



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 8QN2 / pdb_00008qn2
Title : Helical foldamers as selective G-quadruplex ligands
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Deposited on : 2023-09-25
Resolution : 2.51 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 types of validation reports are described at

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

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) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

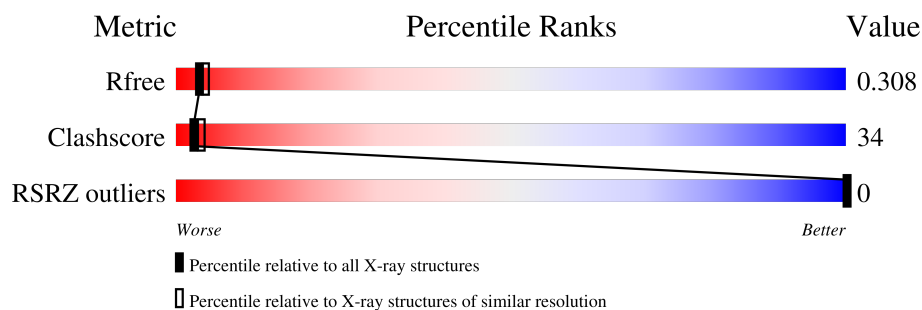
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	20	
2	B	5	
2	C	5	

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
2	QUK	B	108	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 837 atoms, of which 306 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(*TP*GP*GP*GP*TP*TP*GP*GP*GP*TP*TP*GP*GP*GP*TP*TP*GP*GP*GP*TP*TP*GP*GP*GP*T)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	20	Total	C	H	N	O	P	0	0	1
			592	180	204	72	117	19			

- Molecule 2 is a protein called ACE-QUK-QUK-ZY9-QUK.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	5	Total	C	H	N	O	0	0	0
			119	48	51	11	9			
2	C	5	Total	C	H	N	O	0	0	0
			119	48	51	11	9			

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	K	0	0
			3	3		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mg	0	0
			2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	O	0	0
			2	2		

3 Residue-property plots

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(*TP*GP*GP*GP*TP*TP*GP*GP*GP*TP*TP*GP*GP*GP*TP*TP*GP*GP*GP*T)-3')

Chain A: 

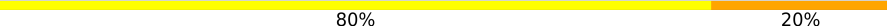


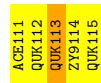
- Molecule 2: ACE-QUK-QUK-ZY9-QUK

Chain B: 



- Molecule 2: ACE-QUK-QUK-ZY9-QUK

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	32.71Å 32.71Å 61.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.50 – 2.51 30.50 – 2.51	Depositor EDS
% Data completeness (in resolution range)	99.1 (30.50-2.51) 100.0 (30.50-2.51)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.29 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.21rc1_5084	Depositor
R, R_{free}	0.272 , 0.308 0.272 , 0.308	Depositor DCC
R_{free} test set	224 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	59.9	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.479 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	837	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.48% 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: MG, ZY9, ACE, QUK, K

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	0.84	0/435	0.93	0/675

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
2	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	108	QUK	Peptide

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	388	204	204	13	0
2	B	68	51	3	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	68	51	3	1	0
3	A	3	0	0	0	0
4	A	2	0	0	0	0
5	A	2	0	0	0	0
All	All	531	306	210	24	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 34.

All (24) 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:B:106:ACE:H2	2:B:108:QUK:C	1.87	1.05
2:B:109:ZY9:N	2:B:110:QUK:OXT	2.10	0.82
2:B:108:QUK:C	2:B:110:QUK:OXT	2.33	0.76
2:B:106:ACE:CH3	2:B:108:QUK:C	2.64	0.75
2:B:106:ACE:H2	2:B:109:ZY9:N	2.08	0.67
1:A:9:DG:H4'	1:A:11:DT:C6	2.33	0.63
1:A:9:DG:C8	2:B:106:ACE:O	2.52	0.63
2:B:106:ACE:H2	2:B:109:ZY9:C2	2.29	0.62
1:A:14:DG:C4	2:B:108:QUK:C4	2.83	0.61
2:B:106:ACE:H2	2:B:108:QUK:O	2.01	0.60
1:A:9:DG:N7	2:B:106:ACE:O	2.35	0.60
2:B:108:QUK:C10	2:B:110:QUK:OXT	2.51	0.59
1:A:9:DG:C5	2:B:106:ACE:O	2.63	0.51
1:A:9:DG:C5	2:B:106:ACE:C	2.88	0.51
2:B:108:QUK:C10	2:B:110:QUK:C	2.92	0.48
2:C:111:ACE:H2	2:C:113:QUK:C	2.43	0.47
1:A:11:DT:H2''	1:A:12:DG:O5'	2.16	0.46
1:A:14:DG:H4'	1:A:16:DT:C6	2.52	0.45
1:A:14:DG:C5	2:B:108:QUK:C3	3.02	0.43
1:A:14:DG:C8	2:B:108:QUK:C5	3.02	0.43
2:B:106:ACE:CH3	2:B:108:QUK:O	2.65	0.42
1:A:10:DT:C5'	1:A:11:DT:OP2	2.68	0.41
1:A:16:DT:O2	1:A:17:DG:H2'	2.21	0.40
1:A:14:DG:C4	2:B:108:QUK:C3	3.05	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report 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](#)

8 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	$\# Z > 2$
2	QUK	B	108	2	19,19,20	0.98	1 (5%)	24,25,27	2.07	4 (16%)
2	QUK	C	113	2	19,19,20	0.64	0	24,25,27	1.40	3 (12%)
2	ZY9	B	109	2	10,10,11	0.78	1 (10%)	11,12,14	1.09	0
2	QUK	C	112	2	19,19,20	0.40	0	24,25,27	0.85	1 (4%)
2	QUK	C	115	2	20,20,20	0.45	0	27,27,27	0.95	2 (7%)
2	ZY9	C	114	2	10,10,11	0.67	0	11,12,14	0.87	1 (9%)
2	QUK	B	107	2	19,19,20	0.47	0	24,25,27	0.83	1 (4%)
2	QUK	B	110	2	20,20,20	0.63	0	27,27,27	0.96	2 (7%)

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
2	QUK	B	108	2	-	6/7/7/9	0/2/2/2
2	QUK	C	113	2	-	2/7/7/9	0/2/2/2
2	ZY9	B	109	2	-	0/4/4/6	0/1/1/1
2	QUK	C	112	2	-	4/7/7/9	0/2/2/2
2	QUK	C	115	2	-	1/9/9/9	0/2/2/2
2	ZY9	C	114	2	-	0/4/4/6	0/1/1/1
2	QUK	B	107	2	-	3/7/7/9	0/2/2/2
2	QUK	B	110	2	-	2/9/9/9	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	108	QUK	C10-C	3.22	1.52	1.48
2	B	109	ZY9	CA-C	2.06	1.50	1.48

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	108	QUK	O-C-C10	-6.50	118.22	124.22
2	B	108	QUK	C-C10-N11	5.60	120.15	114.66
2	B	108	QUK	C9-C10-C	-4.13	113.70	121.11
2	C	113	QUK	O-C-C10	-3.48	121.01	124.22
2	B	110	QUK	C3-CA-N	3.00	126.35	120.39
2	C	113	QUK	C-C10-N11	2.99	117.60	114.66
2	C	113	QUK	CE-CD-CG	2.96	123.73	113.53
2	C	115	QUK	C3-CA-N	2.70	125.74	120.39
2	C	112	QUK	O-C-C10	-2.57	121.85	124.22
2	C	115	QUK	O-C-C10	-2.52	116.23	121.30
2	C	114	ZY9	O-C-CA	-2.36	122.05	124.22
2	B	110	QUK	C7-CA-N	-2.25	113.98	118.16
2	B	107	QUK	O-C-C10	-2.13	122.25	124.22
2	B	108	QUK	C9-C10-N11	2.01	125.49	123.18

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	108	QUK	O-C-C10-N11
2	B	108	QUK	O-C-C10-C9
2	C	113	QUK	O-C-C10-C9
2	C	115	QUK	CE-CD-CG-OB

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Mol	Chain	Res	Type	Atoms
2	C	112	QUK	C6-C8-OB-CG
2	B	110	QUK	CE-CD-CG-OB
2	C	112	QUK	C9-C8-OB-CG
2	C	112	QUK	CE-CD-CG-OB
2	B	108	QUK	C6-C8-OB-CG
2	B	107	QUK	CD-CG-OB-C8
2	B	108	QUK	C9-C8-OB-CG
2	C	113	QUK	CE-CD-CG-OB
2	B	108	QUK	CD-CG-OB-C8
2	C	112	QUK	CG-CD-CE-N1
2	B	108	QUK	CG-CD-CE-N1
2	B	107	QUK	CE-CD-CG-OB
2	B	107	QUK	CG-CD-CE-N1
2	B	110	QUK	CG-CD-CE-N1

There are no ring outliers.

4 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	108	QUK	11	0
2	C	113	QUK	1	0
2	B	109	ZY9	3	0
2	B	110	QUK	4	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 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

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	20/20 (100%)	-1.09	0 100 100	51, 64, 98, 108	0
2	B	0/5	-	-	-	-
2	C	0/5	-	-	-	-
All	All	20/30 (66%)	-1.09	0 100 100	51, 64, 98, 108	0

There are no RSRZ outliers to report.

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

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
2	QUK	C	112	18/19	0.94	0.08	61,100,161,161	0
2	ZY9	B	109	10/11	0.96	0.08	60,74,95,101	0
2	QUK	B	108	18/19	0.97	0.09	61,94,133,141	0
2	QUK	C	113	18/19	0.97	0.07	53,87,126,134	0
2	QUK	B	107	18/19	0.97	0.09	48,77,169,169	0
2	QUK	B	110	19/19	0.97	0.08	63,81,124,124	0
2	ZY9	C	114	10/11	0.98	0.07	67,83,102,108	0
2	QUK	C	115	19/19	0.98	0.08	54,82,118,118	0

6.3 Carbohydrates

There are no oligosaccharides 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(Å ²)	Q<0.9
4	MG	A	104	1/1	0.96	0.17	79,79,79,79	0
3	K	A	101	1/1	0.98	0.02	42,42,42,42	0
4	MG	A	105	1/1	0.98	0.16	78,78,78,78	0
3	K	A	102	1/1	0.99	0.03	41,41,41,41	0
3	K	A	103	1/1	0.99	0.01	28,28,28,28	1

6.5 Other polymers [i](#)

There are no such residues in this entry.