



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 18, 2021 – 11:07 AM GMT

PDB ID : 6QJS
Title : Crystal structure of a DNA dodecamer containing a tetramethylpiperidinoxyl (nitroxide) spin label
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Deposited on : 2019-01-24
Resolution : 1.80 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.17.1.dev1
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) : 2.17.1.dev1

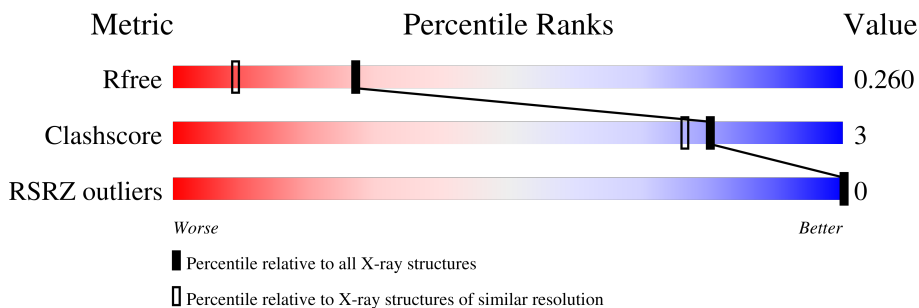
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 $\leq 5\%$. 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	12	 67% 17% 8% 8%
1	B	12	 67% 25% 8%

2 Entry composition [i](#)

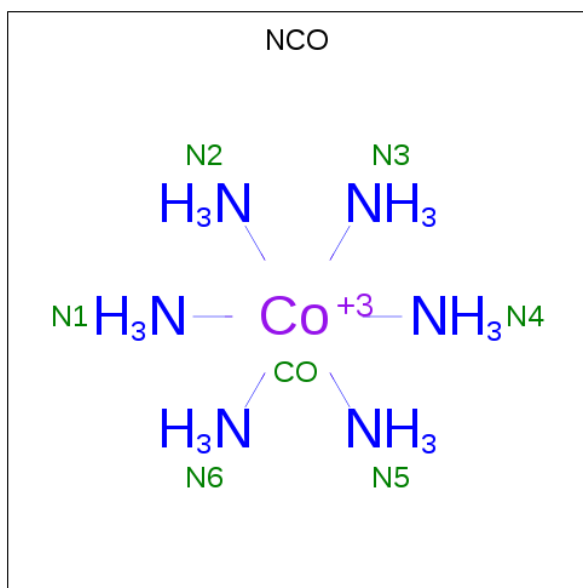
There are 3 unique types of molecules in this entry. The entry contains 519 atoms, of which 0 are hydrogens and 0 are deuteriums.

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(P*GP*CP*AP*AP*AP*TP*TP*(S6M)P*GP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	A	11	Total 242	C 118	N 46	O 67	P 11	0	0	0
1	B	11	Total 242	C 118	N 46	O 67	P 11	0	0	0

- Molecule 2 is COBALT HEXAMMINE(III) (three-letter code: NCO) (formula: CoH₁₈N₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Co	N		
2	A	1	Total 7	Co 1	N 6	0	0
2	A	1	Total 7	Co 1	N 6	0	0
2	A	1	Total 7	Co 1	N 6	0	0
2	B	1	Total 7	Co 1	N 6	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	4	Total O 4 4	0	0
3	B	3	Total O 3 3	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 and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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(P*GP*CP*AP*AP*AP*TP*TP*(S6M)P*GP*CP*G)-3')

Chain A: 



- Molecule 1: DNA (5'-D(P*GP*CP*AP*AP*AP*TP*TP*(S6M)P*GP*CP*G)-3')

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	43.00 Å 43.00 Å 97.49 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.24 – 1.80 37.24 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.6 (37.24-1.80) 96.6 (37.24-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 1.79 Å)	Xtrriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.213 , 0.251 0.218 , 0.260	Depositor DCC
R_{free} test set	555 reflections (5.59%)	wwPDB-VP
Wilson B-factor (Å ²)	29.2	Xtrriage
Anisotropy	0.535	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 13.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	0.476 for -h,-k,l	Xtrriage
Reported twinning fraction	0.481 for H, K, L 0.519 for -h,-k,l	Depositor
Outliers	3 of 9921 reflections (0.030%)	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	519	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.58% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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: NCO, S6M

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	0/231	0.69	0/352
1	B	0.86	1/231 (0.4%)	0.73	0/352
All	All	0.80	1/462 (0.2%)	0.71	0/704

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	17	DA	O3'-P	-5.56	1.54	1.61

There are no bond angle outliers.

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	242	0	113	2	0
1	B	242	0	113	1	0
2	A	21	0	0	0	0
2	B	7	0	0	0	0
3	A	4	0	0	0	0
3	B	3	0	0	0	0
All	All	519	0	226	2	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 3.

All (2) 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:4:DA:H2''	1:A:5:DA:OP2	1.83	0.78
1:A:9:S6M:N3	1:B:16:DA:N1	2.66	0.44

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](#)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	S6M	B	21	1	30,38,39	1.73	6 (20%)	33,59,62	2.59	4 (12%)
1	S6M	A	9	1	30,38,39	2.18	9 (30%)	33,59,62	2.97	5 (15%)

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
1	S6M	B	21	1	-	0/5/55/56	0/4/4/4
1	S6M	A	9	1	-	0/5/55/56	0/4/4/4

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	9	S6M	O4'-C1'	5.82	1.49	1.41
1	A	9	S6M	N15-N14	5.08	1.43	1.34
1	B	21	S6M	C12-N13	-4.86	1.30	1.35
1	A	9	S6M	C6-N1	-4.38	1.30	1.35
1	B	21	S6M	C2-N3	-4.14	1.30	1.38
1	A	9	S6M	C21-N20	3.45	1.53	1.48
1	A	9	S6M	C19-N20	3.20	1.53	1.48
1	B	21	S6M	N15-N14	3.07	1.39	1.34
1	A	9	S6M	C2-N3	-2.64	1.32	1.38
1	B	21	S6M	O4'-C1'	2.46	1.44	1.41
1	B	21	S6M	C6-N1	-2.45	1.32	1.35
1	A	9	S6M	N14-N13	2.24	1.39	1.34
1	A	9	S6M	C12-C11	2.24	1.39	1.36
1	B	21	S6M	C2'-C3'	-2.23	1.51	1.55
1	A	9	S6M	C17-N13	2.07	1.52	1.49

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	S6M	C12-N13-C17	11.98	136.09	125.48
1	B	21	S6M	C12-C11-C2'	-9.60	114.48	129.25
1	B	21	S6M	C2'-C11-N15	7.69	132.81	120.30
1	B	21	S6M	C12-N13-C17	6.78	131.49	125.48
1	A	9	S6M	C12-C11-C2'	-6.47	119.30	129.25
1	A	9	S6M	C22-C21-N20	5.88	115.94	108.69
1	A	9	S6M	C2'-C11-N15	5.81	129.75	120.30
1	A	9	S6M	C18-C19-N20	4.91	114.73	108.69
1	B	21	S6M	C22-C21-N20	3.09	112.50	108.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	9	S6M	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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	$\# Z > 2$
2	NCO	A	103	-	6,6,6	0.83	0	-		
2	NCO	A	102	-	6,6,6	0.70	0	-		
2	NCO	B	101	2	6,6,6	0.44	0	-		
2	NCO	A	101	2	6,6,6	0.60	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.

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

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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	10/12 (83%)	-0.76	0 100 100	26, 34, 42, 51	0
1	B	10/12 (83%)	-0.80	0 100 100	28, 35, 43, 44	0
All	All	20/24 (83%)	-0.78	0 100 100	26, 35, 44, 51	0

There are no RSRZ outliers to report.

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

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. 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	B-factors(Å ²)	Q<0.9
1	S6M	A	9	35/36	0.95	0.10	23,29,38,41	0
1	S6M	B	21	35/36	0.96	0.09	28,31,53,60	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

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. 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	B-factors(\AA^2)	Q<0.9
2	NCO	A	101	7/7	0.99	0.09	29,30,32,32	7
2	NCO	A	102	7/7	0.99	0.15	15,15,17,18	7
2	NCO	A	103	7/7	0.99	0.07	32,38,40,43	0
2	NCO	B	101	7/7	0.99	0.10	26,29,30,31	7

6.5 Other polymers [i](#)

There are no such residues in this entry.