



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 07:12 PM GMT

PDB ID : 1PMQ
Title : The structure of JNK3 in complex with an imidazole-pyrimidine inhibitor
Authors : Scapin, G.; Patel, S.B.; Lisnock, J.; Becker, J.W.; LoGrasso, P.V.
Deposited on : 2003-06-11
Resolution : 2.20 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>

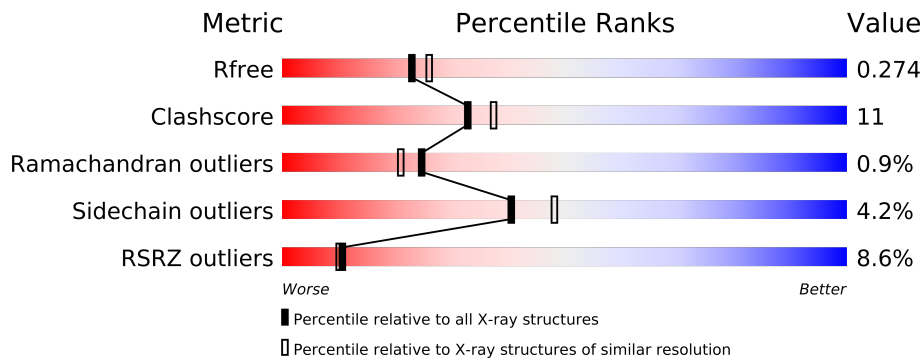
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.20 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	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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	364	

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 2920 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.

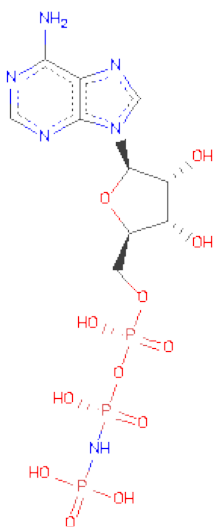
- Molecule 1 is a protein called Mitogen-activated protein kinase 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	348	2800	1800	473	506	21	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	38	MET	-	CLONING ARTIFACT	UNP P53779
A	39	ALA	-	CLONING ARTIFACT	UNP P53779

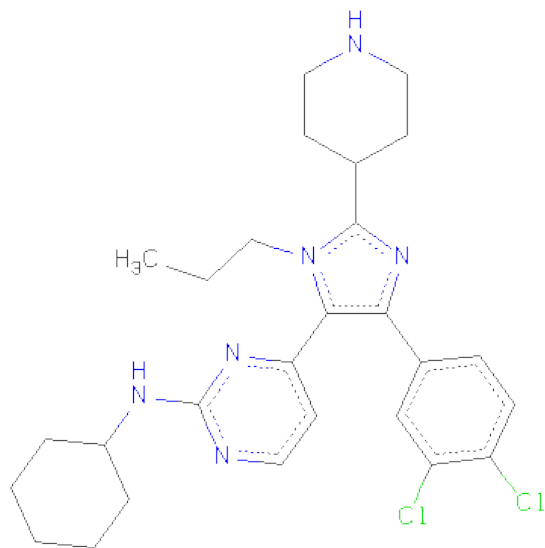
- Molecule 2 is PHOSPHOAMINOPHOSPHONICACID-ADENYLATE ESTER (three-letter code: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	31	10	6	12	3	0	0

- Molecule 3 is CYCLOHEXYL-{4-[5-(3,4-DICHLOROPHENYL)-2-PIPERIDIN-4-YL-3-PR

OPYL-3H-IMIDAZOL-4-YL]-PYRIMIDIN-2-YL}AMINE (three-letter code: 880) (formula: $C_{27}H_{34}Cl_2N_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	Cl	N		
3	A	1	35	27	2	6	0	0

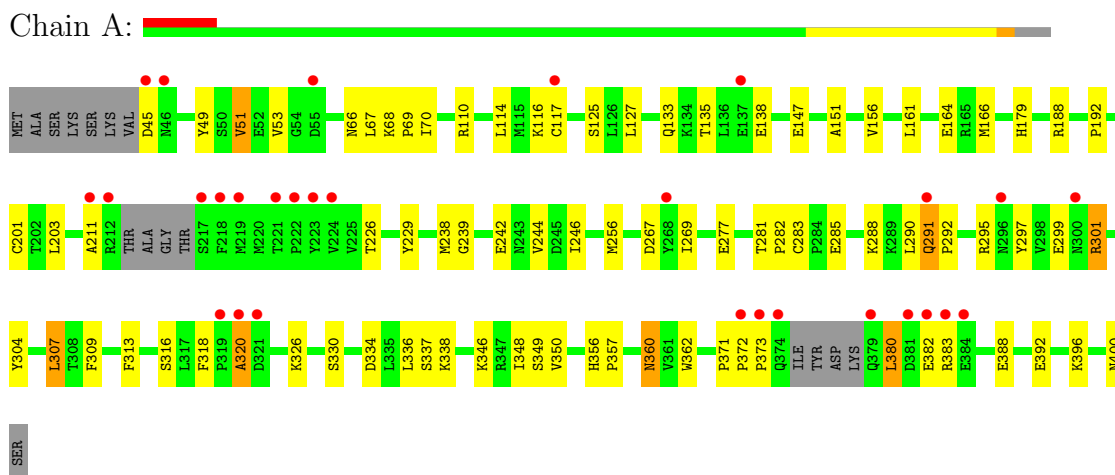
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	54	Total	O	0	0
			54	54		

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: Mitogen-activated protein kinase 10



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	50.95Å 71.53Å 106.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.00 – 2.20 16.94 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.7 (17.00-2.20) 92.9 (16.94-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.14 (at 2.09Å)	Xtrriage
Refinement program	CNX	Depositor
R, R_{free}	0.228 , 0.262 0.229 , 0.274	Depositor DCC
R_{free} test set	1058 reflections (5.12%)	DCC
Wilson B-factor (Å ²)	30.8	Xtrriage
Anisotropy	1.272	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.0	EDS
Estimated twinning fraction	No twinning to report.	Xtrriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Outliers	0 of 21707 reflections	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2920	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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: 880, ANP

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	0.53	0/2864	0.66	0/3879

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 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	2800	0	2801	62	0
2	A	31	0	13	2	0
3	A	35	0	33	4	0
4	A	54	0	0	3	0
All	All	2920	0	2847	64	0

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 11.

All (64) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:110:ARG:CZ	1:A:211:ALA:HB2	1.98	0.93
1:A:291:GLN:HG2	1:A:292:PRO:HD2	1.57	0.83
1:A:151:ALA:HA	3:A:501:880:H572	1.66	0.77
1:A:53:VAL:HG13	1:A:69:PRO:HG3	1.67	0.76
1:A:110:ARG:NE	1:A:211:ALA:HB2	2.09	0.67
1:A:156:VAL:CG1	1:A:161:LEU:HD11	2.31	0.59
1:A:166:MET:HG3	1:A:256:MET:CE	2.33	0.59
1:A:110:ARG:NH2	1:A:211:ALA:HB2	2.17	0.59
1:A:147:GLU:HG2	4:A:624:HOH:O	2.03	0.59
1:A:246:ILE:HD11	1:A:350:VAL:HA	1.86	0.58
1:A:295:ARG:O	1:A:299:GLU:HB2	2.04	0.57
1:A:238:MET:HG2	1:A:239:GLY:N	2.20	0.56
1:A:166:MET:HG3	1:A:256:MET:HE1	1.88	0.56
1:A:320:ALA:HB1	1:A:326:LYS:CG	2.36	0.56
1:A:307:LEU:N	1:A:307:LEU:CD1	2.68	0.55
1:A:290:LEU:HB2	1:A:295:ARG:HB2	1.88	0.55
1:A:283:CYS:SG	1:A:346:LYS:NZ	2.67	0.55
1:A:380:LEU:N	1:A:380:LEU:HD12	2.22	0.55
1:A:291:GLN:HG2	1:A:292:PRO:CD	2.33	0.54
1:A:49:TYR:CE1	1:A:51:VAL:HG13	2.43	0.54
1:A:320:ALA:HB1	1:A:326:LYS:HG2	1.88	0.54
1:A:135:THR:OG1	1:A:138:GLU:HG3	2.07	0.53
1:A:356:HIS:O	1:A:360:ASN:HB2	2.10	0.52
1:A:226:THR:OG1	2:A:502:ANP:H5'2	2.10	0.52
1:A:313:PHE:CE2	1:A:336:LEU:HD13	2.45	0.51
1:A:318:PHE:O	1:A:320:ALA:N	2.39	0.51
1:A:49:TYR:HE1	1:A:51:VAL:HG13	1.75	0.51
3:A:501:880:H062	3:A:501:880:H561	1.92	0.51
1:A:156:VAL:HG12	1:A:161:LEU:HD11	1.92	0.51
1:A:203:LEU:C	1:A:203:LEU:HD23	2.32	0.50
1:A:285:GLU:O	1:A:288:LYS:HB2	2.12	0.50
1:A:179:HIS:CE1	1:A:373:PRO:HD3	2.47	0.49
1:A:288:LYS:HA	1:A:295:ARG:NH1	2.28	0.49
1:A:51:VAL:HG21	1:A:67:LEU:HD22	1.96	0.48
1:A:229:TYR:OH	2:A:502:ANP:N6	2.47	0.48
1:A:164:GLU:HG2	1:A:362:TRP:CZ2	2.48	0.47
1:A:117:CYS:O	1:A:373:PRO:HB2	2.15	0.47
1:A:70:ILE:HG13	3:A:501:880:H101	1.96	0.46
1:A:277:GLU:HG2	1:A:304:TYR:CZ	2.51	0.46
1:A:334:ASP:OD2	1:A:338:LYS:HE2	2.15	0.46
1:A:371:PRO:O	1:A:373:PRO:HD3	2.15	0.46
3:A:501:880:H062	3:A:501:880:N51	2.31	0.45
1:A:392:GLU:HG2	1:A:396:LYS:HE3	1.97	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:192:PRO:HD2	4:A:613:HOH:O	2.17	0.44
1:A:269:ILE:HG23	1:A:297:TYR:OH	2.17	0.44
1:A:309:PHE:CE2	1:A:337:SER:HA	2.52	0.44
1:A:133:GLN:HG2	1:A:138:GLU:O	2.18	0.44
1:A:166:MET:HG3	1:A:256:MET:HE2	2.00	0.44
1:A:307:LEU:HD12	1:A:307:LEU:N	2.32	0.43
1:A:164:GLU:HG2	1:A:362:TRP:HZ2	1.82	0.43
1:A:242:GLU:OE1	1:A:242:GLU:N	2.45	0.43
1:A:238:MET:CG	1:A:239:GLY:N	2.82	0.43
1:A:281:THR:HA	1:A:282:PRO:HD3	1.91	0.43
1:A:114:LEU:C	1:A:116:LYS:H	2.23	0.42
1:A:188:ARG:HG2	1:A:244:VAL:HG11	2.02	0.42
1:A:356:HIS:ND1	1:A:357:PRO:HD2	2.35	0.42
1:A:267:ASP:OD2	1:A:269:ILE:HB	2.20	0.42
1:A:66:ASN:ND2	1:A:68:LYS:HE3	2.35	0.41
1:A:301:ARG:HH11	1:A:301:ARG:HG3	1.85	0.41
1:A:125:SER:HB2	4:A:629:HOH:O	2.21	0.41
1:A:348:ILE:HG12	1:A:349:SER:N	2.36	0.41
1:A:188:ARG:NH2	1:A:211:ALA:O	2.54	0.41
1:A:371:PRO:HA	1:A:372:PRO:HD3	1.87	0.40
1:A:309:PHE:CZ	1:A:337:SER:HA	2.56	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	342/364 (94%)	321 (94%)	18 (5%)	3 (1%)	25 21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	382	GLU
1	A	320	ALA

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Mol	Chain	Res	Type
1	A	383	ARG

5.3.2 Protein sidechains [i](#)

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	309/328 (94%)	296 (96%)	13 (4%)	40 48

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	ASP
1	A	51	VAL
1	A	127	LEU
1	A	201	CYS
1	A	291	GLN
1	A	301	ARG
1	A	307	LEU
1	A	316	SER
1	A	330	SER
1	A	360	ASN
1	A	380	LEU
1	A	388	GLU
1	A	400	ASN

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

Mol	Chain	Res	Type
1	A	65	GLN
1	A	66	ASN
1	A	360	ASN
1	A	400	ASN

5.3.3 RNA [i](#)

There are no RNA chains 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)

2 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
3	880	A	501	-	39,39,39	3.01	11 (28%)	53,54,54	2.83	17 (32%)
2	ANP	A	502	-	33,33,33	3.02	7 (21%)	51,52,52	2.16	14 (27%)

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	880	A	501	-	-	0/13/35/35	0/5/5/5
2	ANP	A	502	-	-	1/18/38/38	0/1/3/3

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	880	C2-C16	-12.10	1.31	1.50
3	A	501	880	C5-N1	8.91	1.52	1.40
2	A	502	ANP	PB-N3B	-8.76	1.56	1.64
2	A	502	ANP	PG-N3B	-7.70	1.57	1.64
2	A	502	ANP	PG-O1G	7.17	1.55	1.46
2	A	502	ANP	PB-O1B	6.02	1.53	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	880	C5-C36	5.04	1.56	1.49
2	A	502	ANP	C4-N3	4.75	1.42	1.35
2	A	502	ANP	O5'-C5'	4.39	1.63	1.44
3	A	501	880	C41-C40	4.30	1.47	1.38
2	A	502	ANP	C5-C4	3.34	1.48	1.40
3	A	501	880	C40-C39	3.27	1.46	1.38
3	A	501	880	C5-C4	2.80	1.49	1.41
3	A	501	880	C60-C55	2.79	1.59	1.51
3	A	501	880	C4-N3	2.57	1.44	1.37
3	A	501	880	C41-C35	2.43	1.44	1.39
3	A	501	880	C38-CL45	-2.32	1.67	1.73
3	A	501	880	C21-C16	2.13	1.59	1.53

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	880	C60-C55-N54	-8.24	97.73	110.55
3	A	501	880	C56-C55-N54	7.81	122.71	110.55
3	A	501	880	C6-N1-C2	-7.13	123.11	127.34
3	A	501	880	C50-N54-C55	6.74	130.59	123.42
2	A	502	ANP	N3-C2-N1	-6.46	123.31	128.71
2	A	502	ANP	C8-N9-C4	5.72	111.27	106.90
3	A	501	880	C16-C2-N3	-5.21	114.28	124.27
2	A	502	ANP	O1G-PG-N3B	-4.61	104.86	111.83
2	A	502	ANP	N3-C4-N9	4.44	133.45	125.43
2	A	502	ANP	PB-N3B-PG	3.94	136.69	130.07
3	A	501	880	C5-C36-N51	3.93	123.60	116.45
3	A	501	880	C36-N51-C50	3.85	118.96	116.44
3	A	501	880	C4-C5-C36	3.81	139.21	130.27
2	A	502	ANP	O4'-C1'-N9	3.60	111.79	108.44
3	A	501	880	C5-N1-C2	-3.60	106.50	109.32
3	A	501	880	C6-N1-C5	3.50	130.37	124.83
3	A	501	880	C4-N3-C2	3.49	110.36	105.87
3	A	501	880	C36-C5-N1	-3.28	115.38	123.93
2	A	502	ANP	C5-C4-N9	-3.24	102.49	107.16
2	A	502	ANP	O2G-PG-O1G	-2.68	106.71	113.60
2	A	502	ANP	O2A-PA-O1A	-2.62	97.57	112.21
3	A	501	880	C17-C16-C2	2.60	114.93	111.44
2	A	502	ANP	PA-O3A-PB	-2.59	122.95	131.81
2	A	502	ANP	C2-N1-C6	2.54	123.35	118.77
3	A	501	880	C47-C36-C5	-2.40	116.17	120.88
2	A	502	ANP	O4'-C1'-C2'	2.38	110.42	106.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	880	C60-C55-C56	2.35	114.95	110.82
3	A	501	880	C7-C6-N1	2.21	117.47	111.67
2	A	502	ANP	C6-C5-C4	-2.15	113.30	117.25
3	A	501	880	C58-C57-C56	2.06	116.03	111.45
2	A	502	ANP	O3G-PG-N3B	2.04	112.13	106.61

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	ANP	O1B-PB-N3B-PG

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 i

6.1 Protein, DNA and RNA chains i

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	348/364 (95%)	0.24	29 (8%) 11 11	26, 44, 68, 84	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	222	PRO	10.2
1	A	223	TYR	7.6
1	A	117	CYS	7.1
1	A	224	VAL	5.6
1	A	221	THR	5.5
1	A	217	SER	4.8
1	A	211	ALA	4.6
1	A	212	ARG	4.6
1	A	321	ASP	4.0
1	A	379	GLN	3.7
1	A	219	MET	3.6
1	A	320	ALA	3.4
1	A	319	PRO	3.4
1	A	382	GLU	3.3
1	A	55	ASP	3.3
1	A	374	GLN	3.1
1	A	373	PRO	3.0
1	A	384	GLU	3.0
1	A	372	PRO	2.9
1	A	268	TYR	2.9
1	A	383	ARG	2.7
1	A	300	ASN	2.6
1	A	45	ASP	2.5
1	A	381	ASP	2.5
1	A	296	ASN	2.5
1	A	137	GLU	2.3
1	A	291	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	218	PHE	2.1
1	A	46	ASN	2.1

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ANP	A	502	31/31	0.28	1.08	40,46,49,51	31
3	880	A	501	35/35	0.14	0.36	27,32,37,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.