



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 08:37 PM UTC

PDB ID : 1GID / pdb\_00001gid  
Title : CRYSTAL STRUCTURE OF A GROUP I RIBOZYME DOMAIN: PRINCIPLES OF RNA PACKING  
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Deposited on : 1996-08-22  
Resolution : 2.50 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

|                                |   |                                                                  |
|--------------------------------|---|------------------------------------------------------------------|
| MolProbity                     | : | 4-5-2 with Phenix2.0                                             |
| Mogul                          | : | 2022.3.0, CSD as543be (2022)                                     |
| Xtriage (Phenix)               | : | NOT EXECUTED                                                     |
| EDS                            | : | NOT EXECUTED                                                     |
| Percentile statistics          | : | 20250101.v01 (using entries in the PDB archive January 1st 2025) |
| Ideal geometry (proteins)      | : | Engh & Huber (2001)                                              |
| Ideal geometry (DNA, RNA)      | : | Parkinson et al. (1996)                                          |
| Validation Pipeline (wwPDB-VP) | : | 2.49                                                             |

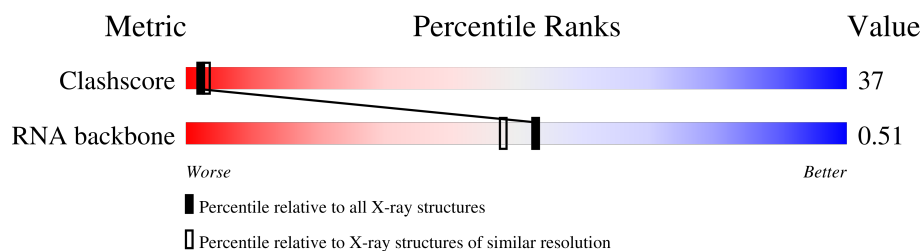
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

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



| Metric       | Whole archive<br>(#Entries) | Similar resolution<br>(#Entries, resolution range(Å)) |
|--------------|-----------------------------|-------------------------------------------------------|
| Clashscore   | 190562                      | 6492 (2.50-2.50)                                      |
| RNA backbone | 3983                        | 1003 (2.78-2.22)                                      |

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ .

Note EDS was not executed.

| Mol | Chain | Length | Quality of chain |
|-----|-------|--------|------------------|
| 1   | A     | 158    |                  |
| 1   | B     | 158    |                  |

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6824 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.

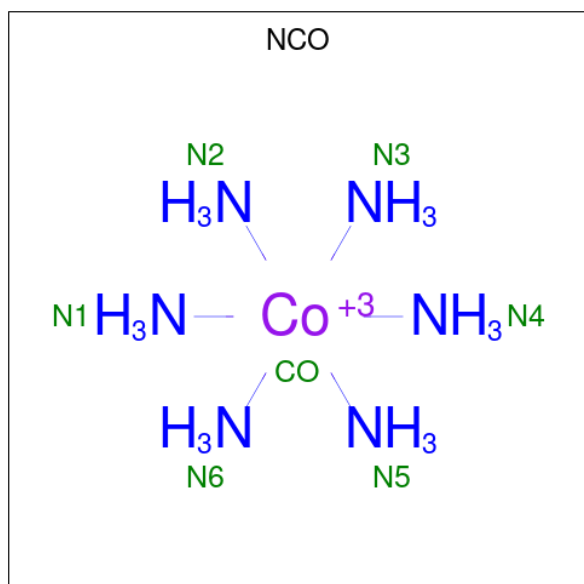
- Molecule 1 is a RNA chain called P4-P6 RNA RIBOZYME DOMAIN.

| Mol | Chain | Residues | Atoms |      |     |      |     | ZeroOcc | AltConf | Trace |
|-----|-------|----------|-------|------|-----|------|-----|---------|---------|-------|
| 1   | A     | 158      | Total | C    | N   | O    | P   | 58      | 0       | 0     |
|     |       |          | 3383  | 1511 | 616 | 1098 | 158 |         |         |       |
| 1   | B     | 158      | Total | C    | N   | O    | P   | 85      | 0       | 0     |
|     |       |          | 3383  | 1511 | 616 | 1098 | 158 |         |         |       |

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

| Mol | Chain | Residues | Atoms |    | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---------|---------|
| 2   | A     | 12       | Total | Mg | 0       | 0       |
|     |       |          | 12    | 12 |         |         |
| 2   | B     | 12       | Total | Mg | 0       | 0       |
|     |       |          | 12    | 12 |         |         |

- Molecule 3 is COBALT HEXAMMINE(III) (CCD ID: NCO) (formula:  $\text{CoH}_{18}\text{N}_6$ ).



| Mol | Chain | Residues | Atoms |    |   | ZeroOcc | AltConf |
|-----|-------|----------|-------|----|---|---------|---------|
| 3   | A     | 1        | Total | Co | N | 0       | 0       |
|     |       |          | 7     | 1  | 6 |         |         |
| 3   | A     | 1        | Total | Co | N | 0       | 0       |
|     |       |          | 7     | 1  | 6 |         |         |
| 3   | B     | 1        | Total | Co | N | 0       | 0       |
|     |       |          | 7     | 1  | 6 |         |         |
| 3   | B     | 1        | Total | Co | N | 0       | 0       |
|     |       |          | 7     | 1  | 6 |         |         |

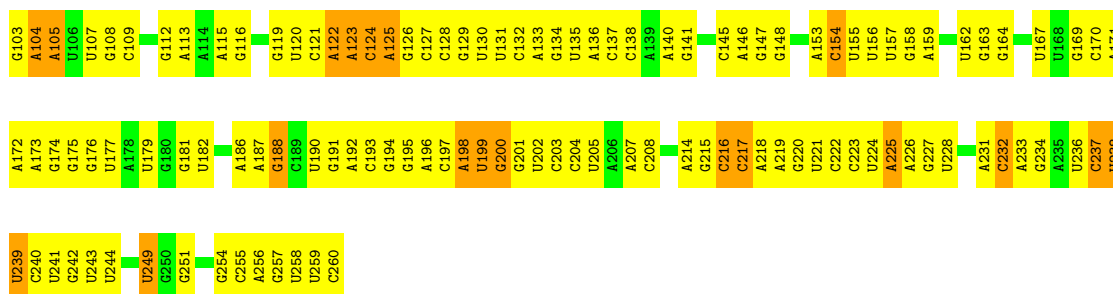
- Molecule 4 is water.

| Mol | Chain | Residues | Atoms |   | ZeroOcc | AltConf |
|-----|-------|----------|-------|---|---------|---------|
| 4   | A     | 3        | Total | O | 0       | 0       |
|     |       |          | 3     | 3 |         |         |
| 4   | B     | 3        | Total | O | 0       | 0       |
|     |       |          | 3     | 3 |         |         |

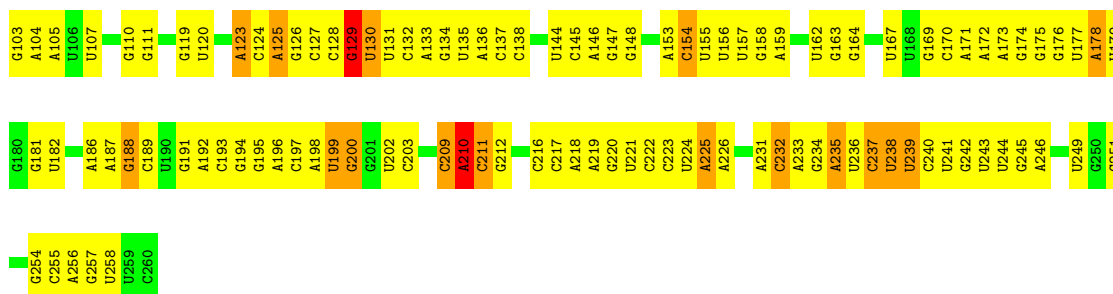


Note EDS was not executed.

Chain A:  26% 62% 12%



Chain B:  32% 56% 10%



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

| Property                                                 | Value                                          | Source    |
|----------------------------------------------------------|------------------------------------------------|-----------|
| Space group                                              | P 21 21 21                                     | Depositor |
| Cell constants<br>a, b, c, $\alpha$ , $\beta$ , $\gamma$ | 74.80Å 128.70Å 145.90Å<br>90.00° 90.00° 90.00° | Depositor |
| Resolution (Å)                                           | 8.00 – 2.50                                    | Depositor |
| % Data completeness<br>(in resolution range)             | 73.5 (8.00-2.50)                               | Depositor |
| $R_{merge}$                                              | (Not available)                                | Depositor |
| $R_{sym}$                                                | 0.04                                           | Depositor |
| Refinement program                                       | X-PLOR 3.843                                   | Depositor |
| R, $R_{free}$                                            | 0.242 , 0.285                                  | Depositor |
| Estimated twinning fraction                              | No twinning to report.                         | Xtriage   |
| Total number of atoms                                    | 6824                                           | wwPDB-VP  |
| Average B, all atoms (Å <sup>2</sup> )                   | 44.0                                           | wwPDB-VP  |

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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     | 0.63         | 0/3787  | 0.85        | 0/5901         |
| 1   | B     | 0.69         | 0/3787  | 0.87        | 4/5901 (0.1%)  |
| All | All   | 0.66         | 0/7574  | 0.86        | 4/11802 (0.0%) |

There are no bond length outliers.

All (4) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms       | Z     | Observed(°) | Ideal(°) |
|-----|-------|-----|------|-------------|-------|-------------|----------|
| 1   | B     | 210 | A    | C2'-C3'-O3' | 6.62  | 119.43      | 109.50   |
| 1   | B     | 178 | A    | C5'-C4'-C3' | -5.52 | 106.92      | 115.20   |
| 1   | B     | 129 | G    | C4'-C3'-O3' | 5.37  | 117.45      | 109.40   |
| 1   | B     | 209 | C    | C2'-C3'-O3' | -5.30 | 105.75      | 113.70   |

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 1   | A     | 3383  | 0        | 1703     | 211     | 0            |
| 1   | B     | 3383  | 0        | 1703     | 155     | 0            |
| 2   | A     | 12    | 0        | 0        | 0       | 0            |
| 2   | B     | 12    | 0        | 0        | 0       | 0            |
| 3   | A     | 14    | 0        | 0        | 3       | 0            |

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| Mol | Chain | Non-H | H(model) | H(added) | Clashes | Symm-Clashes |
|-----|-------|-------|----------|----------|---------|--------------|
| 3   | B     | 14    | 0        | 0        | 3       | 0            |
| 4   | A     | 3     | 0        | 0        | 0       | 0            |
| 4   | B     | 3     | 0        | 0        | 0       | 0            |
| All | All   | 6824  | 0        | 3406     | 365     | 0            |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 37.

The worst 5 of 365 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) |
|---------------|--------------|--------------------------|-------------------|
| 1:A:196:A:H2' | 1:A:197:C:C6 | 1.69                     | 1.27              |
| 1:A:128:C:H2' | 1:A:129:G:H8 | 1.12                     | 1.10              |
| 1:A:128:C:H2' | 1:A:129:G:C8 | 1.93                     | 1.03              |
| 1:A:196:A:H2' | 1:A:197:C:H6 | 0.89                     | 1.03              |
| 1:A:115:A:O5' | 1:A:115:A:H8 | 1.47                     | 0.98              |

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

There are no protein molecules in this entry.

### 5.3.2 Protein sidechains [i](#)

There are no protein molecules in this entry.

### 5.3.3 RNA [i](#)

| Mol | Chain | Analysed      | Backbone Outliers | Pucker Outliers |
|-----|-------|---------------|-------------------|-----------------|
| 1   | A     | 157/158 (99%) | 26 (16%)          | 6 (3%)          |
| 1   | B     | 157/158 (99%) | 23 (14%)          | 5 (3%)          |
| All | All   | 314/316 (99%) | 49 (15%)          | 11 (3%)         |

5 of 49 RNA backbone outliers are listed below:



| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | A     | 104 | A    |
| 1   | A     | 105 | A    |
| 1   | A     | 109 | C    |
| 1   | A     | 122 | A    |
| 1   | A     | 123 | A    |

5 of 11 RNA pucker outliers are listed below:

| Mol | Chain | Res | Type |
|-----|-------|-----|------|
| 1   | B     | 129 | G    |
| 1   | B     | 179 | U    |
| 1   | B     | 236 | U    |
| 1   | B     | 225 | A    |
| 1   | A     | 225 | 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 [i](#)

Of 28 ligands modelled in this entry, 24 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 | $\# Z  > 2$ |
| 3   | NCO  | B     | 264 | -    | 6,6,6        | 0.08 | 0           | -           |      |             |
| 3   | NCO  | A     | 262 | -    | 6,6,6        | 0.08 | 0           | -           |      |             |
| 3   | NCO  | B     | 263 | -    | 6,6,6        | 0.11 | 0           | -           |      |             |
| 3   | NCO  | A     | 261 | -    | 6,6,6        | 0.11 | 0           | -           |      |             |

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.

4 monomers are involved in 6 short contacts:

| Mol | Chain | Res | Type | Clashes | Symm-Clashes |
|-----|-------|-----|------|---------|--------------|
| 3   | B     | 264 | NCO  | 1       | 0            |
| 3   | A     | 262 | NCO  | 1       | 0            |
| 3   | B     | 263 | NCO  | 2       | 0            |
| 3   | A     | 261 | NCO  | 2       | 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.