



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 06:42 PM GMT

PDB ID : 2VQZ
Title : STRUCTURE OF THE CAP-BINDING DOMAIN OF INFLUENZA VIRUS
POLYMERASE SUBUNIT PB2 WITH BOUND M7GTP
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Deposited on : 2008-03-21
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

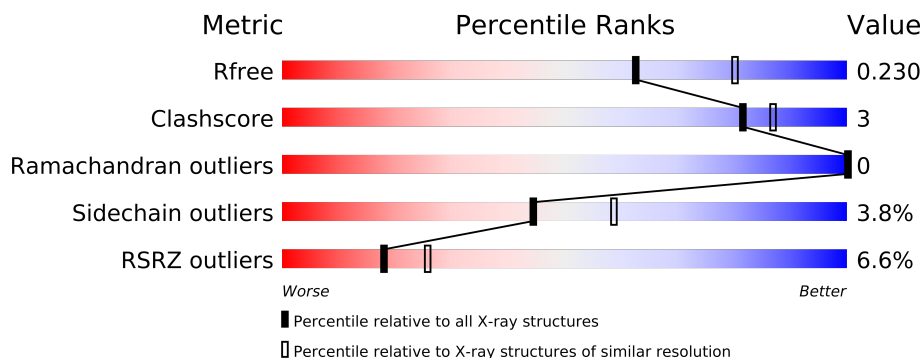
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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}	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	166	
1	B	166	
1	D	166	
1	E	166	
1	F	166	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6920 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

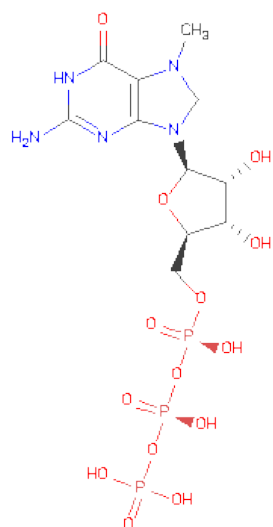
- Molecule 1 is a protein called POLYMERASE BASIC PROTEIN 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	Se	0	2	0
			1303	815	238	238	2	10			
1	B	162	Total	C	N	O	S	Se	0	2	0
			1294	811	235	236	2	10			
1	D	165	Total	C	N	O	S	Se	0	2	0
			1321	827	242	240	2	10			
1	E	160	Total	C	N	O	S	Se	0	1	0
			1270	797	229	232	2	10			
1	F	160	Total	C	N	O	S	Se	0	5	0
			1295	813	235	234	1	12			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	389	LYS	ARG	ENGINEERED MUTATION	UNP P31345
B	389	LYS	ARG	ENGINEERED MUTATION	UNP P31345
D	389	LYS	ARG	ENGINEERED MUTATION	UNP P31345
E	389	LYS	ARG	ENGINEERED MUTATION	UNP P31345
F	389	LYS	ARG	ENGINEERED MUTATION	UNP P31345

- Molecule 2 is 7N-METHYL-8-HYDROGUANOSINE-5'-TRIPHOSPHATE (three-letter code: MGT) (formula: C₁₁H₂₀N₅O₁₄P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			33	11	5	14	3		
2	B	1	Total	C	N	O	P	0	0
			33	11	5	14	3		
2	D	1	Total	C	N	O	P	0	0
			33	11	5	14	3		
2	E	1	Total	C	N	O	P	0	0
			33	11	5	14	3		
2	F	1	Total	C	N	O	P	0	0
			33	11	5	14	3		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	65	Total	O	0	0
			65	65		
3	B	66	Total	O	0	0
			66	66		
3	D	58	Total	O	0	0
			58	58		
3	E	51	Total	O	0	0
			51	51		
3	F	32	Total	O	0	0
			32	32		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

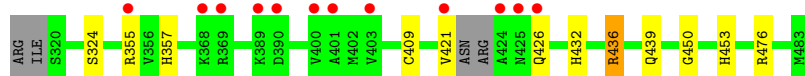
- Molecule 1: POLYMERASE BASIC PROTEIN 2

Chain A: 



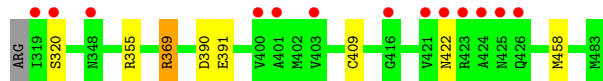
- Molecule 1: POLYMERASE BASIC PROTEIN 2

Chain B: 



- Molecule 1: POLYMERASE BASIC PROTEIN 2

Chain D: 



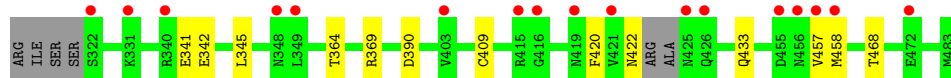
- Molecule 1: POLYMERASE BASIC PROTEIN 2

Chain E: 



- Molecule 1: POLYMERASE BASIC PROTEIN 2

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	92.20Å 94.44Å 220.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.53 – 2.30 29.52 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.53-2.30) 99.6 (29.52-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.186 , 0.235 0.182 , 0.230	Depositor DCC
R_{free} test set	2165 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 31.5	EDS
Estimated twinning fraction	0.047 for -k,-h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 42946 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6920	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MGT

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	0.66	0/1310	0.72	0/1734
1	B	0.66	0/1301	0.70	0/1721
1	D	0.65	0/1329	0.76	1/1760 (0.1%)
1	E	0.62	0/1276	0.70	0/1687
1	F	0.56	0/1303	0.62	0/1724
All	All	0.63	0/6519	0.70	1/8626 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	369	ARG	NE-CZ-NH2	-6.59	117.00	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1303	0	0	3	0
1	B	1294	0	0	7	0
1	D	1321	0	0	3	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1270	0	0	6	0
1	F	1295	0	0	2	0
2	A	33	0	0	1	0
2	B	33	0	0	3	0
2	D	33	0	0	1	0
2	E	33	0	0	1	0
2	F	33	0	0	0	0
3	A	65	0	0	2	0
3	B	66	0	0	2	0
3	D	58	0	0	1	1
3	E	51	0	0	6	0
3	F	32	0	0	0	0
All	All	6920	0	0	21	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (21) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:476:ARG:NH1	3:B:2062:HOH:O	2.24	0.69
1:E:383:GLN:NE2	3:E:2023:HOH:O	2.29	0.66
1:B:436:ARG:NH1	1:B:439:GLN:OE1	2.33	0.62
1:B:453[B]:HIS:CE1	3:E:2022:HOH:O	2.56	0.59
1:E:432:HIS:CE1	2:E:1483:MGT:O2A	2.59	0.56
1:E:453:HIS:NE2	3:E:2042:HOH:O	2.32	0.55
1:A:357:HIS:NE2	2:A:1483:MGT:O2G	2.40	0.55
1:B:355:ARG:CD	2:B:1483:MGT:O1G	2.56	0.54
1:E:436:ARG:NH2	3:E:2034:HOH:O	2.41	0.53
1:B:450:GLY:N	3:B:2050:HOH:O	2.43	0.52
1:A:385:ILE:CG1	3:A:2029:HOH:O	2.58	0.51
1:D:369:ARG:NH2	3:D:2021:HOH:O	2.44	0.51
1:F:369:ARG:NH1	1:F:390:ASP:OD2	2.44	0.51
1:E:383:GLN:CG	3:E:2023:HOH:O	2.59	0.50
1:A:476:ARG:NH1	3:A:2061:HOH:O	2.45	0.49
1:E:476:ARG:NH1	3:E:2047:HOH:O	2.51	0.43
1:D:355:ARG:CD	2:D:1483:MGT:O1G	2.66	0.43
1:D:369:ARG:NH1	1:D:390:ASP:OD2	2.52	0.43
1:B:357:HIS:NE2	2:B:1483:MGT:O2G	2.52	0.42
1:F:420:PHE:N	1:F:433:GLN:OE1	2.53	0.42
1:B:432:HIS:CE1	2:B:1483:MGT:O2A	2.73	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:458:MSE:SE	3:D:2027:HOH:O[4_555]	2.14	0.06

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	163/166 (98%)	159 (98%)	4 (2%)	0	100	100
1	B	160/166 (96%)	158 (99%)	2 (1%)	0	100	100
1	D	165/166 (99%)	161 (98%)	4 (2%)	0	100	100
1	E	157/166 (95%)	153 (98%)	4 (2%)	0	100	100
1	F	159/166 (96%)	156 (98%)	3 (2%)	0	100	100
All	All	804/830 (97%)	787 (98%)	17 (2%)	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	144/135 (107%)	140 (97%)	4 (3%)	56	73
1	B	143/135 (106%)	137 (96%)	6 (4%)	40	53
1	D	146/135 (108%)	141 (97%)	5 (3%)	49	64
1	E	141/135 (104%)	135 (96%)	6 (4%)	40	52
1	F	143/135 (106%)	134 (94%)	9 (6%)	25	32
All	All	717/675 (106%)	687 (96%)	30 (4%)	44	53

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	334	SER
1	A	341	GLU
1	A	369	ARG
1	A	422	ASN
1	B	324	SER
1	B	409[A]	CYS
1	B	409[B]	CYS
1	B	421	VAL
1	B	426	GLN
1	B	436	ARG
1	D	320	SER
1	D	391	GLU
1	D	409[A]	CYS
1	D	409[B]	CYS
1	D	422	ASN
1	E	334	SER
1	E	369	ARG
1	E	409[A]	CYS
1	E	409[B]	CYS
1	E	421	VAL
1	E	422	ASN
1	F	341	GLU
1	F	342	GLU
1	F	345	LEU
1	F	364	THR
1	F	409	CYS
1	F	422	ASN
1	F	457	VAL
1	F	458	MSE
1	F	468	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

5 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	MGT	A	1483	-	35,35,35	3.16	13 (37%)	54,56,56	1.46	7 (12%)
2	MGT	B	1483	-	35,35,35	2.99	11 (31%)	54,56,56	1.35	9 (16%)
2	MGT	D	1483	-	35,35,35	3.06	12 (34%)	54,56,56	1.64	12 (22%)
2	MGT	E	1483	-	35,35,35	3.07	13 (37%)	54,56,56	1.33	9 (16%)
2	MGT	F	1483	-	35,35,35	3.20	16 (45%)	54,56,56	1.41	7 (12%)

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	MGT	A	1483	-	-	0/22/50/50	0/1/3/3
2	MGT	B	1483	-	-	0/22/50/50	0/1/3/3
2	MGT	D	1483	-	-	0/22/50/50	0/1/3/3
2	MGT	E	1483	-	-	0/22/50/50	0/1/3/3
2	MGT	F	1483	-	-	0/22/50/50	0/1/3/3

All (65) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1483	MGT	C8-N9	-14.28	1.35	1.46
2	A	1483	MGT	C8-N9	-14.15	1.35	1.46
2	E	1483	MGT	C8-N9	-13.73	1.35	1.46
2	B	1483	MGT	C8-N9	-13.59	1.35	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1483	MGT	C8-N9	-13.26	1.35	1.46
2	A	1483	MGT	C1'-N9	5.57	1.52	1.45
2	E	1483	MGT	PA-O3A	5.08	1.69	1.59
2	D	1483	MGT	C1'-N9	4.78	1.51	1.45
2	A	1483	MGT	PA-O3A	4.52	1.68	1.59
2	F	1483	MGT	C1'-N9	4.42	1.50	1.45
2	D	1483	MGT	PG-O3B	4.22	1.67	1.60
2	A	1483	MGT	C8-N7	-4.19	1.33	1.45
2	D	1483	MGT	PA-O3A	4.07	1.67	1.59
2	B	1483	MGT	C1'-N9	4.04	1.50	1.45
2	E	1483	MGT	PB-O3A	3.99	1.67	1.59
2	B	1483	MGT	PA-O3A	3.76	1.66	1.59
2	F	1483	MGT	C8-N7	-3.76	1.35	1.45
2	B	1483	MGT	C2-N1	3.76	1.42	1.36
2	D	1483	MGT	C8-N7	-3.75	1.35	1.45
2	F	1483	MGT	PG-O1G	3.68	1.63	1.51
2	E	1483	MGT	C8-N7	-3.58	1.35	1.45
2	E	1483	MGT	PG-O1G	3.58	1.63	1.51
2	B	1483	MGT	C8-N7	-3.46	1.35	1.45
2	E	1483	MGT	PB-O1B	3.44	1.64	1.51
2	A	1483	MGT	C2-N1	3.35	1.42	1.36
2	F	1483	MGT	PB-O1B	3.32	1.64	1.51
2	B	1483	MGT	PG-O1G	3.32	1.62	1.51
2	D	1483	MGT	PB-O3A	3.28	1.65	1.59
2	F	1483	MGT	C2-N1	3.24	1.41	1.36
2	D	1483	MGT	PB-O3B	3.24	1.65	1.59
2	E	1483	MGT	C1'-N9	3.15	1.49	1.45
2	F	1483	MGT	PB-O3A	3.15	1.65	1.59
2	E	1483	MGT	PG-O2G	3.09	1.66	1.54
2	F	1483	MGT	PG-O3B	3.04	1.65	1.60
2	F	1483	MGT	CM7-N7	3.03	1.50	1.46
2	B	1483	MGT	PB-O1B	2.99	1.62	1.51
2	D	1483	MGT	PG-O1G	2.95	1.61	1.51
2	F	1483	MGT	PA-O3A	2.92	1.65	1.59
2	D	1483	MGT	PG-O2G	2.92	1.65	1.54
2	A	1483	MGT	PG-O1G	2.80	1.60	1.51
2	B	1483	MGT	C4-N9	-2.76	1.33	1.37
2	F	1483	MGT	PA-O1A	2.76	1.61	1.51
2	A	1483	MGT	PB-O1B	2.73	1.61	1.51
2	D	1483	MGT	PB-O1B	2.71	1.61	1.51
2	B	1483	MGT	PG-O2G	2.67	1.64	1.54
2	E	1483	MGT	PG-O3B	2.66	1.64	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1483	MGT	PA-O1A	2.60	1.61	1.51
2	F	1483	MGT	PB-O3B	2.60	1.64	1.59
2	F	1483	MGT	C6-N1	2.59	1.41	1.37
2	D	1483	MGT	C4-N9	-2.56	1.33	1.37
2	D	1483	MGT	PA-O1A	2.54	1.61	1.51
2	F	1483	MGT	C4-N9	-2.53	1.33	1.37
2	A	1483	MGT	PA-O1A	2.49	1.60	1.51
2	A	1483	MGT	PB-O3A	2.39	1.64	1.59
2	E	1483	MGT	C5-N7	-2.39	1.34	1.39
2	B	1483	MGT	O4'-C1'	2.37	1.48	1.42
2	E	1483	MGT	CM7-N7	2.23	1.49	1.46
2	E	1483	MGT	C4-N9	-2.15	1.34	1.37
2	A	1483	MGT	C4-N9	-2.13	1.34	1.37
2	F	1483	MGT	C5-N7	-2.12	1.34	1.39
2	F	1483	MGT	PG-O2G	2.12	1.62	1.54
2	A	1483	MGT	O4'-C1'	2.09	1.47	1.42
2	B	1483	MGT	PB-O3A	2.08	1.63	1.59
2	A	1483	MGT	PG-O2G	2.08	1.62	1.54
2	A	1483	MGT	PB-O3B	2.06	1.63	1.59

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1483	MGT	O3A-PB-O3B	4.62	111.06	101.66
2	A	1483	MGT	O4'-C1'-N9	-3.58	102.33	108.53
2	D	1483	MGT	C2-N3-C4	3.52	122.62	117.61
2	D	1483	MGT	C5-C4-N3	-3.35	120.58	126.61
2	A	1483	MGT	C2'-C1'-N9	3.33	122.25	114.59
2	A	1483	MGT	N3-C4-N9	3.29	132.28	127.06
2	A	1483	MGT	N2-C2-N1	3.25	121.43	117.86
2	B	1483	MGT	C2'-C1'-N9	3.22	122.00	114.59
2	E	1483	MGT	C5-C4-N3	-3.17	120.90	126.61
2	F	1483	MGT	N2-C2-N1	3.16	121.34	117.86
2	A	1483	MGT	C2-N3-C4	3.07	121.98	117.61
2	F	1483	MGT	PB-O3B-PG	-3.05	122.73	131.68
2	B	1483	MGT	N2-C2-N1	3.01	121.18	117.86
2	A	1483	MGT	O3G-PG-O3B	2.95	119.12	105.14
2	F	1483	MGT	C2-N3-C4	2.95	121.80	117.61
2	E	1483	MGT	O3A-PB-O3B	2.95	107.65	101.66
2	A	1483	MGT	C5-C4-N3	-2.92	121.36	126.61
2	D	1483	MGT	N3-C4-N9	2.89	131.63	127.06
2	B	1483	MGT	PB-O3B-PG	-2.83	123.38	131.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1483	MGT	C5-C4-N3	-2.82	121.53	126.61
2	F	1483	MGT	C5-C4-N3	-2.76	121.64	126.61
2	B	1483	MGT	C2-N3-C4	2.74	121.50	117.61
2	D	1483	MGT	N2-C2-N1	2.72	120.86	117.86
2	E	1483	MGT	N3-C4-N9	2.68	131.31	127.06
2	B	1483	MGT	N3-C4-N9	2.68	131.30	127.06
2	D	1483	MGT	C6-C5-N7	2.54	137.26	131.87
2	F	1483	MGT	O3G-PG-O3B	2.52	117.10	105.14
2	F	1483	MGT	N3-C4-N9	2.49	131.00	127.06
2	D	1483	MGT	C2'-C1'-N9	2.48	120.28	114.59
2	E	1483	MGT	C2-N3-C4	2.46	121.11	117.61
2	D	1483	MGT	O3G-PG-O3B	2.45	116.75	105.14
2	E	1483	MGT	N2-C2-N1	2.36	120.45	117.86
2	B	1483	MGT	O3G-PG-O3B	2.35	116.28	105.14
2	B	1483	MGT	O4'-C1'-N9	-2.35	104.46	108.53
2	D	1483	MGT	O4'-C1'-N9	-2.32	104.52	108.53
2	D	1483	MGT	O2B-PB-O3A	2.29	115.99	105.14
2	D	1483	MGT	PA-O3A-PB	-2.26	125.06	131.68
2	D	1483	MGT	O2A-PA-O3A	2.25	115.84	105.14
2	E	1483	MGT	PB-O3B-PG	-2.18	125.29	131.68
2	E	1483	MGT	N7-C8-N9	2.15	105.92	103.08
2	E	1483	MGT	C2'-C1'-N9	2.14	119.50	114.59
2	E	1483	MGT	O3G-PG-O3B	2.05	114.86	105.14
2	B	1483	MGT	CM7-N7-C8	2.00	124.19	119.23
2	F	1483	MGT	C6-C5-N7	2.00	136.11	131.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	164/166 (98%)	-0.07	4 (2%) 56 66	15, 26, 43, 64	0
1	B	162/166 (97%)	0.18	12 (7%) 14 21	17, 29, 48, 83	0
1	D	165/166 (99%)	0.20	13 (7%) 13 19	14, 27, 59, 80	0
1	E	160/166 (96%)	0.14	8 (5%) 28 38	17, 31, 54, 70	0
1	F	160/166 (96%)	0.43	17 (10%) 7 11	29, 47, 74, 100	0
All	All	811/830 (97%)	0.17	54 (6%) 18 25	14, 31, 65, 100	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	425	ASN	6.1
1	B	426	GLN	5.6
1	B	424	ALA	5.1
1	D	424	ALA	5.1
1	E	421	VAL	5.0
1	F	456	ASN	5.0
1	D	423	ARG	4.9
1	F	421	VAL	4.6
1	F	349	LEU	4.5
1	A	424	ALA	4.1
1	F	340	ARG	3.6
1	D	425	ASN	3.5
1	F	426	GLN	3.5
1	D	319	ILE	3.5
1	F	455	ASP	3.5
1	A	423	ARG	3.5
1	B	425	ASN	3.4
1	E	422	ASN	3.3
1	F	322	SER	3.2
1	D	422	ASN	3.0

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Mol	Chain	Res	Type	RSRZ
1	F	419	ASN	2.9
1	D	348	ASN	2.9
1	B	421	VAL	2.9
1	E	427	ARG	2.9
1	D	421	VAL	2.8
1	B	355	ARG	2.8
1	F	331	LYS	2.8
1	D	426	GLN	2.7
1	B	401	ALA	2.7
1	A	425	ASN	2.6
1	B	400	VAL	2.6
1	E	389	LYS	2.5
1	F	457	VAL	2.5
1	E	320	SER	2.5
1	B	389	LYS	2.5
1	B	368	LYS	2.4
1	A	426	GLN	2.4
1	B	390	ASP	2.4
1	F	415	ARG	2.3
1	E	355	ARG	2.3
1	F	403	VAL	2.3
1	D	401	ALA	2.3
1	D	403	VAL	2.2
1	B	403	VAL	2.2
1	B	369	ARG	2.1
1	F	416	GLY	2.1
1	F	472[A]	GLU	2.1
1	E	384	LEU	2.1
1	D	416	GLY	2.1
1	D	400	VAL	2.1
1	F	348	ASN	2.1
1	E	349	LEU	2.1
1	F	458	MSE	2.1
1	D	320	SER	2.0

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

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

6.3 Carbohydrates

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MGT	F	1483	33/33	0.18	0.55	54,65,97,100	0
2	MGT	A	1483	33/33	0.14	0.19	25,34,50,51	0
2	MGT	E	1483	33/33	0.15	0.02	30,43,68,70	0
2	MGT	D	1483	33/33	0.13	-0.21	27,41,64,67	0
2	MGT	B	1483	33/33	0.12	-0.23	23,38,52,53	0

6.5 Other polymers

There are no such residues in this entry.