

Full wwPDB X-ray Structure Validation Report (i)

Sep 26, 2022 – 03:51 PM EDT

PDB ID : 7SDG

Title : [C:Ag+:C] Metal-mediated DNA base pair in a self-assembling rhombohedral

lattice

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Deposited on : 2021-09-29

Resolution : 3.63 Å(reported)

This is a Full wwPDB X-ray Structure 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/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

 $\begin{array}{ccc} & Mol Probity & : & 4.02b\text{-}467 \\ Xtriage \left(Phenix\right) & : & 1.13 \end{array}$

EDS : 2.31.2buster-report : 1.1.7 (2018)

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

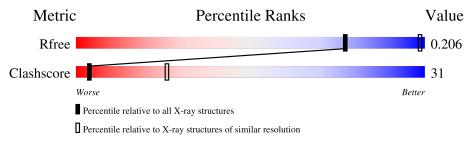
Validation Pipeline (wwPDB-VP) : 2.31.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 3.63 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 $(\# \mathrm{Entries})$ | $\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\AA)}) \end{array}$ |
|------------|---------------------------------------|--|
| R_{free} | 130704 | 1341 (3.78-3.50) |
| Clashscore | 141614 | 1439 (3.78-3.50) |

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 >=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 <=5%

| Mol | Chain | Length | Quality of chain | | | | | | | |
|-----|-------|--------|------------------|-----|-----|--|--|--|--|--|
| 1 | A | 21 | 14% | 67% | 19% | | | | | |
| 2 | В | 7 | | 71% | 29% | | | | | |
| 3 | С | 7 | 14% | 86% | | | | | | |
| 4 | D | 7 | | 86% | 14% | | | | | |



2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 858 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 DNA chain called DNA (5'-D(*GP*AP*GP*CP*AP*GP*CP*TP*GP*TP*GP*TP*GP*AP*TP*CP*A)-3').

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|--------------|----------|---------|----------|---------|---------|---------|-------|
| 1 | A | 21 | Total 428 | C 204 | N 81 | O 123 | P 20 | 0 | 0 | 0 |

• Molecule 2 is a DNA chain called DNA (5'-D(P*CP*CP*AP*CP*AP*CP*A)-3').

| Mol | Chain | Residues | Atoms | | | | | ZeroOcc | AltConf | Trace |
|-----|-------|----------|--------------|---------|---------|---------|--------|---------|---------|-------|
| 2 | В | 7 | Total 139 | C 66 | N 27 | O 39 | P 7 | 0 | 0 | 0 |

• Molecule 3 is a DNA chain called DNA (5'-D(P*GP*GP*CP*TP*GP*CP*T)-3').

| Mol | Chain | Residues | Atoms | | | ZeroOcc | AltConf | Trace | | |
|-----|-------|----------|-------|----|----|---------|---------|-------|---|---|
| 2 | C | 7 | Total | С | N | О | Р | 0 | 0 | 0 |
| 3 | | 1 | 144 | 68 | 25 | 44 | 7 | | U | U |

• Molecule 4 is a DNA chain called DNA (5'-D(P*CP*TP*GP*AP*TP*GP*T)-3').

| Mol | Chain | Residues | Atoms | | | ZeroOcc | AltConf | Trace | | |
|-----|-------|----------|--------------|---------|---------|---------|---------|-------|---|---|
| 4 | D | 7 | Total 145 | C 69 | N 24 | O 45 | P 7 | 0 | 0 | 0 |

• Molecule 5 is SILVER ION (three-letter code: AG) (formula: Ag) (labeled as "Ligand of Interest" by depositor).

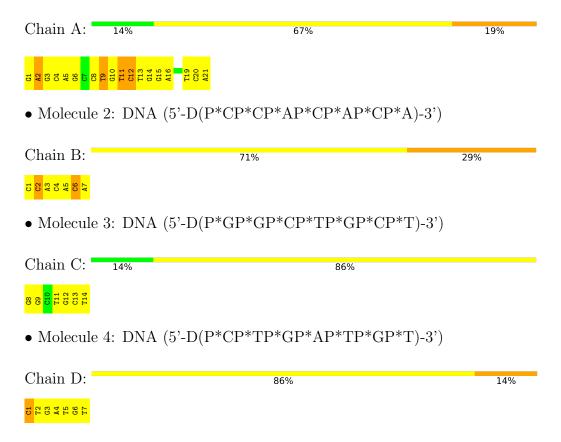
| Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|--------------|---------|---------|
| 5 | В | 2 | Total Ag 2 2 | 0 | 0 |



3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: DNA (5'-D(*GP*AP*GP*CP*AP*GP*CP*CP*TP*GP*TP*CP*TP*GP*AP*AP*CP*AP*TP*CP*A)-3')





4 Data and refinement statistics (i)

| Property | Value | Source |
|--|--|-----------|
| Space group | H 3 | Depositor |
| Cell constants | 105.98Å 105.98Å 90.26Å | Donositon |
| a, b, c, α , β , γ | 90.00° 90.00° 120.00° | Depositor |
| Resolution (Å) | 32.38 - 3.63 | Depositor |
| Resolution (A) | 64.36 - 3.63 | EDS |
| % Data completeness | 60.6 (32.38-3.63) | Depositor |
| (in resolution range) | 60.7 (64.36-3.63) | EDS |
| R_{merge} | (Not available) | Depositor |
| R_{sym} | (Not available) | Depositor |
| $< I/\sigma(I) > 1$ | 0.96 (at 3.67Å) | Xtriage |
| Refinement program | PHENIX 1.18.2_3874 | Depositor |
| D D | 0.188 , 0.207 | Depositor |
| R, R_{free} | 0.187 , 0.206 | DCC |
| R_{free} test set | 127 reflections (4.87%) | wwPDB-VP |
| Wilson B-factor (Å ²) | 281.3 | Xtriage |
| Anisotropy | 0.042 | Xtriage |
| Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$ | 0.21, 209.3 | EDS |
| L-test for twinning ² | $< L > = 0.47, < L^2> = 0.30$ | Xtriage |
| Estimated twinning fraction | 0.090 for h,-h-k,-l | Xtriage |
| F_o, F_c correlation | 0.77 | EDS |
| Total number of atoms | 858 | wwPDB-VP |
| Average B, all atoms (Å ²) | 291.0 | wwPDB-VP |

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.75% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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: AG

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 | Bon | d lengths | Bond angles | | |
|-------|-------|------|--------------|-------------|-------------------|--|
| IVIOI | | RMSZ | # Z > 5 | RMSZ | # Z >5 | |
| 1 | A | 1.21 | 4/480 (0.8%) | 1.25 | $4/739 \ (0.5\%)$ | |
| 2 | В | 1.02 | 0/155 | 1.23 | $2/235 \ (0.9\%)$ | |
| 3 | С | 0.75 | 0/160 | 1.03 | 0/245 | |
| 4 | D | 1.14 | 1/161 (0.6%) | 1.05 | 0/245 | |
| All | All | 1.11 | 5/956~(0.5%) | 1.18 | 6/1464 (0.4%) | |

All (5) bond length outliers are listed below:

| Mol | Chain | Res | Type | Atoms | \mathbf{Z} | Observed(A) | $\operatorname{Ideal}(ext{\AA})$ |
|-----|-------|-----|------|---------|--------------|-------------|-----------------------------------|
| 4 | D | 1 | DC | OP3-P | -10.22 | 1.48 | 1.61 |
| 1 | A | 11 | DT | C3'-O3' | 8.00 | 1.54 | 1.44 |
| 1 | A | 11 | DT | C1'-N1 | 5.86 | 1.56 | 1.49 |
| 1 | A | 8 | DC | C1'-N1 | 5.26 | 1.56 | 1.49 |
| 1 | A | 9 | DT | C1'-N1 | 5.13 | 1.55 | 1.49 |

All (6) bond angle outliers are listed below:

| Mol | Chain | Res | Type | Atoms | \mathbf{Z} | $Observed(^o)$ | $\mathrm{Ideal}(^{o})$ |
|-----|-------|-----|------|-------------|--------------|----------------|------------------------|
| 1 | A | 2 | DA | O4'-C1'-N9 | 6.15 | 112.31 | 108.00 |
| 1 | A | 12 | DC | O5'-P-OP2 | -6.06 | 100.25 | 105.70 |
| 1 | A | 12 | DC | O4'-C1'-N1 | 5.77 | 112.04 | 108.00 |
| 1 | A | 6 | DG | O4'-C1'-N9 | 5.47 | 111.83 | 108.00 |
| 2 | В | 2 | DC | C4'-C3'-C2' | -5.11 | 98.50 | 103.10 |
| 2 | В | 6 | DC | C3'-C2'-C1' | -5.05 | 96.44 | 102.50 |

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 | 428 | 0 | 236 | 22 | 0 |
| 2 | В | 139 | 0 | 78 | 9 | 0 |
| 3 | С | 144 | 0 | 80 | 9 | 0 |
| 4 | D | 145 | 0 | 81 | 5 | 0 |
| 5 | В | 2 | 0 | 0 | 0 | 0 |
| All | All | 858 | 0 | 475 | 41 | 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 31.

All (41) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

| Atom-1 | Atom-2 | $\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ (\rm \mathring{A}) \end{array}$ | Clash overlap (Å) |
|---------------|---------------|---|----------------------|
| 1:A:9:DT:H3' | 1:A:10:DG:C8 | 2.14 | 0.81 |
| 1:A:10:DG:H2' | 1:A:11:DT:C6 | 2.17 | 0.80 |
| 1:A:4:DC:H2" | 1:A:5:DA:H5' | 1.71 | 0.73 |
| 1:A:1:DG:H2' | 1:A:2:DA:C8 | 2.28 | 0.68 |
| 4:D:2:DT:H2' | 4:D:3:DG:C8 | 2.33 | 0.62 |
| 1:A:4:DC:H2' | 1:A:5:DA:C8 | 2.35 | 0.61 |
| 3:C:8:DG:H3' | 3:C:9:DG:C8 | 2.36 | 0.60 |
| 1:A:11:DT:H1' | 1:A:12:DC:OP2 | 2.02 | 0.60 |
| 3:C:8:DG:H3' | 3:C:9:DG:H8 | 1.68 | 0.58 |
| 1:A:14:DG:H2" | 1:A:15:DG:C8 | 2.39 | 0.57 |
| 4:D:1:DC:C6 | 4:D:2:DT:H72 | 2.40 | 0.57 |
| 1:A:3:DG:H2" | 1:A:4:DC:O5' | 2.06 | 0.55 |
| 2:B:5:DA:H2" | 2:B:6:DC:C6 | 2.41 | 0.55 |
| 1:A:13:DT:H1' | 1:A:14:DG:C5 | 2.42 | 0.54 |
| 3:C:13:DC:H2" | 3:C:14:DT:H71 | 1.89 | 0.53 |
| 1:A:19:DT:H2" | 1:A:20:DC:O5' | 2.08 | 0.53 |
| 2:B:7:DA:N9 | 3:C:8:DG:H1' | 2.25 | 0.52 |
| 1:A:13:DT:H4' | 1:A:14:DG:O5' | 2.11 | 0.51 |
| 1:A:20:DC:H2' | 1:A:21:DA:C8 | 2.48 | 0.49 |
| 1:A:2:DA:H1' | 1:A:3:DG:OP2 | 2.13 | 0.48 |
| 2:B:5:DA:H2" | 2:B:6:DC:C5 | 2.49 | 0.48 |
| 3:C:11:DT:H2' | 3:C:12:DG:C8 | 2.48 | 0.47 |

Continued on next page...



| $\alpha \cdots \alpha$ | | |
|------------------------|---------------|------|
| Continued fr | om $previous$ | paae |

| Atom-1 | Atom-2 | $\begin{array}{c} \text{Interatomic} \\ \text{distance } (\text{\AA}) \end{array}$ | Clash overlap (Å) |
|---------------|---------------|--|----------------------|
| 2:B:7:DA:C8 | 3:C:8:DG:H1' | 2.50 | 0.47 |
| 1:A:2:DA:H2" | 1:A:3:DG:OP1 | 2.15 | 0.47 |
| 2:B:2:DC:H2" | 2:B:3:DA:C5 | 2.50 | 0.46 |
| 1:A:13:DT:H3' | 1:A:13:DT:OP2 | 2.15 | 0.45 |
| 4:D:4:DA:C8 | 4:D:5:DT:H72 | 2.52 | 0.45 |
| 3:C:8:DG:C3' | 3:C:9:DG:H8 | 2.28 | 0.44 |
| 1:A:9:DT:H3' | 1:A:10:DG:H8 | 1.76 | 0.43 |
| 1:A:12:DC:P | 1:A:12:DC:O4' | 2.77 | 0.43 |
| 2:B:1:DC:C4 | 4:D:7:DT:C4 | 3.06 | 0.43 |
| 2:B:2:DC:H2" | 2:B:3:DA:N7 | 2.33 | 0.43 |
| 1:A:15:DG:H4' | 1:A:16:DA:OP1 | 2.19 | 0.43 |
| 3:C:8:DG:H2' | 3:C:9:DG:C8 | 2.54 | 0.43 |
| 2:B:4:DC:H2" | 2:B:5:DA:C8 | 2.53 | 0.43 |
| 1:A:10:DG:H2' | 1:A:11:DT:C5 | 2.53 | 0.43 |
| 1:A:1:DG:H2" | 1:A:2:DA:O5' | 2.19 | 0.42 |
| 4:D:5:DT:H2" | 4:D:6:DG:C8 | 2.55 | 0.42 |
| 1:A:4:DC:H2' | 1:A:5:DA:H8 | 1.78 | 0.41 |
| 1:A:11:DT:H2' | 1:A:11:DT:OP2 | 2.20 | 0.41 |
| 2:B:7:DA:C1' | 3:C:8:DG:H1' | 2.52 | 0.40 |

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)

There are no RNA molecules in this entry.

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

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



5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands (i)

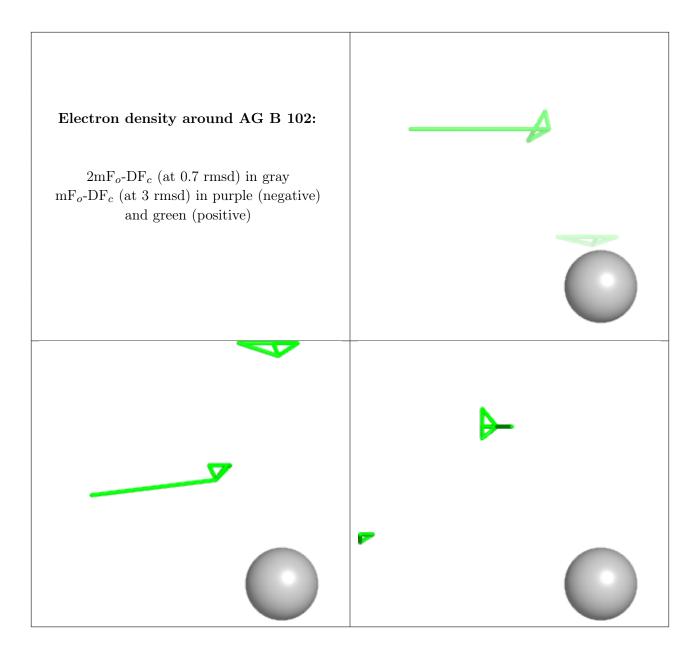
Unable to reproduce the depositors R factor - this section is therefore empty.

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around AG B 101: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)





6.5 Other polymers (i)

Unable to reproduce the depositors R factor - this section is therefore empty.

