



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:01 AM UTC

PDB ID : 5J2M / pdb_00005j2m
Title : HIV-1 reverse transcriptase in complex with DNA and EFdA-triphosphate, a translocation-defective RT inhibitor
Authors : Salie, Z.L.; Kirby, K.A.; Sarafianos, S.G.
Deposited on : 2016-03-29
Resolution : 2.43 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

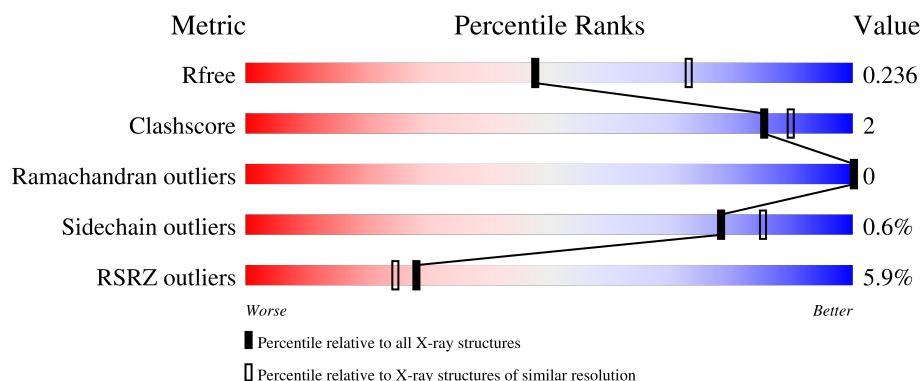
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	560	4% 93% 7%
2	B	440	8% 85% 5% 10%
3	P	21	57% 24% 5% 14%
4	T	27	74% 15% 11%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 9107 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 HIV-1 reverse transcriptase p51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	558	Total	C	N	O	S	7	0	0
			4542	2935	760	839	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	CYS	GLN	engineered mutation	UNP P04585
A	280	SER	CYS	engineered mutation	UNP P04585

- Molecule 2 is a protein called HIV-1 reverse transcriptase p51 subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	398	Total	C	N	O	S	16	0	0
			3281	2132	548	595	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	engineered mutation	UNP P04585

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	18	Total	C	N	O	P	0	0	0
			370	176	64	111	18			

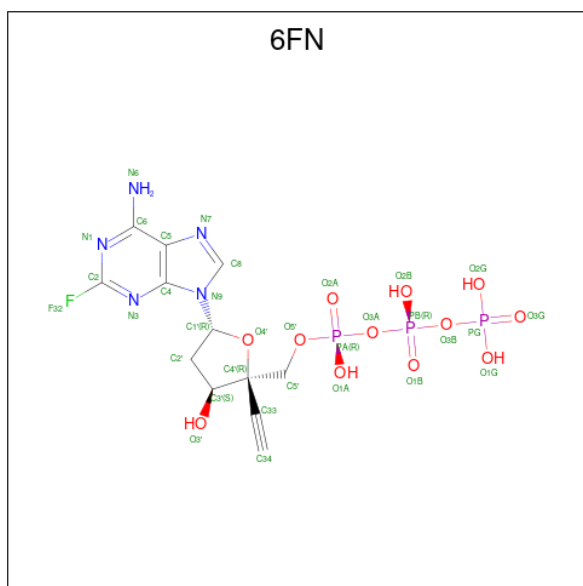
- Molecule 4 is a DNA chain called DNA (27-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	T	24	Total	C	N	O	P	0	0	0
			494	233	97	141	23			

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 6 is 2'-deoxy-4'-ethynyl-2-fluoroadenosine 5'-(tetrahydrogen triphosphate) (CCD ID: 6FN) (formula: C₁₂H₁₅FN₅O₁₂P₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	A	1	Total	C	F	N	O	P	0	0
			33	12	1	5	12	3		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	219	Total	O	0	0
			219	219		
7	B	131	Total	O	0	0
			131	131		
7	P	15	Total	O	0	0
			15	15		
7	T	19	Total	O	0	0
			19	19		

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	167.47Å 169.76Å 101.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.23 – 2.43 65.23 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.5 (65.23-2.43) 99.5 (65.23-2.43)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 2.42Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.199 , 0.233 0.203 , 0.236	Depositor DCC
R_{free} test set	2766 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	44.8	Xtriage
Anisotropy	0.343	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9107	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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: 6FN, MG, MRG

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	0.10	0/4659	0.29	2/6328 (0.0%)
2	B	0.09	0/3370	0.24	0/4575
3	P	0.19	0/382	0.36	0/584
4	T	0.16	0/555	0.30	0/856
All	All	0.11	0/8966	0.28	2/12343 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	419	THR	CA-C-N	5.34	123.56	119.66
1	A	419	THR	C-N-CA	5.34	123.56	119.66

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	4542	0	4599	22	0
2	B	3281	0	3320	10	0
3	P	370	0	206	5	0
4	T	494	0	269	2	0
5	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	33	0	0	0	0
7	A	219	0	0	2	2
7	B	131	0	0	0	0
7	P	15	0	0	0	0
7	T	19	0	0	0	0
All	All	9107	0	8394	37	2

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 2.

All (37) 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:275:LYS:HE3	2:B:302:GLU:HG3	1.80	0.63
1:A:275:LYS:HD2	1:A:332:GLN:HG2	1.82	0.60
3:P:816:DG:H2'	3:P:817:MRG:C8	2.35	0.55
1:A:259:LYS:HA	3:P:819:DG:H5'	1.88	0.55
1:A:220:LYS:NZ	7:A:711:HOH:O	2.38	0.55
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.90	0.54
1:A:60:VAL:HG12	1:A:75:VAL:HG22	1.90	0.54
1:A:184:MET:HG3	3:P:822:DG:H1'	1.89	0.54
1:A:13:LYS:HB2	1:A:16:MET:HG3	1.90	0.54
1:A:556:ILE:HD12	1:A:557:ARG:HG3	1.91	0.52
1:A:16:MET:HE3	1:A:83:ARG:HG2	1.92	0.51
1:A:332:GLN:O	7:A:701:HOH:O	2.20	0.50
4:T:713:DC:H2'	4:T:714:DG:C8	2.48	0.49
2:B:79:GLU:HG3	2:B:83:ARG:HE	1.78	0.47
1:A:503:LEU:HD22	1:A:535:TRP:HB2	1.97	0.46
1:A:255:ASN:HB2	1:A:289:LEU:HB3	1.96	0.46
4:T:703:DG:H2'	4:T:704:DG:C4	2.52	0.45
1:A:557:ARG:HB3	1:A:558:LYS:HD2	1.98	0.45
1:A:317:VAL:HG23	1:A:349:LEU:HD23	1.98	0.44
1:A:325:LEU:HB3	1:A:387:PRO:HB3	1.99	0.43
1:A:134:SER:OG	1:A:135:ILE:N	2.52	0.43
2:B:74:LEU:HD21	2:B:409:THR:HA	2.01	0.43
1:A:58:THR:HG21	1:A:77:PHE:CD1	2.54	0.42
2:B:332:GLN:HB2	2:B:336:GLN:HB2	2.01	0.42
2:B:423:VAL:HA	2:B:426:TRP:CD1	2.55	0.42
3:P:816:DG:H2'	3:P:817:MRG:H8	2.01	0.41
1:A:507:GLN:HE21	1:A:507:GLN:HB3	1.74	0.41
2:B:363:ASN:HB3	2:B:366:LYS:HB3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:ASP:O	1:A:125:ARG:HG3	2.20	0.41
2:B:66:LYS:HG3	2:B:67:ASP:H	1.84	0.41
2:B:261:VAL:HG13	2:B:276:VAL:HG21	2.03	0.41
1:A:405:TYR:CE2	1:A:407:GLN:HB2	2.56	0.41
1:A:438:GLU:OE2	1:A:463:ARG:HD3	2.21	0.41
1:A:111:VAL:HB	1:A:185:ASP:HB2	2.03	0.41
2:B:337:TRP:O	2:B:354:TYR:N	2.49	0.40
1:A:490:GLY:O	1:A:528:LYS:NZ	2.43	0.40
3:P:810:DT:H2"	3:P:811:DG:H8	1.86	0.40

All (2) 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)
7:A:882:HOH:O	7:A:889:HOH:O[7_444]	2.15	0.05
7:A:790:HOH:O	7:A:889:HOH:O[7_444]	2.16	0.04

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	556/560 (99%)	543 (98%)	13 (2%)	0	100	100
2	B	392/440 (89%)	378 (96%)	14 (4%)	0	100	100
All	All	948/1000 (95%)	921 (97%)	27 (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	498/500 (100%)	496 (100%)	2 (0%)	84	89
2	B	360/400 (90%)	357 (99%)	3 (1%)	73	81
All	All	858/900 (95%)	853 (99%)	5 (1%)	78	85

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	LYS
1	A	298	GLU
2	B	185	ASP
2	B	195	ILE
2	B	298	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	207	GLN
1	A	265	ASN
2	B	208	HIS
2	B	428	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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
3	MRG	P	817	1,4,3	25,28,29	2.11	6 (24%)	32,39,42	1.91	9 (28%)

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
3	MRG	P	817	1,4,3	-	2/12/26/27	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	P	817	MRG	O6-C6	6.31	1.35	1.23
3	P	817	MRG	C2-N3	4.87	1.41	1.32
3	P	817	MRG	C2-N2	3.06	1.40	1.34
3	P	817	MRG	C8-N9	2.09	1.42	1.37
3	P	817	MRG	C4-N9	2.04	1.43	1.38
3	P	817	MRG	C5-C4	2.03	1.44	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	P	817	MRG	C2-N1-C6	-6.30	116.94	124.55
3	P	817	MRG	C5-C6-N1	4.02	123.48	113.25
3	P	817	MRG	O6-C6-C5	-3.07	118.43	126.53
3	P	817	MRG	N9-C4-N3	2.85	131.64	125.95
3	P	817	MRG	C5-C4-N3	-2.67	124.14	128.39
3	P	817	MRG	N2-C2-N3	-2.41	116.87	120.19
3	P	817	MRG	C2'-C3'-C4'	-2.22	98.30	102.80
3	P	817	MRG	C6-C5-C4	-2.22	115.50	118.83
3	P	817	MRG	C2-N3-C4	2.04	114.56	112.00

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	P	817	MRG	C21-C22-C23-S24
3	P	817	MRG	N2-C21-C22-C23

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	P	817	MRG	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 3 are 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 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
6	6FN	A	604	5	31,35,35	3.38	16 (51%)	38,56,56	3.78	16 (42%)

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
6	6FN	A	604	5	-	5/23/50/50	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	604	6FN	C8-N9	-8.70	1.22	1.37
6	A	604	6FN	C4-N9	-7.48	1.22	1.37
6	A	604	6FN	PB-O3A	5.25	1.65	1.59
6	A	604	6FN	PA-O3A	5.22	1.65	1.59
6	A	604	6FN	PB-O3B	5.21	1.65	1.59
6	A	604	6FN	O4'-C1'	5.15	1.50	1.42
6	A	604	6FN	C6-N6	4.52	1.45	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	604	6FN	C2'-C3'	-3.56	1.45	1.53
6	A	604	6FN	C2-N1	3.11	1.34	1.31
6	A	604	6FN	C4'-C33	3.02	1.51	1.48
6	A	604	6FN	C5-N7	2.89	1.44	1.39
6	A	604	6FN	PG-O2G	2.81	1.65	1.54
6	A	604	6FN	C1'-N9	-2.43	1.41	1.46
6	A	604	6FN	C2-N3	2.30	1.33	1.31
6	A	604	6FN	PG-O1G	-2.21	1.46	1.54
6	A	604	6FN	C8-N7	2.21	1.35	1.31

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	604	6FN	C4-N9-C8	14.70	121.17	105.74
6	A	604	6FN	N3-C4-N9	6.97	135.83	126.99
6	A	604	6FN	C5-C4-N3	-6.82	119.99	127.18
6	A	604	6FN	N1-C2-N3	-6.55	121.14	130.60
6	A	604	6FN	N9-C8-N7	-5.80	105.71	113.94
6	A	604	6FN	C4-C5-N7	-5.48	104.31	110.58
6	A	604	6FN	F32-C2-N3	4.95	119.45	114.69
6	A	604	6FN	F32-C2-N1	4.92	119.41	114.69
6	A	604	6FN	N6-C6-N1	3.02	121.10	117.03
6	A	604	6FN	C2-N1-C6	2.81	121.59	112.44
6	A	604	6FN	O1A-PA-O2A	-2.42	101.17	112.44
6	A	604	6FN	O1G-PG-O3G	-2.42	101.39	110.83
6	A	604	6FN	O2B-PB-O1B	-2.31	101.69	112.44
6	A	604	6FN	C1'-N9-C8	-2.23	119.13	126.95
6	A	604	6FN	C2'-C1'-N9	-2.23	109.39	114.63
6	A	604	6FN	C6-C5-N7	2.11	136.15	132.09

There are no chirality outliers.

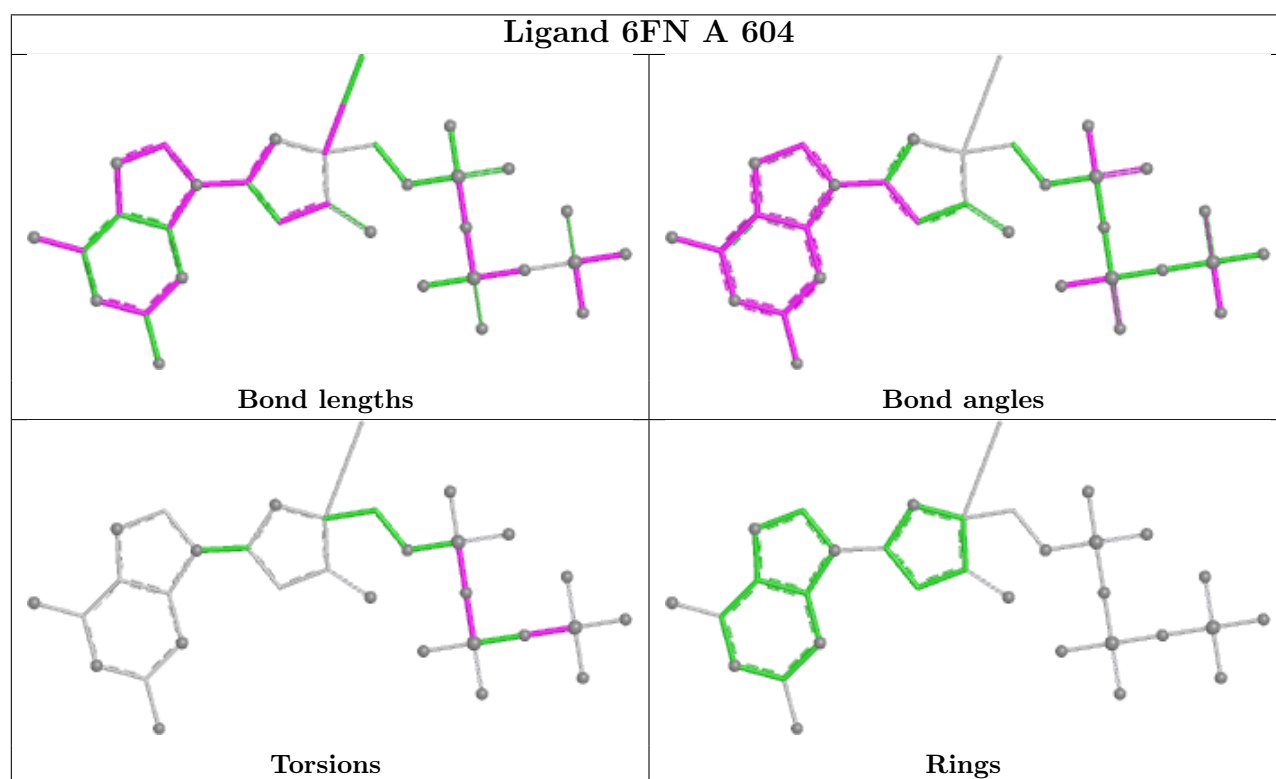
All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	604	6FN	PB-O3B-PG-O2G
6	A	604	6FN	PB-O3B-PG-O1G
6	A	604	6FN	PA-O3A-PB-O2B
6	A	604	6FN	PB-O3A-PA-O2A
6	A	604	6FN	PA-O3A-PB-O1B

There are no ring outliers.

No monomer is involved in short contacts.

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 than 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 any Mogul-identified 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

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	558/560 (99%)	0.13	24 (4%)	40	37	26, 42, 74, 116	2 (0%)
2	B	398/440 (90%)	0.52	35 (8%)	15	13	27, 53, 86, 110	5 (1%)
3	P	17/21 (80%)	-0.00	0	100	100	36, 54, 70, 70	0
4	T	24/27 (88%)	0.18	0	100	100	35, 60, 87, 117	0
All	All	997/1048 (95%)	0.29	59 (5%)	28	25	26, 47, 83, 117	7 (0%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	418	ASN	5.0
1	A	135	ILE	4.9
2	B	419	THR	4.3
2	B	214	LEU	4.1
2	B	430	GLU	4.1
2	B	5	ILE	3.4
1	A	136	ASN	3.3
1	A	134	SER	3.2
2	B	420	PRO	3.2
2	B	425	LEU	3.2
2	B	67	ASP	3.1
2	B	275	LYS	3.1
2	B	195	ILE	3.0
2	B	65	LYS	3.0
1	A	67	ASP	2.9
2	B	212	TRP	2.9
2	B	68	SER	2.9
1	A	543	GLY	2.9
1	A	333	GLY	2.8
1	A	137	ASN	2.8
2	B	429	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	315	HIS	2.8
2	B	7	THR	2.8
1	A	357	MET	2.8
1	A	259	LYS	2.7
2	B	237	ASP	2.7
2	B	240	THR	2.7
1	A	292	VAL	2.6
2	B	4	PRO	2.6
2	B	66	LYS	2.6
2	B	356	ARG	2.6
2	B	245	VAL	2.6
1	A	22	LYS	2.5
1	A	419	THR	2.5
2	B	357	MET	2.5
2	B	386	THR	2.5
2	B	360	ALA	2.4
1	A	558	LYS	2.4
2	B	69	THR	2.3
1	A	297	GLU	2.3
1	A	426	TRP	2.3
1	A	258	CYS	2.3
1	A	140	PRO	2.2
1	A	402	TRP	2.2
2	B	8	VAL	2.2
2	B	301	LEU	2.1
1	A	294	PRO	2.1
1	A	252	TRP	2.1
2	B	361	HIS	2.1
2	B	168	LEU	2.1
1	A	24	TRP	2.1
2	B	211	ARG	2.1
2	B	233	GLU	2.1
1	A	289	LEU	2.1
1	A	404	GLU	2.1
2	B	70	LYS	2.1
1	A	132	ILE	2.0
2	B	421	PRO	2.0
2	B	422	LEU	2.0

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
3	MRG	P	817	26/27	0.91	0.12	45,54,81,91	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

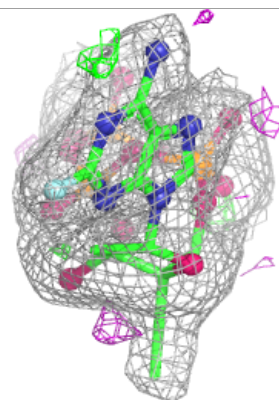
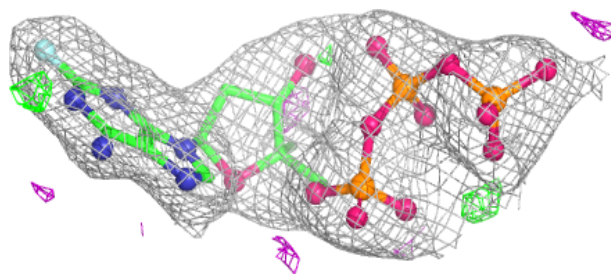
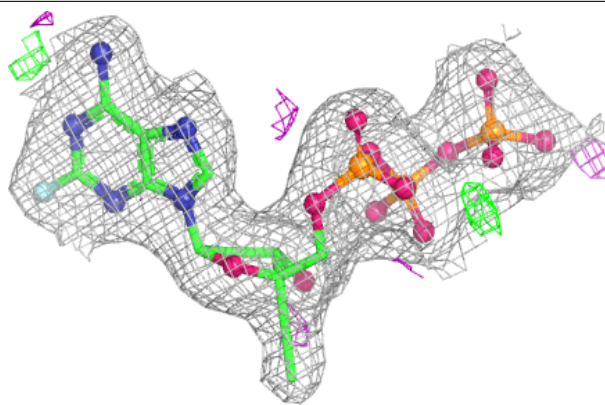
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MG	A	602	1/1	0.90	0.17	57,57,57,57	0
5	MG	A	603	1/1	0.92	0.33	74,74,74,74	0
6	6FN	A	604	33/33	0.97	0.07	27,35,42,45	0
5	MG	A	601	1/1	0.98	0.07	22,22,22,22	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.

Electron density around 6FN A 604:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.