



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 07:44 PM GMT

PDB ID : 4PJO
Title : Minimal U1 snRNP
Authors : Kondo, Y.; Oubridge, C.; van Roon, A.M.; Nagai, K.
Deposited on : 2014-05-12
Resolution : 3.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

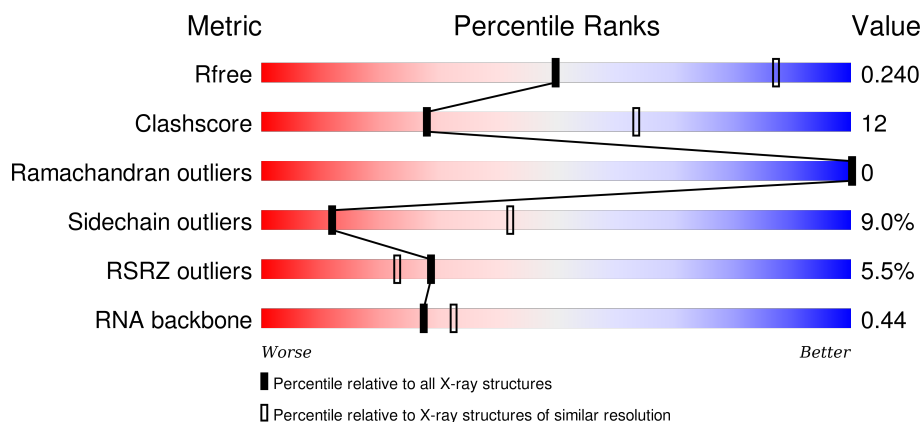
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 Å.

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



| Metric | Whole archive (#Entries) | Similar resolution (#Entries, resolution range(Å)) |
|-----------------------|-----------------------------|---|
| R_{free} | 91344 | 2060 (3.40-3.20) |
| Clashscore | 102246 | 1058 (3.38-3.22) |
| Ramachandran outliers | 100387 | 1038 (3.38-3.22) |
| Sidechain outliers | 100360 | 1037 (3.38-3.22) |
| RSRZ outliers | 91569 | 2070 (3.40-3.20) |
| RNA backbone | 2183 | 1005 (3.82-2.78) |

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

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1 | A | 126 | |
| 1 | O | 126 | |
| 1 | a | 126 | |
| 1 | o | 126 | |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 2 | B | 95 | |
| 2 | P | 95 | |
| 2 | b | 95 | |
| 2 | p | 95 | |
| 3 | C | 91 | |
| 3 | Q | 91 | |
| 3 | c | 91 | |
| 3 | q | 91 | |
| 4 | D | 118 | |
| 4 | R | 118 | |
| 4 | d | 118 | |
| 4 | r | 118 | |
| 5 | E | 92 | |
| 5 | S | 92 | |
| 5 | e | 92 | |
| 5 | s | 92 | |
| 6 | F | 75 | |
| 6 | T | 75 | |
| 6 | f | 75 | |
| 6 | t | 75 | |
| 7 | G | 76 | |
| 7 | U | 76 | |
| 7 | g | 76 | |
| 7 | u | 76 | |
| 8 | K | 60 | |

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| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|--|
| 8 | N | 60 |  |
| 8 | k | 60 |  |
| 8 | n | 60 |  |
| 9 | L | 61 |  |
| 9 | M | 61 |  |
| 9 | l | 61 |  |
| 9 | m | 61 |  |
| 10 | 1 | 60 |  |
| 10 | 2 | 60 |  |
| 10 | 3 | 60 |  |
| 10 | 4 | 60 |  |
| 11 | X | 10 |  |
| 11 | Y | 10 |  |
| 11 | x | 10 |  |
| 11 | y | 10 |  |

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

| Mol | Type | Chain | Res | Chirality | Geometry | Clashes | Electron density |
|-----|------|-------|-----|-----------|----------|---------|------------------|
| 13 | EPE | 2 | 201 | - | - | - | X |
| 14 | MG | 2 | 202 | - | - | - | X |
| 17 | CL | 4 | 203 | - | - | X | - |

2 Entry composition [i](#)

There are 19 unique types of molecules in this entry. The entry contains 26921 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Small nuclear ribonucleoprotein Sm D3.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 1 | A | 84 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 657 | 412 | 116 | 123 | 6 | | | |
| 1 | a | 83 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 652 | 409 | 115 | 122 | 6 | | | |
| 1 | O | 82 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 643 | 403 | 113 | 121 | 6 | | | |
| 1 | o | 81 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 637 | 400 | 112 | 119 | 6 | | | |

- Molecule 2 is a protein called Small nuclear ribonucleoprotein-associated proteins B and B'.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 2 | B | 83 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 673 | 425 | 123 | 118 | 7 | | | |
| 2 | b | 78 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 635 | 402 | 114 | 112 | 7 | | | |
| 2 | P | 83 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 673 | 425 | 123 | 118 | 7 | | | |
| 2 | p | 86 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 692 | 435 | 126 | 124 | 7 | | | |

- Molecule 3 is a protein called Small nuclear ribonucleoprotein Sm D1.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 3 | C | 86 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 668 | 423 | 117 | 125 | 3 | | | |
| 3 | c | 87 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 675 | 428 | 118 | 126 | 3 | | | |
| 3 | Q | 85 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 664 | 421 | 116 | 124 | 3 | | | |
| 3 | q | 87 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 674 | 426 | 118 | 127 | 3 | | | |

There are 28 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|---------|------------|
| C | -5 | GLY | - | linker | UNP P62314 |
| C | -4 | SER | - | linker | UNP P62314 |
| C | -3 | GLY | - | linker | UNP P62314 |
| C | -2 | SER | - | linker | UNP P62314 |
| C | -1 | GLY | - | linker | UNP P62314 |
| C | 0 | SER | - | linker | UNP P62314 |
| C | 1 | VAL | - | linker | UNP P62314 |
| c | -5 | GLY | - | linker | UNP P62314 |
| c | -4 | SER | - | linker | UNP P62314 |
| c | -3 | GLY | - | linker | UNP P62314 |
| c | -2 | SER | - | linker | UNP P62314 |
| c | -1 | GLY | - | linker | UNP P62314 |
| c | 0 | SER | - | linker | UNP P62314 |
| c | 1 | VAL | - | linker | UNP P62314 |
| Q | -5 | GLY | - | linker | UNP P62314 |
| Q | -4 | SER | - | linker | UNP P62314 |
| Q | -3 | GLY | - | linker | UNP P62314 |
| Q | -2 | SER | - | linker | UNP P62314 |
| Q | -1 | GLY | - | linker | UNP P62314 |
| Q | 0 | SER | - | linker | UNP P62314 |
| Q | 1 | VAL | - | linker | UNP P62314 |
| q | -5 | GLY | - | linker | UNP P62314 |
| q | -4 | SER | - | linker | UNP P62314 |
| q | -3 | GLY | - | linker | UNP P62314 |
| q | -2 | SER | - | linker | UNP P62314 |
| q | -1 | GLY | - | linker | UNP P62314 |
| q | 0 | SER | - | linker | UNP P62314 |
| q | 1 | VAL | - | linker | UNP P62314 |

- Molecule 4 is a protein called Small nuclear ribonucleoprotein Sm D2.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 4 | D | 98 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 796 | 498 | 144 | 148 | 6 | | | |
| 4 | d | 92 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 753 | 472 | 138 | 138 | 5 | | | |
| 4 | R | 96 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 790 | 497 | 141 | 146 | 6 | | | |
| 4 | r | 95 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 777 | 486 | 141 | 144 | 6 | | | |

- Molecule 5 is a protein called Small nuclear ribonucleoprotein E.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 5 | E | 77 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 638 | 405 | 113 | 115 | 5 | | | |
| 5 | e | 77 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 638 | 405 | 113 | 115 | 5 | | | |
| 5 | S | 77 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 638 | 405 | 113 | 115 | 5 | | | |
| 5 | s | 77 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 638 | 405 | 113 | 115 | 5 | | | |

- Molecule 6 is a protein called Small nuclear ribonucleoprotein F.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|-----|---|---------|---------|-------|
| 6 | F | 73 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 568 | 367 | 94 | 102 | 5 | | | |
| 6 | f | 74 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 577 | 373 | 95 | 104 | 5 | | | |
| 6 | T | 74 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 577 | 373 | 95 | 104 | 5 | | | |
| 6 | t | 74 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 577 | 373 | 95 | 104 | 5 | | | |

- Molecule 7 is a protein called Small nuclear ribonucleoprotein G.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|---|---------|---------|-------|
| 7 | G | 74 | Total | C | N | O | S | 0 | 1 | 0 |
| | | | 584 | 369 | 106 | 103 | 6 | | | |
| 7 | g | 73 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 568 | 358 | 102 | 102 | 6 | | | |
| 7 | U | 71 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 558 | 353 | 100 | 99 | 6 | | | |
| 7 | u | 73 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 568 | 358 | 102 | 102 | 6 | | | |

- Molecule 8 is a protein called U1 small nuclear ribonucleoprotein 70 kDa.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|---------|-------|
| 8 | K | 58 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 477 | 313 | 82 | 81 | 1 | | | |
| 8 | k | 57 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 470 | 309 | 81 | 79 | 1 | | | |
| 8 | N | 53 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 431 | 283 | 73 | 74 | 1 | | | |

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| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|---------|-------|
| 8 | n | 58 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 477 | 313 | 82 | 81 | 1 | | | |

There are 4 discrepancies between the modelled and reference sequences:

| Chain | Residue | Modelled | Actual | Comment | Reference |
|-------|---------|----------|--------|----------------|------------|
| K | 1 | GLY | - | expression tag | UNP P08621 |
| k | 1 | GLY | - | expression tag | UNP P08621 |
| N | 1 | GLY | - | expression tag | UNP P08621 |
| n | 1 | GLY | - | expression tag | UNP P08621 |

- Molecule 9 is a protein called U1 small nuclear ribonucleoprotein C.

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|----|----|---|---------|---------|-------|
| 9 | L | 53 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 448 | 280 | 77 | 87 | 4 | | | |
| 9 | l | 53 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 448 | 280 | 77 | 87 | 4 | | | |
| 9 | M | 53 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 448 | 280 | 77 | 87 | 4 | | | |
| 9 | m | 50 | Total | C | N | O | S | 0 | 0 | 0 |
| | | | 425 | 266 | 73 | 82 | 4 | | | |

- Molecule 10 is a RNA chain called U1 RNA variant (48-MER) with 4-helix junction replaced by kissing loop (HIV-1 (Mal) DIS) and shorter stem-loop 4..

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|-----|-----|-----|----|---------|---------|-------|
| 10 | 1 | 48 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 1036 | 458 | 183 | 345 | 50 | | | |
| 10 | 2 | 48 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 1036 | 458 | 183 | 345 | 50 | | | |
| 10 | 3 | 47 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 1013 | 448 | 178 | 338 | 49 | | | |
| 10 | 4 | 49 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 1059 | 468 | 188 | 352 | 51 | | | |

- Molecule 11 is a RNA chain called RNA (5'-R(*AP*GP*GP*UP*AP*AP*GP*UP*CP*C)-3').

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|----|----|----|---|---------|---------|-------|
| 11 | X | 10 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 212 | 96 | 40 | 67 | 9 | | | |

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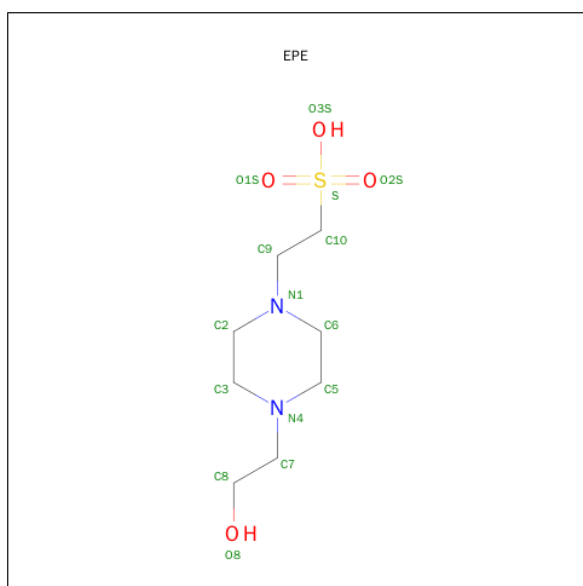
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| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|----|----|----|---|---------|---------|-------|
| 11 | x | 10 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 212 | 96 | 40 | 67 | 9 | | | |
| 11 | Y | 10 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 212 | 96 | 40 | 67 | 9 | | | |
| 11 | y | 10 | Total | C | N | O | P | 0 | 0 | 0 |
| | | | 212 | 96 | 40 | 67 | 9 | | | |

- Molecule 12 is ZINC ION (three-letter code: ZN) (formula: Zn).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 12 | l | 1 | Total | Zn | 0 | 0 |
| | | | 1 | 1 | | |
| 12 | m | 1 | Total | Zn | 0 | 0 |
| | | | 1 | 1 | | |
| 12 | L | 1 | Total | Zn | 0 | 0 |
| | | | 1 | 1 | | |
| 12 | M | 1 | Total | Zn | 0 | 0 |
| | | | 1 | 1 | | |

- Molecule 13 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf |
|-----|-------|----------|-------|---|---|---|---|---------|---------|
| 13 | 1 | 1 | Total | C | N | O | S | 0 | 0 |
| | | | 15 | 8 | 2 | 4 | 1 | | |
| 13 | 2 | 1 | Total | C | N | O | S | 0 | 0 |
| | | | 15 | 8 | 2 | 4 | 1 | | |

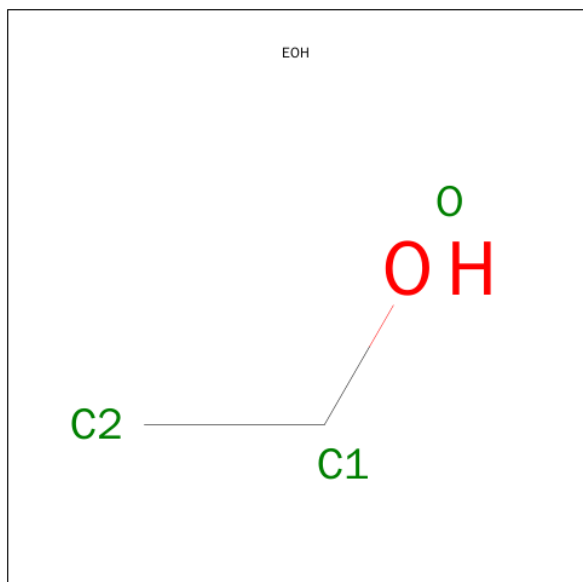
- Molecule 14 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|------------|---------|---------|---------|
| 14 | 1 | 8 | Total 8 | Mg 8 | 0 | 0 |
| 14 | 4 | 1 | Total 1 | Mg 1 | 0 | 0 |
| 14 | 2 | 4 | Total 4 | Mg 4 | 0 | 0 |
| 14 | y | 2 | Total 2 | Mg 2 | 0 | 0 |
| 14 | 1 | 1 | Total 1 | Mg 1 | 0 | 0 |
| 14 | 3 | 4 | Total 4 | Mg 4 | 0 | 0 |

- Molecule 15 is POTASSIUM ION (three-letter code: K) (formula: K).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|------------|--------|---------|---------|
| 15 | b | 2 | Total 2 | K 2 | 0 | 0 |
| 15 | 1 | 1 | Total 1 | K 1 | 0 | 0 |
| 15 | Q | 1 | Total 1 | K 1 | 0 | 0 |
| 15 | 3 | 2 | Total 2 | K 2 | 0 | 0 |

- Molecule 16 is ETHANOL (three-letter code: EOH) (formula: C₂H₆O).

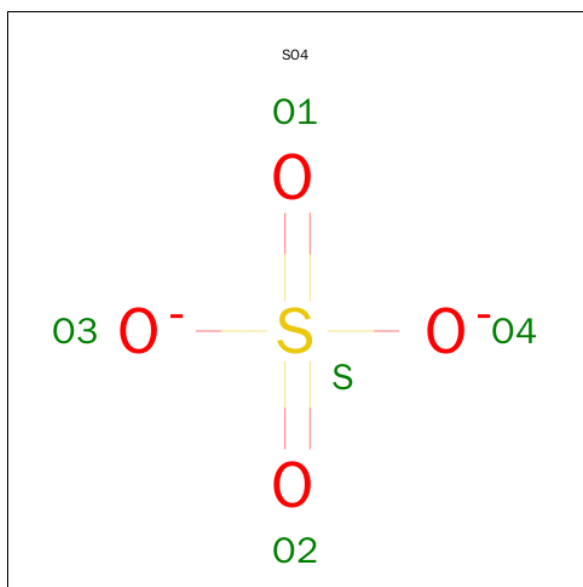


| Mol | Chain | Residues | Atoms | | | ZeroOcc | AltConf |
|-----|-------|----------|-------|---|---|---------|---------|
| 16 | 4 | 1 | Total | C | O | 0 | 0 |
| | | | 3 | 2 | 1 | | |

- Molecule 17 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 17 | 4 | 1 | Total | Cl | 0 | 0 |
| | | | 1 | 1 | | |

- Molecule 18 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



| Mol | Chain | Residues | Atoms | | | ZeroOcc | AltConf |
|-----|-------|----------|-------|---|---|---------|---------|
| 18 | y | 1 | Total | O | S | 0 | 0 |
| | | | 5 | 4 | 1 | | |

- Molecule 19 is water.

| Mol | Chain | Residues | Atoms | | ZeroOcc | AltConf |
|-----|-------|----------|-------|---|---------|---------|
| 19 | D | 5 | Total | O | 0 | 0 |
| | | | 5 | 5 | | |
| 19 | G | 2 | Total | O | 0 | 0 |
| | | | 2 | 2 | | |
| 19 | K | 1 | Total | O | 0 | 0 |
| | | | 1 | 1 | | |
| 19 | 1 | 6 | Total | O | 0 | 0 |
| | | | 6 | 6 | | |

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| Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|----------------|---------|---------|
| 19 | a | 1 | Total O 1 1 | 0 | 0 |
| 19 | b | 1 | Total O 1 1 | 0 | 0 |
| 19 | c | 2 | Total O 2 2 | 0 | 0 |
| 19 | d | 6 | Total O 6 6 | 0 | 0 |
| 19 | g | 1 | Total O 1 1 | 0 | 0 |
| 19 | 2 | 4 | Total O 4 4 | 0 | 0 |
| 19 | P | 2 | Total O 2 2 | 0 | 0 |
| 19 | R | 1 | Total O 1 1 | 0 | 0 |
| 19 | S | 1 | Total O 1 1 | 0 | 0 |
| 19 | T | 1 | Total O 1 1 | 0 | 0 |
| 19 | N | 1 | Total O 1 1 | 0 | 0 |
| 19 | 3 | 2 | Total O 2 2 | 0 | 0 |
| 19 | Y | 1 | Total O 1 1 | 0 | 0 |
| 19 | p | 1 | Total O 1 1 | 0 | 0 |
| 19 | q | 2 | Total O 2 2 | 0 | 0 |
| 19 | r | 1 | Total O 1 1 | 0 | 0 |
| 19 | t | 1 | Total O 1 1 | 0 | 0 |
| 19 | m | 1 | Total O 1 1 | 0 | 0 |
| 19 | 4 | 4 | Total O 4 4 | 0 | 0 |

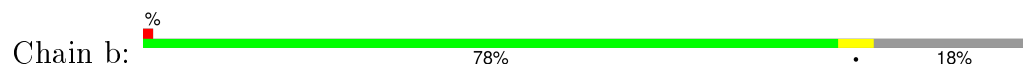
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GLY
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GLY
MET
GLY
ARG
GLY
ASN
ILE
PHE
GLN
LYS
ARG
ARG

- Molecule 2: Small nuclear ribonucleoprotein-associated proteins B and B'



MET THR VAL GLY LYS S6 S7 S8 S9 Q10 Q11 Q12 Q13 D14 Y15 R16 LYS M17 M18 R18 T30 C43 K50 I51 K52 P53 Q58 A59 E60 V78 T81 V82 P86 P87 P88 ASP THR GLY ILE ALA ARG VAL

- Molecule 2: Small nuclear ribonucleoprotein-associated proteins B and B'



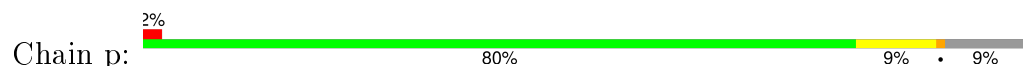
MET THR VAL GLY LYS S6 S7 Q11 R25 P53 LYS ASN SER LYS GLN ALA E60 L71 M80 P89 THR GLY ILE ALA ARG VAL

- Molecule 2: Small nuclear ribonucleoprotein-associated proteins B and B'



MET THR VAL GLY LYS S6 S7 S8 S9 Q10 Q11 Q12 Q13 D14 Y15 R16 LYS M17 M18 R18 T30 C43 K50 I51 K52 P53 Q58 A59 E60 V78 T81 V82 P86 P87 P88 ASP THR GLY ILE ALA ARG VAL

- Molecule 2: Small nuclear ribonucleoprotein-associated proteins B and B'



MET THR VAL GLY LYS S6 S7 H12 I13 D14 R18 K54 N55 S56 K57 Q58 L67 G68 L69 V70 L71 T90 G91 ILE ALA ARG VAL

- Molecule 3: Small nuclear ribonucleoprotein Sm D1



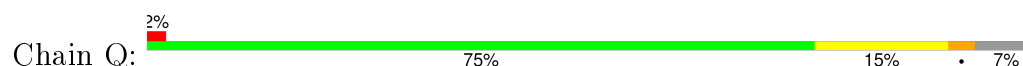
GLY SER G-3 V4 R5 M8 E18 L19 M36 K20 T23 Q24 V25 T30 G31 V32 D33 V34 L40 K41 M45 T46 L47 R50 E51 N63 N64 I65 R66 Y67 V81 D82 VAL GLU PRO

- Molecule 3: Small nuclear ribonucleoprotein Sm D1



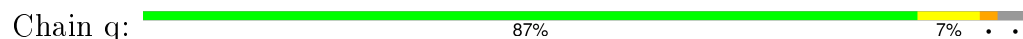
GLY SER G-3 V4 S11 L19 M36 K41 A42 M45 R66 V83 GLU PRO

- Molecule 3: Small nuclear ribonucleoprotein Sm D1

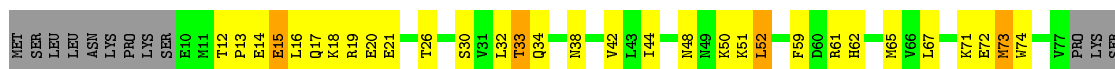




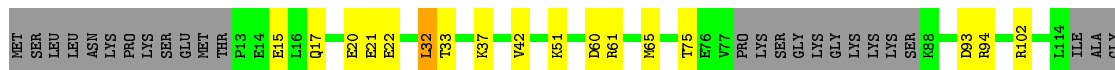
- Molecule 3: Small nuclear ribonucleoprotein Sm D1



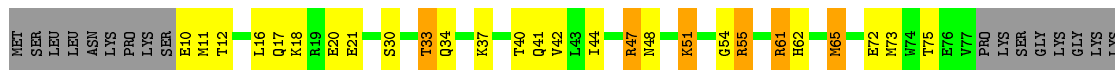
- Molecule 4: Small nuclear ribonucleoprotein Sm D2



- Molecule 4: Small nuclear ribonucleoprotein Sm D2



- Molecule 4: Small nuclear ribonucleoprotein Sm D2

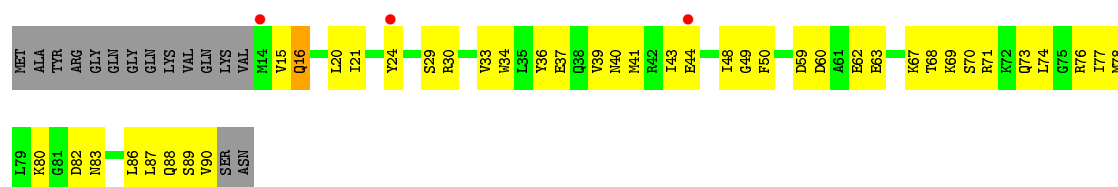


- Molecule 4: Small nuclear ribonucleoprotein Sm D2



- Molecule 5: Small nuclear ribonucleoprotein E





- Molecule 5: Small nuclear ribonucleoprotein E



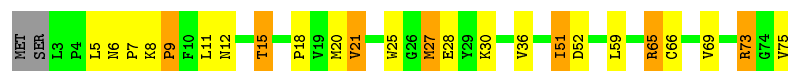
- Molecule 5: Small nuclear ribonucleoprotein E



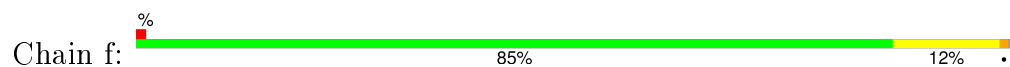
- Molecule 5: Small nuclear ribonucleoprotein E



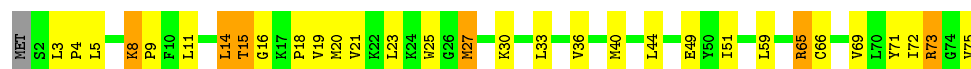
- Molecule 6: Small nuclear ribonucleoprotein F



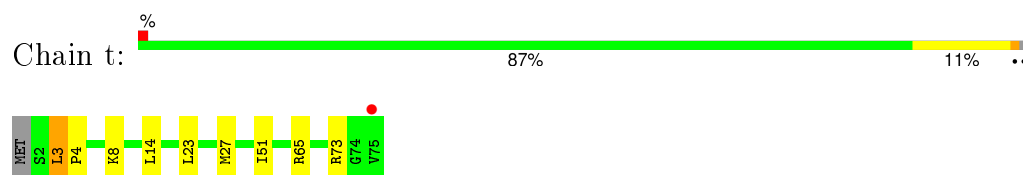
- Molecule 6: Small nuclear ribonucleoprotein F



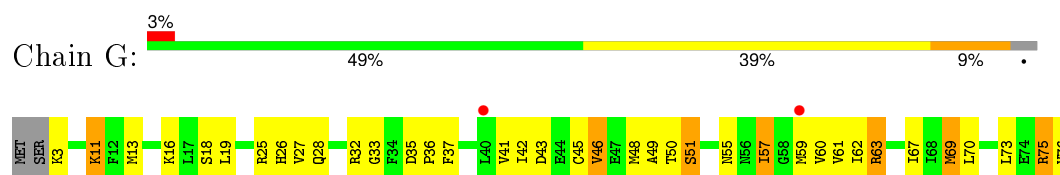
- Molecule 6: Small nuclear ribonucleoprotein F



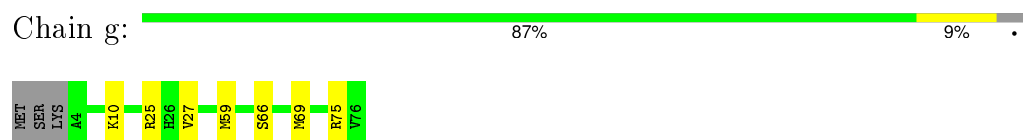
- Molecule 6: Small nuclear ribonucleoprotein F



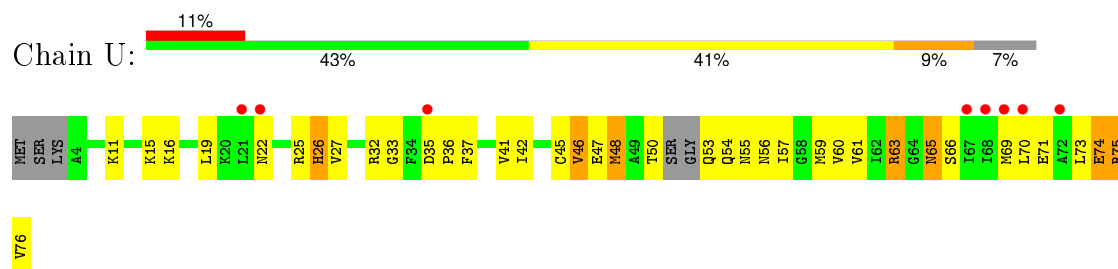
- Molecule 7: Small nuclear ribonucleoprotein G



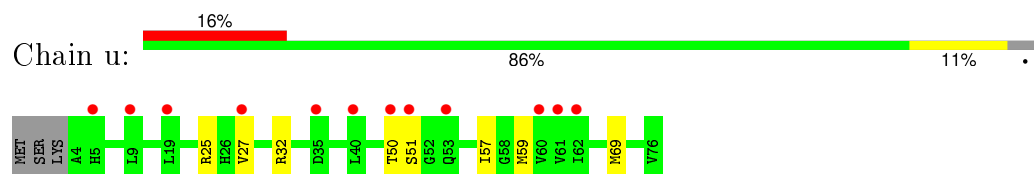
- Molecule 7: Small nuclear ribonucleoprotein G



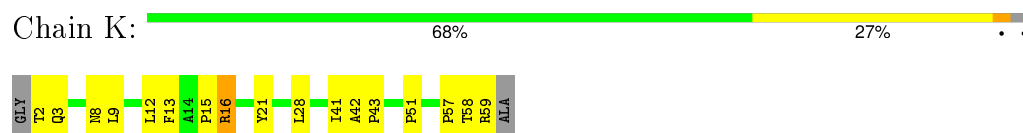
- Molecule 7: Small nuclear ribonucleoprotein G



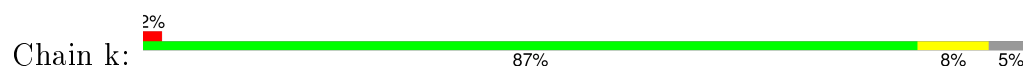
- Molecule 7: Small nuclear ribonucleoprotein G

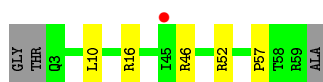


- Molecule 8: U1 small nuclear ribonucleoprotein 70 kDa

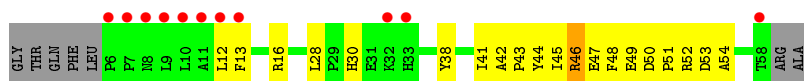


- Molecule 8: U1 small nuclear ribonucleoprotein 70 kDa

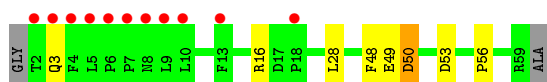
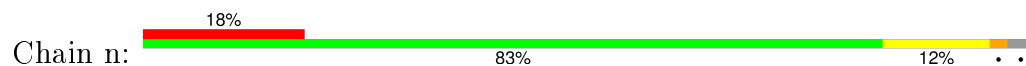




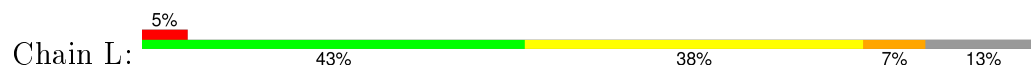
- Molecule 8: U1 small nuclear ribonucleoprotein 70 kDa



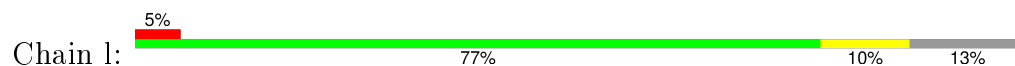
- Molecule 8: U1 small nuclear ribonucleoprotein 70 kDa



- Molecule 9: U1 small nuclear ribonucleoprotein C



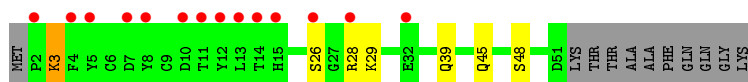
- Molecule 9: U1 small nuclear ribonucleoprotein C



- Molecule 9: U1 small nuclear ribonucleoprotein C



- Molecule 9: U1 small nuclear ribonucleoprotein C



- Molecule 10: U1 RNA variant (48-MER) with 4-helix junction replaced by kissing loop (HIV-1 (Mal) DIS) and shorter stem-loop 4.





- Molecule 10: U1 RNA variant (48-MER) with 4-helix junction replaced by kissing loop (HIV-1 (Mal) DIS) and shorter stem-loop 4.



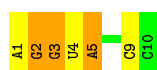
- Molecule 10: U1 RNA variant (48-MER) with 4-helix junction replaced by kissing loop (HIV-1 (Mal) DIS) and shorter stem-loop 4.



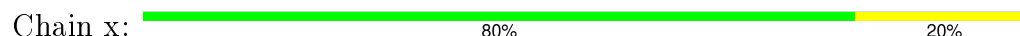
- Molecule 10: U1 RNA variant (48-MER) with 4-helix junction replaced by kissing loop (HIV-1 (Mal) DIS) and shorter stem-loop 4.



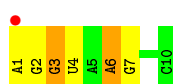
- Molecule 11: RNA (5'-R(*AP*GP*GP*UP*AP*AP*GP*UP*CP*C)-3')



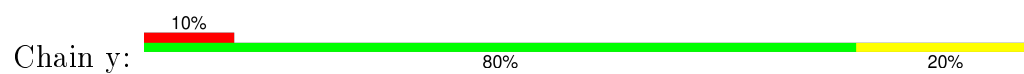
- Molecule 11: RNA (5'-R(*AP*GP*GP*UP*AP*AP*GP*UP*CP*C)-3')



- Molecule 11: RNA (5'-R(*AP*GP*GP*UP*AP*AP*GP*UP*CP*C)-3')



- Molecule 11: RNA (5'-R(*AP*GP*GP*UP*AP*AP*GP*UP*CP*C)-3')



4 Data and refinement statistics

| Property | Value | Source |
|---|---|------------------|
| Space group | P 21 21 21 | Depositor |
| Cell constants a, b, c, α , β , γ | 120.36Å 172.63Å 256.32Å 90.00° 90.00° 90.00° | Depositor |
| Resolution (Å) | 70.00 – 3.30 69.67 – 3.30 | Depositor EDS |
| % Data completeness (in resolution range) | 97.4 (70.00-3.30) 97.4 (69.67-3.30) | Depositor EDS |
| R_{merge} | 0.10 | Depositor |
| R_{sym} | (Not available) | Depositor |
| $\langle I/\sigma(I) \rangle$ ¹ | 3.52 (at 3.33Å) | Xtriage |
| Refinement program | REFMAC | Depositor |
| R, R_{free} | 0.207 , 0.255 0.196 , 0.240 | Depositor DCC |
| R_{free} test set | 4038 reflections (5.39%) | DCC |
| Wilson B-factor (Å ²) | 86.2 | Xtriage |
| Anisotropy | 0.287 | Xtriage |
| Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²) | 0.28 , 72.7 | EDS |
| Estimated twinning fraction | No twinning to report. | Xtriage |
| L-test for twinning ² | $\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$ | Xtriage |
| Outliers | 1 of 78932 reflections (0.001%) | Xtriage |
| F_o, F_c correlation | 0.90 | EDS |
| Total number of atoms | 26921 | wwPDB-VP |
| Average B, all atoms (Å ²) | 102.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.10% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MG, CL, K, EOH, ZN, GTP, SO4, EPE

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 | 0.74 | 1/665 (0.2%) | 1.04 | 4/896 (0.4%) |
| 1 | O | 0.52 | 0/651 | 1.15 | 5/878 (0.6%) |
| 1 | a | 0.71 | 0/660 | 1.18 | 8/889 (0.9%) |
| 1 | o | 0.53 | 0/645 | 1.19 | 6/870 (0.7%) |
| 2 | B | 0.77 | 0/683 | 1.08 | 2/910 (0.2%) |
| 2 | P | 0.65 | 0/683 | 1.08 | 1/910 (0.1%) |
| 2 | b | 0.75 | 0/644 | 1.10 | 1/858 (0.1%) |
| 2 | p | 0.63 | 0/702 | 1.16 | 3/936 (0.3%) |
| 3 | C | 0.87 | 0/676 | 1.19 | 6/914 (0.7%) |
| 3 | Q | 0.74 | 0/672 | 1.19 | 6/909 (0.7%) |
| 3 | c | 0.84 | 0/683 | 1.28 | 4/924 (0.4%) |
| 3 | q | 0.67 | 0/682 | 1.22 | 7/922 (0.8%) |
| 4 | D | 0.75 | 1/805 (0.1%) | 1.15 | 6/1081 (0.6%) |
| 4 | R | 0.80 | 0/803 | 1.32 | 8/1079 (0.7%) |
| 4 | d | 0.76 | 0/762 | 1.18 | 5/1022 (0.5%) |
| 4 | r | 0.69 | 0/786 | 1.14 | 2/1055 (0.2%) |
| 5 | E | 0.67 | 0/646 | 0.98 | 2/867 (0.2%) |
| 5 | S | 0.64 | 0/646 | 1.30 | 6/867 (0.7%) |
| 5 | e | 0.71 | 0/646 | 1.13 | 4/867 (0.5%) |
| 5 | s | 0.62 | 0/646 | 1.27 | 5/867 (0.6%) |
| 6 | F | 0.81 | 1/580 (0.2%) | 1.17 | 3/783 (0.4%) |
| 6 | T | 0.69 | 0/589 | 1.16 | 3/795 (0.4%) |
| 6 | f | 0.77 | 0/589 | 1.24 | 4/795 (0.5%) |
| 6 | t | 0.66 | 1/589 (0.2%) | 1.18 | 4/795 (0.5%) |
| 7 | G | 0.76 | 0/595 | 1.28 | 5/794 (0.6%) |
| 7 | U | 0.57 | 0/564 | 1.06 | 1/752 (0.1%) |
| 7 | g | 0.77 | 1/575 (0.2%) | 1.27 | 4/768 (0.5%) |
| 7 | u | 0.59 | 0/575 | 1.17 | 4/768 (0.5%) |
| 8 | K | 0.76 | 0/499 | 1.10 | 1/688 (0.1%) |
| 8 | N | 0.61 | 0/452 | 1.11 | 1/624 (0.2%) |
| 8 | k | 0.77 | 0/492 | 1.10 | 2/678 (0.3%) |
| 8 | n | 0.63 | 0/499 | 1.09 | 4/688 (0.6%) |

| Mol | Chain | Bond lengths | | Bond angles | |
|-----|-------|--------------|----------------|-------------|------------------|
| | | RMSZ | # Z >5 | RMSZ | # Z >5 |
| 9 | L | 0.60 | 0/460 | 1.00 | 2/618 (0.3%) |
| 9 | M | 0.61 | 0/460 | 1.12 | 6/618 (1.0%) |
| 9 | l | 0.58 | 0/460 | 0.96 | 0/618 |
| 9 | m | 0.56 | 0/437 | 1.16 | 4/587 (0.7%) |
| 10 | 1 | 0.62 | 0/1122 | 1.21 | 12/1747 (0.7%) |
| 10 | 2 | 0.58 | 0/1122 | 1.16 | 7/1747 (0.4%) |
| 10 | 3 | 0.49 | 0/1096 | 1.09 | 6/1706 (0.4%) |
| 10 | 4 | 0.49 | 0/1148 | 1.10 | 7/1788 (0.4%) |
| 11 | X | 0.44 | 0/237 | 1.33 | 5/368 (1.4%) |
| 11 | Y | 0.43 | 0/237 | 1.26 | 4/368 (1.1%) |
| 11 | x | 0.44 | 0/237 | 1.11 | 2/368 (0.5%) |
| 11 | y | 0.40 | 0/237 | 0.94 | 0/368 |
| All | All | 0.67 | 5/27637 (0.0%) | 1.16 | 182/38350 (0.5%) |

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 | 1 |
| 2 | P | 0 | 1 |
| 3 | C | 0 | 1 |
| 3 | c | 0 | 1 |
| 4 | R | 0 | 1 |
| 4 | r | 0 | 1 |
| 6 | T | 0 | 1 |
| 8 | k | 0 | 1 |
| 8 | n | 0 | 1 |
| All | All | 0 | 9 |

All (5) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(Å) | Ideal(Å) |
|-----|-------|-----|------|-------|-------|-------------|----------|
| 4 | D | 89 | PRO | N-CD | 5.34 | 1.55 | 1.47 |
| 1 | A | 81 | PRO | N-CD | 5.18 | 1.55 | 1.47 |
| 6 | F | 9 | PRO | N-CD | 5.13 | 1.55 | 1.47 |
| 7 | g | 66 | SER | CB-OG | -5.13 | 1.35 | 1.42 |
| 6 | t | 4 | PRO | N-CD | 5.04 | 1.54 | 1.47 |

The worst 5 of 182 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed($^{\circ}$) | Ideal($^{\circ}$) |
|-----|-------|-----|------|-------------|--------|------------------------|---------------------|
| 5 | S | 87 | LEU | CB-CG-CD2 | 13.57 | 134.07 | 111.00 |
| 9 | m | 28 | ARG | NE-CZ-NH1 | 12.22 | 126.41 | 120.30 |
| 6 | f | 73 | ARG | NE-CZ-NH1 | -10.81 | 114.89 | 120.30 |
| 4 | d | 32 | LEU | CB-CG-CD2 | 10.65 | 129.10 | 111.00 |
| 10 | 2 | 136 | G | C4'-C3'-O3' | -10.48 | 87.40 | 109.40 |

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

| Mol | Chain | Res | Type | Group |
|-----|-------|-----|------|---------|
| 1 | A | 1 | MET | Peptide |
| 3 | C | 81 | VAL | Peptide |
| 2 | P | 87 | PRO | Peptide |
| 3 | c | 42 | ALA | Peptide |
| 8 | k | 57 | PRO | Peptide |

5.2 Too-close contacts

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

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1 | A | 657 | 0 | 675 | 33 | 0 |
| 1 | O | 643 | 0 | 657 | 31 | 0 |
| 1 | a | 652 | 0 | 670 | 0 | 0 |
| 1 | o | 637 | 0 | 652 | 0 | 0 |
| 2 | B | 673 | 0 | 703 | 31 | 0 |
| 2 | P | 673 | 0 | 703 | 25 | 0 |
| 2 | b | 635 | 0 | 656 | 0 | 0 |
| 2 | p | 692 | 0 | 717 | 0 | 0 |
| 3 | C | 668 | 0 | 709 | 25 | 0 |
| 3 | Q | 664 | 0 | 705 | 14 | 0 |
| 3 | c | 675 | 0 | 718 | 0 | 0 |
| 3 | q | 674 | 0 | 714 | 0 | 0 |
| 4 | D | 796 | 0 | 821 | 68 | 0 |
| 4 | R | 790 | 0 | 813 | 38 | 0 |
| 4 | d | 753 | 0 | 779 | 0 | 0 |
| 4 | r | 777 | 0 | 800 | 0 | 0 |
| 5 | E | 638 | 0 | 657 | 60 | 1 |
| 5 | S | 638 | 0 | 657 | 79 | 0 |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 5 | e | 638 | 0 | 657 | 0 | 1 |
| 5 | s | 638 | 0 | 657 | 0 | 0 |
| 6 | F | 568 | 0 | 575 | 23 | 0 |
| 6 | T | 577 | 0 | 589 | 30 | 0 |
| 6 | f | 577 | 0 | 589 | 0 | 0 |
| 6 | t | 577 | 0 | 589 | 0 | 0 |
| 7 | G | 584 | 0 | 610 | 43 | 0 |
| 7 | U | 558 | 0 | 581 | 52 | 0 |
| 7 | g | 568 | 0 | 590 | 0 | 0 |
| 7 | u | 568 | 0 | 590 | 0 | 0 |
| 8 | K | 477 | 0 | 466 | 23 | 0 |
| 8 | N | 431 | 0 | 419 | 29 | 0 |
| 8 | k | 470 | 0 | 459 | 0 | 0 |
| 8 | n | 477 | 0 | 466 | 0 | 0 |
| 9 | L | 448 | 0 | 420 | 33 | 0 |
| 9 | M | 448 | 0 | 420 | 19 | 0 |
| 9 | l | 448 | 0 | 421 | 0 | 0 |
| 9 | m | 425 | 0 | 393 | 0 | 0 |
| 10 | 1 | 1036 | 0 | 515 | 15 | 0 |
| 10 | 2 | 1036 | 0 | 516 | 11 | 0 |
| 10 | 3 | 1013 | 0 | 505 | 10 | 0 |
| 10 | 4 | 1059 | 0 | 527 | 19 | 0 |
| 11 | X | 212 | 0 | 110 | 7 | 0 |
| 11 | Y | 212 | 0 | 110 | 6 | 0 |
| 11 | x | 212 | 0 | 110 | 0 | 0 |
| 11 | y | 212 | 0 | 110 | 0 | 0 |
| 12 | L | 1 | 0 | 0 | 0 | 0 |
| 12 | M | 1 | 0 | 0 | 0 | 0 |
| 12 | l | 1 | 0 | 0 | 0 | 0 |
| 12 | m | 1 | 0 | 0 | 0 | 0 |
| 13 | 1 | 15 | 0 | 18 | 1 | 0 |
| 13 | 2 | 15 | 0 | 18 | 1 | 0 |
| 14 | 1 | 8 | 0 | 0 | 0 | 0 |
| 14 | 2 | 4 | 0 | 0 | 0 | 0 |
| 14 | 3 | 4 | 0 | 0 | 0 | 0 |
| 14 | 4 | 1 | 0 | 0 | 0 | 0 |
| 14 | l | 1 | 0 | 0 | 0 | 0 |
| 14 | y | 2 | 0 | 0 | 0 | 0 |
| 15 | 1 | 1 | 0 | 0 | 0 | 0 |
| 15 | 3 | 2 | 0 | 0 | 0 | 0 |
| 15 | Q | 1 | 0 | 0 | 0 | 0 |
| 15 | b | 2 | 0 | 0 | 0 | 0 |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 16 | 4 | 3 | 0 | 6 | 0 | 0 |
| 17 | 4 | 1 | 0 | 0 | 4 | 0 |
| 18 | y | 5 | 0 | 0 | 0 | 0 |
| 19 | 1 | 6 | 0 | 0 | 2 | 0 |
| 19 | 2 | 4 | 0 | 0 | 0 | 0 |
| 19 | 3 | 2 | 0 | 0 | 0 | 0 |
| 19 | 4 | 4 | 0 | 0 | 1 | 0 |
| 19 | D | 5 | 0 | 0 | 5 | 0 |
| 19 | G | 2 | 0 | 0 | 0 | 0 |
| 19 | K | 1 | 0 | 0 | 0 | 0 |
| 19 | N | 1 | 0 | 0 | 0 | 0 |
| 19 | P | 2 | 0 | 0 | 3 | 0 |
| 19 | R | 1 | 0 | 0 | 1 | 0 |
| 19 | S | 1 | 0 | 0 | 0 | 0 |
| 19 | T | 1 | 0 | 0 | 0 | 0 |
| 19 | Y | 1 | 0 | 0 | 0 | 0 |
| 19 | a | 1 | 0 | 0 | 0 | 0 |
| 19 | b | 1 | 0 | 0 | 0 | 0 |
| 19 | c | 2 | 0 | 0 | 0 | 0 |
| 19 | d | 6 | 0 | 0 | 0 | 0 |
| 19 | g | 1 | 0 | 0 | 0 | 0 |
| 19 | m | 1 | 0 | 0 | 0 | 0 |
| 19 | p | 1 | 0 | 0 | 0 | 0 |
| 19 | q | 2 | 0 | 0 | 0 | 0 |
| 19 | r | 1 | 0 | 0 | 0 | 0 |
| 19 | t | 1 | 0 | 0 | 0 | 0 |
| All | All | 26921 | 0 | 24842 | 594 | 1 |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

The worst 5 of 594 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|-----------------|-----------------|--------------------------|-------------------|
| 7:U:65:ASN:HD21 | 10:3:127:U:C1' | 1.49 | 1.23 |
| 2:P:8:LYS:O | 2:P:12:HIS:NE2 | 3.81 | 1.23 |
| 2:B:17:MET:CE | 2:B:82:VAL:HA | 1.75 | 1.16 |
| 5:E:69:LYS:HA | 5:E:69:LYS:HE3 | 4.99 | 1.11 |
| 2:B:17:MET:HE3 | 2:B:82:VAL:HG22 | 1.28 | 1.11 |

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

| Atom-1 | Atom-2 | Interatomic distance (Å) | Clash overlap (Å) |
|----------------|-----------------------|--------------------------|-------------------|
| 5:E:62:GLU:OE2 | 5:e:28:ARG:NH1[4_445] | 1.93 | 0.27 |

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Percentiles | |
|-----|-------|--------------|----------|---------|----------|-------------|-----|
| 1 | A | 82/126 (65%) | 80 (98%) | 2 (2%) | 0 | 100 | 100 |
| 1 | O | 80/126 (64%) | 74 (92%) | 6 (8%) | 0 | 100 | 100 |
| 1 | a | 81/126 (64%) | 76 (94%) | 5 (6%) | 0 | 100 | 100 |
| 1 | o | 79/126 (63%) | 75 (95%) | 4 (5%) | 0 | 100 | 100 |
| 2 | B | 81/95 (85%) | 80 (99%) | 1 (1%) | 0 | 100 | 100 |
| 2 | P | 81/95 (85%) | 80 (99%) | 1 (1%) | 0 | 100 | 100 |
| 2 | b | 74/95 (78%) | 73 (99%) | 1 (1%) | 0 | 100 | 100 |
| 2 | p | 84/95 (88%) | 82 (98%) | 2 (2%) | 0 | 100 | 100 |
| 3 | C | 84/91 (92%) | 79 (94%) | 5 (6%) | 0 | 100 | 100 |
| 3 | Q | 83/91 (91%) | 81 (98%) | 2 (2%) | 0 | 100 | 100 |
| 3 | c | 85/91 (93%) | 81 (95%) | 4 (5%) | 0 | 100 | 100 |
| 3 | q | 85/91 (93%) | 83 (98%) | 2 (2%) | 0 | 100 | 100 |
| 4 | D | 94/118 (80%) | 90 (96%) | 4 (4%) | 0 | 100 | 100 |
| 4 | R | 93/118 (79%) | 87 (94%) | 6 (6%) | 0 | 100 | 100 |
| 4 | d | 88/118 (75%) | 86 (98%) | 2 (2%) | 0 | 100 | 100 |
| 4 | r | 91/118 (77%) | 85 (93%) | 6 (7%) | 0 | 100 | 100 |
| 5 | E | 75/92 (82%) | 71 (95%) | 4 (5%) | 0 | 100 | 100 |
| 5 | S | 75/92 (82%) | 72 (96%) | 3 (4%) | 0 | 100 | 100 |
| 5 | e | 75/92 (82%) | 72 (96%) | 3 (4%) | 0 | 100 | 100 |
| 5 | s | 75/92 (82%) | 70 (93%) | 5 (7%) | 0 | 100 | 100 |

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| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Percentiles | |
|-----|-------|-----------------|------------|----------|----------|-------------|-----|
| 6 | F | 71/75 (95%) | 67 (94%) | 4 (6%) | 0 | 100 | 100 |
| 6 | T | 72/75 (96%) | 68 (94%) | 4 (6%) | 0 | 100 | 100 |
| 6 | f | 72/75 (96%) | 69 (96%) | 3 (4%) | 0 | 100 | 100 |
| 6 | t | 72/75 (96%) | 69 (96%) | 3 (4%) | 0 | 100 | 100 |
| 7 | G | 73/76 (96%) | 69 (94%) | 4 (6%) | 0 | 100 | 100 |
| 7 | U | 67/76 (88%) | 64 (96%) | 3 (4%) | 0 | 100 | 100 |
| 7 | g | 71/76 (93%) | 68 (96%) | 3 (4%) | 0 | 100 | 100 |
| 7 | u | 71/76 (93%) | 69 (97%) | 2 (3%) | 0 | 100 | 100 |
| 8 | K | 56/60 (93%) | 52 (93%) | 4 (7%) | 0 | 100 | 100 |
| 8 | N | 51/60 (85%) | 47 (92%) | 4 (8%) | 0 | 100 | 100 |
| 8 | k | 55/60 (92%) | 53 (96%) | 2 (4%) | 0 | 100 | 100 |
| 8 | n | 56/60 (93%) | 47 (84%) | 9 (16%) | 0 | 100 | 100 |
| 9 | L | 51/61 (84%) | 48 (94%) | 3 (6%) | 0 | 100 | 100 |
| 9 | M | 51/61 (84%) | 47 (92%) | 4 (8%) | 0 | 100 | 100 |
| 9 | l | 51/61 (84%) | 48 (94%) | 3 (6%) | 0 | 100 | 100 |
| 9 | m | 48/61 (79%) | 47 (98%) | 1 (2%) | 0 | 100 | 100 |
| All | All | 2633/3176 (83%) | 2509 (95%) | 124 (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 X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|--------------|-----------|----------|-------------|----|
| 1 | A | 73/101 (72%) | 69 (94%) | 4 (6%) | 27 | 66 |
| 1 | O | 72/101 (71%) | 67 (93%) | 5 (7%) | 19 | 57 |
| 1 | a | 73/101 (72%) | 68 (93%) | 5 (7%) | 20 | 57 |
| 1 | o | 71/101 (70%) | 61 (86%) | 10 (14%) | 4 | 20 |
| 2 | B | 76/85 (89%) | 73 (96%) | 3 (4%) | 39 | 75 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|--------------|-----------|----------|-------------|-----|
| 2 | P | 76/85 (89%) | 71 (93%) | 5 (7%) | 21 | 59 |
| 2 | b | 72/85 (85%) | 69 (96%) | 3 (4%) | 36 | 73 |
| 2 | p | 78/85 (92%) | 70 (90%) | 8 (10%) | 9 | 34 |
| 3 | C | 79/83 (95%) | 78 (99%) | 1 (1%) | 76 | 89 |
| 3 | Q | 79/83 (95%) | 77 (98%) | 2 (2%) | 55 | 82 |
| 3 | c | 80/83 (96%) | 77 (96%) | 3 (4%) | 40 | 75 |
| 3 | q | 80/83 (96%) | 76 (95%) | 4 (5%) | 30 | 68 |
| 4 | D | 93/110 (84%) | 88 (95%) | 5 (5%) | 27 | 66 |
| 4 | R | 92/110 (84%) | 82 (89%) | 10 (11%) | 8 | 32 |
| 4 | d | 88/110 (80%) | 75 (85%) | 13 (15%) | 4 | 17 |
| 4 | r | 91/110 (83%) | 83 (91%) | 8 (9%) | 12 | 44 |
| 5 | E | 72/84 (86%) | 68 (94%) | 4 (6%) | 26 | 65 |
| 5 | S | 72/84 (86%) | 61 (85%) | 11 (15%) | 3 | 16 |
| 5 | e | 72/84 (86%) | 57 (79%) | 15 (21%) | 1 | 6 |
| 5 | s | 72/84 (86%) | 61 (85%) | 11 (15%) | 3 | 16 |
| 6 | F | 61/64 (95%) | 55 (90%) | 6 (10%) | 10 | 37 |
| 6 | T | 63/64 (98%) | 56 (89%) | 7 (11%) | 8 | 31 |
| 6 | f | 63/64 (98%) | 54 (86%) | 9 (14%) | 4 | 19 |
| 6 | t | 63/64 (98%) | 56 (89%) | 7 (11%) | 8 | 31 |
| 7 | G | 65/66 (98%) | 57 (88%) | 8 (12%) | 6 | 25 |
| 7 | U | 62/66 (94%) | 53 (86%) | 9 (14%) | 4 | 18 |
| 7 | g | 63/66 (96%) | 59 (94%) | 4 (6%) | 22 | 60 |
| 7 | u | 63/66 (96%) | 57 (90%) | 6 (10%) | 11 | 38 |
| 8 | K | 53/53 (100%) | 53 (100%) | 0 | 100 | 100 |
| 8 | N | 48/53 (91%) | 46 (96%) | 2 (4%) | 36 | 73 |
| 8 | k | 52/53 (98%) | 50 (96%) | 2 (4%) | 40 | 75 |
| 8 | n | 53/53 (100%) | 49 (92%) | 4 (8%) | 17 | 52 |
| 9 | L | 51/56 (91%) | 45 (88%) | 6 (12%) | 6 | 27 |
| 9 | M | 51/56 (91%) | 42 (82%) | 9 (18%) | 2 | 10 |
| 9 | l | 51/56 (91%) | 45 (88%) | 6 (12%) | 6 | 27 |
| 9 | m | 48/56 (86%) | 42 (88%) | 6 (12%) | 6 | 24 |

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| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|-----|-------|-----------------|------------|----------|-------------|----|
| All | All | 2471/2808 (88%) | 2250 (91%) | 221 (9%) | 12 | 43 |

5 of 221 residues with a non-rotameric sidechain are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1 | O | 78 | LYS |
| 5 | S | 60 | ASP |
| 6 | t | 65 | ARG |
| 2 | P | 11 | GLN |
| 4 | R | 33 | THR |

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 5 | E | 73 | GLN |
| 6 | F | 43 | GLN |
| 2 | P | 11 | GLN |
| 7 | U | 26 | HIS |

5.3.3 RNA ⓘ

| Mol | Chain | Analysed | Backbone Outliers | Pucker Outliers |
|-----|-------|---------------|-------------------|-----------------|
| 10 | 1 | 46/60 (76%) | 14 (30%) | 4 (8%) |
| 10 | 2 | 46/60 (76%) | 14 (30%) | 3 (6%) |
| 10 | 3 | 45/60 (75%) | 14 (31%) | 3 (6%) |
| 10 | 4 | 47/60 (78%) | 16 (34%) | 4 (8%) |
| 11 | X | 9/10 (90%) | 1 (11%) | 0 |
| 11 | Y | 9/10 (90%) | 2 (22%) | 0 |
| 11 | x | 9/10 (90%) | 0 | 0 |
| 11 | y | 9/10 (90%) | 2 (22%) | 0 |
| All | All | 220/280 (78%) | 63 (28%) | 14 (6%) |

5 of 63 RNA backbone outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 10 | 1 | 10 | U |
| 10 | 1 | 18 | A |
| 10 | 1 | 19 | G |
| 10 | 1 | 20 | G |
| 10 | 1 | 124 | U |

5 of 14 RNA pucker outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 10 | 2 | 137 | U |
| 10 | 3 | 126 | A |
| 10 | 4 | 128 | U |
| 10 | 2 | 128 | U |
| 10 | 4 | 126 | 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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 35 ligands modelled in this entry, 31 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

| Mol | Type | Chain | Res | Link | Bond lengths | | | Bond angles | | |
|-----|------|-------|-----|------|--------------|------|-------------|-------------|------|-------------|
| | | | | | Counts | RMSZ | $\# Z > 2$ | Counts | RMSZ | $\# Z > 2$ |
| 13 | EPE | 1 | 201 | - | 14,15,15 | 0.75 | 0 | 18,20,20 | 2.56 | 5 (27%) |
| 13 | EPE | 2 | 201 | - | 14,15,15 | 0.88 | 0 | 18,20,20 | 2.46 | 9 (50%) |
| 16 | EOH | 4 | 201 | - | 2,2,2 | 0.54 | 0 | 1,1,1 | 0.16 | 0 |
| 18 | SO4 | y | 103 | - | 4,4,4 | 0.48 | 0 | 6,6,6 | 0.49 | 0 |

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

| Mol | Type | Chain | Res | Link | Chirals | Torsions | Rings |
|-----|------|-------|-----|------|---------|-----------|---------|
| 13 | EPE | 1 | 201 | - | - | 0/9/19/19 | 0/1/1/1 |

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| Mol | Type | Chain | Res | Link | Chirals | Torsions | Rings |
|-----|------|-------|-----|------|---------|-----------|---------|
| 13 | EPE | 2 | 201 | - | - | 0/9/19/19 | 0/1/1/1 |
| 16 | EOH | 4 | 201 | - | - | 0/0/0/0 | 0/0/0/0 |
| 18 | SO4 | y | 103 | - | - | 0/0/0/0 | 0/0/0/0 |

There are no bond length outliers.

The worst 5 of 14 bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | Z | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-----------|-------|-------------|----------|
| 13 | 1 | 201 | EPE | O2S-S-O1S | -3.21 | 101.77 | 113.48 |
| 13 | 2 | 201 | EPE | C2-C3-N4 | -2.42 | 106.29 | 110.63 |
| 13 | 2 | 201 | EPE | O2S-S-O1S | -2.39 | 104.76 | 113.48 |
| 13 | 2 | 201 | EPE | C6-N1-C2 | 2.27 | 113.81 | 108.90 |
| 13 | 2 | 201 | EPE | C5-C6-N1 | 2.40 | 114.93 | 110.63 |

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

| Mol | Chain | Res | Type | Clashes | Symm-Clashes |
|-----|-------|-----|------|---------|--------------|
| 13 | 1 | 201 | EPE | 1 | 0 |
| 13 | 2 | 201 | EPE | 1 | 0 |

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

| Mol | Chain | Analysed | <RSRZ> | #RSRZ>2 | OWAB(Å ²) | Q<0.9 |
|-----|-------|--------------|--------|--------------|-----------------------|-------|
| 1 | A | 84/126 (66%) | 0.21 | 3 (3%) 46 39 | 50, 68, 102, 155 | 0 |
| 1 | O | 82/126 (65%) | 1.33 | 24 (29%) 1 1 | 96, 125, 152, 182 | 0 |
| 1 | a | 83/126 (65%) | -0.20 | 1 (1%) 81 76 | 46, 70, 103, 123 | 0 |
| 1 | o | 81/126 (64%) | 0.87 | 14 (17%) 2 2 | 87, 137, 161, 176 | 0 |
| 2 | B | 83/95 (87%) | 0.66 | 3 (3%) 46 39 | 41, 65, 189, 219 | 0 |
| 2 | P | 83/95 (87%) | 1.01 | 14 (16%) 2 2 | 68, 111, 178, 219 | 0 |
| 2 | b | 78/95 (82%) | 0.08 | 1 (1%) 79 74 | 45, 70, 135, 185 | 0 |
| 2 | p | 86/95 (90%) | 0.10 | 2 (2%) 64 57 | 75, 122, 185, 221 | 0 |
| 3 | C | 86/91 (94%) | -0.08 | 0 100 100 | 39, 56, 120, 139 | 0 |
| 3 | Q | 85/91 (93%) | 0.09 | 2 (2%) 62 55 | 55, 77, 125, 140 | 0 |
| 3 | c | 87/91 (95%) | 0.06 | 0 100 100 | 42, 60, 124, 139 | 0 |
| 3 | q | 87/91 (95%) | -0.20 | 0 100 100 | 57, 82, 131, 150 | 0 |
| 4 | D | 98/118 (83%) | 0.04 | 0 100 100 | 38, 62, 151, 181 | 0 |
| 4 | R | 96/118 (81%) | 0.16 | 0 100 100 | 49, 73, 157, 176 | 0 |
| 4 | d | 92/118 (77%) | -0.03 | 0 100 100 | 41, 63, 154, 176 | 0 |
| 4 | r | 95/118 (80%) | -0.04 | 1 (1%) 82 78 | 63, 85, 173, 233 | 0 |
| 5 | E | 77/92 (83%) | 0.56 | 3 (3%) 43 36 | 49, 75, 113, 134 | 0 |
| 5 | S | 77/92 (83%) | -0.28 | 2 (2%) 59 53 | 75, 117, 159, 177 | 0 |
| 5 | e | 77/92 (83%) | 0.67 | 2 (2%) 59 53 | 52, 81, 126, 150 | 0 |
| 5 | s | 77/92 (83%) | 0.26 | 4 (5%) 31 25 | 86, 115, 155, 171 | 0 |
| 6 | F | 73/75 (97%) | -0.01 | 0 100 100 | 45, 64, 95, 133 | 0 |
| 6 | T | 74/75 (98%) | -0.07 | 0 100 100 | 59, 88, 136, 175 | 0 |
| 6 | f | 74/75 (98%) | 0.39 | 1 (1%) 78 73 | 52, 76, 101, 129 | 0 |
| 6 | t | 74/75 (98%) | 0.31 | 1 (1%) 78 73 | 73, 100, 144, 178 | 0 |

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| Mol | Chain | Analysed | <RSRZ> | #RSRZ>2 | OWAB(Å ²) | Q<0.9 |
|-----|-------|-----------------|--------|----------------|-----------------------|-------|
| 7 | G | 74/76 (97%) | 0.66 | 2 (2%) 58 51 | 51, 80, 120, 137 | 0 |
| 7 | U | 71/76 (93%) | 0.36 | 8 (11%) 7 5 | 95, 130, 165, 178 | 0 |
| 7 | g | 73/76 (96%) | 0.01 | 0 100 100 | 55, 83, 123, 168 | 0 |
| 7 | u | 73/76 (96%) | 0.76 | 12 (16%) 2 2 | 98, 133, 171, 188 | 0 |
| 8 | K | 58/60 (96%) | -0.11 | 0 100 100 | 52, 83, 139, 159 | 0 |
| 8 | N | 53/60 (88%) | 1.08 | 11 (20%) 1 1 | 57, 121, 162, 171 | 0 |
| 8 | k | 57/60 (95%) | -0.16 | 1 (1%) 71 65 | 52, 90, 130, 155 | 0 |
| 8 | n | 58/60 (96%) | 1.01 | 11 (18%) 2 1 | 80, 140, 185, 196 | 0 |
| 9 | L | 53/61 (86%) | 0.29 | 3 (5%) 27 22 | 72, 106, 180, 197 | 0 |
| 9 | M | 53/61 (86%) | 1.53 | 17 (32%) 1 1 | 102, 140, 194, 199 | 0 |
| 9 | l | 53/61 (86%) | 0.30 | 3 (5%) 27 22 | 70, 112, 177, 202 | 0 |
| 9 | m | 50/61 (81%) | 1.36 | 14 (28%) 1 1 | 116, 157, 210, 217 | 0 |
| 10 | 1 | 47/60 (78%) | -0.37 | 0 100 100 | 46, 93, 150, 176 | 0 |
| 10 | 2 | 47/60 (78%) | -0.21 | 0 100 100 | 47, 99, 158, 175 | 0 |
| 10 | 3 | 46/60 (76%) | -0.20 | 0 100 100 | 72, 127, 151, 178 | 0 |
| 10 | 4 | 48/60 (80%) | -0.02 | 0 100 100 | 77, 129, 173, 196 | 0 |
| 11 | X | 10/10 (100%) | -0.75 | 0 100 100 | 88, 119, 157, 164 | 0 |
| 11 | Y | 10/10 (100%) | -0.06 | 1 (10%) 9 8 | 114, 143, 162, 164 | 0 |
| 11 | x | 10/10 (100%) | -0.04 | 0 100 100 | 89, 118, 165, 176 | 0 |
| 11 | y | 10/10 (100%) | 0.61 | 1 (10%) 9 8 | 130, 152, 171, 177 | 0 |
| All | All | 2943/3456 (85%) | 0.29 | 162 (5%) 29 23 | 38, 95, 165, 233 | 0 |

The worst 5 of 162 RSRZ outliers are listed below:

| Mol | Chain | Res | Type | RSRZ |
|-----|-------|-----|------|------|
| 2 | P | 56 | SER | 18.1 |
| 8 | n | 2 | THR | 9.7 |
| 2 | P | 55 | ASN | 8.7 |
| 8 | n | 3 | GLN | 7.5 |
| 2 | P | 57 | LYS | 7.1 |

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

| Mol | Type | Chain | Res | Atoms | RSCC | RSR | LLDF | B-factors(Å ²) | Q<0.9 |
|-----|------|-------|-----|-------|------|------|-------|----------------------------|-------|
| 14 | MG | 2 | 202 | 1/1 | 0.92 | 0.36 | 12.31 | 69,69,69,69 | 0 |
| 13 | EPE | 2 | 201 | 15/15 | 0.81 | 0.43 | 3.02 | 92,116,141,146 | 0 |
| 13 | EPE | 1 | 201 | 15/15 | 0.77 | 0.35 | 1.93 | 131,146,150,162 | 0 |
| 14 | MG | 2 | 203 | 1/1 | 0.35 | 0.18 | 1.18 | 87,87,87,87 | 0 |
| 14 | MG | y | 102 | 1/1 | 0.88 | 0.21 | -0.72 | 100,100,100,100 | 0 |
| 12 | ZN | l | 201 | 1/1 | 0.99 | 0.11 | -1.11 | 98,98,98,98 | 0 |
| 12 | ZN | L | 201 | 1/1 | 0.99 | 0.13 | -1.54 | 83,83,83,83 | 0 |
| 14 | MG | 3 | 201 | 1/1 | 0.89 | 0.15 | -1.82 | 75,75,75,75 | 0 |
| 15 | K | Q | 101 | 1/1 | 0.93 | 0.13 | -2.41 | 97,97,97,97 | 0 |
| 17 | CL | 4 | 203 | 1/1 | 0.95 | 0.11 | -3.16 | 111,111,111,111 | 0 |
| 14 | MG | 1 | 204 | 1/1 | 0.95 | 0.15 | -5.05 | 64,64,64,64 | 0 |
| 14 | MG | 3 | 203 | 1/1 | 0.90 | 0.12 | - | 80,80,80,80 | 0 |
| 16 | EOH | 4 | 201 | 3/3 | 0.43 | 1.09 | - | 112,112,115,119 | 0 |
| 14 | MG | l | 202 | 1/1 | 0.44 | 0.94 | - | 106,106,106,106 | 0 |
| 14 | MG | 1 | 203 | 1/1 | 0.87 | 0.20 | - | 52,52,52,52 | 0 |
| 14 | MG | 1 | 209 | 1/1 | 0.65 | 0.27 | - | 95,95,95,95 | 0 |
| 14 | MG | 2 | 204 | 1/1 | 0.60 | 0.27 | - | 94,94,94,94 | 0 |
| 14 | MG | y | 101 | 1/1 | 0.89 | 0.30 | - | 100,100,100,100 | 0 |
| 12 | ZN | M | 201 | 1/1 | 0.97 | 0.10 | - | 126,126,126,126 | 0 |
| 15 | K | 3 | 206 | 1/1 | 0.88 | 0.41 | - | 125,125,125,125 | 0 |
| 15 | K | 1 | 210 | 1/1 | 0.95 | 0.09 | - | 91,91,91,91 | 0 |
| 14 | MG | 1 | 205 | 1/1 | 0.70 | 0.30 | - | 71,71,71,71 | 0 |
| 15 | K | b | 101 | 1/1 | 0.88 | 0.33 | - | 105,105,105,105 | 0 |
| 14 | MG | 3 | 204 | 1/1 | 0.81 | 0.25 | - | 97,97,97,97 | 0 |
| 18 | SO4 | y | 103 | 5/5 | 0.39 | 0.59 | - | 192,193,205,208 | 0 |
| 14 | MG | 1 | 207 | 1/1 | 0.80 | 0.18 | - | 82,82,82,82 | 0 |
| 14 | MG | 1 | 202 | 1/1 | 0.80 | 0.27 | - | 100,100,100,100 | 0 |
| 14 | MG | 4 | 202 | 1/1 | 0.78 | 0.31 | - | 86,86,86,86 | 0 |
| 15 | K | b | 102 | 1/1 | 0.89 | 0.24 | - | 124,124,124,124 | 0 |
| 12 | ZN | m | 201 | 1/1 | 0.98 | 0.09 | - | 136,136,136,136 | 0 |

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| Mol | Type | Chain | Res | Atoms | RSCC | RSR | LLDF | B-factors(\AA^2) | Q<0.9 |
|-----|------|-------|-----|-------|------|------|------|-----------------------------|-------|
| 14 | MG | 3 | 202 | 1/1 | 0.48 | 0.13 | - | 91,91,91,91 | 0 |
| 14 | MG | 1 | 206 | 1/1 | 0.86 | 0.27 | - | 86,86,86,86 | 0 |
| 14 | MG | 2 | 205 | 1/1 | 0.87 | 0.26 | - | 61,61,61,61 | 0 |
| 14 | MG | 1 | 208 | 1/1 | 0.72 | 0.27 | - | 75,75,75,75 | 0 |
| 15 | K | 3 | 205 | 1/1 | 0.44 | 0.23 | - | 122,122,122,122 | 0 |

6.5 Other polymers [i](#)

There are no such residues in this entry.