

wwPDB X-ray Structure Validation Summary Report (i)

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PDB ID	:	268D
Title	:	STRUCTURAL STUDIES ON NUCLEIC ACIDS
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Deposited on		
Resolution	:	2.02 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 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:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-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)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\;DIFFRACTION$

The reported resolution of this entry is 2.02 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	Percentile Ranks	Value
Clashscore		44
RSRZ outliers		0
	Worse	Better
	Percentile relative to all X-ray structures	
	Percentile relative to X-ray structures of similar resolution	

Metric	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
Clashscore	141614	11643 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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% 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	А	12		42%		50%	8%
1	В	12	17%		83%	6	

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
1	5CM	А	9	-	-	Х	-
2	BRN	А	25	-	-	Х	-



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2 Entry composition (i)

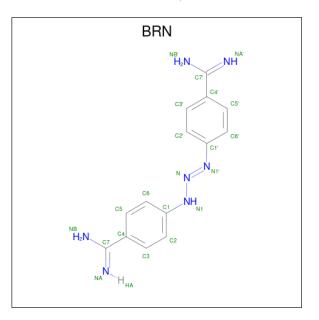
There are 4 unique types of molecules in this entry. The entry contains 621 atoms, of which 0 are hydrogens and 0 are deuteriums.

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(*CP*GP*CP*GP*AP*AP*TP*TP*(5CM)P* GP*CP*G)-3').

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	19	Total	С	Ν	Ο	Р	0	0	0
		12	244	117	46	70	11	0	0	0
1	В	12	Total	С	Ν	Ο	Р	0	0	0
	D	12	244	117	46	70	11	0		0

• Molecule 2 is BERENIL (three-letter code: BRN) (formula: $C_{14}H_{15}N_7$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	А	1	Total 21	C 14	N 7	0	0

• Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).



]	Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
	3	А	1	Total 1	Mg 1	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	62	$\begin{array}{cc} \text{Total} & \text{O} \\ 62 & 62 \end{array}$	0	0
4	В	49	Total O 49 49	0	0



C13 G14 G14 C15 C15 C15 A17 A18 A17 T19 T19 T19 C21 C21 G22 C23 C23 C23

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(*CP*GP*CP*GP*AP*AP*TP*TP*(5CM)P*GP*CP*G)-3')

Chain A:	42%	50%	8%	
C1 C2 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3 C3	69 611 611 611 611 611			
• Molecule	e 1: DNA (5'-D(*CP*GP*C	CP*GP*AP*AP*TP*TP*(5	6CM)P*GP*CP*G)-3')
Chain B:	17%	83%		



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	24.74Å 40.19Å 65.32Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) - 2.02	Depositor
Resolution (A)	40.19 - 2.02	EDS
% Data completeness	90.5 ((Not available)- 2.02)	Depositor
(in resolution range)	98.9 (40.19-2.02)	EDS
R _{merge}	0.06	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$2.21 (at 2.01 \text{\AA})$	Xtriage
Refinement program	NUCLSQ	Depositor
D D.	0.160 , (Not available)	Depositor
R, R_{free}	0.172 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor $(Å^2)$	30.9	Xtriage
Anisotropy	0.288	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32, 134.8	EDS
L-test for twinning ²	$ \langle L \rangle = 0.47, \langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	621	wwPDB-VP
Average B, all atoms $(Å^2)$	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 13.01% of the height of the origin peak. No significant pseudotranslation is detected.

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.



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: MG, BRN, $5\mathrm{CM}$

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	Bo	nd lengths	Bond angles		
MOI C	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	3.72	44/250~(17.6%)	4.70	82/382~(21.5%)	
1	В	3.62	44/250~(17.6%)	5.02	107/382~(28.0%)	
All	All	3.67	88/500~(17.6%)	4.86	189/764~(24.7%)	

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	А	0	1

The worst 5 of 88 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	В	17	DA	C6-N6	12.01	1.43	1.33
1	А	7	DT	C4'-O4'	-11.68	1.33	1.45
1	В	18	DA	N3-C4	11.31	1.41	1.34
1	А	6	DA	N9-C8	-10.65	1.29	1.37
1	В	18	DA	N7-C5	10.29	1.45	1.39

The worst 5 of 189 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	В	18	DA	N1-C2-N3	-25.29	116.65	129.30
1	А	1	DC	P-O3'-C3'	19.63	143.26	119.70
1	А	2	DG	P-O3'-C3'	18.15	141.47	119.70
1	В	16	DG	P-O3'-C3'	17.00	140.09	119.70
1	В	18	DA	C6-N1-C2	16.48	128.49	118.60



There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group	
1	А	2	DG	Sidechain	

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	А	244	0	134	14	0
1	В	244	0	138	19	1
2	А	21	0	15	13	0
3	А	1	0	0	0	0
4	А	62	0	0	8	1
4	В	49	0	0	3	0
All	All	621	0	287	34	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 44.

The worst 5 of 34 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)
2:A:25:BRN:H5	1:B:18:DA:C2	1.94	1.02
2:A:25:BRN:H5	1:B:18:DA:N3	1.77	0.99
1:A:9:5CM:H5'	2:A:25:BRN:C3	2.10	0.79
1:B:17:DA:O5'	1:B:17:DA:H2'	1.83	0.78
1:A:9:5CM:H5'	2:A:25:BRN:H3	1.67	0.77

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:DC:OP1	4:A:55:HOH:O[2_664]	2.16	0.04



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).

Mal	Turne	Chain	Res	Tinle	Bo	Bond lengths			Bond angles		
IVIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2	
1	5CM	А	9	1	17,21,22	1.30	2 (11%)	24,30,33	2.05	6 (25%)	
1	5CM	В	21	1	17,21,22	1.36	3 (17%)	24,30,33	2.38	8 (33%)	

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	5CM	А	9	1	-	3/7/21/22	0/2/2/2
1	5CM	В	21	1	-	0/7/21/22	0/2/2/2

All (5) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
1	В	21	5CM	C6-C5	3.12	1.39	1.34
1	В	21	5CM	O4'-C4'	-2.77	1.38	1.45
1	А	9	5CM	O4'-C1'	2.67	1.48	1.42
1	В	21	5CM	C2-N3	2.15	1.40	1.36
1	А	9	5CM	C6-N1	2.01	1.41	1.38

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

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	В	21	5CM	C5-C6-N1	-5.78	117.39	123.34
1	А	9	5CM	C5-C6-N1	-5.22	117.97	123.34
1	В	21	5CM	C5A-C5-C6	-5.12	116.01	122.85
1	А	9	5CM	C5A-C5-C6	-4.90	116.30	122.85
1	В	21	5CM	C5-C4-N4	4.03	127.51	121.48

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	А	9	5CM	O4'-C4'-C5'-O5'
1	А	9	5CM	C3'-C4'-C5'-O5'
1	А	9	5CM	C4'-C5'-O5'-P

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	9	5CM	8	0
1	В	21	5CM	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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



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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).

Mo	Type	Chain	Dog	Link	Bond lengths			Bond angles		
Mol	Type		nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BRN	А	25	-	22,22,22	2.03	6 (27%)	21,29,29	1.12	2 (9%)

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.

Μ	ol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2		BRN	А	25	-	-	0/14/14/14	0/2/2/2

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
2	А	25	BRN	C4-C7	-5.19	1.37	1.47
2	А	25	BRN	C4'-C7'	-5.13	1.38	1.47
2	А	25	BRN	C7-NA	2.96	1.40	1.28
2	А	25	BRN	N1-N	2.75	1.39	1.32
2	А	25	BRN	C7'-NA'	2.31	1.38	1.28

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	А	25	BRN	C4-C7-NB	2.86	122.36	118.05
2	А	25	BRN	C4'-C7'-NB'	2.31	121.53	118.05

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

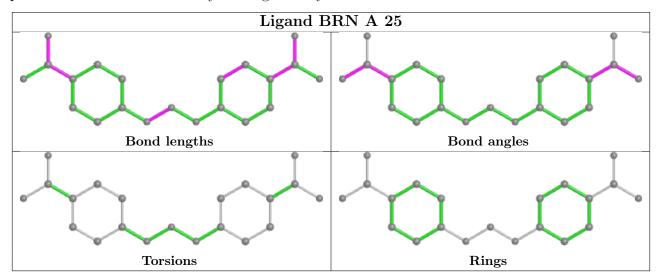
1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes	
2	А	25	BRN	13	0	

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will



also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and similar rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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)

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, 95^{th} 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	$\langle RSRZ \rangle$	#	#RSRZ>2		$\mathbf{OWAB}(\mathbf{A}^2)$	Q<0.9
1	А	11/12~(91%)	-0.84	0	100	100	19, 24, 34, 35	0
1	В	11/12 (91%)	-0.86	0	100	100	23, 26, 32, 33	0
All	All	22/24~(91%)	-0.85	0	100	100	19, 25, 33, 35	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, 95^{th} 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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
1	5CM	А	9	20/21	0.92	0.12	16,34,60,60	0
1	5CM	В	21	20/21	0.96	0.09	11,25,37,47	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

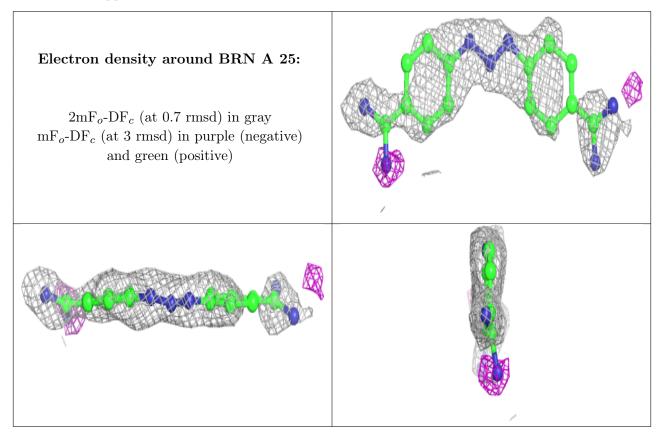
6.4 Ligands (i)

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, 95^{th} 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
2	BRN	А	25	21/21	0.84	0.28	$36,\!53,\!65,\!92$	0
3	MG	А	26	1/1	0.97	0.06	23,23,23,23	0

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.



6.5 Other polymers (i)

There are no such residues in this entry.

