLE | Depositor |
| Imposed symmetry | POINT, C1 | Depositor |
| Number of particles used | 46904 | Depositor |
| Resolution determination method | FSC 0.143 CUT-OFF | Depositor |
| CTF correction method | PHASE FLIPPING AND AMPLITUDE CORRECTION | Depositor |
| Microscope | FEI TECNAI F30 | Depositor |
| Voltage (kV) | 300 | Depositor |
| Electron dose ($e^-/\text{\AA}^2$) | 70.9 | Depositor |
| Minimum defocus (nm) | 1000 | Depositor |
| Maximum defocus (nm) | 2000 | Depositor |
| Magnification | 52000 | Depositor |
| Image detector | GATAN K3 (6k x 4k) | Depositor |
| Maximum map value | 0.117 | Depositor |
| Minimum map value | -0.050 | Depositor |
| Average map value | 0.001 | Depositor |
| Map value standard deviation | 0.006 | Depositor |
| Recommended contour level | 0.0175 | Depositor |
| Map size (Å) | 380.0, 380.0, 380.0 | wwPDB |
| Map dimensions | 400, 400, 400 | wwPDB |
| Map angles (°) | 90.0, 90.0, 90.0 | wwPDB |
| Pixel spacing (Å) | 0.95, 0.95, 0.95 | 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 | 2 | 0.27 | 0/40506 | 1.01 | 174/63123 (0.3%) |
| 2 | A | 0.29 | 0/809 | 0.65 | 0/1083 |
| 3 | B | 0.27 | 0/1747 | 0.61 | 0/2374 |
| 4 | C | 0.27 | 0/1756 | 0.59 | 0/2350 |
| 5 | D | 0.29 | 0/1753 | 0.61 | 1/2369 (0.0%) |
| 6 | E | 0.30 | 0/1796 | 0.67 | 1/2417 (0.0%) |
| 7 | F | 0.29 | 0/2118 | 0.66 | 2/2849 (0.1%) |
| 8 | G | 0.28 | 0/1531 | 0.69 | 1/2059 (0.0%) |
| 9 | H | 0.29 | 0/1946 | 0.67 | 0/2590 |
| 10 | I | 0.31 | 0/1510 | 0.66 | 0/2022 |
| 11 | J | 0.29 | 0/1723 | 0.69 | 0/2298 |
| 12 | K | 0.28 | 0/1550 | 0.64 | 1/2069 (0.0%) |
| 13 | L | 0.33 | 0/834 | 0.66 | 0/1125 |
| 14 | M | 0.28 | 0/1254 | 0.63 | 0/1677 |
| 15 | N | 0.27 | 0/918 | 0.61 | 0/1233 |
| 16 | O | 0.29 | 0/1226 | 0.64 | 1/1649 (0.1%) |
| 17 | P | 0.29 | 0/1029 | 0.64 | 1/1380 (0.1%) |
| 18 | Q | 0.29 | 0/1017 | 0.70 | 1/1358 (0.1%) |
| 19 | R | 0.30 | 0/1146 | 0.62 | 0/1534 |
| 20 | S | 0.28 | 0/1082 | 0.64 | 0/1452 |
| 21 | T | 0.27 | 0/1208 | 0.66 | 0/1618 |
| 22 | U | 0.30 | 0/1115 | 0.67 | 0/1493 |
| 23 | V | 0.25 | 0/805 | 0.61 | 0/1081 |
| 24 | W | 0.27 | 0/643 | 0.62 | 0/860 |
| 25 | X | 0.30 | 0/1051 | 0.64 | 0/1406 |
| 26 | Y | 0.30 | 0/1116 | 0.63 | 0/1490 |
| 27 | Z | 0.27 | 0/1028 | 0.61 | 0/1366 |
| 28 | a | 0.32 | 0/620 | 0.68 | 0/831 |
| 29 | b | 0.26 | 0/828 | 0.64 | 1/1109 (0.1%) |
| 30 | c | 0.31 | 0/665 | 0.66 | 1/891 (0.1%) |
| 31 | d | 0.29 | 0/532 | 0.80 | 0/712 |
| 32 | e | 0.34 | 0/470 | 0.76 | 0/623 |

| Mol | Chain | Bond lengths | | Bond angles | |
|-----|-------|--------------|----------------|-------------|-------------------|
| | | RMSZ | # Z >5 | RMSZ | # Z >5 |
| 33 | f | 0.27 | 0/462 | 0.62 | 0/607 |
| 34 | g | 0.26 | 0/567 | 0.63 | 0/753 |
| 35 | h | 0.28 | 0/2493 | 0.61 | 0/3394 |
| 36 | i | 0.24 | 0/1795 | 0.89 | 1/2798 (0.0%) |
| 37 | j | 0.34 | 1/1263 (0.1%) | 0.63 | 2/1708 (0.1%) |
| 38 | n | 0.27 | 0/240 | 0.80 | 0/305 |
| 39 | z | 0.26 | 0/4488 | 0.95 | 3/6995 (0.0%) |
| All | All | 0.28 | 1/88640 (0.0%) | 0.86 | 191/129051 (0.1%) |

All (1) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|---------|------|-------------|----------|
| 37 | j | 136 | TRP | CD2-CE2 | 6.11 | 1.48 | 1.41 |

All (191) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|------|------|-------------|--------|-------------|----------|
| 1 | 2 | 1852 | C | N3-C2-O2 | -11.46 | 113.88 | 121.90 |
| 1 | 2 | 1063 | C | N1-C2-O2 | 9.61 | 124.67 | 118.90 |
| 1 | 2 | 1591 | C | N1-C2-O2 | 9.59 | 124.65 | 118.90 |
| 1 | 2 | 853 | C | N1-C2-O2 | 9.33 | 124.50 | 118.90 |
| 1 | 2 | 1139 | C | N1-C2-O2 | 9.29 | 124.47 | 118.90 |
| 1 | 2 | 1139 | C | C2-N1-C1' | 9.06 | 128.76 | 118.80 |
| 1 | 2 | 179 | C | N1-C2-O2 | 8.60 | 124.06 | 118.90 |
| 1 | 2 | 1700 | C | C2-N1-C1' | 8.60 | 128.26 | 118.80 |
| 1 | 2 | 457 | C | N3-C2-O2 | -8.56 | 115.91 | 121.90 |
| 1 | 2 | 1453 | C | C2-N1-C1' | 8.53 | 128.18 | 118.80 |
| 1 | 2 | 1624 | U | C2-N1-C1' | 8.41 | 127.79 | 117.70 |
| 1 | 2 | 1453 | C | N1-C2-O2 | 8.35 | 123.91 | 118.90 |
| 1 | 2 | 1852 | C | C6-N1-C2 | -8.35 | 116.96 | 120.30 |
| 1 | 2 | 1206 | C | C2-N1-C1' | 8.32 | 127.95 | 118.80 |
| 1 | 2 | 853 | C | C2-N1-C1' | 8.19 | 127.81 | 118.80 |
| 37 | j | 136 | TRP | CD2-CE3-CZ3 | 8.12 | 129.36 | 118.80 |
| 1 | 2 | 1206 | C | N1-C2-O2 | 8.12 | 123.77 | 118.90 |
| 1 | 2 | 1591 | C | C2-N1-C1' | 8.09 | 127.70 | 118.80 |
| 1 | 2 | 1078 | C | N1-C2-O2 | 8.06 | 123.74 | 118.90 |
| 1 | 2 | 1700 | C | N1-C2-O2 | 7.95 | 123.67 | 118.90 |
| 1 | 2 | 1520 | G | C4-N9-C1' | 7.92 | 136.80 | 126.50 |
| 1 | 2 | 1852 | C | N1-C2-O2 | 7.88 | 123.63 | 118.90 |
| 1 | 2 | 853 | C | N3-C2-O2 | -7.78 | 116.46 | 121.90 |
| 1 | 2 | 1442 | U | N1-C2-O2 | 7.76 | 128.23 | 122.80 |
| 1 | 2 | 894 | G | C4-N9-C1' | 7.71 | 136.52 | 126.50 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|------|------|-----------|-------|-------------|----------|
| 1 | 2 | 1591 | C | N3-C2-O2 | -7.69 | 116.52 | 121.90 |
| 1 | 2 | 1063 | C | N3-C2-O2 | -7.68 | 116.53 | 121.90 |
| 1 | 2 | 1303 | C | N1-C2-O2 | 7.66 | 123.49 | 118.90 |
| 1 | 2 | 884 | C | N1-C2-O2 | 7.59 | 123.45 | 118.90 |
| 1 | 2 | 1139 | C | N3-C2-O2 | -7.55 | 116.61 | 121.90 |
| 1 | 2 | 1123 | C | N1-C2-O2 | 7.54 | 123.42 | 118.90 |
| 1 | 2 | 1442 | U | N3-C2-O2 | -7.53 | 116.93 | 122.20 |
| 1 | 2 | 585 | C | N1-C2-O2 | 7.45 | 123.37 | 118.90 |
| 1 | 2 | 1078 | C | N3-C2-O2 | -7.36 | 116.75 | 121.90 |
| 1 | 2 | 1298 | G | C4-N9-C1' | 7.36 | 136.06 | 126.50 |
| 1 | 2 | 1520 | G | N3-C4-N9 | 7.35 | 130.41 | 126.00 |
| 18 | Q | 82 | ASP | CB-CG-OD1 | 7.27 | 124.84 | 118.30 |
| 1 | 2 | 1303 | C | C2-N1-C1' | 7.23 | 126.76 | 118.80 |
| 1 | 2 | 1063 | C | C2-N1-C1' | 7.23 | 126.75 | 118.80 |
| 1 | 2 | 1298 | G | N3-C4-N9 | 7.21 | 130.33 | 126.00 |
| 1 | 2 | 1520 | G | N3-C4-C5 | -7.20 | 125.00 | 128.60 |
| 6 | E | 154 | ASP | CB-CG-OD1 | 7.11 | 124.69 | 118.30 |
| 1 | 2 | 578 | C | N1-C2-O2 | 7.09 | 123.16 | 118.90 |
| 1 | 2 | 1311 | C | N1-C2-O2 | 7.08 | 123.15 | 118.90 |
| 1 | 2 | 183 | G | N3-C4-C5 | -7.08 | 125.06 | 128.60 |
| 1 | 2 | 894 | G | N3-C4-N9 | 7.04 | 130.22 | 126.00 |
| 1 | 2 | 853 | C | C6-N1-C2 | -7.01 | 117.49 | 120.30 |
| 1 | 2 | 894 | G | N3-C4-C5 | -7.00 | 125.10 | 128.60 |
| 1 | 2 | 179 | C | N3-C2-O2 | -6.93 | 117.05 | 121.90 |
| 1 | 2 | 1298 | G | N3-C4-C5 | -6.91 | 125.14 | 128.60 |
| 1 | 2 | 179 | C | C2-N1-C1' | 6.90 | 126.39 | 118.80 |
| 1 | 2 | 1063 | C | C6-N1-C2 | -6.87 | 117.55 | 120.30 |
| 1 | 2 | 356 | C | N1-C2-O2 | 6.87 | 123.02 | 118.90 |
| 1 | 2 | 1591 | C | C6-N1-C2 | -6.84 | 117.56 | 120.30 |
| 1 | 2 | 1624 | U | N1-C2-O2 | 6.84 | 127.59 | 122.80 |
| 29 | b | 60 | ASP | CB-CG-OD2 | 6.79 | 124.41 | 118.30 |
| 1 | 2 | 1311 | C | C2-N1-C1' | 6.75 | 126.23 | 118.80 |
| 1 | 2 | 902 | G | C5-C6-O6 | 6.73 | 132.64 | 128.60 |
| 1 | 2 | 688 | U | P-O3'-C3' | 6.72 | 127.77 | 119.70 |
| 1 | 2 | 1442 | U | C2-N1-C1' | 6.69 | 125.73 | 117.70 |
| 1 | 2 | 1520 | G | C8-N9-C1' | -6.66 | 118.34 | 127.00 |
| 1 | 2 | 539 | C | N1-C2-O2 | 6.63 | 122.88 | 118.90 |
| 1 | 2 | 183 | G | C4-N9-C1' | 6.63 | 135.11 | 126.50 |
| 1 | 2 | 1443 | C | N1-C2-O2 | 6.57 | 122.84 | 118.90 |
| 1 | 2 | 1022 | U | C2-N1-C1' | 6.52 | 125.52 | 117.70 |
| 1 | 2 | 894 | G | C8-N9-C1' | -6.50 | 118.54 | 127.00 |
| 1 | 2 | 553 | U | P-O3'-C3' | 6.48 | 127.48 | 119.70 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|------|------|-----------|-------|-------------|----------|
| 1 | 2 | 1298 | G | C8-N9-C1' | -6.46 | 118.60 | 127.00 |
| 1 | 2 | 1624 | U | N3-C2-O2 | -6.43 | 117.70 | 122.20 |
| 1 | 2 | 183 | G | C2-N3-C4 | 6.42 | 115.11 | 111.90 |
| 1 | 2 | 118 | C | N1-C2-O2 | 6.42 | 122.75 | 118.90 |
| 1 | 2 | 4 | C | C2-N1-C1' | 6.41 | 125.85 | 118.80 |
| 1 | 2 | 356 | C | C2-N1-C1' | 6.40 | 125.84 | 118.80 |
| 1 | 2 | 752 | G | P-O3'-C3' | 6.37 | 127.34 | 119.70 |
| 1 | 2 | 750 | C | C2-N1-C1' | 6.34 | 125.77 | 118.80 |
| 36 | i | 73 | C | P-O3'-C3' | 6.31 | 127.27 | 119.70 |
| 1 | 2 | 1700 | C | N3-C2-O2 | -6.30 | 117.49 | 121.90 |
| 1 | 2 | 884 | C | N3-C2-O2 | -6.30 | 117.49 | 121.90 |
| 1 | 2 | 1395 | C | P-O3'-C3' | 6.29 | 127.25 | 119.70 |
| 1 | 2 | 1453 | C | N3-C2-O2 | -6.27 | 117.51 | 121.90 |
| 1 | 2 | 1019 | C | N1-C2-O2 | 6.24 | 122.64 | 118.90 |
| 1 | 2 | 1303 | C | N3-C2-O2 | -6.23 | 117.54 | 121.90 |
| 1 | 2 | 1057 | C | C2-N1-C1' | 6.23 | 125.65 | 118.80 |
| 1 | 2 | 1139 | C | C6-N1-C2 | -6.22 | 117.81 | 120.30 |
| 1 | 2 | 1206 | C | N3-C2-O2 | -6.18 | 117.57 | 121.90 |
| 1 | 2 | 1637 | A | P-O3'-C3' | 6.17 | 127.10 | 119.70 |
| 1 | 2 | 1139 | C | C6-N1-C1' | -6.16 | 113.40 | 120.80 |
| 1 | 2 | 902 | G | N1-C6-O6 | -6.16 | 116.20 | 119.90 |
| 1 | 2 | 188 | C | N1-C2-O2 | 6.16 | 122.59 | 118.90 |
| 1 | 2 | 1078 | C | C6-N1-C2 | -6.14 | 117.84 | 120.30 |
| 17 | P | 131 | ASP | CB-CG-OD2 | 6.14 | 123.82 | 118.30 |
| 1 | 2 | 1453 | C | C6-N1-C1' | -6.11 | 113.47 | 120.80 |
| 39 | z | 124 | C | N1-C2-O2 | 6.11 | 122.56 | 118.90 |
| 1 | 2 | 1123 | C | N3-C2-O2 | -6.10 | 117.63 | 121.90 |
| 1 | 2 | 1590 | C | N1-C2-O2 | 6.09 | 122.56 | 118.90 |
| 1 | 2 | 179 | C | C6-N1-C2 | -6.08 | 117.87 | 120.30 |
| 1 | 2 | 659 | G | C4-N9-C1' | 6.08 | 134.41 | 126.50 |
| 7 | F | 23 | LEU | CA-CB-CG | 6.07 | 129.26 | 115.30 |
| 1 | 2 | 1700 | C | C6-N1-C1' | -6.07 | 113.52 | 120.80 |
| 1 | 2 | 188 | C | C6-N1-C2 | -6.02 | 117.89 | 120.30 |
| 1 | 2 | 1311 | C | N3-C2-O2 | -6.02 | 117.69 | 121.90 |
| 1 | 2 | 578 | C | N3-C2-O2 | -6.01 | 117.69 | 121.90 |
| 1 | 2 | 541 | U | N1-C2-O2 | 6.01 | 127.00 | 122.80 |
| 39 | z | 160 | U | N3-C2-O2 | -5.98 | 118.01 | 122.20 |
| 1 | 2 | 1019 | C | N3-C2-O2 | -5.95 | 117.73 | 121.90 |
| 1 | 2 | 585 | C | N3-C2-O2 | -5.92 | 117.76 | 121.90 |
| 1 | 2 | 445 | A | P-O3'-C3' | 5.90 | 126.78 | 119.70 |
| 1 | 2 | 1556 | A | C2-N3-C4 | 5.90 | 113.55 | 110.60 |
| 1 | 2 | 183 | G | N3-C4-N9 | 5.89 | 129.53 | 126.00 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|------|------|------------|-------|-------------|----------|
| 1 | 2 | 383 | G | P-O3'-C3' | 5.89 | 126.77 | 119.70 |
| 1 | 2 | 973 | C | N1-C2-O2 | 5.86 | 122.41 | 118.90 |
| 1 | 2 | 1660 | C | C2-N1-C1' | 5.85 | 125.24 | 118.80 |
| 8 | G | 27 | ASP | CB-CG-OD1 | 5.85 | 123.56 | 118.30 |
| 16 | O | 110 | ASP | CB-CG-OD1 | 5.82 | 123.54 | 118.30 |
| 7 | F | 19 | MET | CA-CB-CG | 5.82 | 123.19 | 113.30 |
| 1 | 2 | 1551 | U | N3-C2-O2 | -5.81 | 118.13 | 122.20 |
| 1 | 2 | 457 | C | C6-N1-C2 | -5.80 | 117.98 | 120.30 |
| 1 | 2 | 823 | U | C2-N1-C1' | 5.80 | 124.66 | 117.70 |
| 1 | 2 | 1206 | C | C6-N1-C1' | -5.80 | 113.84 | 120.80 |
| 1 | 2 | 895 | G | P-O3'-C3' | 5.79 | 126.65 | 119.70 |
| 1 | 2 | 293 | C | N1-C2-O2 | 5.76 | 122.36 | 118.90 |
| 1 | 2 | 1518 | C | C2-N1-C1' | 5.75 | 125.12 | 118.80 |
| 1 | 2 | 1753 | C | N3-C2-O2 | -5.70 | 117.91 | 121.90 |
| 1 | 2 | 1311 | C | C6-N1-C2 | -5.69 | 118.03 | 120.30 |
| 30 | c | 6 | ASP | CB-CG-OD1 | 5.69 | 123.42 | 118.30 |
| 1 | 2 | 874 | G | P-O3'-C3' | 5.67 | 126.50 | 119.70 |
| 1 | 2 | 1551 | U | N1-C2-O2 | 5.66 | 126.76 | 122.80 |
| 1 | 2 | 630 | U | C2-N1-C1' | 5.63 | 124.46 | 117.70 |
| 1 | 2 | 4 | C | C6-N1-C2 | -5.63 | 118.05 | 120.30 |
| 1 | 2 | 541 | U | N3-C2-O2 | -5.62 | 118.26 | 122.20 |
| 1 | 2 | 570 | C | N1-C2-O2 | 5.61 | 122.27 | 118.90 |
| 1 | 2 | 356 | C | N3-C2-O2 | -5.61 | 117.97 | 121.90 |
| 1 | 2 | 884 | C | N3-C4-N4 | -5.59 | 114.09 | 118.00 |
| 1 | 2 | 457 | C | N1-C2-N3 | 5.57 | 123.10 | 119.20 |
| 1 | 2 | 1624 | U | C6-N1-C1' | -5.56 | 113.42 | 121.20 |
| 1 | 2 | 1078 | C | C2-N1-C1' | 5.55 | 124.91 | 118.80 |
| 1 | 2 | 1865 | C | C2-N1-C1' | 5.55 | 124.90 | 118.80 |
| 1 | 2 | 295 | C | N1-C2-O2 | 5.54 | 122.22 | 118.90 |
| 1 | 2 | 571 | U | N1-C2-O2 | 5.54 | 126.67 | 122.80 |
| 1 | 2 | 568 | C | N1-C2-O2 | 5.53 | 122.22 | 118.90 |
| 1 | 2 | 1742 | C | N1-C2-O2 | 5.53 | 122.22 | 118.90 |
| 1 | 2 | 1314 | U | C2-N1-C1' | 5.52 | 124.33 | 117.70 |
| 1 | 2 | 1591 | C | C5-C6-N1 | 5.49 | 123.75 | 121.00 |
| 1 | 2 | 188 | C | C2-N1-C1' | 5.47 | 124.82 | 118.80 |
| 1 | 2 | 1063 | C | C5-C6-N1 | 5.47 | 123.73 | 121.00 |
| 1 | 2 | 1395 | C | OP1-P-O3' | 5.45 | 117.19 | 105.20 |
| 1 | 2 | 1123 | C | C2-N1-C1' | 5.45 | 124.79 | 118.80 |
| 1 | 2 | 1624 | U | O4'-C1'-N1 | 5.42 | 112.54 | 108.20 |
| 5 | D | 255 | LEU | CA-CB-CG | 5.37 | 127.65 | 115.30 |
| 1 | 2 | 884 | C | C5-C4-N4 | 5.34 | 123.94 | 120.20 |
| 1 | 2 | 1057 | C | N1-C2-O2 | 5.34 | 122.10 | 118.90 |

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| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|------|------|-------------|-------|-------------|----------|
| 1 | 2 | 571 | U | N3-C2-O2 | -5.33 | 118.47 | 122.20 |
| 1 | 2 | 853 | C | C5-C6-N1 | 5.33 | 123.67 | 121.00 |
| 1 | 2 | 1442 | U | C5-C6-N1 | 5.33 | 125.37 | 122.70 |
| 1 | 2 | 142 | C | N1-C2-O2 | 5.32 | 122.09 | 118.90 |
| 1 | 2 | 541 | U | C2-N1-C1' | 5.31 | 124.07 | 117.70 |
| 1 | 2 | 1865 | C | N1-C2-O2 | 5.29 | 122.07 | 118.90 |
| 1 | 2 | 1206 | C | C6-N1-C2 | -5.25 | 118.20 | 120.30 |
| 1 | 2 | 585 | C | C2-N1-C1' | 5.25 | 124.57 | 118.80 |
| 1 | 2 | 579 | C | N1-C2-O2 | 5.23 | 122.03 | 118.90 |
| 1 | 2 | 1700 | C | C6-N1-C2 | -5.20 | 118.22 | 120.30 |
| 1 | 2 | 750 | C | N1-C2-O2 | 5.19 | 122.02 | 118.90 |
| 1 | 2 | 1389 | C | N1-C2-O2 | 5.19 | 122.02 | 118.90 |
| 12 | K | 29 | LEU | CA-CB-CG | 5.19 | 127.24 | 115.30 |
| 1 | 2 | 1443 | C | N3-C2-O2 | -5.17 | 118.28 | 121.90 |
| 39 | z | 330 | A | C2-N3-C4 | 5.17 | 113.19 | 110.60 |
| 1 | 2 | 457 | C | C2-N3-C4 | -5.17 | 117.32 | 119.90 |
| 1 | 2 | 293 | C | C2-N1-C1' | 5.16 | 124.48 | 118.80 |
| 1 | 2 | 183 | G | C8-N9-C1' | -5.16 | 120.29 | 127.00 |
| 37 | j | 136 | TRP | CE2-CD2-CE3 | 5.16 | 124.89 | 118.70 |
| 1 | 2 | 118 | C | N3-C2-O2 | -5.13 | 118.31 | 121.90 |
| 1 | 2 | 659 | G | C8-N9-C1' | -5.12 | 120.34 | 127.00 |
| 1 | 2 | 853 | C | C6-N1-C1' | -5.12 | 114.66 | 120.80 |
| 1 | 2 | 539 | C | N3-C2-O2 | -5.11 | 118.32 | 121.90 |
| 1 | 2 | 659 | G | N3-C4-N9 | 5.11 | 129.06 | 126.00 |
| 1 | 2 | 1389 | C | C2-N1-C1' | 5.09 | 124.41 | 118.80 |
| 1 | 2 | 1591 | C | C6-N1-C1' | -5.09 | 114.69 | 120.80 |
| 1 | 2 | 1389 | C | C6-N1-C2 | -5.09 | 118.26 | 120.30 |
| 1 | 2 | 119 | U | N3-C2-O2 | -5.08 | 118.65 | 122.20 |
| 1 | 2 | 531 | A | C4-N9-C1' | 5.06 | 135.41 | 126.30 |
| 1 | 2 | 1123 | C | C6-N1-C2 | -5.06 | 118.28 | 120.30 |
| 1 | 2 | 1303 | C | C6-N1-C1' | -5.06 | 114.73 | 120.80 |
| 1 | 2 | 341 | C | N1-C2-O2 | 5.05 | 121.93 | 118.90 |
| 1 | 2 | 630 | U | N3-C2-O2 | -5.04 | 118.67 | 122.20 |
| 1 | 2 | 688 | U | OP2-P-O3' | 5.03 | 116.26 | 105.20 |
| 1 | 2 | 1660 | C | N1-C2-O2 | 5.03 | 121.92 | 118.90 |
| 1 | 2 | 1590 | C | N3-C2-O2 | -5.02 | 118.39 | 121.90 |
| 1 | 2 | 1016 | U | C2-N1-C1' | 5.01 | 123.71 | 117.70 |
| 1 | 2 | 437 | G | N1-C2-N3 | 5.00 | 126.90 | 123.90 |
| 1 | 2 | 394 | G | C4-N9-C1' | 5.00 | 133.00 | 126.50 |
| 1 | 2 | 1590 | C | C6-N1-C2 | -5.00 | 118.30 | 120.30 |

There are no chirality outliers.

There are no planarity outliers.

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 | |
|-----|-------|---------------|-----------|----------|----------|-------------|-----|
| 2 | A | 97/144 (67%) | 84 (87%) | 13 (13%) | 0 | 100 | 100 |
| 3 | B | 215/295 (73%) | 201 (94%) | 14 (6%) | 0 | 100 | 100 |
| 4 | C | 211/264 (80%) | 198 (94%) | 13 (6%) | 0 | 100 | 100 |
| 5 | D | 219/221 (99%) | 214 (98%) | 5 (2%) | 0 | 100 | 100 |
| 6 | E | 226/281 (80%) | 219 (97%) | 7 (3%) | 0 | 100 | 100 |
| 7 | F | 260/263 (99%) | 250 (96%) | 10 (4%) | 0 | 100 | 100 |
| 8 | G | 189/204 (93%) | 179 (95%) | 10 (5%) | 0 | 100 | 100 |
| 9 | H | 235/249 (94%) | 229 (97%) | 6 (3%) | 0 | 100 | 100 |
| 10 | I | 181/432 (42%) | 175 (97%) | 6 (3%) | 0 | 100 | 100 |
| 11 | J | 205/208 (99%) | 192 (94%) | 13 (6%) | 0 | 100 | 100 |
| 12 | K | 183/194 (94%) | 175 (96%) | 8 (4%) | 0 | 100 | 100 |
| 13 | L | 94/149 (63%) | 88 (94%) | 6 (6%) | 0 | 100 | 100 |
| 14 | M | 149/158 (94%) | 137 (92%) | 12 (8%) | 0 | 100 | 100 |
| 15 | N | 115/132 (87%) | 110 (96%) | 5 (4%) | 0 | 100 | 100 |
| 16 | O | 147/151 (97%) | 143 (97%) | 4 (3%) | 0 | 100 | 100 |
| 17 | P | 134/168 (80%) | 123 (92%) | 11 (8%) | 0 | 100 | 100 |
| 18 | Q | 118/145 (81%) | 109 (92%) | 9 (8%) | 0 | 100 | 100 |
| 19 | R | 140/172 (81%) | 136 (97%) | 4 (3%) | 0 | 100 | 100 |
| 20 | S | 130/135 (96%) | 124 (95%) | 6 (5%) | 0 | 100 | 100 |

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| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Percentiles | |
|-----|-------|-----------------|------------|----------|----------|-------------|-----|
| 21 | T | 142/152 (93%) | 137 (96%) | 5 (4%) | 0 | 100 | 100 |
| 22 | U | 139/145 (96%) | 133 (96%) | 6 (4%) | 0 | 100 | 100 |
| 23 | V | 98/119 (82%) | 95 (97%) | 3 (3%) | 0 | 100 | 100 |
| 24 | W | 81/83 (98%) | 81 (100%) | 0 | 0 | 100 | 100 |
| 25 | X | 127/130 (98%) | 121 (95%) | 6 (5%) | 0 | 100 | 100 |
| 26 | Y | 139/143 (97%) | 135 (97%) | 4 (3%) | 0 | 100 | 100 |
| 27 | Z | 122/131 (93%) | 117 (96%) | 5 (4%) | 0 | 100 | 100 |
| 28 | a | 75/124 (60%) | 72 (96%) | 3 (4%) | 0 | 100 | 100 |
| 29 | b | 99/101 (98%) | 95 (96%) | 4 (4%) | 0 | 100 | 100 |
| 30 | c | 81/84 (96%) | 77 (95%) | 4 (5%) | 0 | 100 | 100 |
| 31 | d | 65/69 (94%) | 62 (95%) | 3 (5%) | 0 | 100 | 100 |
| 32 | e | 53/56 (95%) | 50 (94%) | 3 (6%) | 0 | 100 | 100 |
| 33 | f | 55/133 (41%) | 53 (96%) | 2 (4%) | 0 | 100 | 100 |
| 34 | g | 66/188 (35%) | 62 (94%) | 4 (6%) | 0 | 100 | 100 |
| 35 | h | 311/317 (98%) | 287 (92%) | 24 (8%) | 0 | 100 | 100 |
| 37 | j | 180/315 (57%) | 174 (97%) | 6 (3%) | 0 | 100 | 100 |
| 38 | n | 23/25 (92%) | 23 (100%) | 0 | 0 | 100 | 100 |
| All | All | 5104/6280 (81%) | 4860 (95%) | 244 (5%) | 0 | 100 | 100 |

There are no Ramachandran outliers to report.

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 | |
|-----|-------|----------------|------------|----------|-------------|-----|
| 2 | A | 84/123 (68%) | 84 (100%) | 0 | 100 | 100 |
| 3 | B | 180/245 (74%) | 180 (100%) | 0 | 100 | 100 |
| 4 | C | 194/231 (84%) | 194 (100%) | 0 | 100 | 100 |
| 5 | D | 187/187 (100%) | 187 (100%) | 0 | 100 | 100 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|----------------|------------|----------|-------------|-----|
| 6 | E | 190/232 (82%) | 188 (99%) | 2 (1%) | 70 | 83 |
| 7 | F | 224/225 (100%) | 222 (99%) | 2 (1%) | 75 | 87 |
| 8 | G | 161/170 (95%) | 161 (100%) | 0 | 100 | 100 |
| 9 | H | 207/218 (95%) | 205 (99%) | 2 (1%) | 73 | 85 |
| 10 | I | 165/360 (46%) | 164 (99%) | 1 (1%) | 84 | 92 |
| 11 | J | 179/180 (99%) | 179 (100%) | 0 | 100 | 100 |
| 12 | K | 161/168 (96%) | 160 (99%) | 1 (1%) | 84 | 92 |
| 13 | L | 87/125 (70%) | 87 (100%) | 0 | 100 | 100 |
| 14 | M | 136/142 (96%) | 135 (99%) | 1 (1%) | 81 | 90 |
| 15 | N | 99/108 (92%) | 99 (100%) | 0 | 100 | 100 |
| 16 | O | 130/131 (99%) | 129 (99%) | 1 (1%) | 79 | 88 |
| 17 | P | 106/130 (82%) | 105 (99%) | 1 (1%) | 75 | 87 |
| 18 | Q | 109/130 (84%) | 109 (100%) | 0 | 100 | 100 |
| 19 | R | 117/140 (84%) | 116 (99%) | 1 (1%) | 75 | 87 |
| 20 | S | 119/121 (98%) | 118 (99%) | 1 (1%) | 79 | 88 |
| 21 | T | 125/132 (95%) | 125 (100%) | 0 | 100 | 100 |
| 22 | U | 111/116 (96%) | 111 (100%) | 0 | 100 | 100 |
| 23 | V | 92/107 (86%) | 92 (100%) | 0 | 100 | 100 |
| 24 | W | 67/67 (100%) | 67 (100%) | 0 | 100 | 100 |
| 25 | X | 112/113 (99%) | 112 (100%) | 0 | 100 | 100 |
| 26 | Y | 113/115 (98%) | 112 (99%) | 1 (1%) | 75 | 87 |
| 27 | Z | 107/113 (95%) | 107 (100%) | 0 | 100 | 100 |
| 28 | a | 68/102 (67%) | 68 (100%) | 0 | 100 | 100 |
| 29 | b | 88/88 (100%) | 88 (100%) | 0 | 100 | 100 |
| 30 | c | 75/76 (99%) | 75 (100%) | 0 | 100 | 100 |
| 31 | d | 60/62 (97%) | 60 (100%) | 0 | 100 | 100 |
| 32 | e | 48/49 (98%) | 47 (98%) | 1 (2%) | 48 | 71 |
| 33 | f | 47/106 (44%) | 47 (100%) | 0 | 100 | 100 |
| 34 | g | 61/154 (40%) | 60 (98%) | 1 (2%) | 58 | 76 |
| 35 | h | 272/275 (99%) | 272 (100%) | 0 | 100 | 100 |
| 37 | j | 90/280 (32%) | 88 (98%) | 2 (2%) | 47 | 69 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|-----------------|-------------|----------|-------------|-----|
| 38 | n | 24/24 (100%) | 24 (100%) | 0 | 100 | 100 |
| All | All | 4395/5345 (82%) | 4377 (100%) | 18 (0%) | 88 | 95 |

All (18) residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 6 | E | 76 | ARG |
| 6 | E | 227 | LYS |
| 7 | F | 62 | LYS |
| 7 | F | 200 | ARG |
| 9 | H | 98 | ARG |
| 9 | H | 132 | ARG |
| 10 | I | 81 | ARG |
| 12 | K | 150 | ARG |
| 14 | M | 69 | ARG |
| 16 | O | 36 | GLN |
| 17 | P | 26 | ASN |
| 19 | R | 71 | ARG |
| 20 | S | 48 | ASN |
| 26 | Y | 127 | ASN |
| 32 | e | 26 | ASN |
| 34 | g | 93 | HIS |
| 37 | j | 147 | TYR |
| 37 | j | 182 | ARG |

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 5 | D | 113 | GLN |
| 8 | G | 148 | ASN |
| 8 | G | 203 | ASN |
| 17 | P | 26 | ASN |
| 20 | S | 26 | ASN |
| 34 | g | 93 | HIS |

5.3.3 RNA ⓘ

| Mol | Chain | Analysed | Backbone Outliers | Pucker Outliers |
|-----|-------|-----------------|-------------------|-----------------|
| 1 | 2 | 1685/1870 (90%) | 317 (18%) | 12 (0%) |

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| Mol | Chain | Analysed | Backbone Outliers | Pucker Outliers |
|-----|-------|-----------------|-------------------|-----------------|
| 36 | i | 74/75 (98%) | 15 (20%) | 0 |
| 39 | z | 186/400 (46%) | 54 (29%) | 0 |
| All | All | 1945/2345 (82%) | 386 (19%) | 12 (0%) |

All (386) RNA backbone outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | 2 | 2 | A |
| 1 | 2 | 4 | C |
| 1 | 2 | 11 | A |
| 1 | 2 | 26 | U |
| 1 | 2 | 33 | G |
| 1 | 2 | 41 | G |
| 1 | 2 | 42 | A |
| 1 | 2 | 44 | U |
| 1 | 2 | 46 | A |
| 1 | 2 | 56 | G |
| 1 | 2 | 59 | U |
| 1 | 2 | 62 | G |
| 1 | 2 | 67 | C |
| 1 | 2 | 68 | A |
| 1 | 2 | 69 | C |
| 1 | 2 | 73 | C |
| 1 | 2 | 74 | G |
| 1 | 2 | 76 | U |
| 1 | 2 | 102 | A |
| 1 | 2 | 103 | A |
| 1 | 2 | 113 | G |
| 1 | 2 | 115 | U |
| 1 | 2 | 126 | G |
| 1 | 2 | 130 | G |
| 1 | 2 | 142 | C |
| 1 | 2 | 143 | U |
| 1 | 2 | 149 | A |
| 1 | 2 | 155 | G |
| 1 | 2 | 162 | C |
| 1 | 2 | 170 | A |
| 1 | 2 | 171 | A |
| 1 | 2 | 177 | G |
| 1 | 2 | 178 | C |
| 1 | 2 | 180 | G |
| 1 | 2 | 182 | C |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | 2 | 183 | G |
| 1 | 2 | 184 | G |
| 1 | 2 | 185 | G |
| 1 | 2 | 188 | C |
| 1 | 2 | 192 | C |
| 1 | 2 | 199 | C |
| 1 | 2 | 292 | A |
| 1 | 2 | 293 | C |
| 1 | 2 | 305 | U |
| 1 | 2 | 306 | C |
| 1 | 2 | 307 | G |
| 1 | 2 | 309 | G |
| 1 | 2 | 312 | G |
| 1 | 2 | 313 | A |
| 1 | 2 | 314 | U |
| 1 | 2 | 319 | C |
| 1 | 2 | 320 | G |
| 1 | 2 | 323 | C |
| 1 | 2 | 347 | G |
| 1 | 2 | 364 | A |
| 1 | 2 | 368 | U |
| 1 | 2 | 370 | G |
| 1 | 2 | 383 | G |
| 1 | 2 | 384 | U |
| 1 | 2 | 385 | G |
| 1 | 2 | 386 | C |
| 1 | 2 | 398 | A |
| 1 | 2 | 399 | C |
| 1 | 2 | 400 | C |
| 1 | 2 | 408 | A |
| 1 | 2 | 409 | C |
| 1 | 2 | 418 | A |
| 1 | 2 | 429 | C |
| 1 | 2 | 435 | A |
| 1 | 2 | 438 | G |
| 1 | 2 | 446 | G |
| 1 | 2 | 448 | A |
| 1 | 2 | 449 | A |
| 1 | 2 | 450 | C |
| 1 | 2 | 465 | A |
| 1 | 2 | 471 | G |
| 1 | 2 | 472 | C |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | 2 | 474 | G |
| 1 | 2 | 487 | U |
| 1 | 2 | 488 | U |
| 1 | 2 | 492 | C |
| 1 | 2 | 493 | A |
| 1 | 2 | 500 | A |
| 1 | 2 | 502 | C |
| 1 | 2 | 508 | A |
| 1 | 2 | 525 | A |
| 1 | 2 | 531 | A |
| 1 | 2 | 532 | C |
| 1 | 2 | 536 | A |
| 1 | 2 | 539 | C |
| 1 | 2 | 541 | U |
| 1 | 2 | 544 | G |
| 1 | 2 | 546 | G |
| 1 | 2 | 549 | C |
| 1 | 2 | 550 | C |
| 1 | 2 | 551 | U |
| 1 | 2 | 554 | A |
| 1 | 2 | 555 | A |
| 1 | 2 | 556 | U |
| 1 | 2 | 562 | U |
| 1 | 2 | 576 | A |
| 1 | 2 | 583 | A |
| 1 | 2 | 587 | A |
| 1 | 2 | 588 | G |
| 1 | 2 | 589 | G |
| 1 | 2 | 590 | A |
| 1 | 2 | 591 | U |
| 1 | 2 | 593 | C |
| 1 | 2 | 603 | C |
| 1 | 2 | 604 | A |
| 1 | 2 | 606 | G |
| 1 | 2 | 607 | U |
| 1 | 2 | 608 | C |
| 1 | 2 | 614 | C |
| 1 | 2 | 617 | G |
| 1 | 2 | 628 | A |
| 1 | 2 | 643 | A |
| 1 | 2 | 644 | G |
| 1 | 2 | 660 | C |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | 2 | 662 | G |
| 1 | 2 | 668 | A |
| 1 | 2 | 669 | A |
| 1 | 2 | 671 | A |
| 1 | 2 | 672 | A |
| 1 | 2 | 673 | G |
| 1 | 2 | 688 | U |
| 1 | 2 | 689 | U |
| 1 | 2 | 696 | G |
| 1 | 2 | 732 | U |
| 1 | 2 | 752 | G |
| 1 | 2 | 753 | C |
| 1 | 2 | 754 | G |
| 1 | 2 | 799 | U |
| 1 | 2 | 811 | A |
| 1 | 2 | 821 | G |
| 1 | 2 | 822 | U |
| 1 | 2 | 830 | A |
| 1 | 2 | 834 | C |
| 1 | 2 | 844 | U |
| 1 | 2 | 847 | A |
| 1 | 2 | 870 | A |
| 1 | 2 | 872 | A |
| 1 | 2 | 873 | G |
| 1 | 2 | 874 | G |
| 1 | 2 | 875 | A |
| 1 | 2 | 878 | G |
| 1 | 2 | 883 | U |
| 1 | 2 | 885 | U |
| 1 | 2 | 886 | A |
| 1 | 2 | 887 | U |
| 1 | 2 | 888 | U |
| 1 | 2 | 890 | U |
| 1 | 2 | 891 | G |
| 1 | 2 | 892 | U |
| 1 | 2 | 895 | G |
| 1 | 2 | 896 | U |
| 1 | 2 | 901 | G |
| 1 | 2 | 913 | A |
| 1 | 2 | 914 | U |
| 1 | 2 | 919 | A |
| 1 | 2 | 920 | A |

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| Mol | Chain | Res | Type |
|-----|-------|------|------|
| 1 | 2 | 921 | G |
| 1 | 2 | 933 | G |
| 1 | 2 | 943 | U |
| 1 | 2 | 970 | G |
| 1 | 2 | 971 | G |
| 1 | 2 | 985 | G |
| 1 | 2 | 986 | G |
| 1 | 2 | 990 | A |
| 1 | 2 | 992 | A |
| 1 | 2 | 999 | G |
| 1 | 2 | 1002 | U |
| 1 | 2 | 1017 | U |
| 1 | 2 | 1022 | U |
| 1 | 2 | 1023 | A |
| 1 | 2 | 1045 | U |
| 1 | 2 | 1061 | U |
| 1 | 2 | 1070 | A |
| 1 | 2 | 1078 | C |
| 1 | 2 | 1083 | A |
| 1 | 2 | 1085 | C |
| 1 | 2 | 1089 | G |
| 1 | 2 | 1109 | C |
| 1 | 2 | 1114 | U |
| 1 | 2 | 1121 | G |
| 1 | 2 | 1133 | A |
| 1 | 2 | 1138 | C |
| 1 | 2 | 1149 | A |
| 1 | 2 | 1150 | A |
| 1 | 2 | 1153 | C |
| 1 | 2 | 1154 | U |
| 1 | 2 | 1155 | U |
| 1 | 2 | 1157 | G |
| 1 | 2 | 1165 | G |
| 1 | 2 | 1170 | A |
| 1 | 2 | 1195 | A |
| 1 | 2 | 1215 | C |
| 1 | 2 | 1217 | A |
| 1 | 2 | 1224 | G |
| 1 | 2 | 1242 | U |
| 1 | 2 | 1251 | A |
| 1 | 2 | 1253 | A |
| 1 | 2 | 1256 | G |

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| Mol | Chain | Res | Type |
|-----|-------|------|------|
| 1 | 2 | 1257 | G |
| 1 | 2 | 1259 | A |
| 1 | 2 | 1264 | C |
| 1 | 2 | 1266 | C |
| 1 | 2 | 1274 | G |
| 1 | 2 | 1275 | G |
| 1 | 2 | 1284 | A |
| 1 | 2 | 1285 | G |
| 1 | 2 | 1286 | G |
| 1 | 2 | 1293 | A |
| 1 | 2 | 1294 | G |
| 1 | 2 | 1297 | U |
| 1 | 2 | 1299 | A |
| 1 | 2 | 1300 | U |
| 1 | 2 | 1301 | A |
| 1 | 2 | 1302 | G |
| 1 | 2 | 1303 | C |
| 1 | 2 | 1313 | A |
| 1 | 2 | 1318 | G |
| 1 | 2 | 1333 | U |
| 1 | 2 | 1341 | C |
| 1 | 2 | 1342 | U |
| 1 | 2 | 1358 | U |
| 1 | 2 | 1361 | G |
| 1 | 2 | 1362 | U |
| 1 | 2 | 1363 | C |
| 1 | 2 | 1364 | U |
| 1 | 2 | 1371 | U |
| 1 | 2 | 1372 | U |
| 1 | 2 | 1378 | A |
| 1 | 2 | 1396 | A |
| 1 | 2 | 1401 | A |
| 1 | 2 | 1402 | A |
| 1 | 2 | 1406 | G |
| 1 | 2 | 1428 | G |
| 1 | 2 | 1439 | A |
| 1 | 2 | 1442 | U |
| 1 | 2 | 1454 | A |
| 1 | 2 | 1455 | A |
| 1 | 2 | 1462 | U |
| 1 | 2 | 1463 | U |
| 1 | 2 | 1476 | A |

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| Mol | Chain | Res | Type |
|-----|-------|------|------|
| 1 | 2 | 1477 | U |
| 1 | 2 | 1480 | A |
| 1 | 2 | 1489 | A |
| 1 | 2 | 1490 | G |
| 1 | 2 | 1494 | U |
| 1 | 2 | 1497 | G |
| 1 | 2 | 1498 | A |
| 1 | 2 | 1509 | U |
| 1 | 2 | 1521 | C |
| 1 | 2 | 1523 | C |
| 1 | 2 | 1533 | A |
| 1 | 2 | 1535 | U |
| 1 | 2 | 1536 | G |
| 1 | 2 | 1548 | G |
| 1 | 2 | 1551 | U |
| 1 | 2 | 1552 | G |
| 1 | 2 | 1554 | C |
| 1 | 2 | 1555 | U |
| 1 | 2 | 1556 | A |
| 1 | 2 | 1567 | G |
| 1 | 2 | 1573 | G |
| 1 | 2 | 1575 | G |
| 1 | 2 | 1578 | U |
| 1 | 2 | 1580 | A |
| 1 | 2 | 1582 | C |
| 1 | 2 | 1585 | U |
| 1 | 2 | 1588 | A |
| 1 | 2 | 1601 | A |
| 1 | 2 | 1604 | G |
| 1 | 2 | 1606 | G |
| 1 | 2 | 1621 | U |
| 1 | 2 | 1622 | U |
| 1 | 2 | 1623 | A |
| 1 | 2 | 1637 | A |
| 1 | 2 | 1638 | G |
| 1 | 2 | 1646 | C |
| 1 | 2 | 1648 | G |
| 1 | 2 | 1654 | G |
| 1 | 2 | 1659 | U |
| 1 | 2 | 1660 | C |
| 1 | 2 | 1665 | G |
| 1 | 2 | 1695 | A |

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| Mol | Chain | Res | Type |
|-----|-------|------|------|
| 1 | 2 | 1699 | A |
| 1 | 2 | 1715 | A |
| 1 | 2 | 1721 | U |
| 1 | 2 | 1722 | G |
| 1 | 2 | 1744 | G |
| 1 | 2 | 1746 | U |
| 1 | 2 | 1751 | C |
| 1 | 2 | 1754 | G |
| 1 | 2 | 1755 | C |
| 1 | 2 | 1756 | C |
| 1 | 2 | 1757 | G |
| 1 | 2 | 1760 | G |
| 1 | 2 | 1774 | C |
| 1 | 2 | 1776 | G |
| 1 | 2 | 1779 | G |
| 1 | 2 | 1781 | A |
| 1 | 2 | 1782 | G |
| 1 | 2 | 1783 | C |
| 1 | 2 | 1813 | A |
| 1 | 2 | 1823 | A |
| 1 | 2 | 1824 | A |
| 1 | 2 | 1826 | G |
| 1 | 2 | 1834 | A |
| 1 | 2 | 1836 | G |
| 1 | 2 | 1838 | U |
| 1 | 2 | 1849 | G |
| 1 | 2 | 1861 | G |
| 1 | 2 | 1862 | G |
| 1 | 2 | 1863 | A |
| 1 | 2 | 1865 | C |
| 36 | i | 4 | A |
| 36 | i | 16 | C |
| 36 | i | 17 | G |
| 36 | i | 18 | G |
| 36 | i | 19 | A |
| 36 | i | 20 | A |
| 36 | i | 47 | C |
| 36 | i | 48 | G |
| 36 | i | 57 | A |
| 36 | i | 58 | A |
| 36 | i | 59 | A |
| 36 | i | 60 | C |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 36 | i | 73 | C |
| 36 | i | 74 | C |
| 36 | i | 75 | A |
| 39 | z | 121 | C |
| 39 | z | 124 | C |
| 39 | z | 136 | A |
| 39 | z | 137 | G |
| 39 | z | 144 | U |
| 39 | z | 154 | A |
| 39 | z | 157 | C |
| 39 | z | 160 | U |
| 39 | z | 161 | G |
| 39 | z | 162 | A |
| 39 | z | 163 | G |
| 39 | z | 164 | U |
| 39 | z | 165 | A |
| 39 | z | 166 | C |
| 39 | z | 168 | C |
| 39 | z | 169 | C |
| 39 | z | 229 | G |
| 39 | z | 232 | C |
| 39 | z | 234 | U |
| 39 | z | 237 | C |
| 39 | z | 244 | A |
| 39 | z | 253 | G |
| 39 | z | 258 | G |
| 39 | z | 259 | U |
| 39 | z | 265 | U |
| 39 | z | 266 | G |
| 39 | z | 279 | C |
| 39 | z | 280 | C |
| 39 | z | 281 | U |
| 39 | z | 282 | U |
| 39 | z | 288 | A |
| 39 | z | 296 | A |
| 39 | z | 297 | U |
| 39 | z | 303 | G |
| 39 | z | 306 | U |
| 39 | z | 324 | U |
| 39 | z | 328 | G |
| 39 | z | 330 | A |
| 39 | z | 331 | G |

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| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 39 | z | 332 | A |
| 39 | z | 333 | C |
| 39 | z | 334 | C |
| 39 | z | 335 | G |
| 39 | z | 337 | G |
| 39 | z | 339 | A |
| 39 | z | 345 | A |
| 39 | z | 346 | G |
| 39 | z | 348 | A |
| 39 | z | 349 | C |
| 39 | z | 351 | A |
| 39 | z | 352 | A |
| 39 | z | 353 | U |
| 39 | z | 355 | C |
| 39 | z | 356 | U |

All (12) RNA pucker outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|------|------|
| 1 | 2 | 383 | G |
| 1 | 2 | 445 | A |
| 1 | 2 | 553 | U |
| 1 | 2 | 561 | A |
| 1 | 2 | 688 | U |
| 1 | 2 | 752 | G |
| 1 | 2 | 874 | G |
| 1 | 2 | 886 | A |
| 1 | 2 | 895 | G |
| 1 | 2 | 1137 | U |
| 1 | 2 | 1395 | C |
| 1 | 2 | 1637 | A |

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 oligosaccharides in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is 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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

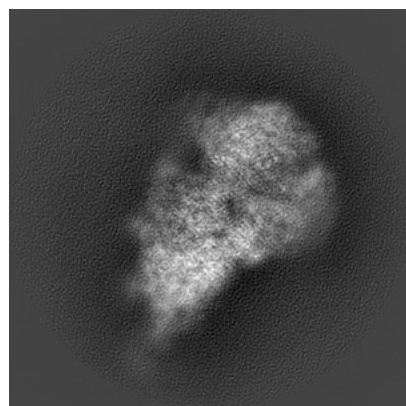
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-25538. These allow visual inspection of the internal detail of the map and identification of artifacts.

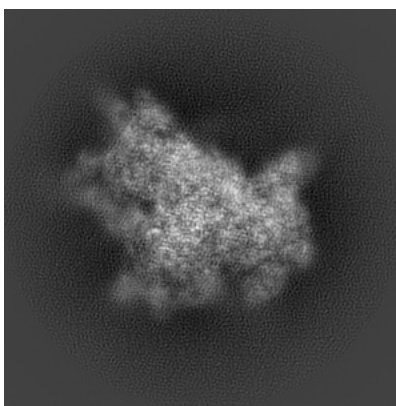
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

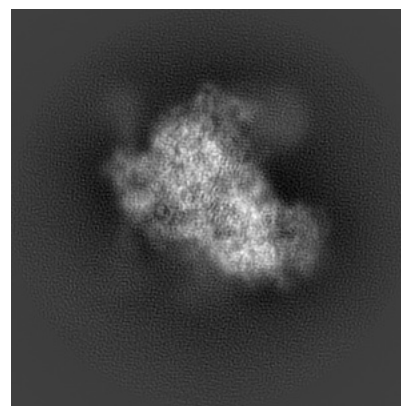
6.1.1 Primary map



X

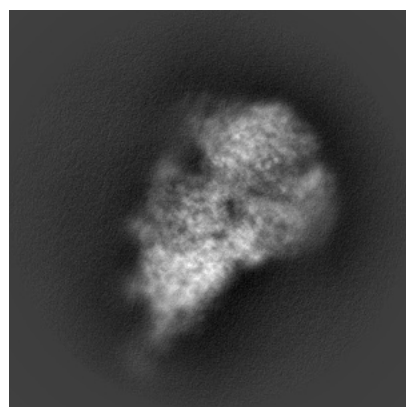


Y

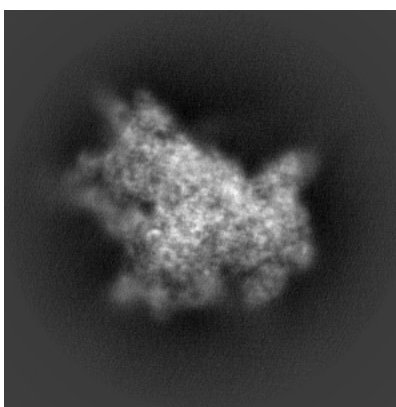


Z

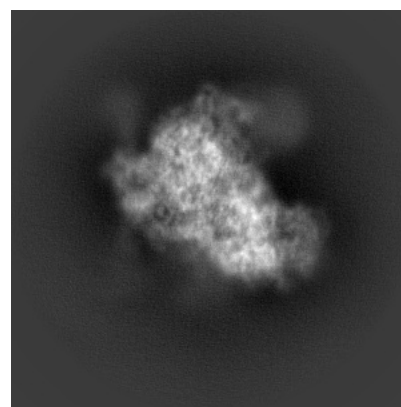
6.1.2 Raw map



X



Y

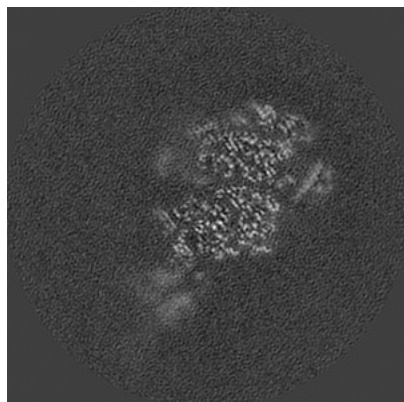


Z

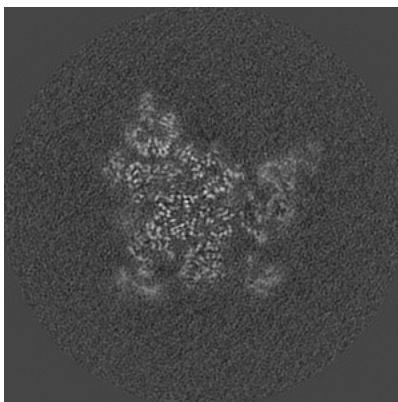
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

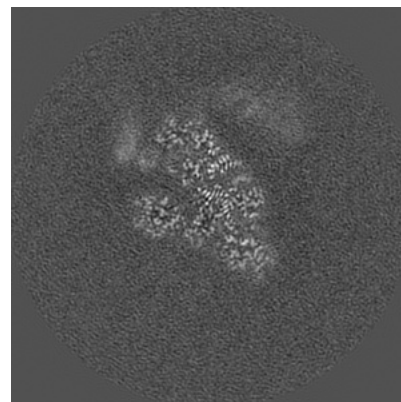
6.2.1 Primary map



X Index: 200

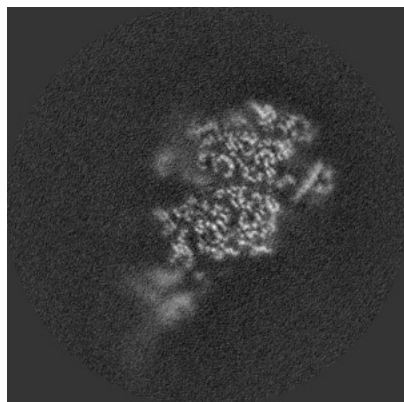


Y Index: 200

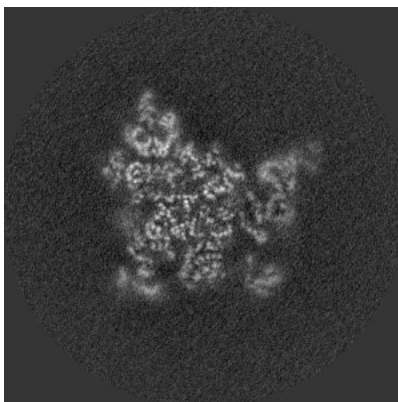


Z Index: 200

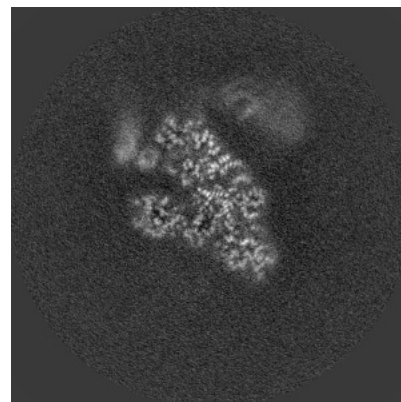
6.2.2 Raw map



X Index: 200



Y Index: 200

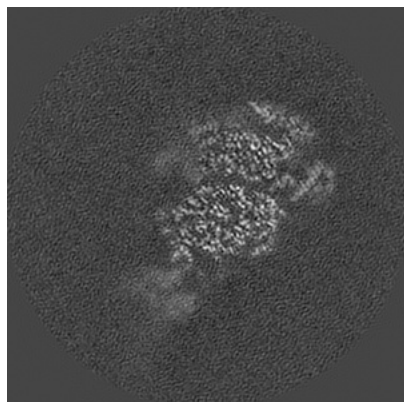


Z Index: 200

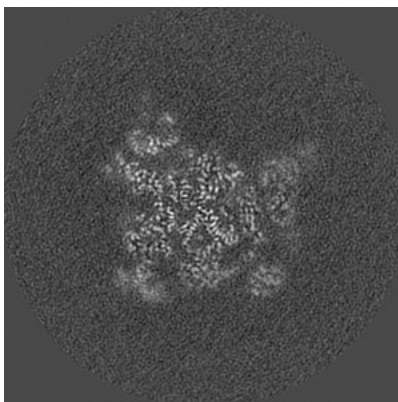
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

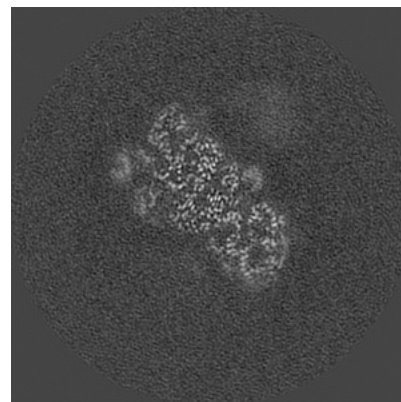
6.3.1 Primary map



X Index: 198

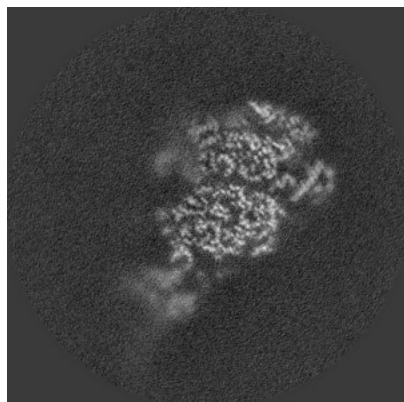


Y Index: 204

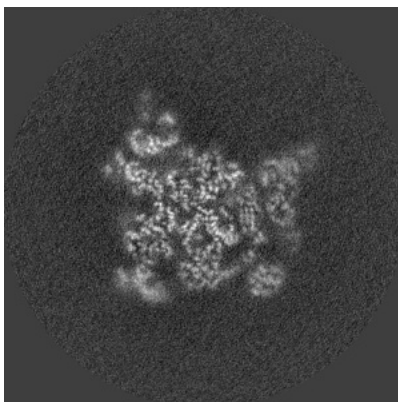


Z Index: 179

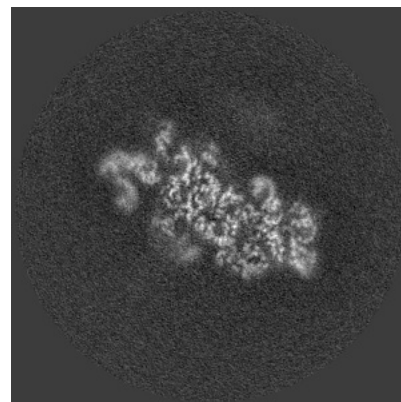
6.3.2 Raw map



X Index: 198



Y Index: 204

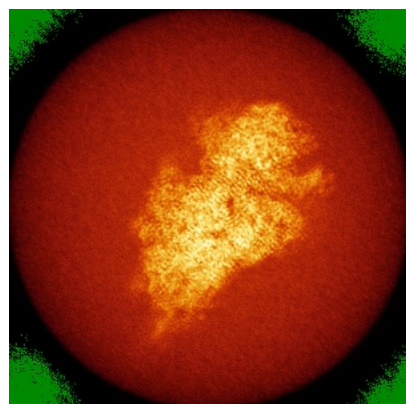


Z Index: 154

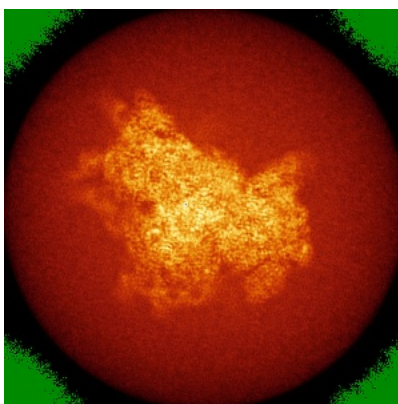
The images above show the largest variance slices of the map in three orthogonal directions.

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

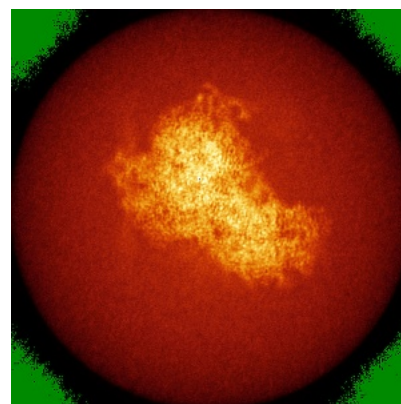
6.4.1 Primary map



X

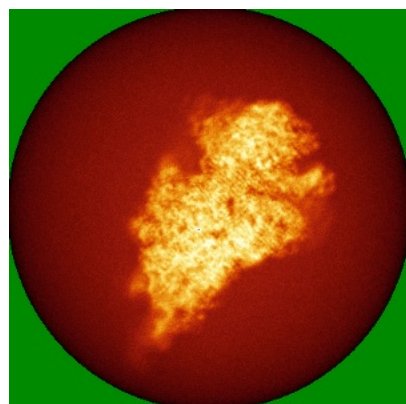


Y

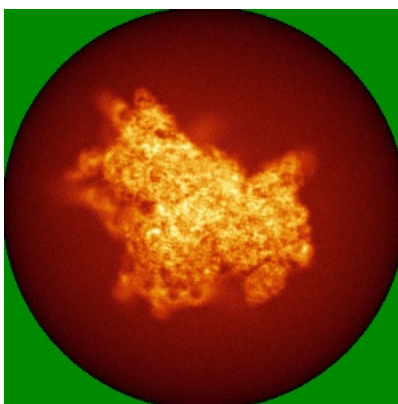


Z

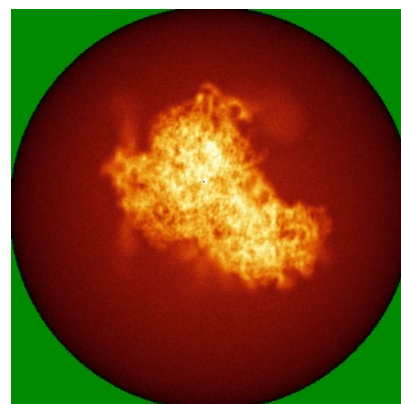
6.4.2 Raw map



X



Y

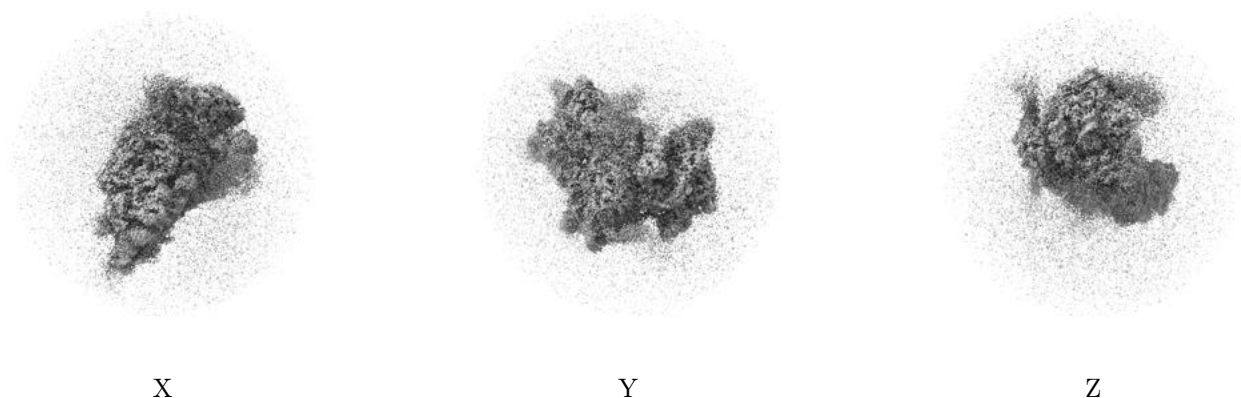


Z

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.

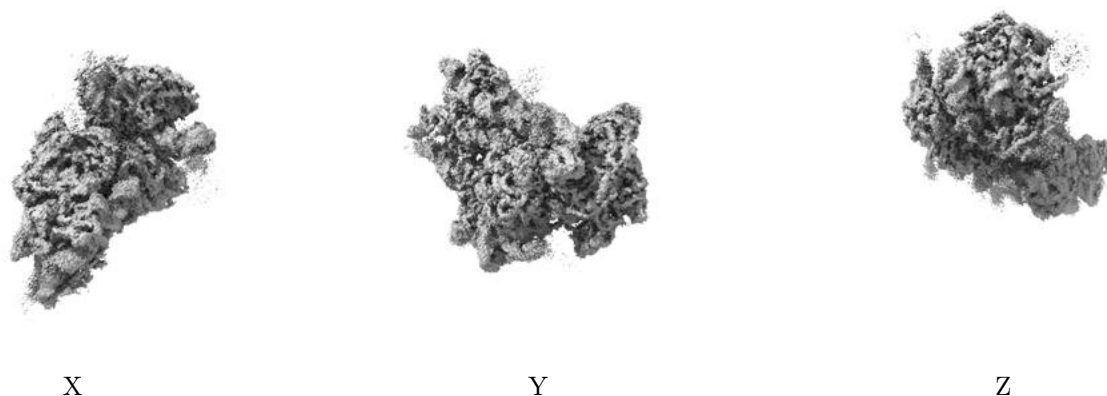
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0175. 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



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.

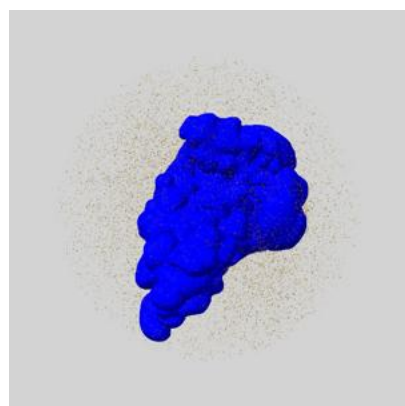
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

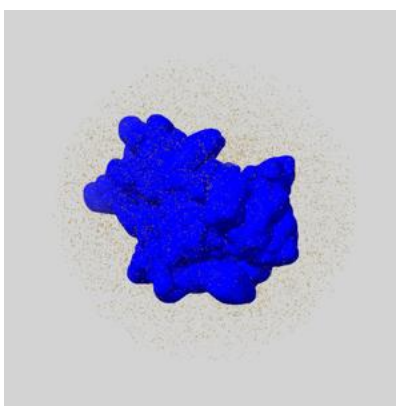
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

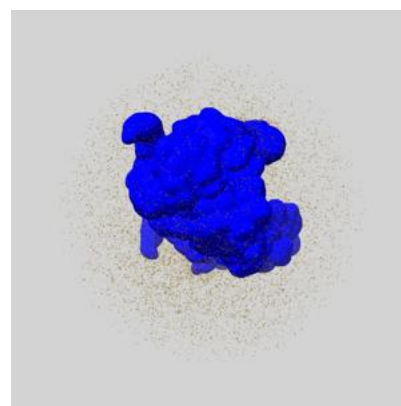
6.6.1 emd_25538_msk_1.map [i](#)



X



Y

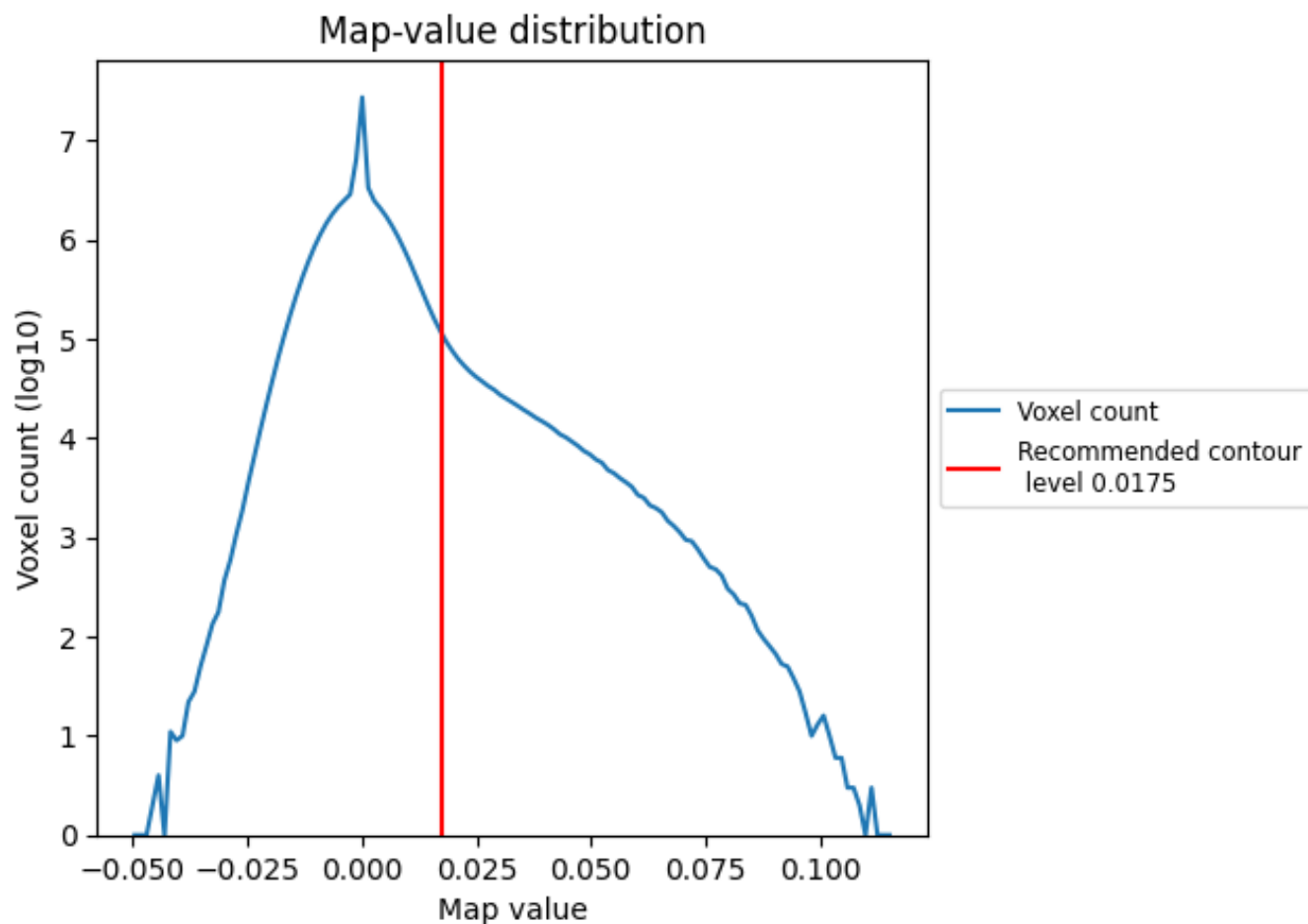


Z

7 Map analysis [i](#)

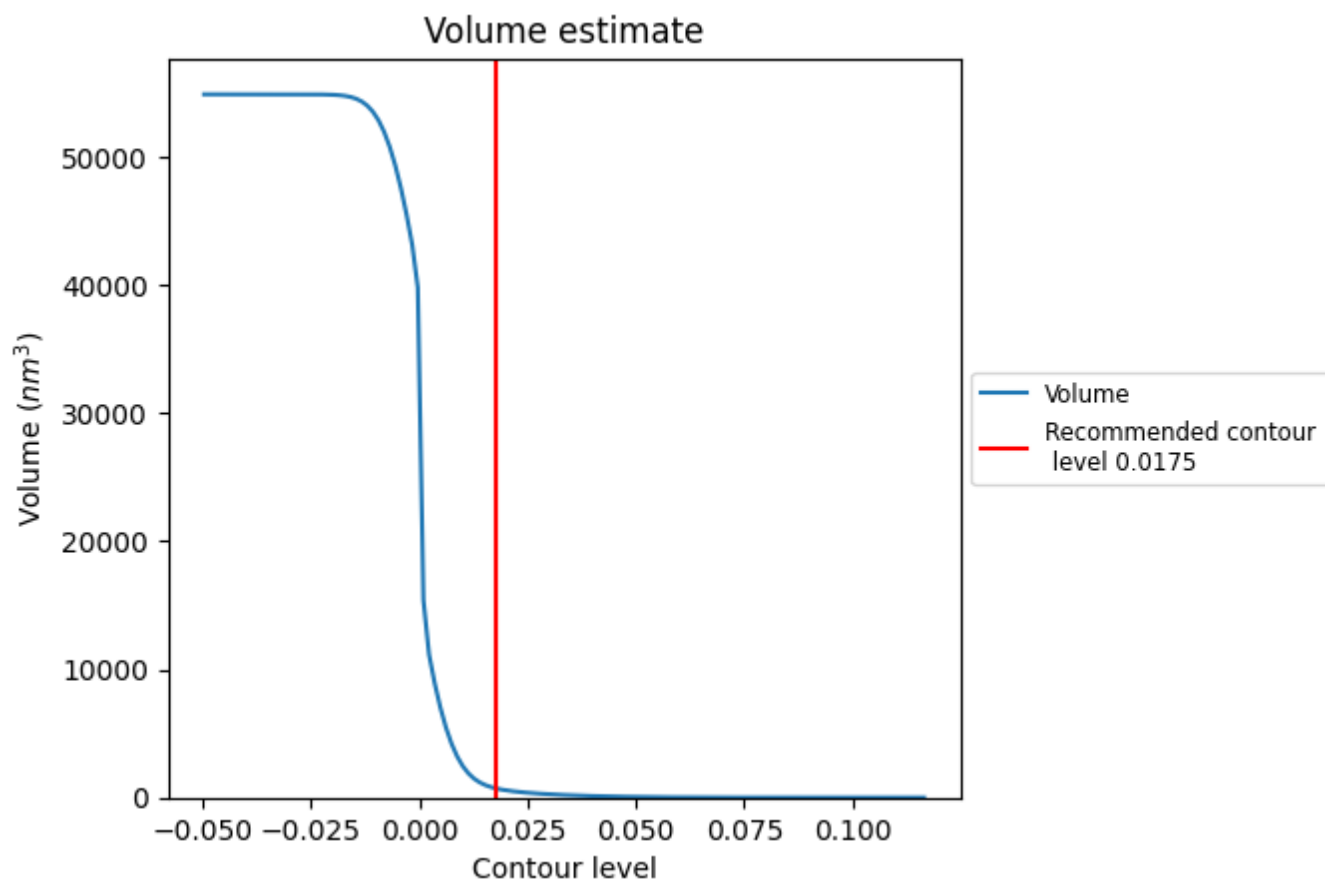
This section contains the results of statistical analysis of the map.

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.

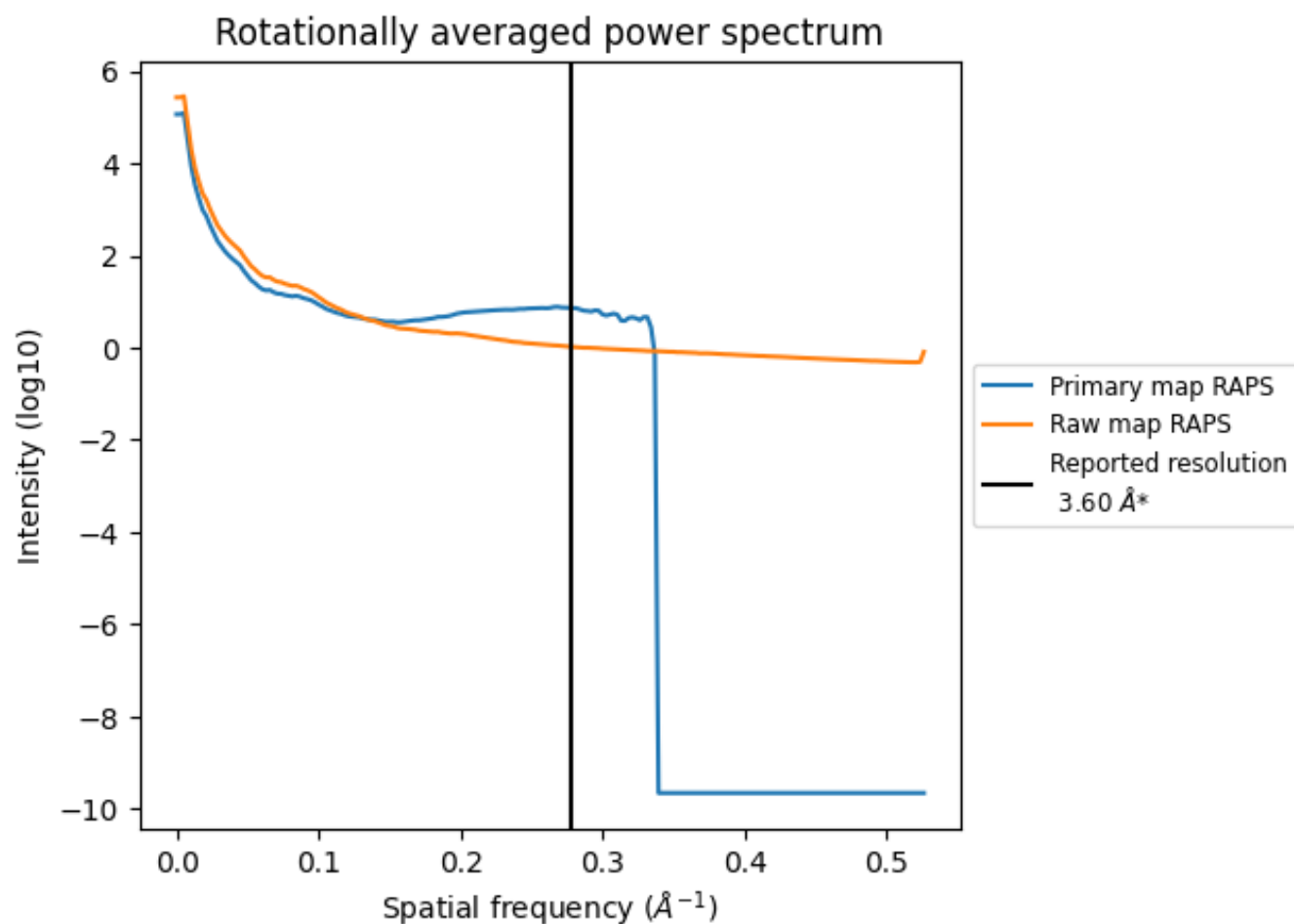
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 738 nm³; this corresponds to an approximate mass of 666 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 ⓘ

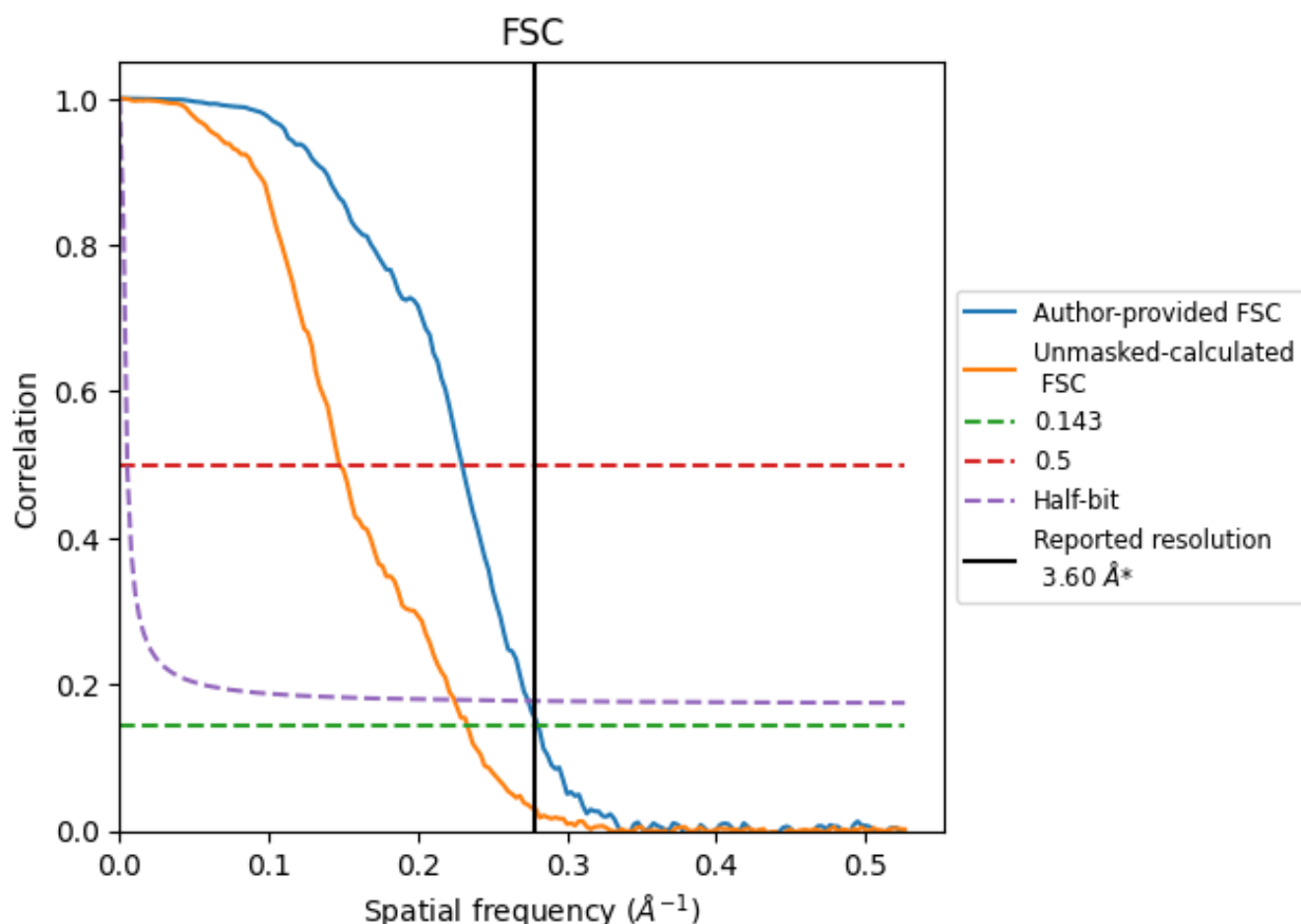


*Reported resolution corresponds to spatial frequency of 0.278 Å⁻¹

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 0.278 Å⁻¹

8.2 Resolution estimates [i](#)

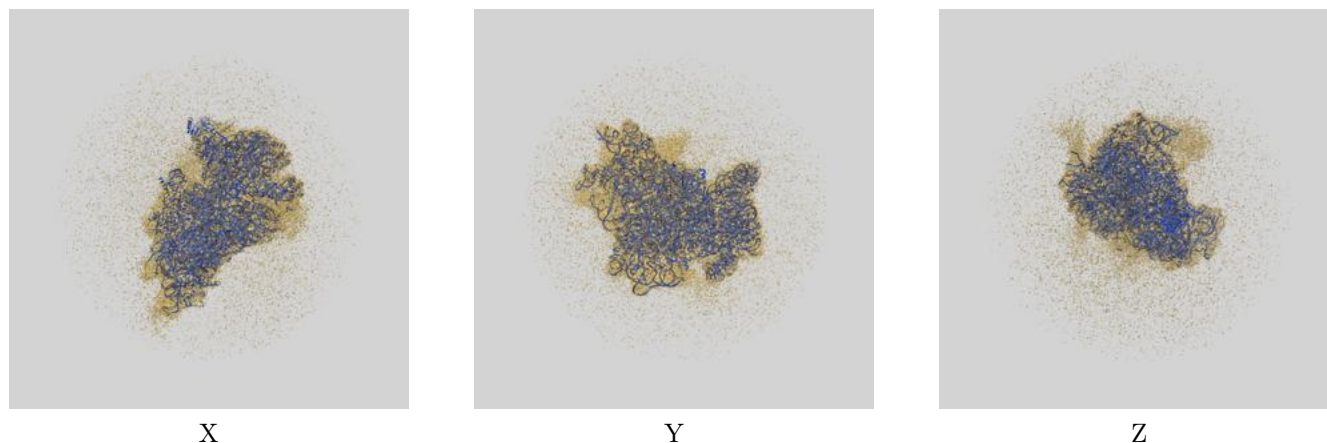
| Resolution estimate (Å) | Estimation criterion (FSC cut-off) | | |
|---------------------------|------------------------------------|------|----------|
| | 0.143 | 0.5 | Half-bit |
| Reported by author | 3.60 | - | - |
| Author-provided FSC curve | 3.57 | 4.36 | 3.65 |
| Unmasked-calculated* | 4.29 | 6.79 | 4.46 |

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.29 differs from the reported value 3.6 by more than 10 %

9 Map-model fit [i](#)

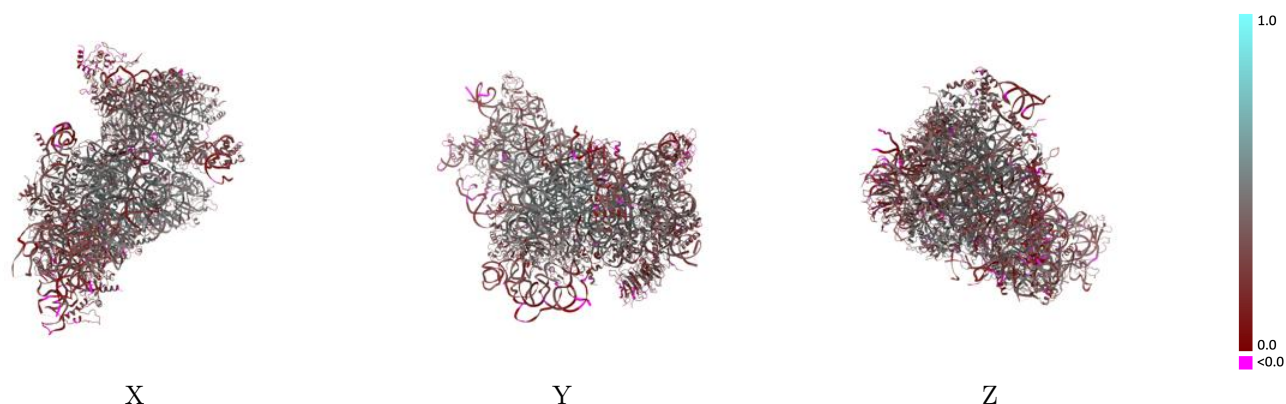
This section contains information regarding the fit between EMDB map EMD-25538 and PDB model 7SYR. Per-residue inclusion information can be found in [section 3](#) on [page 12](#).

9.1 Map-model overlay [i](#)



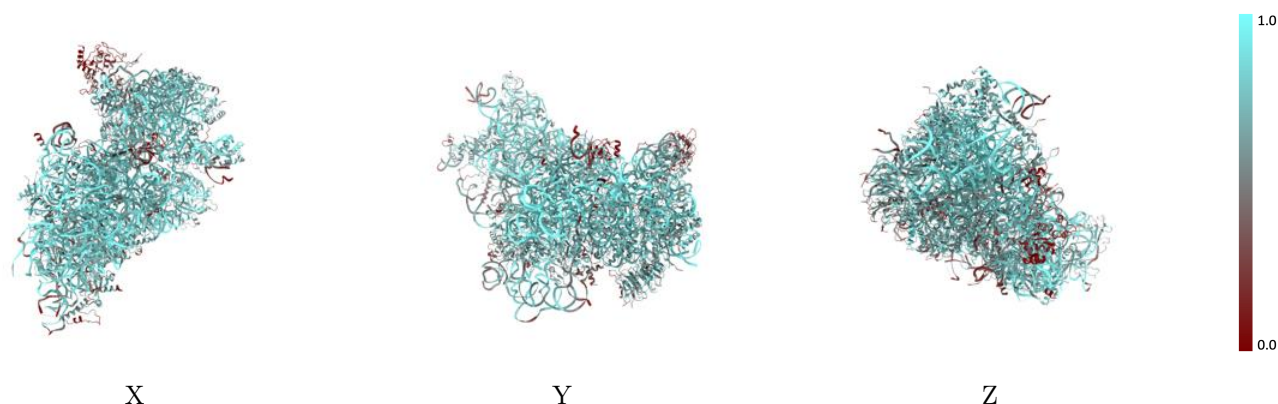
The images above show the 3D surface view of the map at the recommended contour level 0.0175 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



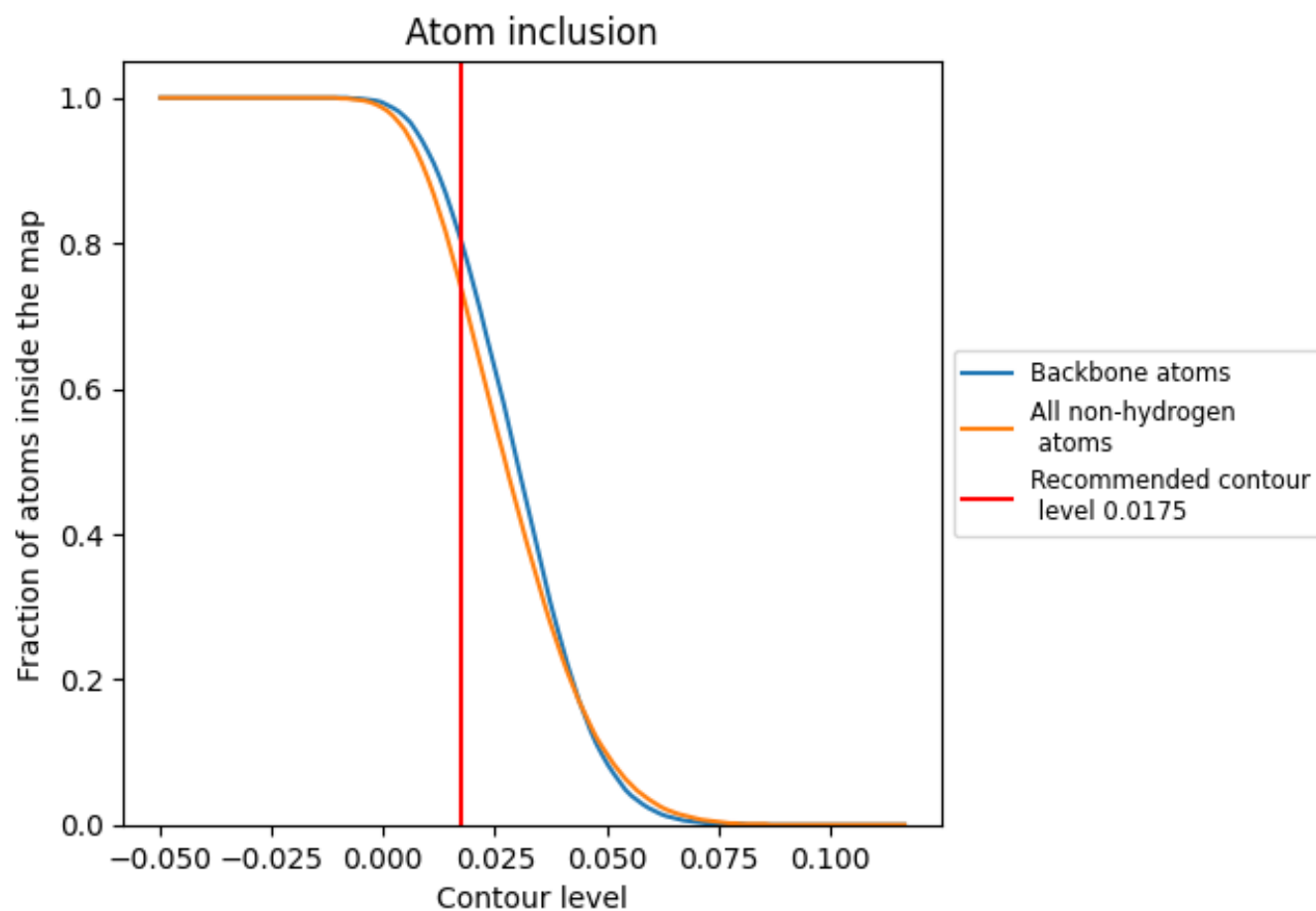
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.0175).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 80% of all backbone atoms, 74% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ











The table lists the average atom inclusion at the recommended contour level (0.0175) and Q-score for the entire model and for each chain.

| Chain | Atom inclusion | Q-score |
|-------|--|--|
| All |  0.7380 |  0.3470 |
| 2 |  0.8500 |  0.3810 |
| A |  0.1010 |  0.1500 |
| B |  0.7090 |  0.3860 |
| C |  0.6900 |  0.3620 |
| D |  0.7440 |  0.4260 |
| E |  0.5860 |  0.3100 |
| F |  0.7520 |  0.3750 |
| G |  0.7200 |  0.3950 |
| H |  0.6590 |  0.2560 |
| I |  0.5520 |  0.3240 |
| J |  0.6400 |  0.2860 |
| K |  0.7300 |  0.3900 |
| L |  0.5540 |  0.3040 |
| M |  0.6770 |  0.3620 |
| N |  0.1250 |  0.1340 |
| O |  0.7290 |  0.3640 |
| P |  0.7600 |  0.4110 |
| Q |  0.6490 |  0.3090 |
| R |  0.7180 |  0.3450 |
| S |  0.6070 |  0.3120 |
| T |  0.6930 |  0.3500 |
| U |  0.7440 |  0.3500 |
| V |  0.5150 |  0.2770 |
| W |  0.7190 |  0.3950 |
| X |  0.7700 |  0.4470 |
| Y |  0.7470 |  0.4060 |
| Z |  0.7120 |  0.3020 |
| a |  0.6480 |  0.3480 |
| b |  0.7890 |  0.4340 |
| c |  0.6420 |  0.3260 |
| d |  0.6260 |  0.3380 |
| e |  0.7190 |  0.3750 |
| f |  0.5270 |  0.2860 |
| g |  0.1960 |  0.1600 |



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| Chain | Atom inclusion | Q-score |
|-------|--|--|
| h |  0.6160 |  0.2670 |
| i |  0.7310 |  0.2760 |
| j |  0.7330 |  0.3140 |
| n |  0.6970 |  0.3790 |
| z |  0.6050 |  0.1890 |