



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:50 PM UTC

PDB ID : 3X03 / pdb_00003x03
Title : Crystal structure of PIP4KIIBETA complex with AMPPNP
Authors : Takeuchi, K.; Lo, Y.H.; Sumita, K.; Senda, M.; Terakawa, J.; Dimitoris, A.; Locasale, J.W.; Sasaki, M.; Yoshino, H.; Zhang, Y.; Kahoud, E.R.; Takano, T.; Yokota, T.; Emerling, B.; Asara, J.A.; Ishida, T.; Shimada, I.; Daikoku, T.; Cantley, L.C.; Senda, T.; Sasaki, A.T.
Deposited on : 2014-10-09
Resolution : 2.70 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](#) ①) 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

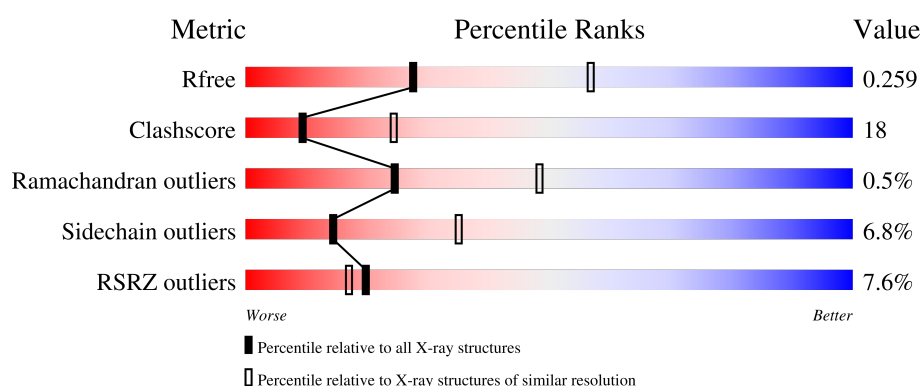
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	393	6% 54% 25% • 19%
1	B	393	6% 46% 28% • 23%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5237 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 Phosphatidylinositol 5-phosphate 4-kinase type-2 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	0	0
			2629	1672	448	495	14			
1	B	302	Total	C	N	O	S	0	0	0
			2490	1595	426	456	13			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	GLY	-	expression tag	UNP P78356
A	25	PRO	-	expression tag	UNP P78356
A	26	ASN	-	expression tag	UNP P78356
A	27	CYS	-	expression tag	UNP P78356
A	28	ALA	-	expression tag	UNP P78356
A	29	PRO	-	expression tag	UNP P78356
A	30	GLY	-	expression tag	UNP P78356
B	24	GLY	-	expression tag	UNP P78356
B	25	PRO	-	expression tag	UNP P78356
B	26	ASN	-	expression tag	UNP P78356
B	27	CYS	-	expression tag	UNP P78356
B	28	ALA	-	expression tag	UNP P78356
B	29	PRO	-	expression tag	UNP P78356
B	30	GLY	-	expression tag	UNP P78356

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



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

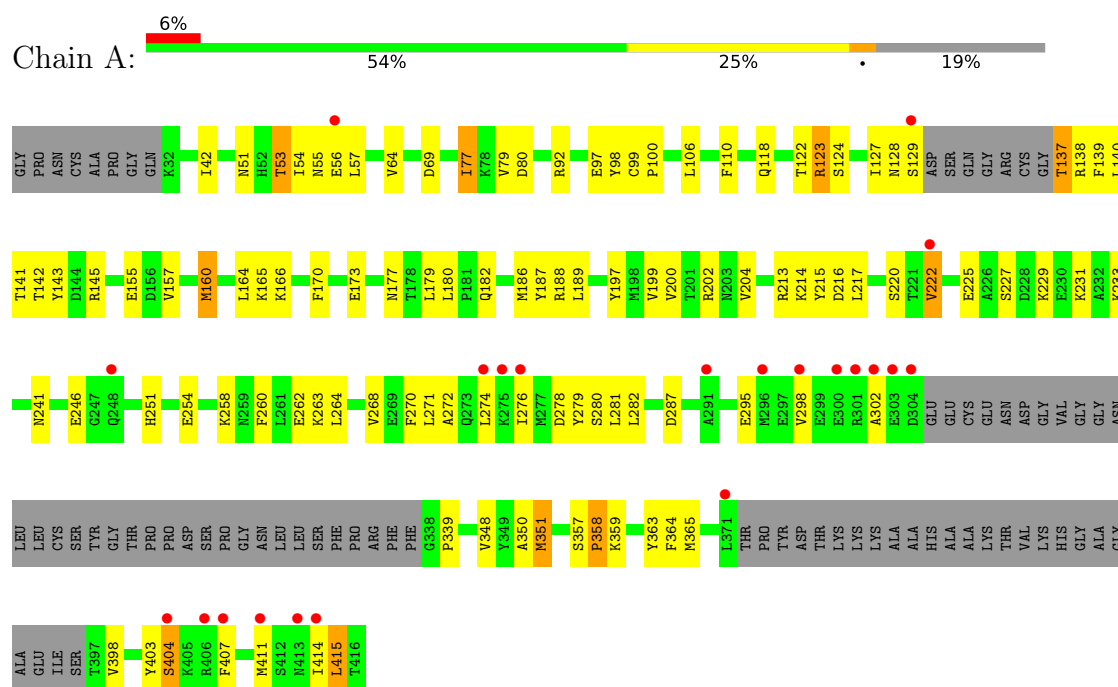
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	11	Total O 11 11	0	0
3	B	14	Total O 14 14	0	0

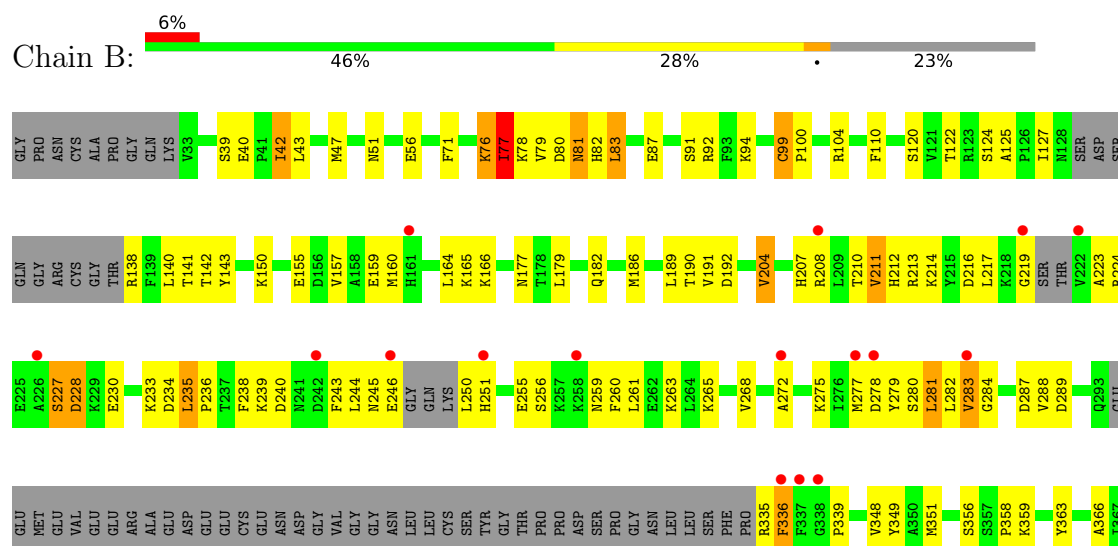
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Phosphatidylinositol 5-phosphate 4-kinase type-2 beta



- Molecule 1: Phosphatidylinositol 5-phosphate 4-kinase type-2 beta



I368	D389	L371	T372	PRO	TYR	ASP	THR	LYS	LYS	ALA	ALA	HIS	ALA	ALA	LYS	THR	VAL	LYS	HIS	GLY	ALA	GLY	ALA	GLU	ILE	SER	THR	VAL	ASN	F400	E401	Q402	Y403	S404	F405	F406	F407	F410	M411	I414	L415	T416
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4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	108.64Å 182.98Å 106.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	53.18 – 2.70 53.18 – 2.70	Depositor EDS
% Data completeness (in resolution range)	98.6 (53.18-2.70) 98.7 (53.18-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.3_928)	Depositor
R, R_{free}	0.234 , 0.280 0.218 , 0.259	Depositor DCC
R_{free} test set	1451 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	75.6	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 76.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.005 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.012 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5237	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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 ⓘ

5.1 Standard geometry ⓘ

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	0.96	3/2682 (0.1%)	1.10	5/3608 (0.1%)
1	B	1.06	12/2542 (0.5%)	1.14	11/3418 (0.3%)
All	All	1.01	15/5224 (0.3%)	1.12	16/7026 (0.2%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	99	CYS	CA-C	-10.03	1.47	1.53
1	B	207	HIS	C-N	-8.51	1.22	1.33
1	B	208	ARG	N-CA	7.64	1.55	1.46
1	B	401	GLU	C-N	-7.41	1.24	1.33
1	B	83	LEU	N-CA	6.92	1.56	1.46
1	B	82	HIS	C-N	-6.91	1.23	1.33
1	B	244	LEU	C-N	-6.61	1.25	1.34
1	B	245	ASN	C-N	6.22	1.42	1.33
1	A	251	HIS	CA-C	-6.16	1.45	1.52
1	B	207	HIS	CG-ND1	-6.06	1.31	1.38
1	B	81	ASN	C-O	6.04	1.31	1.24
1	B	77	ILE	CA-CB	5.89	1.62	1.54
1	A	271	LEU	C-N	-5.54	1.26	1.33
1	A	200	VAL	CA-C	5.25	1.59	1.52
1	B	283	VAL	CA-C	-5.10	1.46	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	271	LEU	O-C-N	8.84	131.49	122.12
1	A	271	LEU	CA-C-O	-7.61	112.48	120.55
1	B	401	GLU	CA-C-N	6.64	129.07	120.44
1	B	401	GLU	C-N-CA	6.64	129.07	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	254	GLU	N-CA-C	6.53	118.99	111.02
1	B	235	LEU	CA-C-N	6.21	126.15	119.76
1	B	235	LEU	C-N-CA	6.21	126.15	119.76
1	B	401	GLU	O-C-N	-6.15	115.60	122.12
1	A	53	THR	N-CA-C	5.82	117.29	111.07
1	B	402	GLN	O-C-N	5.67	127.91	122.07
1	B	179	LEU	N-CA-C	-5.57	105.83	113.30
1	B	207	HIS	CA-C-O	5.50	125.73	119.18
1	B	208	ARG	N-CA-CB	-5.06	100.81	110.63
1	B	204	VAL	CB-CA-C	-5.05	105.47	112.24
1	A	214	LYS	N-CA-C	5.01	117.00	109.23
1	B	164	LEU	N-CA-C	5.01	116.74	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	2629	0	2601	90	0
1	B	2490	0	2472	94	0
2	A	62	0	22	10	0
2	B	31	0	11	2	0
3	A	11	0	0	0	0
3	B	14	0	0	2	0
All	All	5237	0	5106	186	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 18.

All (186) 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:281:LEU:HD22	1:A:411:MET:SD	1.87	1.12
1:B:411:MET:O	1:B:414:ILE:HG12	1.51	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:LEU:CD2	1:A:411:MET:SD	2.47	1.03
1:A:56:GLU:HG3	1:A:57:LEU:N	1.75	1.02
1:B:251:HIS:ND1	1:B:416:THR:HG21	1.78	0.98
1:A:217:LEU:HB2	1:A:411:MET:HE1	1.47	0.94
1:A:53:THR:O	1:A:56:GLU:HG2	1.66	0.93
1:B:281:LEU:HD22	1:B:411:MET:SD	2.10	0.92
2:A:501:ANP:PG	2:A:501:ANP:O1A	2.30	0.89
1:B:281:LEU:HB2	1:B:407:PHE:HZ	1.43	0.83
1:A:160:MET:HE1	1:A:164:LEU:HD13	1.63	0.81
1:B:251:HIS:ND1	1:B:416:THR:CG2	2.44	0.81
1:A:407:PHE:CE2	1:A:411:MET:HG3	2.16	0.80
1:B:335:ARG:HG3	1:B:336:PHE:N	1.96	0.79
1:A:56:GLU:HG3	1:A:57:LEU:H	1.47	0.79
1:A:217:LEU:CB	1:A:411:MET:HE1	2.15	0.76
1:A:53:THR:O	1:A:56:GLU:CG	2.34	0.75
1:A:128:ASN:HB3	1:A:140:LEU:HD23	1.69	0.75
1:B:411:MET:O	1:B:414:ILE:CG1	2.32	0.74
1:A:357:SER:OG	1:A:358:PRO:HD2	1.87	0.74
1:A:77:ILE:HG13	1:B:77:ILE:HD11	1.72	0.72
1:B:268:VAL:CG1	1:B:404:SER:HB2	2.21	0.71
1:A:415:LEU:HD23	1:A:415:LEU:O	1.91	0.70
1:A:258:LYS:NZ	1:A:262:GLU:OE2	2.24	0.70
1:A:53:THR:HA	1:A:56:GLU:HG2	1.73	0.69
1:A:268:VAL:HG12	1:A:404:SER:HB2	1.76	0.68
1:B:99:CYS:N	1:B:100:PRO:CD	2.57	0.68
1:A:281:LEU:HD23	1:A:411:MET:SD	2.35	0.65
1:A:56:GLU:CG	1:A:57:LEU:N	2.56	0.65
1:A:141:THR:CG2	1:A:145:ARG:HA	2.27	0.64
1:B:47:MET:HE3	1:B:125:ALA:HA	1.79	0.64
1:B:210:THR:HG22	1:B:211:VAL:N	2.12	0.64
1:B:217:LEU:CB	1:B:411:MET:HE1	2.28	0.63
1:A:53:THR:C	1:A:56:GLU:HG2	2.24	0.62
1:B:268:VAL:HG12	1:B:404:SER:HB2	1.81	0.62
1:B:212:HIS:O	1:B:213:ARG:HG2	2.00	0.61
1:B:94:LYS:HB2	1:B:190:THR:HB	1.83	0.60
1:B:219:GLY:N	1:B:240:ASP:OD2	2.35	0.59
1:B:47:MET:CE	1:B:125:ALA:HA	2.31	0.59
1:B:228:ASP:N	1:B:228:ASP:OD1	2.33	0.59
1:B:281:LEU:HB2	1:B:407:PHE:CZ	2.33	0.58
1:A:155:GLU:N	1:A:155:GLU:OE1	2.33	0.58
1:A:188:ARG:NH1	2:A:502:ANP:O4'	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:VAL:HG23	2:A:501:ANP:N6	2.19	0.58
1:B:279:TYR:O	1:B:403:TYR:OH	2.22	0.58
1:B:287:ASP:OD2	1:B:359:LYS:HD3	2.04	0.58
1:A:51:ASN:HB2	1:A:122:THR:HG21	1.86	0.57
1:A:281:LEU:CB	1:A:411:MET:SD	2.92	0.57
1:B:78:LYS:HD2	1:B:94:LYS:HE2	1.85	0.57
1:A:270:PHE:CZ	1:A:274:LEU:HD11	2.39	0.57
1:B:277:MET:HG3	1:B:278:ASP:OD2	2.06	0.56
1:A:69:ASP:HB3	1:B:83:LEU:HD13	1.87	0.56
1:B:240:ASP:HB3	1:B:410:PHE:CZ	2.40	0.56
1:A:53:THR:CA	1:A:56:GLU:HG2	2.36	0.55
1:A:268:VAL:HG13	1:A:279:TYR:OH	2.06	0.55
2:A:501:ANP:O1A	2:A:501:ANP:N3B	2.30	0.55
1:B:217:LEU:HB2	1:B:411:MET:HE1	1.88	0.55
1:A:264:LEU:HD11	1:A:407:PHE:HE2	1.72	0.55
1:A:216:ASP:C	1:A:217:LEU:HD23	2.31	0.55
1:A:229:LYS:HG2	1:A:233:LYS:NZ	2.22	0.55
1:B:288:VAL:CG1	1:B:339:PRO:HB2	2.37	0.55
1:B:236:PRO:HB2	1:B:238:PHE:CZ	2.42	0.54
1:B:217:LEU:HB3	1:B:411:MET:HE1	1.89	0.54
1:B:250:LEU:HD22	1:B:363:TYR:CD2	2.43	0.53
1:A:350:ALA:HB2	1:A:364:PHE:CE2	2.43	0.53
1:A:287:ASP:OD2	1:A:359:LYS:HD3	2.08	0.53
1:B:336:PHE:O	1:B:336:PHE:CD2	2.62	0.53
1:B:277:MET:O	1:B:278:ASP:HB2	2.09	0.53
1:A:229:LYS:HG2	1:A:233:LYS:HZ2	1.74	0.53
1:B:160:MET:CB	1:B:186:MET:HE1	2.38	0.53
1:B:251:HIS:CE1	1:B:416:THR:HG21	2.44	0.53
1:B:43:LEU:HD21	1:B:140:LEU:HD11	1.91	0.53
1:A:157:VAL:HG21	1:A:197:TYR:CD2	2.43	0.52
1:B:40:GLU:HB3	1:B:42:ILE:HG22	1.90	0.52
1:B:177:ASN:ND2	1:B:263:LYS:HD2	2.24	0.52
1:B:204:VAL:HG23	2:B:501:ANP:N6	2.24	0.52
1:B:282:LEU:HB2	1:B:368:ILE:HD13	1.90	0.52
1:A:186:MET:HG3	1:A:199:VAL:HG22	1.92	0.52
1:B:371:LEU:C	1:B:372:THR:HG23	2.34	0.52
1:A:80:ASP:O	1:B:76:LYS:HB2	2.09	0.52
1:A:295:GLU:O	1:A:295:GLU:CD	2.53	0.52
1:A:129:SER:HG	1:A:139:PHE:HD1	1.57	0.52
1:B:411:MET:C	1:B:414:ILE:HG12	2.30	0.51
1:A:56:GLU:CG	1:A:57:LEU:H	2.17	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:VAL:HG12	1:B:284:GLY:N	2.25	0.51
1:B:191:VAL:O	1:B:192:ASP:HB2	2.10	0.51
2:A:502:ANP:C3'	2:A:502:ANP:PA	2.99	0.50
1:A:77:ILE:HD12	3:B:614:HOH:O	2.11	0.50
1:B:236:PRO:CB	1:B:238:PHE:CZ	2.94	0.50
2:A:501:ANP:O1A	2:A:501:ANP:O3G	2.30	0.50
2:A:501:ANP:O1G	2:A:501:ANP:O1B	2.30	0.50
1:B:233:LYS:HG2	1:B:234:ASP:N	2.26	0.50
1:B:160:MET:HB2	1:B:186:MET:HE1	1.93	0.49
1:A:216:ASP:O	1:A:217:LEU:HD23	2.12	0.49
1:B:214:LYS:HD2	1:B:235:LEU:CD1	2.43	0.49
1:B:43:LEU:HD21	1:B:140:LEU:CD1	2.43	0.49
1:B:120:SER:O	1:B:142:THR:HB	2.12	0.48
1:B:213:ARG:HD2	1:B:238:PHE:HE1	1.76	0.48
1:B:79:VAL:O	1:B:92:ARG:HA	2.13	0.48
1:A:99:CYS:N	1:A:100:PRO:CD	2.76	0.48
1:A:225:GLU:CG	1:A:241:ASN:HB2	2.44	0.48
1:B:81:ASN:O	1:B:91:SER:HB3	2.13	0.48
1:B:110:PHE:CD1	1:B:182:GLN:HB3	2.47	0.48
1:B:127:ILE:HG13	1:B:141:THR:HG22	1.94	0.48
1:A:227:SER:O	1:A:231:LYS:HG3	2.13	0.48
1:A:77:ILE:CG1	1:B:77:ILE:HD11	2.42	0.47
1:A:213:ARG:NH1	1:A:246:GLU:OE1	2.47	0.47
1:B:250:LEU:CD2	1:B:363:TYR:CD2	2.97	0.47
1:A:141:THR:HG21	1:A:145:ARG:HA	1.95	0.47
1:B:210:THR:CG2	1:B:211:VAL:N	2.77	0.47
1:B:227:SER:OG	1:B:230:GLU:N	2.44	0.47
1:B:124:SER:HB3	1:B:143:TYR:CG	2.50	0.47
1:A:98:TYR:O	1:A:99:CYS:C	2.57	0.47
1:B:216:ASP:OD1	1:B:280:SER:CB	2.63	0.47
1:B:335:ARG:CG	1:B:336:PHE:N	2.69	0.47
1:A:139:PHE:CE2	2:A:501:ANP:H1'	2.49	0.47
1:B:110:PHE:CE1	1:B:182:GLN:HB3	2.50	0.47
1:B:272:ALA:O	1:B:275:LYS:N	2.47	0.46
1:B:214:LYS:HG3	1:B:235:LEU:HD11	1.97	0.46
1:A:137:THR:HB	1:A:138:ARG:HB3	1.98	0.46
1:B:189:LEU:HD12	1:B:189:LEU:N	2.30	0.46
1:A:264:LEU:O	1:A:268:VAL:HG23	2.16	0.46
1:A:160:MET:O	1:A:164:LEU:HB2	2.16	0.46
1:B:243:PHE:C	1:B:243:PHE:CD2	2.94	0.46
1:A:97:GLU:HG3	1:A:187:TYR:CE2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:407:PHE:CD2	1:A:407:PHE:C	2.93	0.45
1:B:349:TYR:CD1	1:B:366:ALA:HB2	2.51	0.45
1:B:336:PHE:CD2	1:B:336:PHE:C	2.93	0.45
1:B:236:PRO:HB2	1:B:238:PHE:CE1	2.51	0.45
1:A:110:PHE:CD1	1:A:182:GLN:HB3	2.52	0.45
1:B:235:LEU:O	1:B:235:LEU:HD13	2.17	0.45
1:A:274:LEU:O	1:A:276:ILE:HG13	2.16	0.45
1:A:166:LYS:HD3	1:A:274:LEU:CD2	2.47	0.45
1:B:99:CYS:N	1:B:100:PRO:HD3	2.32	0.44
1:A:189:LEU:N	1:A:189:LEU:HD12	2.32	0.44
1:B:223:ALA:O	1:B:224:ARG:HG3	2.17	0.44
1:A:128:ASN:OD1	1:A:128:ASN:C	2.60	0.44
1:A:220:SER:HB3	1:A:222:VAL:HG23	2.00	0.44
1:B:260:PHE:HB2	1:B:351:MET:HE1	2.00	0.44
1:A:270:PHE:CE1	1:A:274:LEU:HD11	2.52	0.44
1:A:80:ASP:HB3	1:A:92:ARG:HE	1.82	0.44
1:B:213:ARG:HH22	1:B:246:GLU:HG3	1.83	0.44
1:A:215:TYR:HB3	1:A:217:LEU:HD21	1.99	0.44
1:B:214:LYS:HB2	1:B:236:PRO:O	2.18	0.44
1:B:261:LEU:HB3	1:B:265:LYS:HE2	2.00	0.43
1:A:268:VAL:CG1	1:A:404:SER:HA	2.48	0.43
1:A:106:LEU:HD23	1:A:106:LEU:HA	1.89	0.43
1:A:282:LEU:O	1:A:365:MET:HA	2.19	0.43
1:A:278:ASP:N	1:A:403:TYR:CE1	2.86	0.43
1:B:177:ASN:HD21	1:B:263:LYS:HD2	1.82	0.43
1:B:371:LEU:C	1:B:372:THR:CG2	2.92	0.43
1:B:255:GLU:O	1:B:259:ASN:OD1	2.37	0.42
1:A:179:LEU:HG	1:A:263:LYS:HB3	2.00	0.42
2:B:501:ANP:O1G	2:B:501:ANP:O1B	2.37	0.42
1:A:42:ILE:H	1:A:42:ILE:HD12	1.83	0.42
1:B:283:VAL:CG1	1:B:284:GLY:N	2.83	0.42
1:A:272:ALA:HB2	1:A:404:SER:HB3	2.01	0.42
1:A:129:SER:OG	1:A:139:PHE:HD1	2.02	0.42
1:B:142:THR:O	1:B:143:TYR:C	2.62	0.42
1:A:139:PHE:CZ	2:A:501:ANP:Hi'	2.54	0.42
1:A:141:THR:HG22	1:A:142:THR:O	2.19	0.42
1:A:225:GLU:HG2	1:A:241:ASN:HB2	2.00	0.42
1:B:124:SER:HB3	1:B:143:TYR:CB	2.49	0.42
1:B:51:ASN:HB2	1:B:122:THR:HG21	2.01	0.42
1:B:250:LEU:HD23	1:B:363:TYR:CE2	2.54	0.42
1:B:71:PHE:CE2	1:B:165:LYS:HA	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:LEU:HB2	1:A:411:MET:SD	2.60	0.41
1:B:79:VAL:HG13	3:B:614:HOH:O	2.19	0.41
1:A:77:ILE:HG13	1:B:77:ILE:CD1	2.45	0.41
1:A:260:PHE:HB2	1:A:351:MET:HE1	2.01	0.41
1:B:216:ASP:O	1:B:239:LYS:HA	2.20	0.41
1:A:141:THR:HG22	1:A:142:THR:N	2.34	0.41
1:A:272:ALA:HB2	1:A:404:SER:CB	2.51	0.41
1:A:281:LEU:HB3	1:A:411:MET:SD	2.60	0.41
1:A:123:ARG:HG3	1:A:123:ARG:NH1	2.36	0.41
1:A:124:SER:HB3	1:A:143:TYR:CG	2.56	0.41
1:B:40:GLU:OE1	1:B:138:ARG:NH2	2.53	0.41
1:A:350:ALA:HA	1:A:363:TYR:O	2.21	0.41
2:A:501:ANP:O1A	2:A:501:ANP:O1G	2.39	0.41
1:B:159:GLU:HA	1:B:159:GLU:OE2	2.20	0.41
1:A:54:ILE:HG21	1:A:118:GLN:HB2	2.03	0.40
1:A:298:VAL:O	1:A:302:ALA:HB2	2.21	0.40
1:A:79:VAL:O	1:A:92:ARG:HA	2.22	0.40
1:A:170:PHE:CE1	1:A:180:LEU:HD12	2.57	0.40
1:B:40:GLU:O	1:B:43:LEU:HB2	2.22	0.40
1:B:157:VAL:HA	1:B:186:MET:HE1	2.04	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	312/393 (79%)	295 (95%)	15 (5%)	2 (1%)	21	44
1	B	290/393 (74%)	272 (94%)	17 (6%)	1 (0%)	36	60
All	All	602/786 (77%)	567 (94%)	32 (5%)	3 (0%)	24	48

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	358	PRO
1	B	358	PRO
1	A	339	PRO

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	297/352 (84%)	278 (94%)	19 (6%)	16	38
1	B	280/352 (80%)	260 (93%)	20 (7%)	13	33
All	All	577/704 (82%)	538 (93%)	39 (7%)	14	35

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	64	VAL
1	A	77	ILE
1	A	123	ARG
1	A	127	ILE
1	A	137	THR
1	A	160	MET
1	A	165	LYS
1	A	173	GLU
1	A	177	ASN
1	A	202	ARG
1	A	222	VAL
1	A	280	SER
1	A	348	VAL
1	A	351	MET
1	A	398	VAL
1	A	404	SER
1	A	414	ILE
1	A	415	LEU
1	B	39	SER
1	B	42	ILE

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Mol	Chain	Res	Type
1	B	56	GLU
1	B	76	LYS
1	B	77	ILE
1	B	80	ASP
1	B	87	GLU
1	B	104	ARG
1	B	150	LYS
1	B	155	GLU
1	B	166	LYS
1	B	211	VAL
1	B	227	SER
1	B	228	ASP
1	B	256	SER
1	B	281	LEU
1	B	289	ASP
1	B	336	PHE
1	B	348	VAL
1	B	356	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	ASN
1	B	177	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry

3 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	ANP	B	501	-	33,33,33	2.64	13 (39%)	45,52,52	3.28	16 (35%)
2	ANP	A	501	-	33,33,33	2.48	13 (39%)	45,52,52	3.33	18 (40%)
2	ANP	A	502	-	33,33,33	2.79	17 (51%)	45,52,52	3.58	24 (53%)

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	ANP	B	501	-	-	5/18/38/38	0/3/3/3
2	ANP	A	501	-	-	7/18/38/38	0/3/3/3
2	ANP	A	502	-	-	8/18/38/38	0/3/3/3

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	ANP	PA-O3A	8.54	1.68	1.59
2	A	502	ANP	PA-O3A	7.87	1.68	1.59
2	A	501	ANP	PA-O3A	6.95	1.67	1.59
2	B	501	ANP	PB-O1B	6.88	1.56	1.46
2	A	502	ANP	PB-O1B	5.86	1.55	1.46
2	A	501	ANP	PB-O1B	5.81	1.55	1.46
2	A	502	ANP	C6-N6	4.58	1.45	1.34
2	A	501	ANP	C5-N7	-4.02	1.31	1.39
2	A	502	ANP	PA-O2A	3.91	1.73	1.55
2	A	501	ANP	C8-N9	-3.89	1.30	1.37
2	B	501	ANP	C5'-C4'	3.88	1.63	1.51
2	A	501	ANP	C5'-C4'	3.81	1.63	1.51
2	B	501	ANP	C6-N6	3.80	1.43	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	502	ANP	PA-O5'	3.65	1.73	1.59
2	A	502	ANP	PG-O1G	3.52	1.51	1.46
2	A	502	ANP	C2'-C3'	-3.45	1.44	1.53
2	A	502	ANP	PB-O3A	-3.37	1.54	1.59
2	A	501	ANP	O3'-C3'	-3.21	1.35	1.43
2	B	501	ANP	PG-O1G	3.10	1.50	1.46
2	B	501	ANP	C5-N7	-3.03	1.33	1.39
2	A	502	ANP	C3'-C4'	2.97	1.60	1.53
2	A	502	ANP	O3'-C3'	-2.90	1.35	1.43
2	A	502	ANP	PB-N3B	2.82	1.70	1.63
2	B	501	ANP	PA-O2A	2.78	1.68	1.55
2	B	501	ANP	C2'-C3'	-2.76	1.45	1.53
2	A	502	ANP	PA-O1A	2.74	1.60	1.50
2	A	501	ANP	PB-O3A	-2.60	1.55	1.59
2	B	501	ANP	O3'-C3'	-2.47	1.36	1.43
2	B	501	ANP	PA-O5'	2.47	1.69	1.59
2	A	501	ANP	C2'-C3'	-2.45	1.46	1.53
2	A	502	ANP	C5-N7	-2.40	1.34	1.39
2	A	501	ANP	O2'-C2'	-2.38	1.37	1.43
2	A	501	ANP	C6-N6	2.37	1.40	1.34
2	A	501	ANP	PA-O2A	2.25	1.65	1.55
2	B	501	ANP	C8-N9	-2.24	1.33	1.37
2	A	502	ANP	C8-N9	-2.23	1.33	1.37
2	A	502	ANP	C1'-N9	-2.22	1.40	1.46
2	A	501	ANP	O4'-C1'	-2.21	1.36	1.42
2	A	502	ANP	PG-N3B	2.12	1.68	1.63
2	B	501	ANP	PA-O1A	2.10	1.58	1.50
2	A	501	ANP	C5-C6	-2.07	1.35	1.41
2	A	502	ANP	C2-N1	2.06	1.37	1.33
2	B	501	ANP	O2'-C2'	-2.05	1.37	1.43

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	ANP	O3A-PA-O1A	-12.09	74.32	110.70
2	A	501	ANP	O3A-PA-O1A	-11.15	77.17	110.70
2	A	502	ANP	O3A-PA-O1A	-9.48	82.18	110.70
2	A	501	ANP	O5'-PA-O1A	-8.71	74.43	108.94
2	A	502	ANP	O2A-PA-O1A	-7.70	76.63	112.44
2	B	501	ANP	O2A-PA-O1A	-7.68	76.72	112.44
2	B	501	ANP	O5'-PA-O1A	-7.53	79.08	108.94
2	A	502	ANP	O5'-PA-O1A	-7.42	79.53	108.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ANP	O2A-PA-O1A	-7.32	78.42	112.44
2	A	502	ANP	C4-N9-C8	7.00	113.09	105.74
2	A	502	ANP	N3-C2-N1	-6.92	118.11	128.58
2	B	501	ANP	N3-C2-N1	-6.43	118.85	128.58
2	A	501	ANP	N3-C2-N1	-6.07	119.39	128.58
2	A	501	ANP	C5-C4-N3	-5.72	118.84	126.72
2	A	502	ANP	O2A-PA-O5'	5.58	132.85	107.57
2	A	501	ANP	N3-C4-N9	5.27	136.13	127.17
2	A	502	ANP	C5'-C4'-C3'	5.24	134.09	115.21
2	A	502	ANP	N9-C8-N7	-5.21	106.55	113.94
2	B	501	ANP	C5-C4-N3	-4.75	120.17	126.72
2	B	501	ANP	N3-C4-N9	4.34	134.55	127.17
2	A	502	ANP	N3-C4-N9	4.10	134.14	127.17
2	A	502	ANP	N6-C6-N1	4.03	127.34	118.38
2	A	502	ANP	C5-C4-N9	-3.97	101.48	105.81
2	A	501	ANP	C5-N7-C8	3.93	109.63	103.45
2	B	501	ANP	C5'-C4'-C3'	3.91	129.27	115.21
2	A	501	ANP	C2-N3-C4	3.90	121.36	111.83
2	B	501	ANP	C5-N7-C8	3.88	109.54	103.45
2	A	501	ANP	O4'-C4'-C5'	3.75	121.34	109.33
2	A	501	ANP	N9-C8-N7	-3.63	108.79	113.94
2	A	502	ANP	C5-C6-N6	-3.60	114.37	123.29
2	B	501	ANP	O2A-PA-O3A	3.57	116.92	107.27
2	A	502	ANP	O3'-C3'-C4'	3.57	121.33	111.08
2	B	501	ANP	O4'-C4'-C5'	3.57	120.76	109.33
2	B	501	ANP	C2-N3-C4	3.56	120.52	111.83
2	A	502	ANP	C2-N1-C6	3.51	124.49	118.73
2	A	502	ANP	C5-N7-C8	3.44	108.85	103.45
2	B	501	ANP	N9-C8-N7	-3.40	109.12	113.94
2	A	501	ANP	C3'-C2'-C1'	3.22	107.55	101.46
2	A	502	ANP	O2'-C2'-C1'	-3.21	99.04	110.10
2	A	501	ANP	C4-N9-C8	3.02	108.91	105.74
2	A	501	ANP	C5'-C4'-C3'	2.96	125.86	115.21
2	A	502	ANP	C4'-O4'-C1'	-2.91	103.04	109.47
2	A	502	ANP	O4'-C1'-C2'	2.80	112.60	106.62
2	B	501	ANP	C2-N1-C6	2.78	123.29	118.73
2	A	502	ANP	C4-N9-C1'	-2.67	120.39	126.63
2	A	502	ANP	O3G-PG-O1G	-2.64	106.83	113.45
2	A	502	ANP	O5'-C5'-C4'	-2.64	100.01	108.99
2	B	501	ANP	C4-C5-N7	-2.62	107.59	110.58
2	B	501	ANP	C4-N9-C8	2.61	108.48	105.74
2	A	501	ANP	C4-C5-N7	-2.57	107.64	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	502	ANP	C2-N3-C4	2.46	117.84	111.83
2	A	501	ANP	O2A-PA-O3A	2.45	113.89	107.27
2	A	501	ANP	O4'-C1'-C2'	2.42	111.80	106.62
2	A	502	ANP	O1B-PB-N3B	-2.36	108.29	111.77
2	A	502	ANP	C3'-C2'-C1'	2.09	105.43	101.46
2	A	501	ANP	C4'-O4'-C1'	-2.02	105.00	109.47
2	B	501	ANP	C4'-O4'-C1'	-2.00	105.04	109.47
2	A	501	ANP	C5-C6-N6	-2.00	118.33	123.29

There are no chirality outliers.

All (20) torsion outliers are listed below:

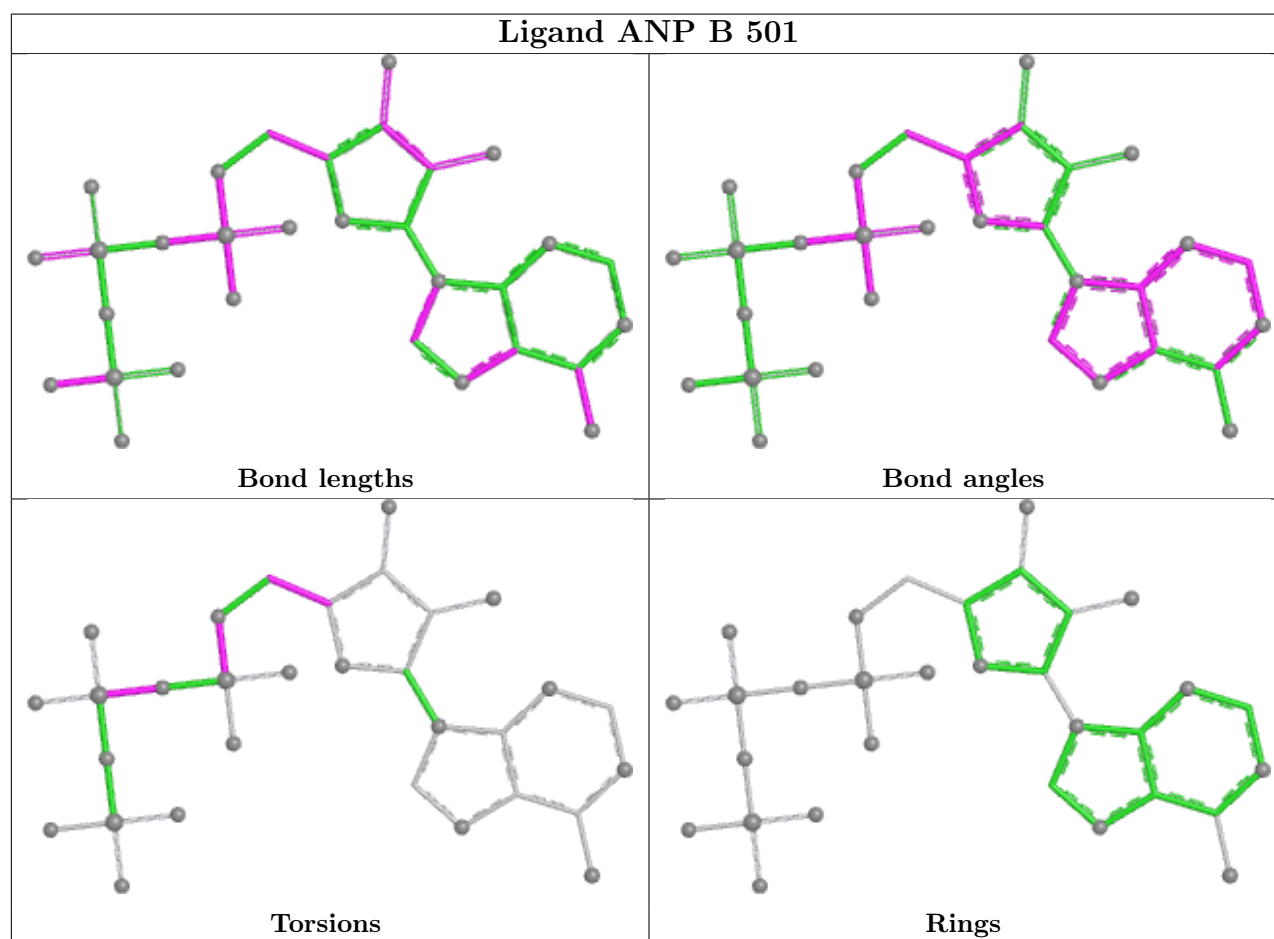
Mol	Chain	Res	Type	Atoms
2	A	501	ANP	C5'-O5'-PA-O1A
2	A	501	ANP	C5'-O5'-PA-O2A
2	A	501	ANP	C5'-O5'-PA-O3A
2	A	501	ANP	C4'-C5'-O5'-PA
2	A	501	ANP	O4'-C4'-C5'-O5'
2	A	502	ANP	PB-N3B-PG-O1G
2	A	502	ANP	PG-N3B-PB-O1B
2	A	502	ANP	PG-N3B-PB-O3A
2	A	502	ANP	PA-O3A-PB-O2B
2	A	502	ANP	C5'-O5'-PA-O1A
2	B	501	ANP	C5'-O5'-PA-O3A
2	A	502	ANP	C4'-C5'-O5'-PA
2	A	502	ANP	O4'-C4'-C5'-O5'
2	B	501	ANP	C3'-C4'-C5'-O5'
2	B	501	ANP	C5'-O5'-PA-O1A
2	A	501	ANP	C2'-C1'-N9-C8
2	A	501	ANP	C2'-C1'-N9-C4
2	A	502	ANP	PB-O3A-PA-O1A
2	B	501	ANP	PA-O3A-PB-O1B
2	B	501	ANP	O4'-C4'-C5'-O5'

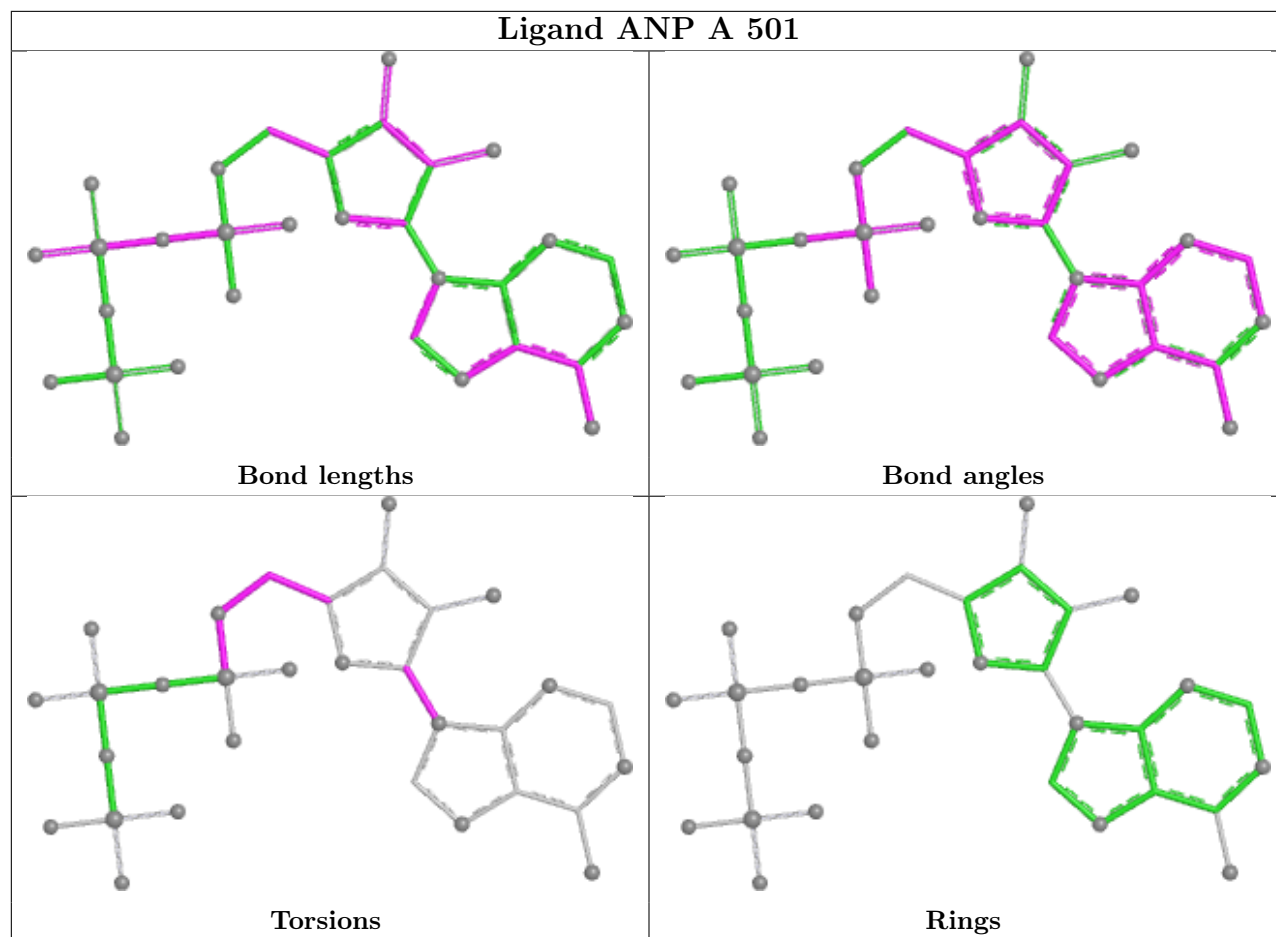
There are no ring outliers.

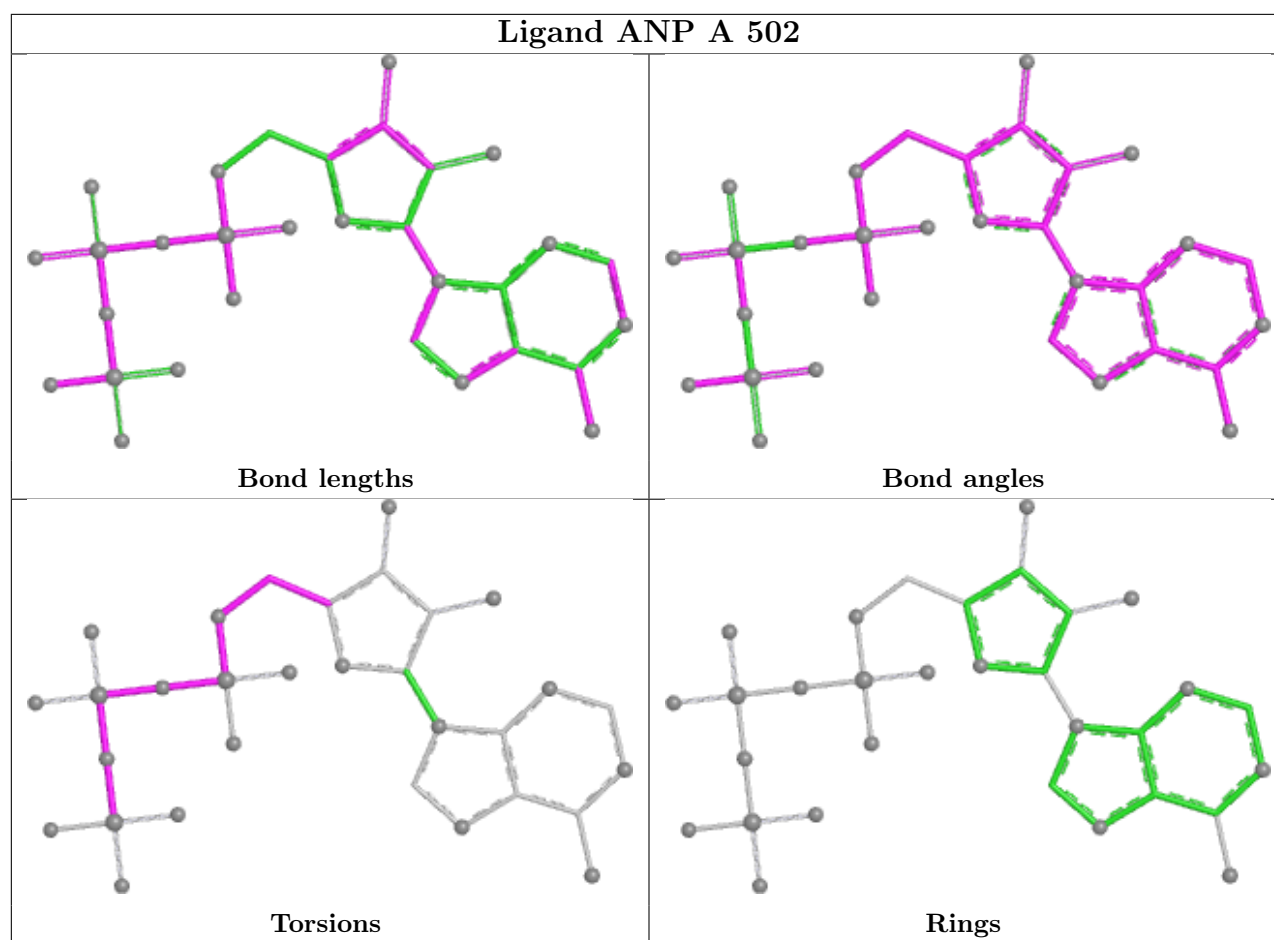
3 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	ANP	2	0
2	A	501	ANP	8	0
2	A	502	ANP	2	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

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	320/393 (81%)	0.41	22 (6%)	23 20	39, 78, 132, 181	0
1	B	302/393 (76%)	0.56	25 (8%)	17 14	28, 86, 158, 173	0
All	All	622/786 (79%)	0.48	47 (7%)	20 17	28, 81, 151, 181	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	416	THR	7.2
1	B	338	GLY	6.8
1	B	372	THR	5.6
1	A	56	GLU	5.0
1	B	337	PHE	4.6
1	B	219	GLY	4.4
1	B	251	HIS	4.4
1	A	291	ALA	4.3
1	B	400	PRO	3.9
1	B	411	MET	3.9
1	B	407	PHE	3.7
1	A	274	LEU	3.6
1	A	298	VAL	3.4
1	B	222	VAL	3.4
1	A	296	MET	3.3
1	A	304	ASP	3.3
1	A	411	MET	3.2
1	B	242	ASP	3.1
1	B	246	GLU	3.0
1	A	371	LEU	3.0
1	B	208	ARG	2.9
1	A	248	GLN	2.8
1	B	258	LYS	2.7
1	A	129	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	413	ASN	2.7
1	A	302	ALA	2.6
1	A	222	VAL	2.5
1	A	407	PHE	2.5
1	B	336	PHE	2.5
1	A	303	GLU	2.5
1	B	415	LEU	2.4
1	B	278	ASP	2.4
1	A	406	ARG	2.4
1	A	300	GLU	2.4
1	A	404	SER	2.4
1	A	275	LYS	2.4
1	B	272	ALA	2.2
1	B	277	MET	2.2
1	B	226	ALA	2.2
1	A	301	ARG	2.2
1	B	406	ARG	2.1
1	B	161	HIS	2.1
1	A	414	ILE	2.1
1	B	410	PHE	2.1
1	A	276	ILE	2.0
1	B	369	ASP	2.0
1	B	283	VAL	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.

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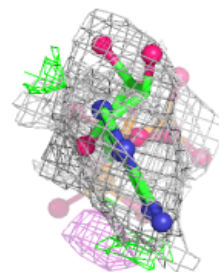
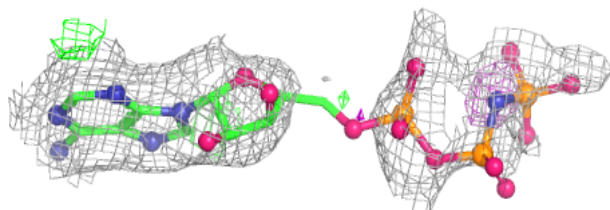
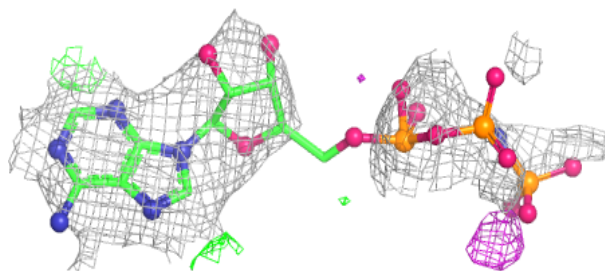
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ANP	B	501	31/31	0.71	0.14	110,129,249,301	0
2	ANP	A	501	31/31	0.77	0.14	69,117,221,271	0
2	ANP	A	502	31/31	0.82	0.17	41,87,161,228	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.

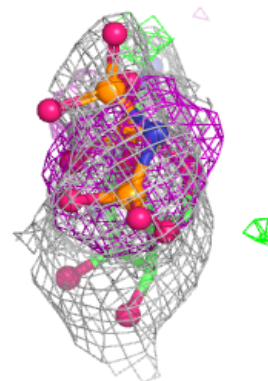
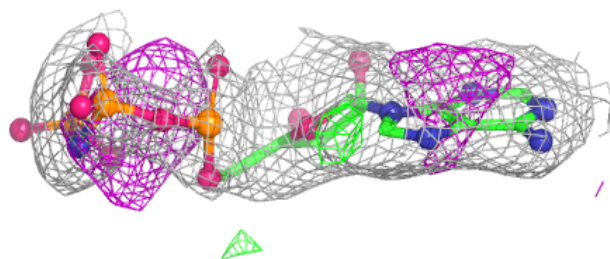
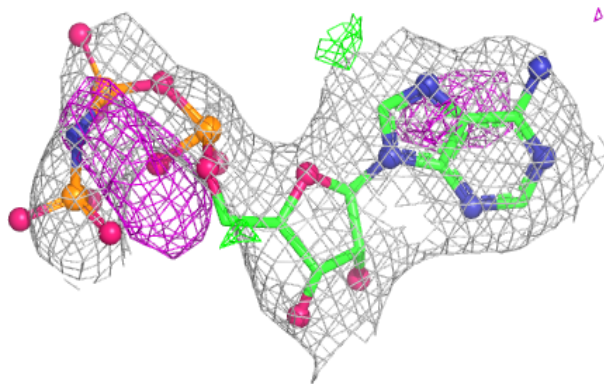
Electron density around ANP B 501:

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

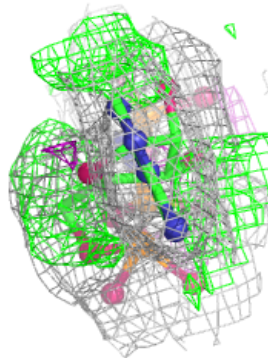
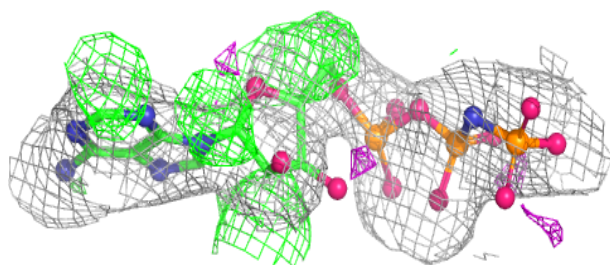
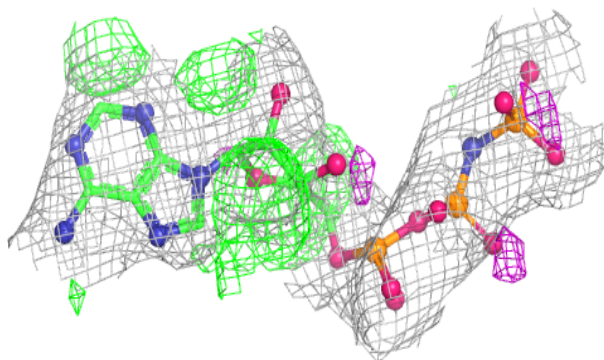


Electron density around ANP A 501:

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

**Electron density around ANP A 502:**

$2mF_o-DF_c$ (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.