



Full wwPDB X-ray Structure Validation Report

Mar 8, 2026 – 06:55 AM UTC

PDB ID : 2EB3 / pdb_00002eb3
Title : Crystal structure of mutated EGFR kinase domain (L858R) in complex with AMPPNP
Authors : Yoshikawa, S.; Kukimoto-Niino, M.; Shirouzu, M.; Senba, K.; Yamamoto, T.; Yokoyama, S.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2007-02-06
Resolution : 2.84 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)
Xtrriage (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

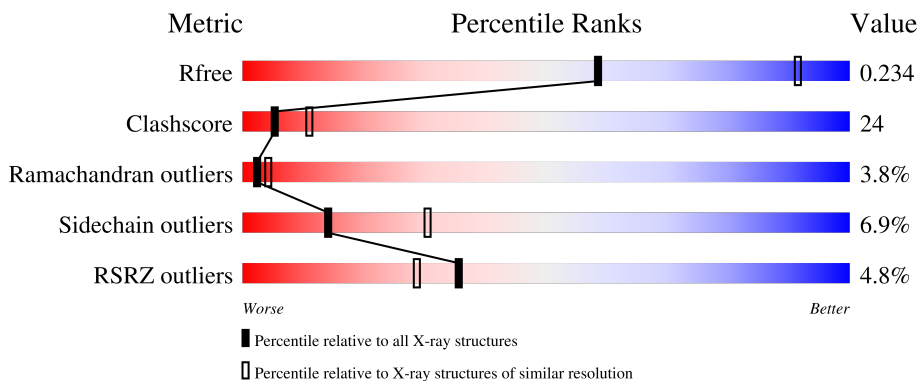
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	334	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 2397 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.

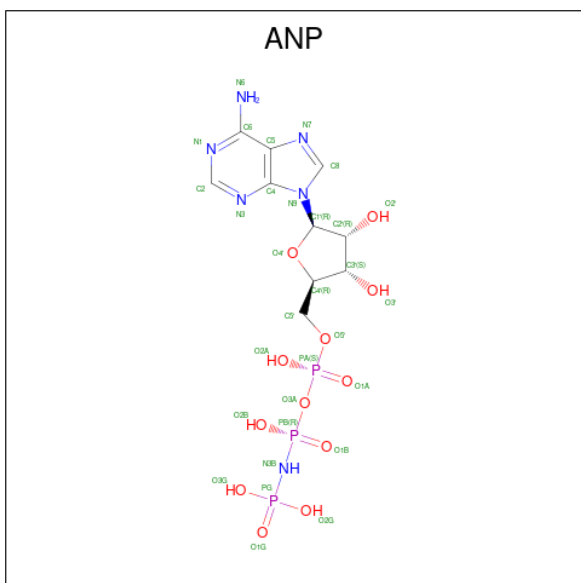
- Molecule 1 is a protein called Epidermal growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	294	2365	1520	400	428	17	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	689	GLY	-	expression tag	UNP P00533
A	690	ALA	-	expression tag	UNP P00533
A	691	MET	-	expression tag	UNP P00533
A	692	GLY	-	expression tag	UNP P00533
A	693	ILE	-	expression tag	UNP P00533
A	694	ARG	-	expression tag	UNP P00533
A	858	ARG	LEU	SEE REMARK 999	UNP P00533

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	144.69Å 144.69Å 144.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.17 – 2.84 36.17 – 2.84	Depositor EDS
% Data completeness (in resolution range)	97.4 (36.17-2.84) 97.3 (36.17-2.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.77 (at 2.86Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.190 , 0.236 0.188 , 0.234	Depositor DCC
R_{free} test set	1248 reflections (10.36%)	wwPDB-VP
Wilson B-factor (Å ²)	65.9	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 62.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.035 for -l,-k,-h	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2397	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: 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	0.59	0/2414	1.09	17/3265 (0.5%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	939	CYS	N-CA-C	8.60	122.64	108.96
1	A	835	HIS	N-CA-C	-7.41	99.79	110.24
1	A	834	VAL	N-CA-C	6.91	117.44	106.32
1	A	974	ASP	CA-C-N	6.02	126.19	119.32
1	A	974	ASP	C-N-CA	6.02	126.19	119.32
1	A	895	SER	N-CA-C	-5.79	105.72	112.89
1	A	881	MET	N-CA-C	5.73	119.59	109.96
1	A	909	THR	N-CA-C	-5.46	106.40	112.57
1	A	896	ASP	N-CA-C	-5.35	105.53	111.36
1	A	724	GLY	N-CA-C	-5.19	100.88	113.18
1	A	765	VAL	N-CA-C	5.15	115.30	110.30
1	A	882	ALA	N-CA-C	-5.14	104.08	110.41
1	A	740	ILE	N-CA-C	5.14	115.83	107.71
1	A	933	LEU	N-CA-C	-5.14	102.74	110.24
1	A	900	TYR	N-CA-C	-5.12	105.59	111.07
1	A	935	GLN	CA-C-N	-5.07	115.16	120.38
1	A	935	GLN	C-N-CA	-5.07	115.16	120.38

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	2365	0	2409	115	0
2	A	31	0	13	3	0
3	A	1	0	0	0	0
All	All	2397	0	2422	117	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 24.

All (117) 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:1005:GLU:HB3	1:A:1007:MET:HE3	1.49	0.91
1:A:952:MET:HE2	1:A:952:MET:HA	1.67	0.76
1:A:932:ARG:NH1	1:A:951:TRP:O	2.19	0.75
1:A:773:HIS:CE1	1:A:820:GLN:HG2	2.21	0.75
1:A:835:HIS:CD2	1:A:856:PHE:HB3	2.23	0.74
1:A:713:LYS:HE3	1:A:715:ILE:HD11	1.71	0.73
1:A:720:SER:O	1:A:722:ALA:HB3	1.89	0.72
1:A:825:MET:HE2	1:A:838:LEU:HD22	1.70	0.71
1:A:747:LEU:HD21	1:A:759:ILE:HG21	1.73	0.70
1:A:893:HIS:O	1:A:896:ASP:HB2	1.92	0.69
1:A:848:PRO:HD2	1:A:849:GLN:HE22	1.58	0.69
1:A:756:ASN:O	1:A:760:LEU:HD23	1.95	0.67
1:A:748:ARG:C	1:A:750:ALA:H	2.03	0.67
1:A:858:ARG:O	1:A:858:ARG:HG2	1.96	0.64
1:A:931:GLU:O	1:A:932:ARG:NE	2.27	0.64
1:A:747:LEU:HD21	1:A:759:ILE:CG2	2.28	0.64
1:A:888:HIS:O	1:A:890:ILE:HG13	1.98	0.64
1:A:877:PRO:O	1:A:881:MET:HG3	1.99	0.62
1:A:748:ARG:O	1:A:750:ALA:N	2.33	0.62
1:A:882:ALA:HB1	1:A:884:GLU:OE2	2.00	0.61
1:A:985:GLU:H	1:A:985:GLU:CD	2.08	0.61
1:A:791:GLN:NE2	1:A:1012:ASP:OD2	2.34	0.61
1:A:1006:ASP:O	1:A:1007:MET:HE2	2.01	0.60
1:A:857:GLY:C	1:A:859:ALA:H	2.10	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:834:VAL:HG12	1:A:836:ARG:HG3	1.83	0.59
1:A:825:MET:SD	1:A:853:ILE:HD13	2.42	0.59
1:A:707:LEU:HD13	1:A:779:GLY:HA3	1.85	0.59
1:A:829:GLU:HG3	1:A:893:HIS:CD2	2.39	0.58
1:A:707:LEU:CD1	1:A:779:GLY:HA3	2.34	0.58
1:A:856:PHE:O	1:A:859:ALA:HB3	2.05	0.57
1:A:960:LYS:HZ2	1:A:960:LYS:HA	1.70	0.57
1:A:960:LYS:HA	1:A:960:LYS:NZ	2.20	0.56
1:A:733:PRO:O	1:A:734:GLU:C	2.49	0.56
1:A:759:ILE:HD11	1:A:788:LEU:HG	1.88	0.56
1:A:714:LYS:HD3	1:A:727:TYR:CD2	2.40	0.56
1:A:748:ARG:HG2	1:A:749:GLU:N	2.21	0.56
1:A:923:ILE:H	1:A:923:ILE:HD12	1.71	0.55
1:A:748:ARG:C	1:A:750:ALA:N	2.63	0.55
1:A:932:ARG:HH11	1:A:951:TRP:HB3	1.72	0.55
1:A:879:LYS:HD3	1:A:914:PRO:O	2.06	0.55
1:A:834:VAL:O	1:A:836:ARG:HG3	2.07	0.55
1:A:1006:ASP:C	1:A:1007:MET:HE2	2.32	0.54
1:A:726:VAL:HG22	1:A:745:LYS:HG3	1.88	0.54
1:A:888:HIS:HB2	1:A:890:ILE:HG13	1.90	0.54
1:A:833:LEU:HD12	1:A:861:LEU:HA	1.89	0.53
1:A:751:THR:O	1:A:751:THR:HG23	2.07	0.53
1:A:752:SER:O	1:A:754:LYS:N	2.42	0.53
1:A:759:ILE:HG23	1:A:760:LEU:HD22	1.91	0.52
1:A:805:HIS:O	1:A:806:LYS:C	2.52	0.52
1:A:857:GLY:C	1:A:859:ALA:N	2.67	0.52
1:A:780:ILE:HG22	1:A:788:LEU:HD22	1.92	0.52
1:A:783:THR:O	1:A:784:SER:C	2.53	0.51
1:A:882:ALA:HA	1:A:898:TRP:CD2	2.45	0.51
1:A:961:PHE:O	1:A:965:ILE:HG13	2.11	0.51
1:A:707:LEU:HD22	1:A:789:ILE:HG13	1.92	0.51
1:A:835:HIS:CE1	1:A:837:ASP:C	2.89	0.51
1:A:831:ARG:HH11	1:A:831:ARG:HG3	1.75	0.50
1:A:848:PRO:HD2	1:A:849:GLN:NE2	2.25	0.50
1:A:920:ALA:HA	1:A:923:ILE:HD13	1.92	0.50
1:A:830:ASP:C	1:A:832:ARG:H	2.20	0.49
1:A:740:ILE:O	1:A:742:VAL:HG13	2.12	0.49
1:A:935:GLN:HG3	1:A:944:TYR:CG	2.48	0.49
1:A:905:TRP:HA	1:A:947:MET:HE3	1.95	0.49
1:A:856:PHE:HB2	1:A:859:ALA:HB2	1.94	0.49
1:A:829:GLU:HA	1:A:893:HIS:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:825:MET:CE	1:A:838:LEU:HD22	2.39	0.48
2:A:2001:ANP:O1G	2:A:2001:ANP:O2A	2.32	0.48
1:A:792:LEU:O	1:A:794:PRO:HD3	2.14	0.48
1:A:835:HIS:O	1:A:896:ASP:OD2	2.32	0.48
1:A:745:LYS:HE3	2:A:2001:ANP:O1A	2.14	0.47
1:A:726:VAL:CG2	1:A:745:LYS:HG3	2.45	0.47
1:A:825:MET:HE2	1:A:838:LEU:CD2	2.41	0.47
1:A:808:ASN:ND2	1:A:988:HIS:O	2.40	0.47
2:A:2001:ANP:O5'	2:A:2001:ANP:H2'	2.15	0.46
1:A:806:LYS:O	1:A:808:ASN:N	2.49	0.46
1:A:795:PHE:HB2	1:A:845:VAL:HB	1.97	0.46
1:A:857:GLY:O	1:A:859:ALA:N	2.48	0.46
1:A:902:VAL:O	1:A:906:GLU:HG3	2.16	0.45
1:A:887:LEU:HB3	1:A:888:HIS:CE1	2.52	0.45
1:A:764:TYR:CD1	1:A:764:TYR:C	2.94	0.45
1:A:713:LYS:O	1:A:713:LYS:HG3	2.17	0.45
1:A:734:GLU:HB3	1:A:735:GLY:H	1.65	0.45
1:A:949:LYS:O	1:A:952:MET:HG2	2.17	0.44
1:A:888:HIS:O	1:A:889:ARG:C	2.60	0.44
1:A:793:MET:HE1	1:A:846:LYS:HB2	2.00	0.44
1:A:778:LEU:HB2	1:A:789:ILE:O	2.18	0.44
1:A:766:MET:HB3	1:A:777:LEU:HB2	1.99	0.44
1:A:750:ALA:O	1:A:751:THR:HB	2.19	0.43
1:A:884:GLU:H	1:A:884:GLU:CD	2.19	0.43
1:A:974:ASP:N	1:A:975:PRO:HD3	2.33	0.43
1:A:849:GLN:NE2	1:A:849:GLN:H	2.17	0.43
1:A:962:ARG:O	1:A:966:ILE:HG13	2.19	0.43
1:A:747:LEU:O	1:A:748:ARG:C	2.59	0.43
1:A:878:ILE:HA	1:A:881:MET:SD	2.59	0.43
1:A:878:ILE:HG21	1:A:920:ALA:HB1	2.00	0.42
1:A:918:ILE:HG21	1:A:926:ILE:HD12	2.00	0.42
1:A:924:SER:O	1:A:928:GLU:HG3	2.19	0.42
1:A:752:SER:C	1:A:754:LYS:N	2.78	0.42
1:A:894:GLN:OE1	1:A:960:LYS:NZ	2.50	0.42
1:A:923:ILE:O	1:A:924:SER:C	2.63	0.42
1:A:897:VAL:O	1:A:900:TYR:HB3	2.19	0.42
1:A:983:GLY:O	1:A:984:ASP:C	2.63	0.42
1:A:829:GLU:HG3	1:A:893:HIS:CG	2.55	0.41
1:A:968:PHE:HA	1:A:971:MET:HE3	2.01	0.41
1:A:834:VAL:HG12	1:A:834:VAL:O	2.20	0.41
1:A:894:GLN:O	1:A:897:VAL:HB	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:745:LYS:O	1:A:787:GLN:HA	2.21	0.41
1:A:831:ARG:O	1:A:832:ARG:C	2.64	0.41
1:A:736:GLU:O	1:A:737:LYS:C	2.64	0.41
1:A:954:ASP:OD2	1:A:954:ASP:C	2.64	0.41
1:A:932:ARG:HD3	1:A:932:ARG:HA	1.86	0.41
1:A:713:LYS:CE	1:A:715:ILE:HD11	2.46	0.41
1:A:780:ILE:CG2	1:A:788:LEU:HD22	2.51	0.40
1:A:905:TRP:HD1	1:A:947:MET:CE	2.35	0.40
1:A:923:ILE:HD12	1:A:923:ILE:N	2.34	0.40
1:A:905:TRP:CZ2	1:A:909:THR:HG21	2.57	0.40
1:A:905:TRP:CA	1:A:947:MET:HE3	2.51	0.40

There are no symmetry-related clashes.

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	286/334 (86%)	248 (87%)	27 (9%)	11 (4%)	2 4

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	734	GLU
1	A	751	THR
1	A	749	GLU
1	A	807	ASP
1	A	889	ARG
1	A	858	ARG
1	A	753	PRO
1	A	806	LYS
1	A	923	ILE
1	A	926	ILE

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Mol	Chain	Res	Type
1	A	733	PRO

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	262/291 (90%)	244 (93%)	18 (7%)	14 30

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	703	LEU
1	A	704	LEU
1	A	718	LEU
1	A	747	LEU
1	A	753	PRO
1	A	808	ASN
1	A	884	GLU
1	A	888	HIS
1	A	891	TYR
1	A	921	SER
1	A	932	ARG
1	A	937	PRO
1	A	941	ILE
1	A	953	ILE
1	A	960	LYS
1	A	962	ARG
1	A	964	LEU
1	A	985	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	773	HIS
1	A	816	ASN

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Mol	Chain	Res	Type
1	A	826	ASN
1	A	849	GLN
1	A	893	HIS

5.3.3 RNA [i](#)

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand 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
2	ANP	A	2001	-	33,33,33	1.22	5 (15%)	45,52,52	1.71	7 (15%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ANP	A	2001	-	-	5/18/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2001	ANP	C5-N7	-3.11	1.33	1.39
2	A	2001	ANP	C8-N9	-2.34	1.33	1.37
2	A	2001	ANP	PA-O3A	2.08	1.61	1.59
2	A	2001	ANP	PB-O3A	2.06	1.61	1.59
2	A	2001	ANP	C4-N9	-2.01	1.33	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2001	ANP	C5-C4-N3	-5.68	118.89	126.72
2	A	2001	ANP	N3-C2-N1	-4.24	122.16	128.58
2	A	2001	ANP	N3-C4-N9	3.89	133.78	127.17
2	A	2001	ANP	C2-N3-C4	3.57	120.54	111.83
2	A	2001	ANP	N9-C8-N7	-2.79	109.98	113.94
2	A	2001	ANP	C4-C5-N7	-2.78	107.40	110.58
2	A	2001	ANP	C5-N7-C8	2.67	107.65	103.45

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2001	ANP	PB-N3B-PG-O1G
2	A	2001	ANP	PG-N3B-PB-O1B
2	A	2001	ANP	PB-O3A-PA-O1A
2	A	2001	ANP	PA-O3A-PB-O2B
2	A	2001	ANP	PA-O3A-PB-O1B

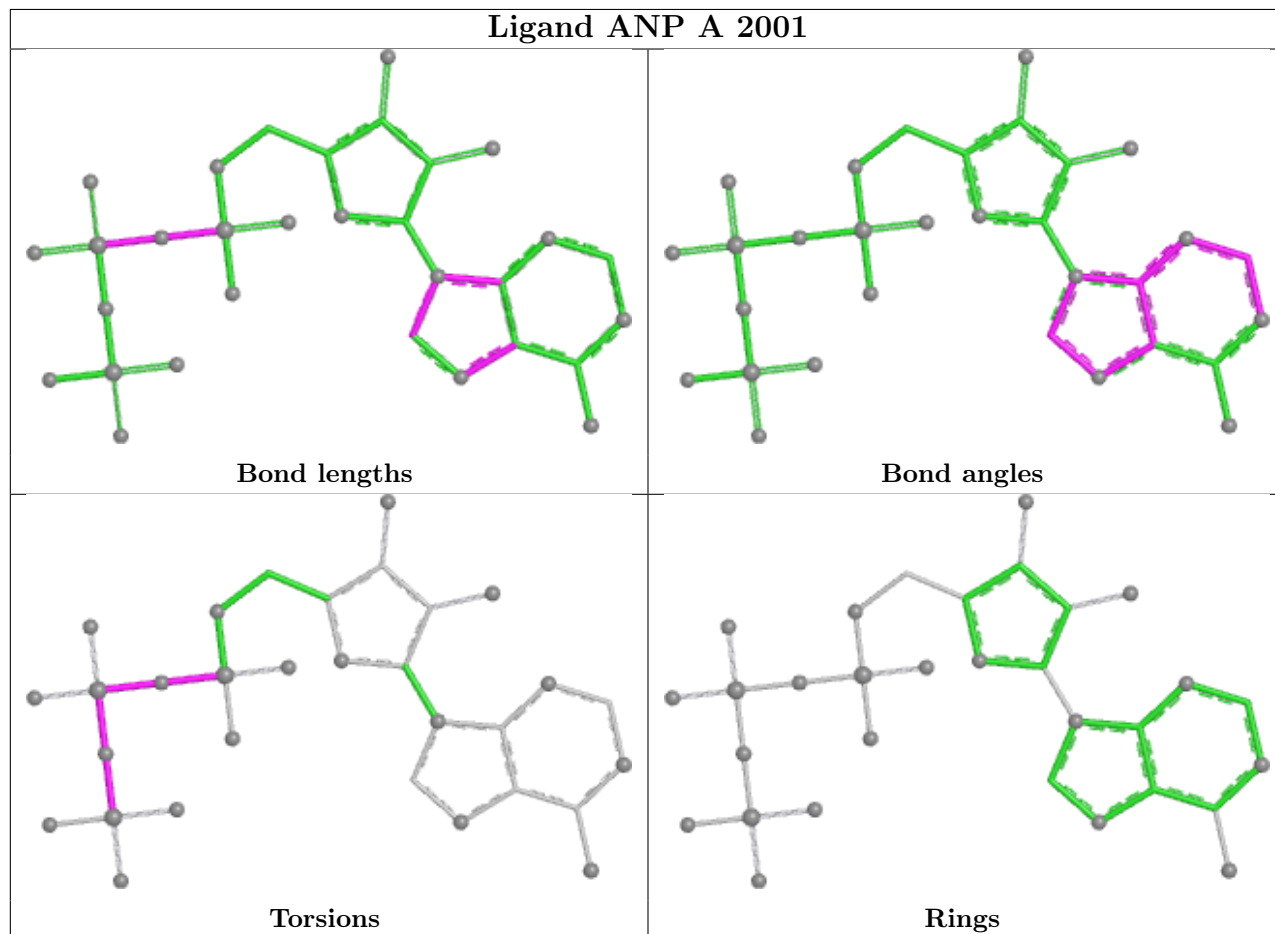
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	2001	ANP	3	0

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 [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	294/334 (88%)	-0.20	14 (4%) 35 29	29, 58, 119, 175	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1018	ILE	3.8
1	A	723	PHE	3.6
1	A	861	LEU	3.5
1	A	990	PRO	3.5
1	A	753	PRO	3.4
1	A	750	ALA	3.0
1	A	857	GLY	2.9
1	A	722	ALA	2.8
1	A	720	SER	2.7
1	A	747	LEU	2.5
1	A	858	ARG	2.3
1	A	860	LYS	2.1
1	A	751	THR	2.1
1	A	724	GLY	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](#)

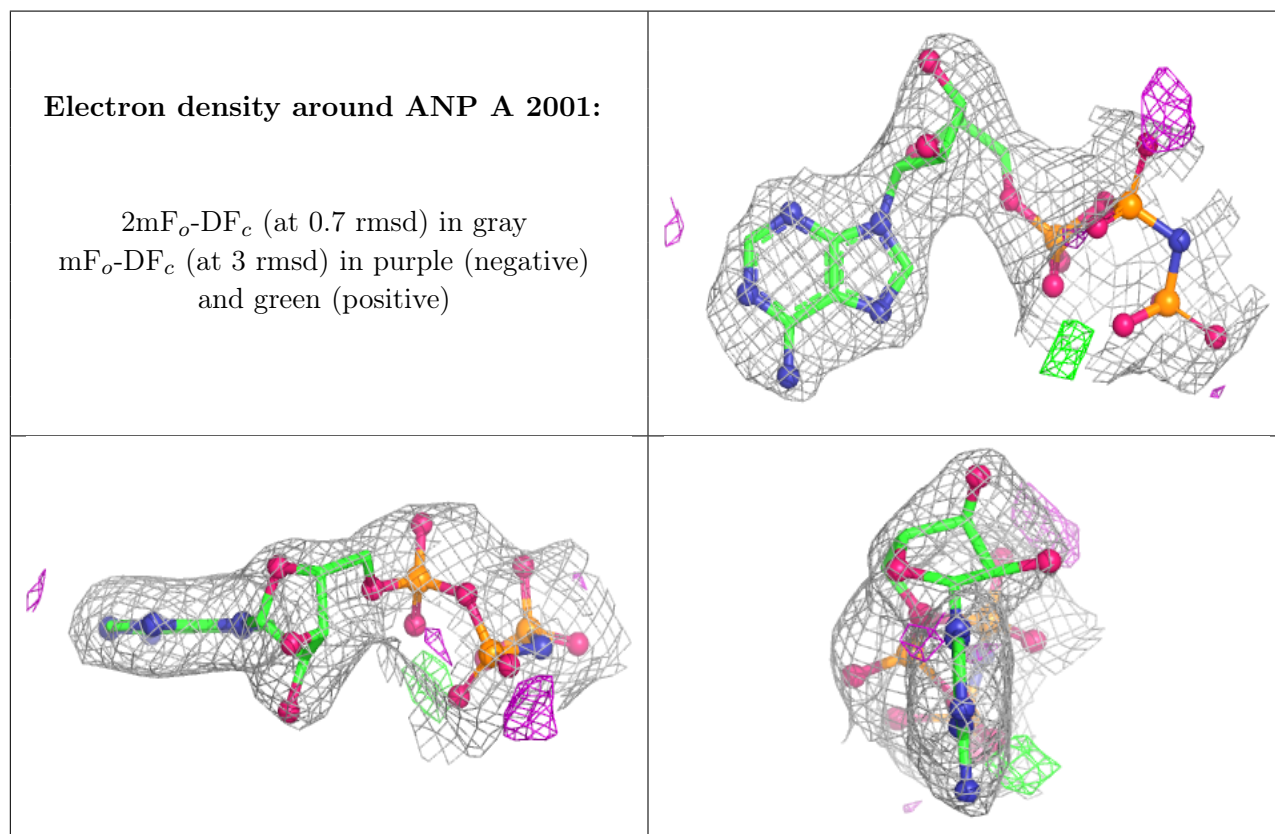
There are no oligosaccharides in this entry.

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
2	ANP	A	2001	31/31	0.90	0.09	44,92,124,132	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.