



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 15, 2017 – 07:34 AM EDT

PDB ID : 3B2T
Title : Structure of phosphotransferase
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Deposited on : unknown
Resolution : 1.80 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

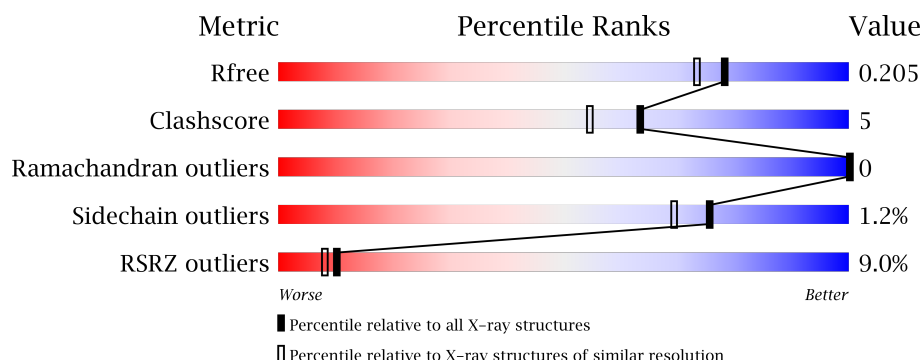
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	100719	4827 (1.80-1.80)
Clashscore	112137	5742 (1.80-1.80)
Ramachandran outliers	110173	5676 (1.80-1.80)
Sidechain outliers	110143	5675 (1.80-1.80)
RSRZ outliers	101464	4906 (1.80-1.80)

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. 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	311	<div> <div>9%</div> <div>81%</div> <div>11%</div> <div>7%</div> </div>
1	B	311	<div> <div>8%</div> <div>84%</div> <div>8%</div> <div>8%</div> </div>

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	PO4	A	1	-	-	-	X
2	PO4	B	2	-	-	-	X
3	M33	B	1996	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5191 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.

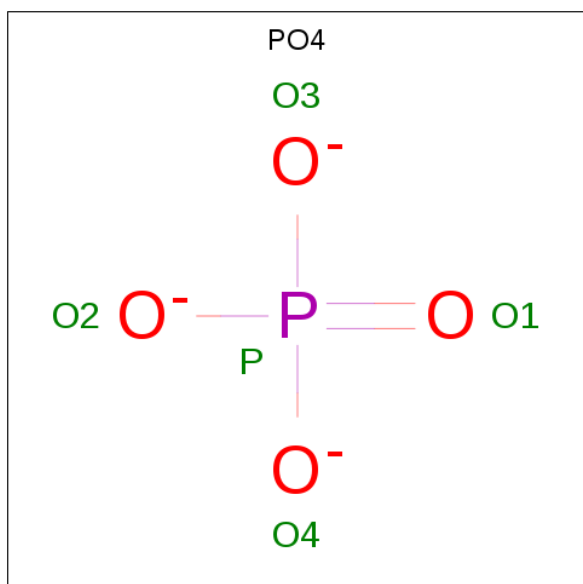
- Molecule 1 is a protein called Fibroblast growth factor receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2284	1452	386	425	21			
1	B	287	Total	C	N	O	S	0	0	0
			2272	1454	381	416	21			

There are 4 discrepancies between the modelled and reference sequences:

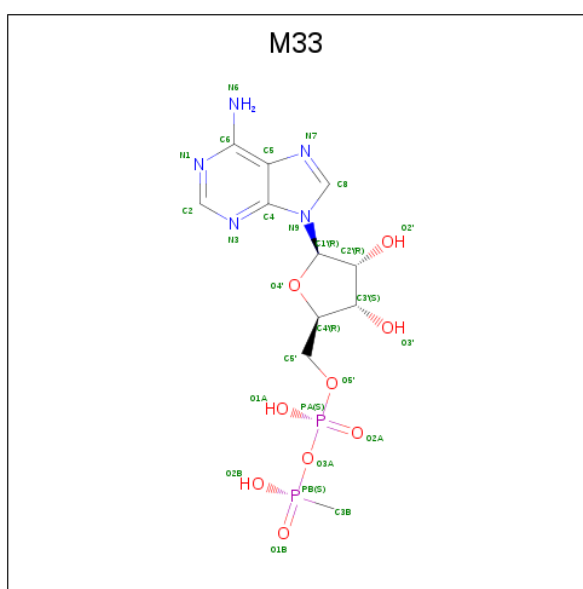
Chain	Residue	Modelled	Actual	Comment	Reference
A	628	THR	ALA	ENGINEERED	UNP P21802
A	767	GLN	GLU	ENGINEERED	UNP P21802
B	628	THR	ALA	ENGINEERED	UNP P21802
B	767	GLN	GLU	ENGINEERED	UNP P21802

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O P 5 4 1	0	0
2	A	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0
2	B	1	Total O P 5 4 1	0	0

- Molecule 3 is 5'-O-[(S)-hydroxy{[(S)-hydroxy(methyl)phosphoryl]oxy}phosphoryl]adenosine (three-letter code: M33) (formula: C₁₁H₁₇N₅O₉P₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P 27 11 5 9 2	0	0
3	B	1	Total C N O P 27 11 5 9 2	0	0

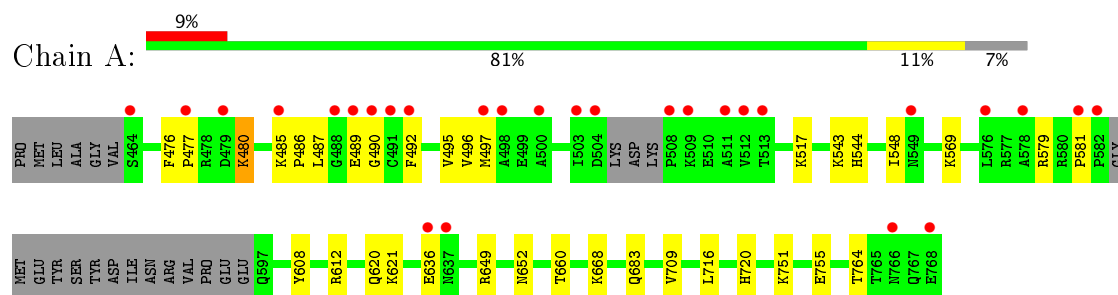
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	292	Total O 292 292	0	0
4	B	269	Total O 269 269	0	0

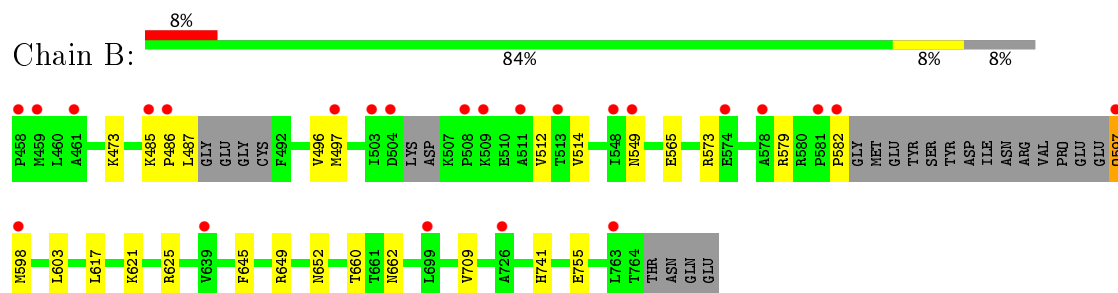
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Fibroblast growth factor receptor 2



• Molecule 1: Fibroblast growth factor receptor 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	67.70 Å 80.34 Å 118.60 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80 47.71 – 1.44	Depositor EDS
% Data completeness (in resolution range)	99.0 (50.00-1.80) 97.2 (47.71-1.44)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 1.44 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.201 , 0.233 0.204 , 0.205	Depositor DCC
R_{free} test set	3023 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.648	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5191	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 69.94 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.4008e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: PO4, M33

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	0/2330	0.66	1/3150 (0.0%)
1	B	0.44	0/2318	0.66	1/3132 (0.0%)
All	All	0.45	0/4648	0.66	2/6282 (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	581	PRO	N-CA-CB	5.60	110.03	103.30
1	B	582	PRO	N-CA-CB	5.47	109.86	103.30

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	2284	0	2261	27	0
1	B	2272	0	2265	15	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	27	0	15	1	0
3	B	27	0	15	2	0
4	A	292	0	0	7	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	269	0	0	3	1
All	All	5191	0	4556	45	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 5.

All (45) 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:B:487:LEU:HD21	1:B:497:MET:HG3	1.63	0.81
3:B:1996:M33:H5'1	3:B:1996:M33:H8	1.63	0.79
1:A:485:LYS:HD2	1:A:486:PRO:HD2	1.73	0.71
1:A:543:LYS:HD3	1:A:544:HIS:N	2.06	0.71
1:A:764:THR:HG23	4:A:2066:HOH:O	1.92	0.68
1:A:720:HIS:HE1	4:A:2072:HOH:O	1.78	0.67
1:A:720:HIS:HD2	4:A:2206:HOH:O	1.82	0.62
1:B:549:ASN:OD1	1:B:565:GLU:HB2	2.00	0.61
1:B:621:LYS:HD2	1:B:652:ASN:ND2	2.16	0.60
1:A:486:PRO:HA	1:A:496:VAL:HG12	1.84	0.59
1:A:517:LYS:NZ	1:A:517:LYS:HB3	2.18	0.58
1:A:480:LYS:HE3	1:A:480:LYS:HA	1.87	0.56
1:A:490:GLY:C	1:A:492:PHE:H	2.10	0.55
1:A:649:ARG:HD2	4:A:2189:HOH:O	2.06	0.55
1:A:485:LYS:HD2	1:A:486:PRO:CD	2.37	0.54
1:B:755:GLU:HG3	4:B:2236:HOH:O	2.07	0.54
1:A:620:GLN:NE2	4:A:2149:HOH:O	2.40	0.51
1:B:649:ARG:HD2	4:B:2203:HOH:O	2.13	0.49
1:B:486:PRO:HA	1:B:496:VAL:HG12	1.94	0.49
1:A:569:LYS:HE3	1:A:636:GLU:O	2.15	0.47
1:A:660:THR:HG23	1:A:660:THR:O	2.15	0.47
1:B:741:HIS:HE1	4:B:2261:HOH:O	1.98	0.47
3:B:1996:M33:C8	3:B:1996:M33:H5'1	2.38	0.47
1:B:512:VAL:HG23	1:B:514:VAL:HG13	1.97	0.46
1:A:608:TYR:CZ	1:A:612:ARG:HD3	2.51	0.45
1:A:683:GLN:HG3	4:A:2244:HOH:O	2.17	0.45
1:A:517:LYS:HZ2	1:A:517:LYS:HB3	1.81	0.45
1:A:487:LEU:HD12	1:A:495:VAL:HG12	1.98	0.44
1:B:579:ARG:HG3	1:B:579:ARG:HH11	1.83	0.43
1:A:621:LYS:HG3	1:A:652:ASN:ND2	2.34	0.43
3:A:1996:M33:H2'	3:A:1996:M33:O5'	2.19	0.43
1:A:476:PHE:CD1	1:A:477:PRO:HD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:617:LEU:CD1	1:B:645:PHE:HE2	2.32	0.42
1:B:597:GLN:NE2	1:B:597:GLN:O	2.52	0.42
1:A:548:ILE:HD12	1:A:548:ILE:HA	1.81	0.42
1:A:716:LEU:HD11	4:A:1998:HOH:O	2.19	0.42
1:A:751:LYS:O	1:A:755:GLU:HG3	2.19	0.42
1:B:598:MET:HE2	1:B:603:LEU:HD23	2.02	0.42
1:A:487:LEU:HD21	1:A:497:MET:HG3	2.02	0.41
1:B:660:THR:O	1:B:660:THR:HG23	2.20	0.41
1:A:579:ARG:HG3	1:A:579:ARG:HH11	1.86	0.41
1:B:485:LYS:HA	1:B:486:PRO:HD3	1.89	0.41
1:A:543:LYS:C	1:A:543:LYS:HD3	2.41	0.41
1:A:668:LYS:HG3	1:A:709:VAL:HG22	2.04	0.40
1:B:625:ARG:NH2	1:B:662:ASN:HD22	2.20	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)
4:A:2051:HOH:O	4:A:2051:HOH:O[2_655]	2.06	0.14
4:A:2287:HOH:O	4:B:2264:HOH:O[2_655]	2.13	0.07

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	282/311 (91%)	276 (98%)	6 (2%)	0	100	100
1	B	279/311 (90%)	273 (98%)	6 (2%)	0	100	100
All	All	561/622 (90%)	549 (98%)	12 (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	247/274 (90%)	245 (99%)	2 (1%)	85	81
1	B	243/274 (89%)	239 (98%)	4 (2%)	68	58
All	All	490/548 (89%)	484 (99%)	6 (1%)	75	69

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	480	LYS
1	A	489	GLU
1	B	473	LYS
1	B	573	ARG
1	B	597	GLN
1	B	709	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	620	GLN
1	A	631	ASN
1	A	652	ASN
1	A	662	ASN
1	A	720	HIS
1	A	741	HIS
1	B	597	GLN
1	B	631	ASN
1	B	652	ASN
1	B	662	ASN
1	B	683	GLN
1	B	741	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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	PO4	A	1	-	4,4,4	1.21	0	6,6,6	0.41	0
3	M33	A	1996	-	25,29,29	1.20	3 (12%)	23,45,45	2.22	4 (17%)
2	PO4	A	2	-	4,4,4	1.12	0	6,6,6	0.34	0
2	PO4	B	1	-	4,4,4	1.28	0	6,6,6	0.43	0
3	M33	B	1996	-	25,29,29	1.20	3 (12%)	23,45,45	2.30	5 (21%)
2	PO4	B	2	-	4,4,4	1.23	0	6,6,6	0.37	0

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	PO4	A	1	-	-	0/0/0/0	0/0/0/0
3	M33	A	1996	-	-	0/9/32/32	0/3/3/3
2	PO4	A	2	-	-	0/0/0/0	0/0/0/0
2	PO4	B	1	-	-	0/0/0/0	0/0/0/0
3	M33	B	1996	-	-	0/9/32/32	0/3/3/3
2	PO4	B	2	-	-	0/0/0/0	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1996	M33	PB-O2B	-3.18	1.48	1.54
3	B	1996	M33	PB-O2B	-3.12	1.48	1.54
3	A	1996	M33	O4'-C1'	2.24	1.44	1.41
3	B	1996	M33	O4'-C1'	2.27	1.44	1.41
3	A	1996	M33	PB-O3A	2.79	1.61	1.58
3	B	1996	M33	PB-O3A	2.81	1.61	1.58

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1996	M33	N3-C2-N1	-8.17	121.74	128.86
3	B	1996	M33	N3-C2-N1	-7.91	121.97	128.86
3	B	1996	M33	O2B-PB-O3A	-4.02	98.04	107.45
3	A	1996	M33	O2B-PB-O3A	-3.85	98.44	107.45
3	B	1996	M33	PA-O3A-PB	-3.66	120.60	132.39
3	A	1996	M33	PA-O3A-PB	-3.43	121.34	132.39
3	B	1996	M33	C2'-C3'-C4'	-2.65	97.46	102.62
3	B	1996	M33	C4'-O4'-C1'	-2.47	107.14	109.77
3	A	1996	M33	C4-C5-N7	-2.13	107.35	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1996	M33	1	0
3	B	1996	M33	2	0

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	288/311 (92%)	0.63	28 (9%) 8 7	3, 14, 34, 50	0
1	B	287/311 (92%)	0.48	24 (8%) 12 10	4, 15, 30, 45	0
All	All	575/622 (92%)	0.56	52 (9%) 10 8	3, 14, 31, 50	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	490	GLY	8.7
1	B	504	ASP	5.9
1	A	512	VAL	5.3
1	A	503	ILE	5.2
1	A	488	GLY	5.1
1	B	582	PRO	5.1
1	A	491	CYS	4.9
1	A	504	ASP	4.6
1	A	582	PRO	4.6
1	A	489	GLU	4.5
1	A	581	PRO	4.1
1	B	597	GLN	3.8
1	B	458	PRO	3.8
1	B	581	PRO	3.6
1	A	637	ASN	3.5
1	A	464	SER	3.5
1	A	513	THR	3.4
1	A	511	ALA	3.3
1	B	508	PRO	3.3
1	B	509	LYS	3.1
1	B	549	ASN	3.1
1	B	578	ALA	3.0
1	A	578	ALA	2.9
1	A	498	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	461	ALA	2.8
1	A	509	LYS	2.7
1	A	508	PRO	2.7
1	B	726	ALA	2.7
1	B	503	ILE	2.6
1	A	549	ASN	2.6
1	B	485	LYS	2.5
1	B	459	MET	2.5
1	A	492	PHE	2.5
1	B	598	MET	2.5
1	A	766	ASN	2.5
1	A	485	LYS	2.5
1	A	768	GLU	2.5
1	A	479	ASP	2.4
1	A	497	MET	2.4
1	A	477	PRO	2.4
1	B	548	ILE	2.4
1	B	763	LEU	2.4
1	B	513	THR	2.4
1	B	497	MET	2.3
1	B	486	PRO	2.3
1	A	576	LEU	2.3
1	B	574	GLU	2.2
1	A	500	ALA	2.1
1	A	636	GLU	2.1
1	B	639	VAL	2.0
1	B	511	ALA	2.0
1	B	699	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	PO4	A	1	5/5	0.80	0.21	12.20	21,28,29,29	0
2	PO4	B	2	5/5	0.66	0.21	6.33	35,39,40,41	0
3	M33	B	1996	27/27	0.67	0.23	2.94	16,33,58,59	0
3	M33	A	1996	27/27	0.71	0.21	0.12	16,37,57,57	0
2	PO4	B	1	5/5	0.98	0.12	-3.75	4,4,6,9	0
2	PO4	A	2	5/5	0.97	0.12	-5.91	4,5,6,7	0

6.5 Other polymers [i](#)

There are no such residues in this entry.