Full wwPDB X-ray Structure Validation Report ⓘ

May 2, 2024 – 12:32 PM EDT

PDB ID	:	8VF9
Title	:	Binary DNA Polymerase Beta bound to DNA containing primer terminal dA base-paired with FapydG
Authors	:	Oden, P.N.; Ryan, B.J.; Freudenthal, B.D.
Deposited on	:	2023-12-21
Resolution	:	1.90 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:

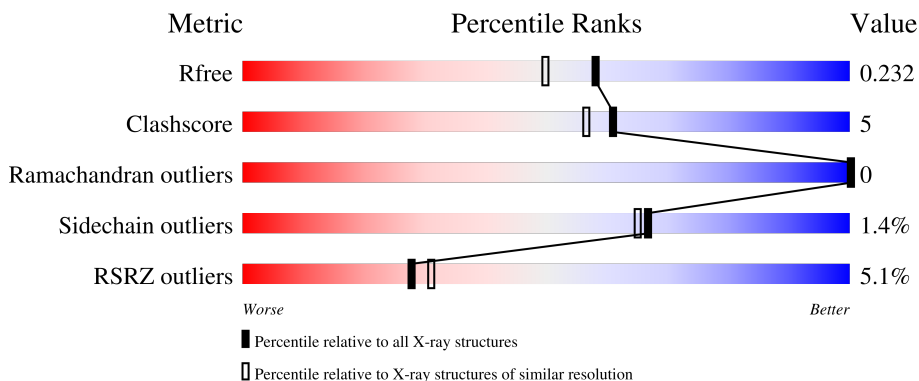|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.90 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
Rfree		0.232
Clashscore		5
Ramachandran outliers		0
Sidechain outliers		1.4%
RSRZ outliers		5.1%

Worse   *Better*

▮ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R$_{free}$	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	A	335	5% / 85% / 11% · ·
2	T	16	94% / 6%
3	P	10	70% / 30%
4	D	5	100%

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	8NI	T	7[A]	X	-	-	-
2	8NI	T	7[B]	X	-	-	-

2 Entry composition ⓘ

There are 6 unique types of molecules in this entry. The entry contains 3503 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 protein called DNA polymerase beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total 2586	C 1636	N 453	O 488	S 9	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*CP*GP*AP*CP*CP*(FAP)P*CP*GP* CP*AP*TP*CP*AP*GP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	16	Total 344	C 162	N 66	O 100	P 16	0	1	0

- Molecule 3 is a DNA chain called DNA (5'-D(*GP*CP*TP*GP*AP*TP*GP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	10	Total 205	C 98	N 40	O 58	P 9	0	0	0

- Molecule 4 is a DNA chain called DNA (5'-D(P*GP*TP*CP*GP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	5	Total 106	C 49	N 20	O 32	P 5	0	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total 3	Na 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	177	Total 177	O 177	0	0
6	T	41	Total 41	O 41	0	0
6	P	31	Total 31	O 31	0	0
6	D	10	Total 10	O 10	0	0

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 polymerase beta

Chain A:
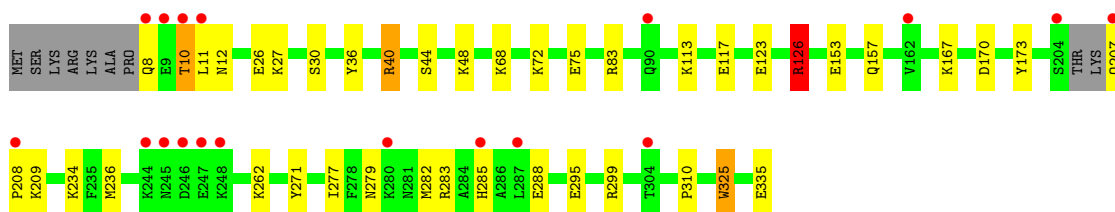
• Molecule 2: DNA (5'-D(*CP*CP*GP*AP*CP*CP*(FAP)P*CP*GP*CP*AP*TP*CP*AP*GP*C)-3')

Chain T:

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

Chain P:

• Molecule 4: DNA (5'-D(P*GP*TP*CP*GP*G)-3')

Chain D:

There are no outlier residues recorded for this chain.

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	52.44Å 78.63Å 54.52Å 90.00° 105.89° 90.00°	Depositor
Resolution (Å)	24.92 – 1.90 24.92 – 1.90	Depositor EDS
% Data completeness (in resolution range)	91.7 (24.92-1.90) 86.9 (24.92-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.44 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.208 , 0.232 0.208 , 0.232	Depositor DCC
R_{free} test set	1992 reflections (6.42%)	wwPDB-VP
Wilson B-factor (Å^2)	18.2	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 44.0	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.026 for l,-k,h	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3503	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.21% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical 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.

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: 8NI, NA

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	0.46	0/2634	0.68	0/3542
2	T	0.86	0/332	1.08	0/506
3	P	0.97	0/230	1.12	0/354
4	D	0.84	0/118	0.93	0/179
All	All	0.57	0/3314	0.78	0/4581

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	A	0	2
2	T	3	0
All	All	3	2

There are no bond length outliers.

There are no bond angle outliers.

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	T	7[A]	8NI	C5
2	T	7[B]	8NI	C5,C1'

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	126	ARG	Sidechain
1	A	40	ARG	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	A	2586	0	2582	25	0
2	T	344	0	168	2	0
3	P	205	0	114	3	0
4	D	106	0	57	0	0
5	A	3	0	0	0	0
6	A	177	0	0	4	0
6	D	10	0	0	0	0
6	P	31	0	0	0	0
6	T	41	0	0	0	0
All	All	3503	0	2921	28	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 5.

All (28) 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:167:LYS:HE2	6:A:551:HOH:O	1.59	1.01
1:A:271:TYR:CD2	1:A:295:GLU:HB2	2.31	0.65
2:T:7[A]:8NI:N1	3:P:10:DA:N1	2.46	0.63
1:A:10:THR:O	1:A:11:LEU:HB2	2.06	0.56
1:A:299:ARG:HG2	1:A:310:PRO:HA	1.91	0.53
1:A:36:TYR:CZ	1:A:40:ARG:HD3	2.50	0.47
1:A:234:LYS:HE2	1:A:236:MET:SD	2.55	0.46
2:T:7[B]:8NI:N1	3:P:10:DA:N1	2.64	0.46
1:A:277:ILE:HG22	1:A:335:GLU:HG3	1.97	0.45
1:A:27:LYS:HB2	1:A:36:TYR:CG	2.52	0.45
1:A:285:HIS:HA	1:A:288:GLU:HG2	1.99	0.45
1:A:170:ASP:HB3	1:A:173:TYR:CD2	2.53	0.44
1:A:279:ASN:O	1:A:283:ARG:HG3	2.18	0.44
3:P:1:DG:H2'	3:P:2:DC:C6	2.52	0.44
1:A:123:GLU:HA	1:A:126:ARG:NH1	2.33	0.43
1:A:27:LYS:HD2	6:A:662:HOH:O	2.18	0.43
1:A:113:LYS:O	1:A:117:GLU:HG3	2.19	0.43
1:A:234:LYS:HG2	6:A:547:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:SER:O	1:A:48:LYS:HE2	2.20	0.42
1:A:68:LYS:O	1:A:72:LYS:HD2	2.20	0.42
1:A:75:GLU:OE2	1:A:83:ARG:HG3	2.20	0.42
1:A:282:MET:HG3	1:A:325:TRP:CZ3	2.55	0.42
1:A:153:GLU:O	1:A:157:GLN:HG3	2.20	0.41
1:A:207:GLN:N	1:A:208:PRO:CD	2.83	0.41
1:A:208:PRO:O	1:A:209:LYS:HB2	2.19	0.41
1:A:208:PRO:HA	6:A:590:HOH:O	2.21	0.41
1:A:8:GLN:O	1:A:12:ASN:N	2.55	0.40
1:A:26:GLU:HA	1:A:30:SER:OG	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	322/335 (96%)	311 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	279/295 (95%)	275 (99%)	4 (1%)	67	65

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	THR
1	A	126	ARG
1	A	262	LYS
1	A	325	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	217	GLN

5.3.3   RNA ⓘ

There are no RNA molecules in this entry.

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

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	$\#|Z| > 2$
2	8NI	T	7[B]	2	18,24,25	1.94	6 (33%)	19,33,36	2.38	5 (26%)
2	8NI	T	7[A]	2	18,24,25	1.98	6 (33%)	19,33,36	1.80	6 (31%)

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	8NI	T	7[B]	2	2/2/8/11	4/7/40/41	0/1/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	8NI	T	7[A]	2	1/1/8/11	2/7/40/41	0/1/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	T	7[B]	8NI	C2-N2	3.87	1.45	1.34
2	T	7[A]	8NI	C2-N2	3.84	1.45	1.34
2	T	7[A]	8NI	C8-N7	3.61	1.45	1.33
2	T	7[B]	8NI	C8-N7	3.60	1.45	1.33
2	T	7[B]	8NI	C5-C6	-3.25	1.35	1.49
2	T	7[A]	8NI	C5-C6	-3.22	1.35	1.49
2	T	7[A]	8NI	C6-N1	-2.96	1.32	1.39
2	T	7[B]	8NI	C6-N1	-2.95	1.32	1.39
2	T	7[A]	8NI	O6-C6	-2.69	1.18	1.23
2	T	7[B]	8NI	O6-C6	-2.66	1.18	1.23
2	T	7[A]	8NI	O5'-C5'	-2.30	1.39	1.44
2	T	7[B]	8NI	O5'-C5'	-2.12	1.39	1.44

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	T	7[B]	8NI	O4'-C1'-N9	7.48	117.59	110.04
2	T	7[A]	8NI	N3-C2-N1	-4.00	119.94	126.43
2	T	7[B]	8NI	N3-C2-N1	-3.49	120.77	126.43
2	T	7[A]	8NI	C6-C5-N7	3.25	120.84	110.72
2	T	7[B]	8NI	C6-C5-N7	3.08	120.33	110.72
2	T	7[B]	8NI	C2'-C3'-C4'	2.49	107.95	102.76
2	T	7[B]	8NI	N2-C2-N3	2.32	120.34	116.57
2	T	7[A]	8NI	O4'-C1'-N9	2.18	112.24	110.04
2	T	7[A]	8NI	N2-C2-N1	2.08	120.33	117.06
2	T	7[A]	8NI	C5-N7-C8	-2.07	119.50	122.62
2	T	7[A]	8NI	C2'-C3'-C4'	2.01	106.94	102.76

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	T	7[A]	8NI	C5
2	T	7[B]	8NI	C5
2	T	7[B]	8NI	C1'

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	T	7[A]	8NI	C3'-C4'-C5'-O5'
2	T	7[A]	8NI	O4'-C4'-C5'-O5'
2	T	7[B]	8NI	O4'-C1'-N9-C4
2	T	7[B]	8NI	C3'-C4'-C5'-O5'
2	T	7[B]	8NI	O4'-C4'-C5'-O5'
2	T	7[B]	8NI	C5-C4-N9-C1'

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	T	7[B]	8NI	1	0
2	T	7[A]	8NI	1	0

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 3 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 ⓘ

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 ⓘ

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	326/335 (97%)	0.25	18 (5%)	25	28	7, 22, 46, 70	0
2	T	15/16 (93%)	-0.22	0	100	100	11, 26, 40, 47	0
3	P	10/10 (100%)	-0.29	0	100	100	12, 19, 27, 41	0
4	D	5/5 (100%)	-0.58	0	100	100	18, 21, 26, 28	0
All	All	356/366 (97%)	0.20	18 (5%)	28	31	7, 22, 46, 70	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	GLN	6.6
1	A	245	ASN	5.7
1	A	10	THR	5.3
1	A	9	GLU	4.9
1	A	208	PRO	4.4
1	A	287	LEU	3.8
1	A	11	LEU	3.6
1	A	207	GLN	3.0
1	A	248	LYS	2.9
1	A	246	ASP	2.8
1	A	244	LYS	2.8
1	A	280	LYS	2.6
1	A	247	GLU	2.5
1	A	162	VAL	2.5
1	A	90	GLN	2.5
1	A	304	THR	2.3
1	A	285	HIS	2.1
1	A	204	SER	2.0

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	8NI	T	7[A]	23/24	0.80	0.25	36,41,51,55	23
2	8NI	T	7[B]	23/24	0.80	0.25	37,41,50,54	23

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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
5	NA	A	402	1/1	0.97	0.04	20,20,20,20	0
5	NA	A	403	1/1	0.98	0.12	17,17,17,17	0
5	NA	A	401	1/1	0.99	0.07	6,6,6,6	0

6.5 Other polymers ⓘ

There are no such residues in this entry.