



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2022 – 12:10 PM JST

PDB ID : 7CTV
Title : Crystal structure of Arabidopsis thaliana SOBIR1 kinase domain D489A mutant in complex with AMP-PNP and magnesium
Authors : Wei, X.; Wang, Y.L.; Gu, T.Y.; Xin, F.J.
Deposited on : 2020-08-20
Resolution : 2.89 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.27
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.27

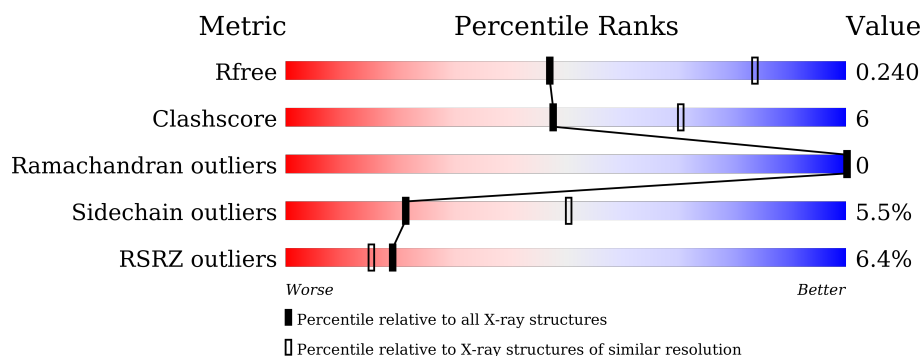
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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}	130704	2691 (2.90-2.86)
Clashscore	141614	2947 (2.90-2.86)
Ramachandran outliers	138981	2868 (2.90-2.86)
Sidechain outliers	138945	2871 (2.90-2.86)
RSRZ outliers	127900	2629 (2.90-2.86)

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	336	<div> <div>6%</div> <div> <div></div> <div>75%</div> <div>12%</div> <div>•</div> <div>11%</div> </div> </div>
1	B	336	<div> <div>5%</div> <div> <div></div> <div>71%</div> <div>15%</div> <div>•</div> <div>13%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4507 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 Leucine-rich repeat receptor-like serine/threonine/tyrosine-protein kinase SOBIR1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	0	0
			2264	1444	384	420	16			
1	B	293	Total	C	N	O	S	0	0	0
			2179	1389	370	406	14			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	305	MET	-	initiating methionine	UNP Q9SKB2
A	306	GLY	-	expression tag	UNP Q9SKB2
A	307	SER	-	expression tag	UNP Q9SKB2
A	308	SER	-	expression tag	UNP Q9SKB2
A	309	HIS	-	expression tag	UNP Q9SKB2
A	310	HIS	-	expression tag	UNP Q9SKB2
A	311	HIS	-	expression tag	UNP Q9SKB2
A	312	HIS	-	expression tag	UNP Q9SKB2
A	313	HIS	-	expression tag	UNP Q9SKB2
A	314	HIS	-	expression tag	UNP Q9SKB2
A	315	SER	-	expression tag	UNP Q9SKB2
A	316	SER	-	expression tag	UNP Q9SKB2
A	317	GLY	-	expression tag	UNP Q9SKB2
A	318	LEU	-	expression tag	UNP Q9SKB2
A	319	VAL	-	expression tag	UNP Q9SKB2
A	320	PRO	-	expression tag	UNP Q9SKB2
A	321	ARG	-	expression tag	UNP Q9SKB2
A	322	GLY	-	expression tag	UNP Q9SKB2
A	323	SER	-	expression tag	UNP Q9SKB2
A	324	HIS	-	expression tag	UNP Q9SKB2
A	325	MET	-	expression tag	UNP Q9SKB2
A	489	ALA	ASP	engineered mutation	UNP Q9SKB2
B	305	MET	-	initiating methionine	UNP Q9SKB2
B	306	GLY	-	expression tag	UNP Q9SKB2

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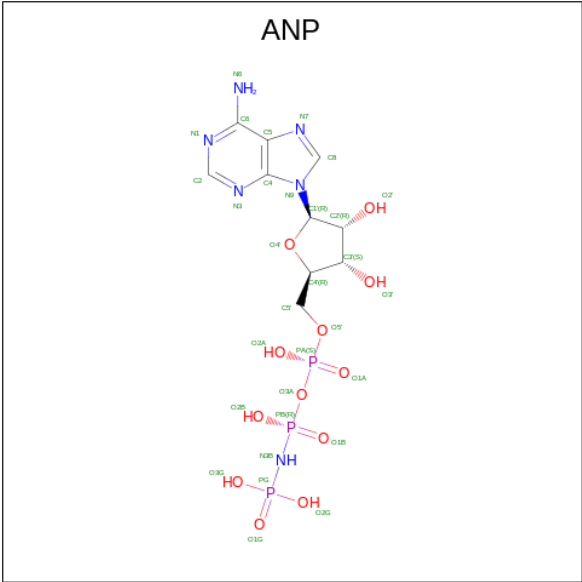
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Chain	Residue	Modelled	Actual	Comment	Reference
B	307	SER	-	expression tag	UNP Q9SKB2
B	308	SER	-	expression tag	UNP Q9SKB2
B	309	HIS	-	expression tag	UNP Q9SKB2
B	310	HIS	-	expression tag	UNP Q9SKB2
B	311	HIS	-	expression tag	UNP Q9SKB2
B	312	HIS	-	expression tag	UNP Q9SKB2
B	313	HIS	-	expression tag	UNP Q9SKB2
B	314	HIS	-	expression tag	UNP Q9SKB2
B	315	SER	-	expression tag	UNP Q9SKB2
B	316	SER	-	expression tag	UNP Q9SKB2
B	317	GLY	-	expression tag	UNP Q9SKB2
B	318	LEU	-	expression tag	UNP Q9SKB2
B	319	VAL	-	expression tag	UNP Q9SKB2
B	320	PRO	-	expression tag	UNP Q9SKB2
B	321	ARG	-	expression tag	UNP Q9SKB2
B	322	GLY	-	expression tag	UNP Q9SKB2
B	323	SER	-	expression tag	UNP Q9SKB2
B	324	HIS	-	expression tag	UNP Q9SKB2
B	325	MET	-	expression tag	UNP Q9SKB2
B	489	ALA	ASP	engineered mutation	UNP Q9SKB2

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Mg 1 1	0	0
2	B	1	Total Mg 1 1	0	0

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (three-letter code: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).

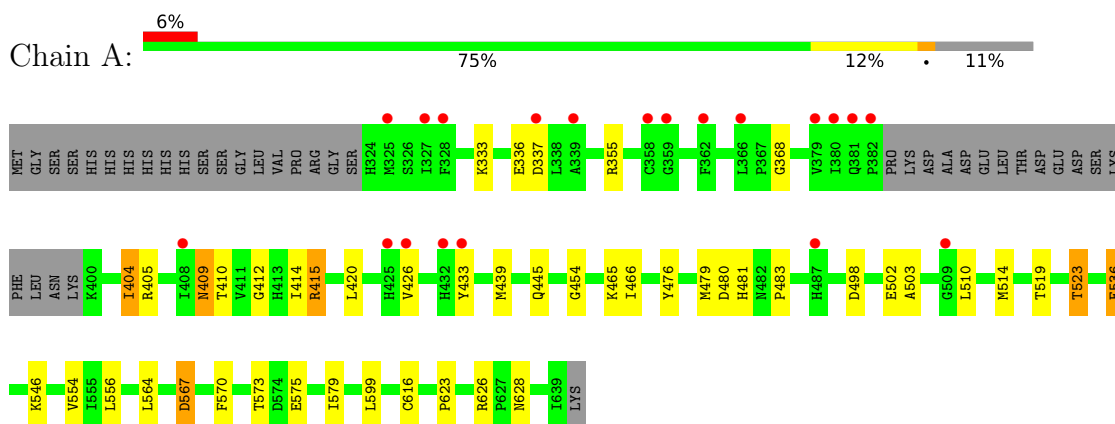


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			31	10	6	12	3		
3	B	1	Total	C	N	O	P	0	0
			31	10	6	12	3		

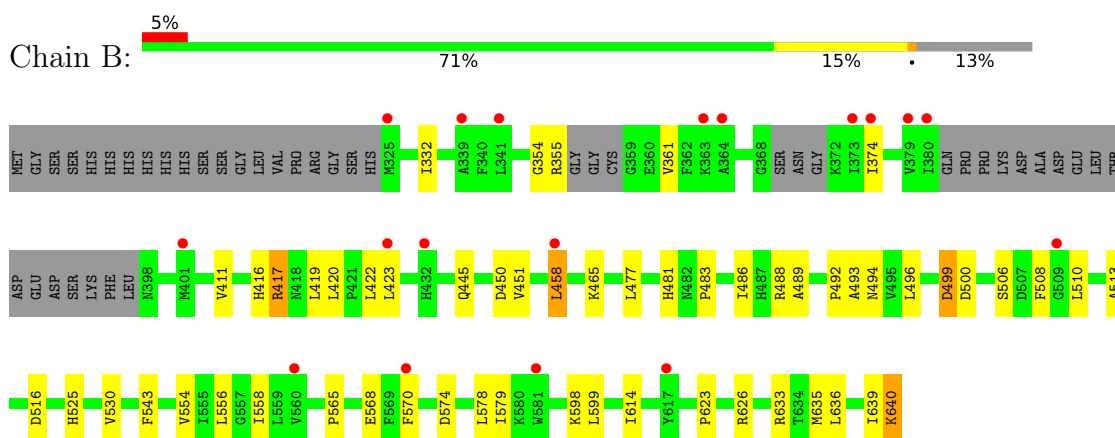
3 Residue-property plots

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: Leucine-rich repeat receptor-like serine/threonine/tyrosine-protein kinase SO-BIR1



- Molecule 1: Leucine-rich repeat receptor-like serine/threonine/tyrosine-protein kinase SO-BIR1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.57Å 50.89Å 105.42Å 90.00° 95.35° 90.00°	Depositor
Resolution (Å)	43.39 – 2.89 43.39 – 2.89	Depositor EDS
% Data completeness (in resolution range)	98.2 (43.39-2.89) 98.3 (43.39-2.89)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.217 , 0.239 0.217 , 0.240	Depositor DCC
R_{free} test set	775 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	82.3	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 47.6	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	0.94	EDS
Total number of atoms	4507	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.61% 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, MG

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	0.24	0/2310	0.41	0/3139
1	B	0.23	0/2221	0.40	0/3022
All	All	0.24	0/4531	0.40	0/6161

There are no bond length outliers.

There are no bond angle outliers.

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	2264	0	2194	24	0
1	B	2179	0	2076	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	31	0	13	0	0
3	B	31	0	13	0	0
All	All	4507	0	4296	54	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 6.

All (54) 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:416:HIS:HB3	1:B:419:LEU:HB2	1.72	0.71
1:B:417:ARG:HH12	1:B:633:ARG:HH12	1.42	0.67
1:A:479:MET:HE1	1:A:628:ASN:HB2	1.78	0.65
1:B:639:ILE:HG22	1:B:640:LYS:HG3	1.78	0.65
1:A:412:GLY:O	1:A:415:ARG:NH2	2.31	0.64
1:B:445:GLN:HB2	1:B:492:PRO:HB2	1.81	0.61
1:B:477:LEU:HD21	1:B:508:PHE:HE2	1.67	0.60
1:B:451:VAL:O	1:B:598:LYS:NZ	2.36	0.58
1:B:419:LEU:HD21	1:B:477:LEU:HD11	1.85	0.57
1:B:499:ASP:OD2	1:B:499:ASP:N	2.38	0.56
1:B:574:ASP:OD1	1:B:574:ASP:N	2.38	0.56
1:B:558:ILE:HD11	1:B:565:PRO:HD3	1.89	0.54
1:B:623:PRO:HA	1:B:626:ARG:HG3	1.92	0.51
1:A:466:ILE:HG23	1:A:503:ALA:HB2	1.91	0.51
1:B:489:ALA:O	1:B:494:ASN:ND2	2.43	0.51
1:B:486:ILE:HG22	1:B:488:ARG:HG3	1.93	0.50
1:B:510:LEU:HB2	1:B:513:ALA:HB3	1.94	0.50
1:B:556:LEU:HD21	1:B:636:LEU:HD21	1.93	0.50
1:A:476:TYR:HA	1:A:480:ASP:HB2	1.93	0.49
1:A:554:VAL:HG22	1:A:579:ILE:HD11	1.93	0.48
1:A:404:ILE:H	1:A:404:ILE:HG12	1.44	0.48
1:A:405:ARG:O	1:A:409:ASN:HB2	2.13	0.47
1:B:354:GLY:HA3	1:B:361:VAL:H	1.80	0.47
1:A:575:GLU:N	1:A:575:GLU:OE1	2.47	0.47
1:A:510:LEU:O	1:A:514:MET:HG3	2.15	0.47
1:B:355:ARG:NH2	1:B:493:ALA:HB2	2.31	0.46
1:A:454:GLY:HA3	1:B:332:ILE:O	2.15	0.46
1:A:410:THR:O	1:A:414:ILE:HG13	2.16	0.46
1:A:439:MET:HE2	1:A:498:ASP:HA	1.97	0.46
1:A:564:LEU:O	1:A:567:ASP:HB2	2.16	0.45
1:B:614:ILE:HG12	1:B:635:MET:HB3	1.97	0.45
1:B:516:ASP:N	1:B:516:ASP:OD1	2.46	0.45
1:A:333:LYS:O	1:A:336:GLU:HG2	2.17	0.45
1:B:374:ILE:HD11	1:B:423:LEU:HD12	1.99	0.45
1:B:570:PHE:CZ	1:B:578:LEU:HD13	2.52	0.45
1:B:355:ARG:CZ	1:B:493:ALA:HB2	2.47	0.45
1:B:465:LYS:NZ	1:B:500:ASP:O	2.50	0.44
1:A:519:THR:O	1:A:523:THR:HG22	2.19	0.43
1:A:426:VAL:HB	1:A:433:TYR:HB2	2.00	0.43
1:A:414:ILE:HG12	1:A:481:HIS:NE2	2.33	0.43
1:B:554:VAL:HG22	1:B:579:ILE:HD11	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:420:LEU:HD22	1:B:496:LEU:HD12	2.00	0.43
1:A:570:PHE:O	1:A:573:THR:HG22	2.18	0.42
1:A:414:ILE:HG23	1:A:476:TYR:HE2	1.85	0.42
1:A:481:HIS:ND1	1:A:483:PRO:HD2	2.34	0.42
1:A:536:GLU:OE1	1:A:626:ARG:NH1	2.41	0.42
1:B:481:HIS:CE1	1:B:483:PRO:HG2	2.55	0.41
1:B:494:ASN:HB3	1:B:506:SER:O	2.20	0.41
1:A:623:PRO:HA	1:A:626:ARG:HG3	2.03	0.41
1:A:336:GLU:CD	1:A:368:GLY:HA3	2.41	0.40
1:B:374:ILE:HD13	1:B:374:ILE:HA	1.85	0.40
1:A:465:LYS:HE2	1:A:502:GLU:HG2	2.04	0.40
1:B:488:ARG:HG2	1:B:543:PHE:HB2	2.02	0.40
1:B:458:LEU:HD23	1:B:458:LEU:HA	1.82	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	295/336 (88%)	289 (98%)	6 (2%)	0	100	100
1	B	285/336 (85%)	277 (97%)	8 (3%)	0	100	100
All	All	580/672 (86%)	566 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	235/292 (80%)	221 (94%)	14 (6%)	19	46
1	B	220/292 (75%)	209 (95%)	11 (5%)	24	54
All	All	455/584 (78%)	430 (94%)	25 (6%)	21	50

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	337	ASP
1	A	355	ARG
1	A	404	ILE
1	A	409	ASN
1	A	415	ARG
1	A	420	LEU
1	A	445	GLN
1	A	523	THR
1	A	536	GLU
1	A	546	LYS
1	A	556	LEU
1	A	567	ASP
1	A	599	LEU
1	A	616	CYS
1	B	411	VAL
1	B	417	ARG
1	B	422	LEU
1	B	450	ASP
1	B	458	LEU
1	B	499	ASP
1	B	525	HIS
1	B	530	VAL
1	B	568	GLU
1	B	599	LEU
1	B	640	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ANP	A	702	2	29,33,33	4.82	14 (48%)	31,52,52	2.57	6 (19%)
3	ANP	B	702	2	29,33,33	4.82	15 (51%)	31,52,52	2.63	5 (16%)

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	ANP	A	702	2	-	6/14/38/38	0/3/3/3
3	ANP	B	702	2	-	4/14/38/38	0/3/3/3

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	ANP	C2'-C1'	-16.11	1.29	1.53
3	B	702	ANP	C2'-C1'	-16.11	1.29	1.53
3	B	702	ANP	O4'-C1'	14.56	1.61	1.41
3	A	702	ANP	O4'-C1'	14.47	1.61	1.41
3	B	702	ANP	PB-O2B	-5.26	1.42	1.56
3	A	702	ANP	PB-N3B	5.26	1.77	1.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	ANP	PB-O2B	-5.22	1.42	1.56
3	B	702	ANP	PB-N3B	5.13	1.76	1.63
3	A	702	ANP	O4'-C4'	-4.85	1.34	1.45
3	B	702	ANP	O4'-C4'	-4.84	1.34	1.45
3	A	702	ANP	PB-O3A	4.41	1.64	1.59
3	A	702	ANP	C6-N6	4.29	1.49	1.34
3	B	702	ANP	C6-N6	4.29	1.49	1.34
3	A	702	ANP	PG-O1G	4.18	1.52	1.46
3	B	702	ANP	PG-O1G	4.17	1.52	1.46
3	B	702	ANP	PB-O1B	4.13	1.52	1.46
3	B	702	ANP	PB-O3A	4.12	1.64	1.59
3	A	702	ANP	PB-O1B	4.02	1.52	1.46
3	A	702	ANP	O2'-C2'	2.71	1.49	1.43
3	B	702	ANP	O2'-C2'	2.66	1.49	1.43
3	B	702	ANP	C5-C4	-2.61	1.34	1.40
3	A	702	ANP	C2-N3	2.61	1.36	1.32
3	A	702	ANP	C5-C4	-2.60	1.34	1.40
3	B	702	ANP	PA-O5'	2.58	1.69	1.59
3	B	702	ANP	C2-N3	2.55	1.36	1.32
3	A	702	ANP	PA-O5'	2.47	1.69	1.59
3	A	702	ANP	PG-N3B	2.38	1.69	1.63
3	B	702	ANP	PG-N3B	2.30	1.69	1.63
3	B	702	ANP	C2'-C3'	2.00	1.58	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	702	ANP	C5-C6-N6	9.69	135.08	120.35
3	A	702	ANP	C5-C6-N6	9.50	134.79	120.35
3	B	702	ANP	N6-C6-N1	-6.58	104.91	118.57
3	A	702	ANP	N6-C6-N1	-6.44	105.21	118.57
3	B	702	ANP	N3-C2-N1	-6.06	119.20	128.68
3	A	702	ANP	N3-C2-N1	-5.96	119.37	128.68
3	B	702	ANP	C3'-C2'-C1'	3.63	106.45	100.98
3	A	702	ANP	C3'-C2'-C1'	3.27	105.91	100.98
3	B	702	ANP	PB-O3A-PA	-2.45	123.99	132.62
3	A	702	ANP	PB-O3A-PA	-2.24	124.72	132.62
3	A	702	ANP	O1B-PB-N3B	-2.04	108.77	111.77

There are no chirality outliers.

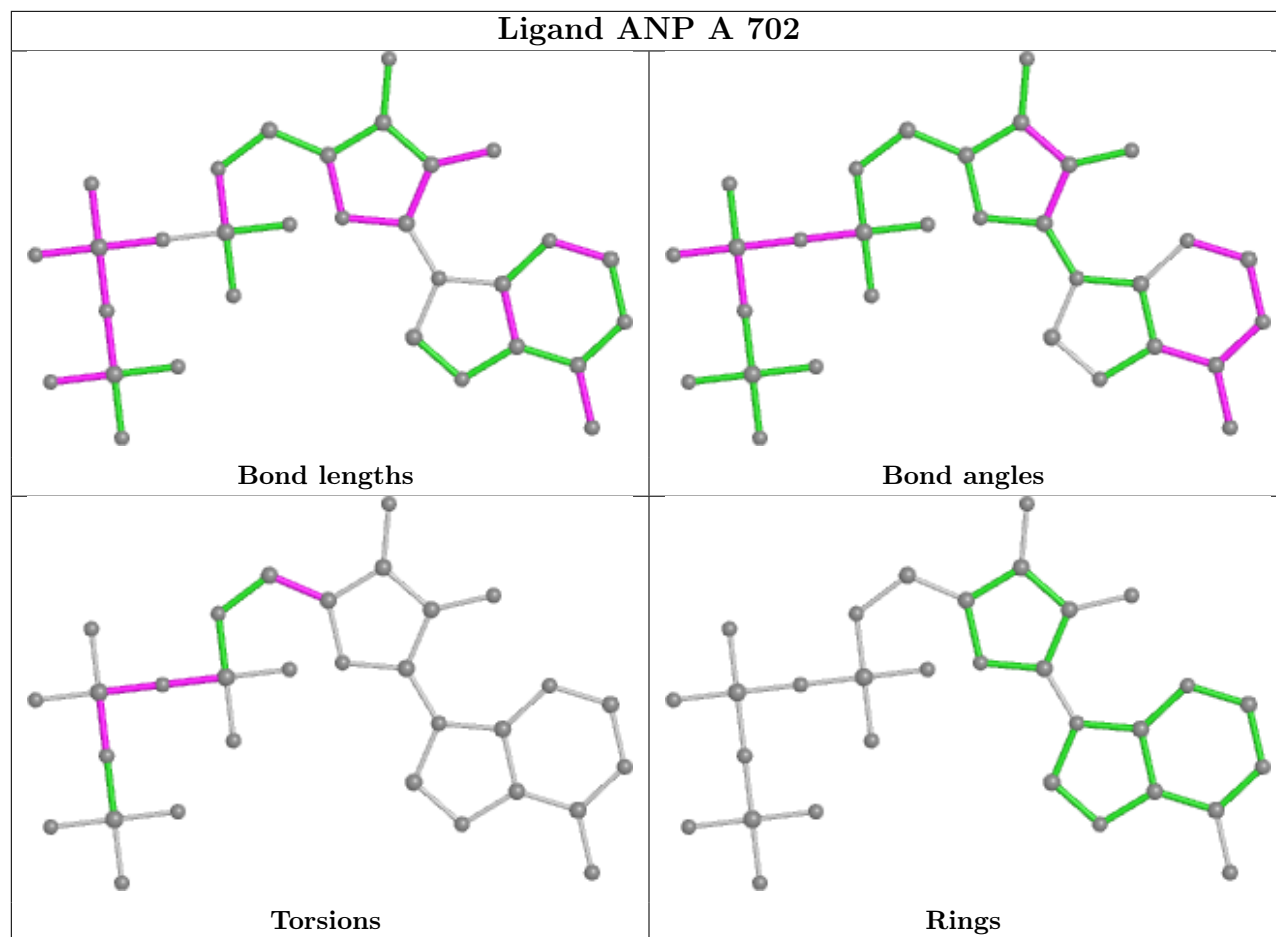
All (10) torsion outliers are listed below:

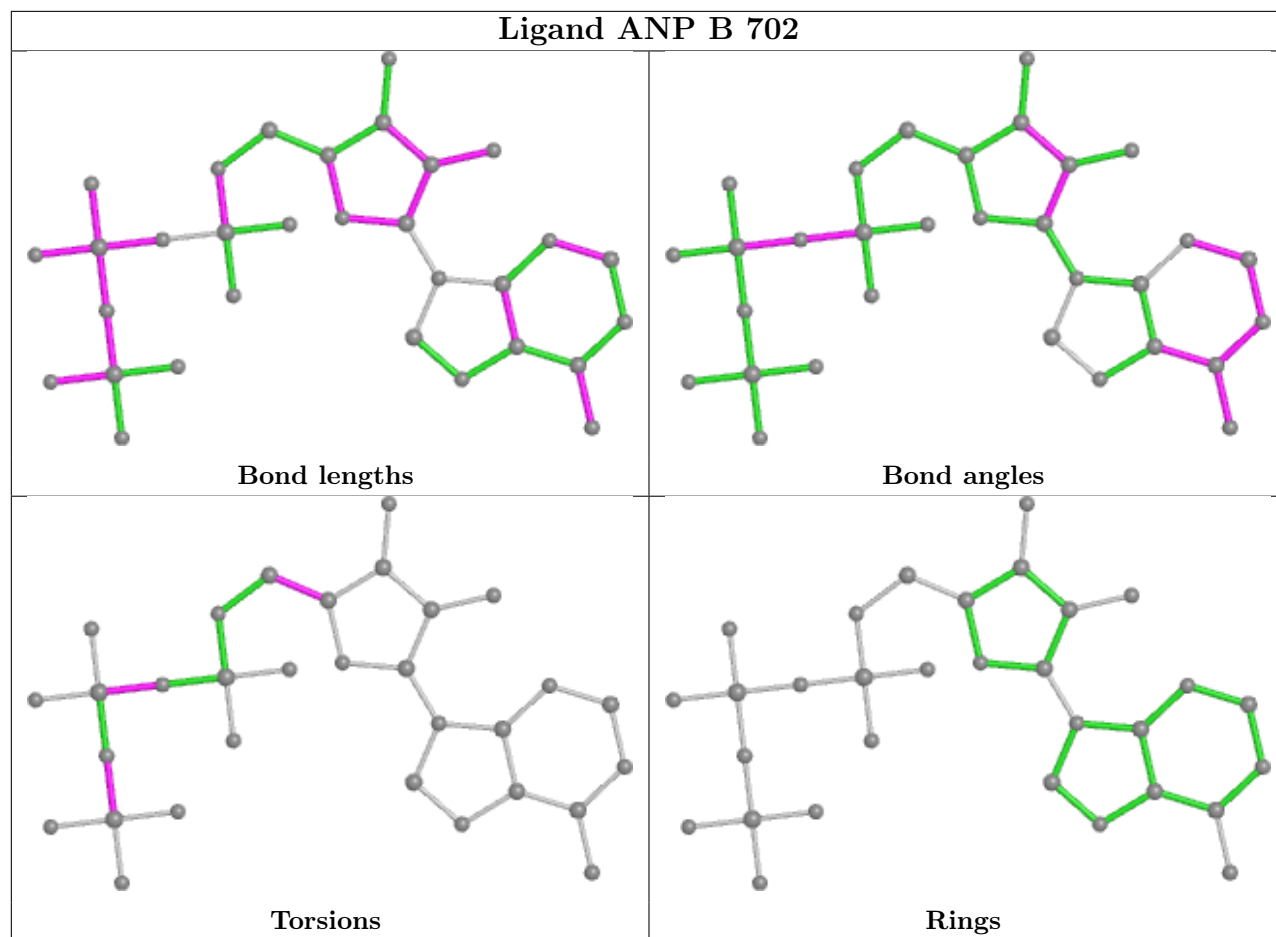
Mol	Chain	Res	Type	Atoms
3	A	702	ANP	PA-O3A-PB-O1B
3	A	702	ANP	PA-O3A-PB-O2B
3	B	702	ANP	PA-O3A-PB-O1B
3	B	702	ANP	PA-O3A-PB-O2B
3	A	702	ANP	PB-O3A-PA-O1A
3	A	702	ANP	C3'-C4'-C5'-O5'
3	A	702	ANP	PG-N3B-PB-O3A
3	B	702	ANP	C3'-C4'-C5'-O5'
3	A	702	ANP	O4'-C4'-C5'-O5'
3	B	702	ANP	PB-N3B-PG-O1G

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	299/336 (88%)	0.25	20 (6%) 17 13	49, 74, 126, 178	0
1	B	293/336 (87%)	0.34	18 (6%) 21 17	74, 104, 149, 185	0
All	All	592/672 (88%)	0.29	38 (6%) 19 15	49, 93, 143, 185	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	364	ALA	7.5
1	A	381	GLN	5.2
1	A	337	ASP	4.4
1	B	374	ILE	4.2
1	A	382	PRO	3.9
1	B	339	ALA	3.8
1	B	363	LYS	3.7
1	B	341	LEU	3.5
1	B	373	ILE	3.5
1	A	325	MET	3.4
1	A	509	GLY	3.4
1	A	379	VAL	3.2
1	B	423	LEU	3.2
1	B	570	PHE	3.1
1	A	327	ILE	3.0
1	A	432	HIS	2.9
1	A	362	PHE	2.8
1	B	380	ILE	2.7
1	B	509	GLY	2.7
1	B	458	LEU	2.6
1	A	339	ALA	2.5
1	A	358	CYS	2.5
1	A	433	TYR	2.5
1	B	379	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	617	TYR	2.4
1	A	359	GLY	2.3
1	A	366	LEU	2.3
1	A	328	PHE	2.3
1	B	581	TRP	2.3
1	A	380	ILE	2.3
1	B	432	HIS	2.2
1	A	408	ILE	2.2
1	A	426	VAL	2.1
1	A	425	HIS	2.1
1	B	325	MET	2.1
1	B	401	MET	2.1
1	A	487	HIS	2.0
1	B	560	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 monosaccharides in this entry.

6.4 Ligands [i](#)

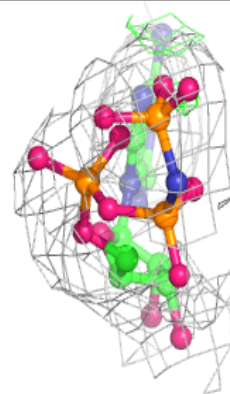
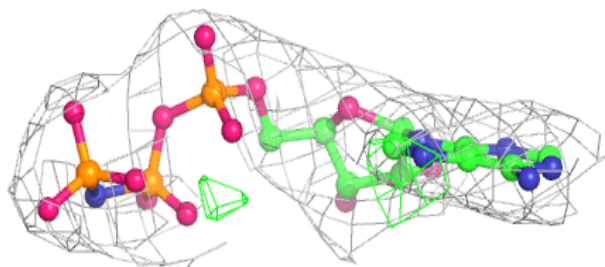
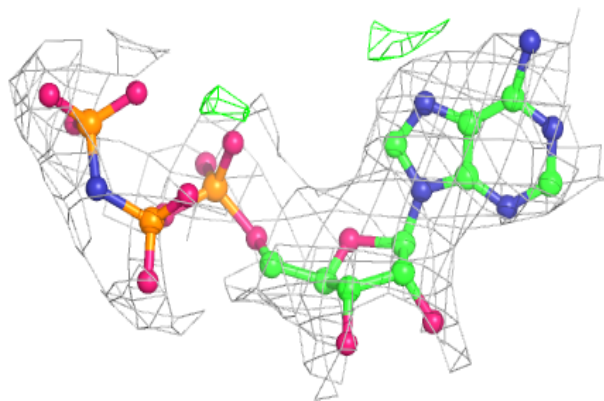
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ANP	B	702	31/31	0.90	0.17	81,95,108,112	0
2	MG	A	701	1/1	0.94	0.36	103,103,103,103	0
3	ANP	A	702	31/31	0.96	0.19	44,61,77,82	0
2	MG	B	701	1/1	0.96	0.35	201,201,201,201	0

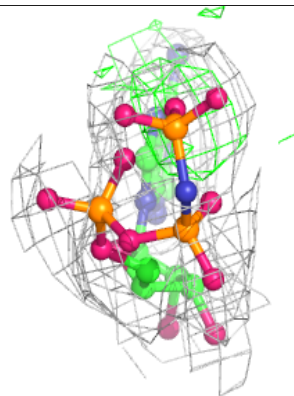
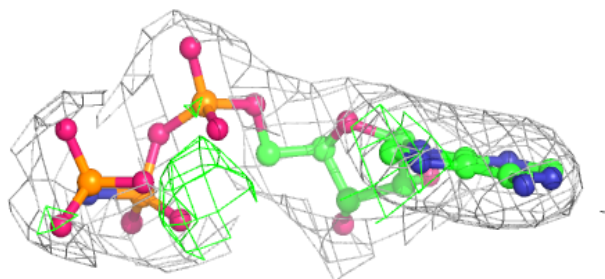
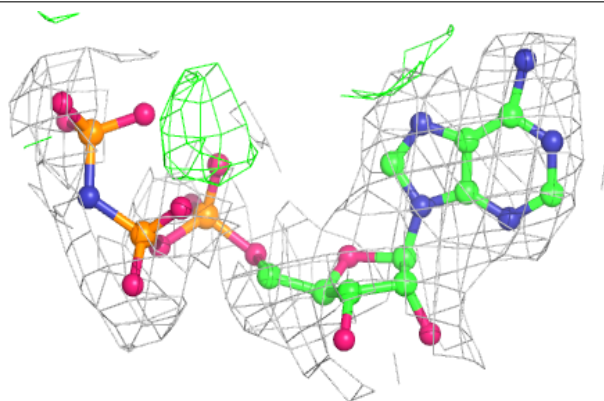
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ANP B 702:

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

**Electron density around ANP A 702:**

$2mF_o-DF_c$ (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.