



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:40 PM UTC

PDB ID : 5C1O / pdb_00005c1o
Title : Crystal structure of AMP-PNP complexed D-alanine-D-alanine ligase(DDL)
from *Yersinia pestis*
Authors : Tran, H.T.; Kang, L.W.; Hong, M.K.
Deposited on : 2015-06-15
Resolution : 2.50 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)
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

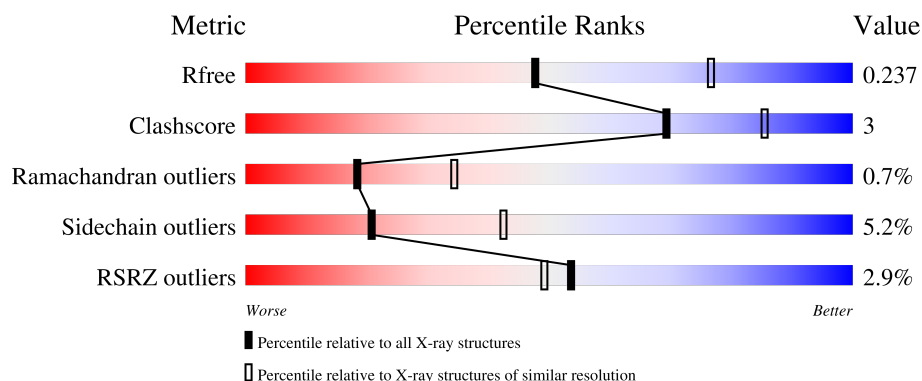
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	306	4% 81% 15% ..
1	B	306	2% 81% 13% 5%
1	C	306	0% 91% 8% .
1	D	306	4% 83% 13% ..

2 Entry composition [i](#)

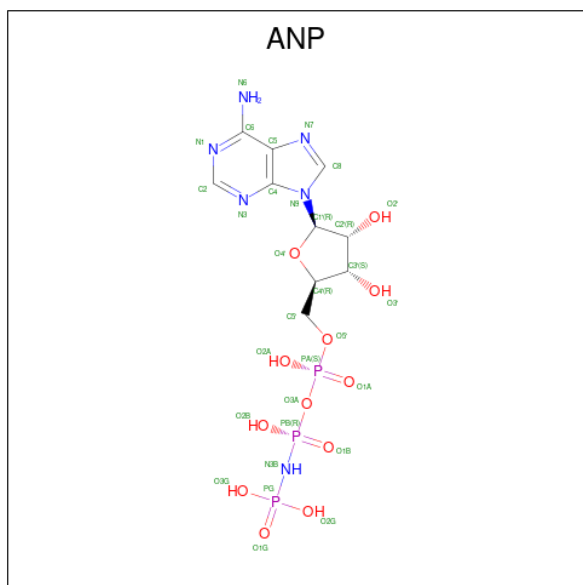
There are 5 unique types of molecules in this entry. The entry contains 9210 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 D-alanine–D-alanine ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2265	1446	374	434	11			
1	B	290	Total	C	N	O	S	0	0	0
			2203	1403	366	423	11			
1	C	306	Total	C	N	O	S	0	0	0
			2332	1487	384	450	11			
1	D	295	Total	C	N	O	S	0	0	0
			2241	1429	372	429	11			

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$).



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

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

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	2	Total	Na	0	0
			2	2		
3	C	1	Total	Na	0	0
			1	1		
3	D	1	Total	Na	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

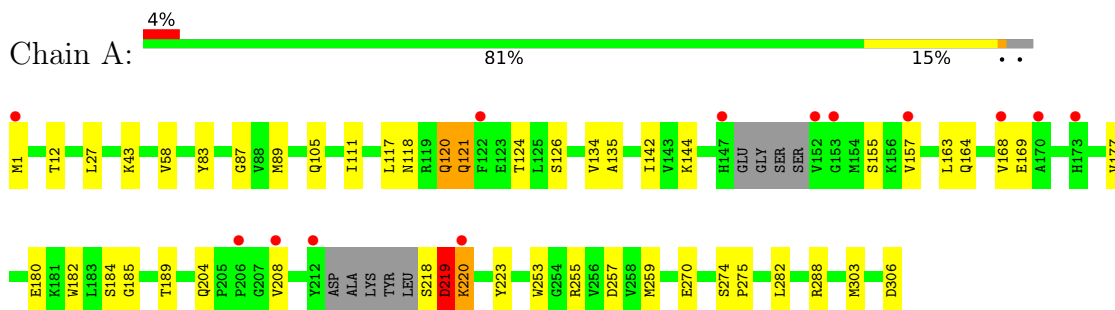
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	8	Total	O	0	0
			8	8		
5	B	7	Total	O	0	0
			7	7		
5	C	17	Total	O	0	0
			17	17		
5	D	4	Total	O	0	0
			4	4		

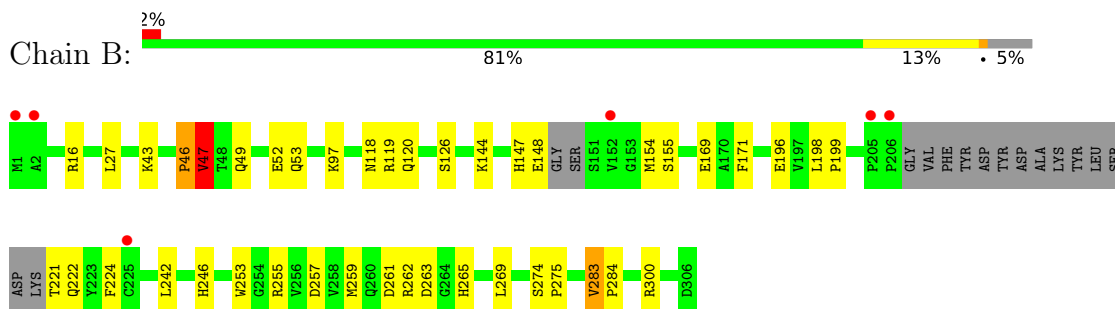
3 Residue-property plots [i](#)

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.

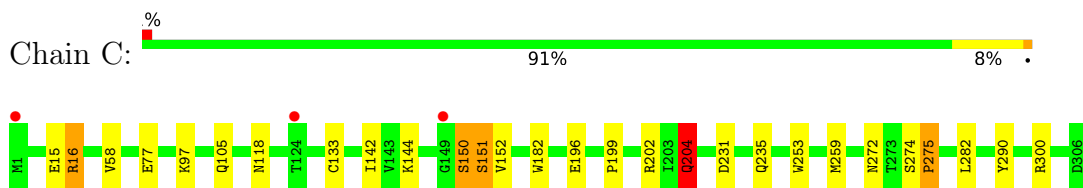
- Molecule 1: D-alanine–D-alanine ligase



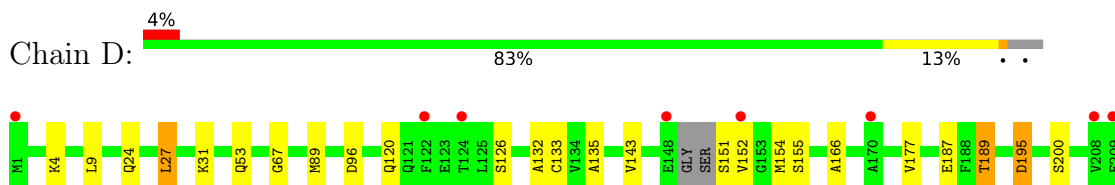
- Molecule 1: D-alanine–D-alanine ligase



- Molecule 1: D-alanine–D-alanine ligase



- Molecule 1: D-alanine–D-alanine ligase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.03Å 106.80Å 212.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.69 – 2.50 47.69 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.69-2.50) 99.8 (47.69-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	17.07 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.188 , 0.243 (Not available) , 0.237	Depositor DCC
R_{free} test set	2567 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	47.7	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 30.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9210	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.99% 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, MG, NA

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	1.01	2/2309 (0.1%)	1.01	4/3129 (0.1%)
1	B	1.06	3/2244 (0.1%)	1.09	5/3041 (0.2%)
1	C	1.03	1/2379 (0.0%)	1.08	11/3226 (0.3%)
1	D	0.94	1/2283 (0.0%)	1.01	0/3094
All	All	1.01	7/9215 (0.1%)	1.05	20/12490 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	274	SER	N-CA	7.74	1.55	1.46
1	A	274	SER	N-CA	5.98	1.54	1.46
1	B	246	HIS	C-O	-5.87	1.16	1.24
1	D	274	SER	N-CA	5.72	1.53	1.46
1	A	87	GLY	C-O	-5.64	1.18	1.23
1	C	274	SER	N-CA	5.34	1.52	1.46
1	B	274	SER	CA-C	5.02	1.58	1.52

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	ASP	N-CA-C	7.08	125.87	110.80
1	B	47	VAL	N-CA-CB	6.94	122.68	111.23
1	A	288	ARG	N-CA-C	-6.93	103.66	111.14
1	B	126	SER	CA-C-N	6.21	127.61	119.84
1	B	126	SER	C-N-CA	6.21	127.61	119.84
1	C	58	VAL	N-CA-C	6.19	116.77	108.11
1	C	204	GLN	CA-C-N	-5.96	114.24	120.38
1	C	204	GLN	C-N-CA	-5.96	114.24	120.38
1	C	77	GLU	N-CA-C	-5.85	104.98	111.36
1	A	126	SER	CA-C-N	5.80	127.09	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	SER	C-N-CA	5.80	127.09	119.84
1	C	274	SER	CA-C-N	-5.69	114.75	120.21
1	C	274	SER	C-N-CA	-5.69	114.75	120.21
1	C	282	LEU	CA-C-N	-5.54	115.90	120.33
1	C	282	LEU	C-N-CA	-5.54	115.90	120.33
1	C	275	PRO	CA-C-N	5.51	125.21	119.92
1	C	275	PRO	C-N-CA	5.51	125.21	119.92
1	C	144	LYS	N-CA-C	5.36	116.91	108.23
1	B	147	HIS	N-CA-C	5.15	117.10	107.99
1	B	283	VAL	N-CA-CB	5.10	114.43	110.45

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	2265	0	2260	18	0
1	B	2203	0	2206	20	0
1	C	2332	0	2323	10	0
1	D	2241	0	2244	14	0
2	A	31	0	13	0	0
2	B	31	0	13	0	0
2	C	31	0	13	3	0
2	D	31	0	13	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	8	0	0	0	0
5	B	7	0	0	0	0
5	C	17	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	4	0	0	0	0
All	All	9210	0	9085	63	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 3.

All (63) 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:C:151:SER:HB3	2:C:401:ANP:O3G	1.87	0.75
1:C:150:SER:O	1:C:152:VAL:N	2.26	0.68
1:A:253:TRP:CH2	1:A:275:PRO:HD2	2.31	0.65
1:B:259:MET:HG3	1:B:269:LEU:HD11	1.77	0.65
1:B:263:ASP:OD1	1:B:265:HIS:CD2	2.52	0.63
1:C:202:ARG:NH1	1:C:204:GLN:OE1	2.33	0.62
1:A:144:LYS:NZ	1:A:180:GLU:OE1	2.29	0.61
1:B:196:GLU:OE2	1:B:300:ARG:NH2	2.33	0.60
1:A:142:ILE:HD11	1:A:182:TRP:HD1	1.69	0.58
1:C:196:GLU:OE1	1:C:290:TYR:OH	2.19	0.57
1:B:255:ARG:NH1	1:B:257:ASP:OD1	2.38	0.55
1:C:253:TRP:CZ3	1:C:275:PRO:HG2	2.42	0.54
1:C:142:ILE:HD11	1:C:182:TRP:HD1	1.75	0.51
1:B:253:TRP:CH2	1:B:275:PRO:HD2	2.46	0.50
1:A:255:ARG:NH1	1:A:257:ASP:OD1	2.45	0.50
1:D:189:THR:OG1	1:D:255:ARG:HD3	2.11	0.50
1:A:223:TYR:CE2	1:A:282:LEU:HD21	2.47	0.49
1:A:58:VAL:O	1:A:83:TYR:HA	2.12	0.49
1:A:118:ASN:OD1	1:A:120:GLN:N	2.45	0.49
1:A:182:TRP:CH2	1:A:184:SER:HA	2.48	0.49
1:D:27:LEU:HD22	1:D:31:LYS:HD3	1.95	0.48
2:C:401:ANP:H8	2:C:401:ANP:O5'	2.13	0.48
1:A:184:SER:OG	1:A:185:GLY:O	2.30	0.47
1:B:97:LYS:NZ	1:B:148:GLU:OE1	2.43	0.47
1:C:272:ASN:C	1:C:272:ASN:OD1	2.58	0.47
1:A:218:SER:OG	1:A:219:ASP:N	2.48	0.46
1:D:225:CYS:SG	1:D:289:GLN:HG2	2.55	0.46
1:B:261:ASP:OD1	1:B:263:ASP:OD1	2.32	0.46
1:C:150:SER:C	1:C:152:VAL:H	2.23	0.46
1:A:218:SER:O	1:A:220:LYS:N	2.43	0.46
1:D:240:LEU:HG	1:D:268:LEU:HD21	1.98	0.46
1:A:117:LEU:HA	1:A:121:GLN:OE1	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:VAL:O	1:A:135:ALA:C	2.59	0.45
1:B:283:VAL:HB	1:B:284:PRO:HD3	1.98	0.45
1:C:97:LYS:HE2	2:C:401:ANP:PB	2.57	0.45
1:B:144:LYS:HB3	1:B:154:MET:HG3	1.98	0.45
1:B:119:ARG:NH2	1:B:171:PHE:O	2.46	0.44
1:D:132:ALA:HA	1:D:135:ALA:HB2	1.98	0.44
1:D:230:SER:O	1:D:231:ASP:C	2.60	0.44
1:B:46:PRO:O	1:B:47:VAL:HG22	2.17	0.44
1:B:46:PRO:O	1:B:47:VAL:HG13	2.18	0.44
1:A:12:THR:HG22	1:A:43:LYS:HG3	2.00	0.43
1:B:49:GLN:HG2	1:B:52:GLU:OE2	2.17	0.43
1:A:163:LEU:O	1:A:164:GLN:C	2.62	0.43
1:B:242:LEU:HD12	1:B:242:LEU:HA	1.83	0.43
1:D:143:VAL:O	1:D:154:MET:HA	2.18	0.43
1:D:285:MET:HE3	1:D:285:MET:HB2	1.88	0.43
1:D:195:ASP:OD1	1:D:195:ASP:N	2.51	0.42
1:A:105:GLN:HB2	1:A:111:ILE:HD11	2.00	0.42
1:D:242:LEU:HD12	1:D:242:LEU:HA	1.91	0.42
1:D:96:ASP:OD1	1:D:96:ASP:C	2.63	0.42
1:D:283:VAL:HB	1:D:284:PRO:HD3	2.01	0.42
1:A:89:MET:HE3	1:A:89:MET:HB3	1.94	0.42
1:D:155:SER:HB2	1:D:166:ALA:HB1	2.01	0.42
1:B:53:GLN:O	1:B:53:GLN:HG3	2.20	0.42
1:B:118:ASN:OD1	1:B:120:GLN:HG3	2.20	0.41
1:D:253:TRP:CH2	1:D:275:PRO:HD2	2.56	0.41
1:B:16:ARG:NH2	1:B:43:LYS:HE3	2.36	0.41
1:B:198:LEU:HB3	1:B:199:PRO:CD	2.50	0.41
1:B:255:ARG:HH11	1:B:255:ARG:HG2	1.86	0.41
1:A:253:TRP:CZ3	1:A:275:PRO:HD2	2.56	0.40
1:C:15:GLU:O	1:C:16:ARG:C	2.65	0.40
1:B:224:PHE:CD1	1:B:224:PHE:N	2.89	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	291/306 (95%)	276 (95%)	12 (4%)	3 (1%)	12	24
1	B	284/306 (93%)	277 (98%)	5 (2%)	2 (1%)	18	34
1	C	304/306 (99%)	291 (96%)	11 (4%)	2 (1%)	18	34
1	D	289/306 (94%)	282 (98%)	6 (2%)	1 (0%)	36	55
All	All	1168/1224 (95%)	1126 (96%)	34 (3%)	8 (1%)	18	34

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	121	GLN
1	A	219	ASP
1	B	47	VAL
1	C	151	SER
1	C	16	ARG
1	B	46	PRO
1	D	67	GLY

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	244/251 (97%)	228 (93%)	16 (7%)	15	32
1	B	238/251 (95%)	232 (98%)	6 (2%)	42	69
1	C	251/251 (100%)	241 (96%)	10 (4%)	28	54
1	D	242/251 (96%)	223 (92%)	19 (8%)	11	24
All	All	975/1004 (97%)	924 (95%)	51 (5%)	21	42

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	27	LEU
1	A	124	THR
1	A	155	SER
1	A	157	VAL
1	A	168	VAL
1	A	169	GLU
1	A	177	VAL
1	A	189	THR
1	A	204	GLN
1	A	208	VAL
1	A	220	LYS
1	A	259	MET
1	A	270	GLU
1	A	303	MET
1	A	306	ASP
1	B	27	LEU
1	B	155	SER
1	B	169	GLU
1	B	221	THR
1	B	222	GLN
1	B	262	ARG
1	C	105	GLN
1	C	118	ASN
1	C	133	CYS
1	C	150	SER
1	C	199	PRO
1	C	204	GLN
1	C	231	ASP
1	C	235	GLN
1	C	259	MET
1	C	300	ARG
1	D	4	LYS
1	D	9	LEU
1	D	24	GLN
1	D	27	LEU
1	D	53	GLN
1	D	89	MET
1	D	120	GLN
1	D	126	SER
1	D	133	CYS
1	D	151	SER
1	D	152	VAL

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Mol	Chain	Res	Type
1	D	177	VAL
1	D	187	GLU
1	D	189	THR
1	D	195	ASP
1	D	200	SER
1	D	219	ASP
1	D	263	ASP
1	D	306	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	105	GLN
1	A	147	HIS
1	A	159	HIS
1	A	246	HIS
1	B	49	GLN
1	B	235	GLN
1	B	265	HIS
1	B	272	ASN
1	C	24	GLN
1	C	147	HIS
1	C	173	HIS
1	C	235	GLN
1	D	24	GLN
1	D	222	GLN
1	D	243	GLN

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

Of 13 ligands modelled in this entry, 9 are monoatomic - leaving 4 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
2	ANP	A	401	4	33,33,33	2.14	12 (36%)	45,52,52	2.07	15 (33%)
2	ANP	D	401	4	33,33,33	2.03	10 (30%)	45,52,52	2.27	13 (28%)
2	ANP	B	401	4	33,33,33	2.19	10 (30%)	45,52,52	2.44	11 (24%)
2	ANP	C	401	4	33,33,33	1.99	9 (27%)	45,52,52	2.45	17 (37%)

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	A	401	4	-	5/18/38/38	0/3/3/3
2	ANP	D	401	4	-	2/18/38/38	0/3/3/3
2	ANP	B	401	4	-	3/18/38/38	0/3/3/3
2	ANP	C	401	4	-	4/18/38/38	0/3/3/3

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	ANP	C5-C4	5.41	1.48	1.39
2	C	401	ANP	PG-O1G	5.41	1.54	1.46
2	B	401	ANP	C5-C4	5.35	1.48	1.39
2	A	401	ANP	C5-C4	5.24	1.48	1.39
2	C	401	ANP	C5-C4	4.70	1.47	1.39
2	A	401	ANP	PB-O1B	4.70	1.53	1.46
2	B	401	ANP	PB-N3B	4.19	1.74	1.63
2	B	401	ANP	PA-O3A	4.18	1.64	1.59
2	A	401	ANP	PG-N3B	4.10	1.74	1.63
2	B	401	ANP	PB-O3A	4.09	1.64	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ANP	PG-N3B	4.06	1.74	1.63
2	A	401	ANP	PB-N3B	3.98	1.73	1.63
2	D	401	ANP	PB-N3B	3.93	1.73	1.63
2	D	401	ANP	PG-N3B	3.78	1.73	1.63
2	B	401	ANP	PG-O1G	3.67	1.51	1.46
2	D	401	ANP	PB-O3A	3.59	1.63	1.59
2	B	401	ANP	PB-O1B	3.59	1.51	1.46
2	C	401	ANP	PB-O1B	3.50	1.51	1.46
2	A	401	ANP	PG-O1G	3.49	1.51	1.46
2	C	401	ANP	PG-N3B	3.23	1.71	1.63
2	B	401	ANP	C8-N7	3.09	1.37	1.31
2	D	401	ANP	PA-O3A	2.98	1.62	1.59
2	C	401	ANP	C5-N7	-2.98	1.33	1.39
2	D	401	ANP	C8-N7	2.86	1.37	1.31
2	D	401	ANP	PB-O1B	2.84	1.50	1.46
2	D	401	ANP	C5-C6	2.63	1.48	1.41
2	A	401	ANP	C4-N9	-2.61	1.32	1.37
2	A	401	ANP	C8-N7	2.60	1.36	1.31
2	A	401	ANP	PB-O3A	2.60	1.62	1.59
2	C	401	ANP	PB-N3B	2.58	1.70	1.63
2	A	401	ANP	PA-O3A	2.52	1.62	1.59
2	A	401	ANP	C5-N7	-2.49	1.34	1.39
2	D	401	ANP	PG-O1G	2.46	1.49	1.46
2	C	401	ANP	C4-N9	-2.45	1.32	1.37
2	C	401	ANP	C5-C6	2.37	1.47	1.41
2	A	401	ANP	PB-O2B	-2.36	1.50	1.56
2	C	401	ANP	C8-N7	2.28	1.36	1.31
2	A	401	ANP	C5-C6	2.27	1.47	1.41
2	B	401	ANP	C4-N9	-2.26	1.33	1.37
2	B	401	ANP	C5-N7	-2.15	1.35	1.39
2	D	401	ANP	C5-N7	-2.09	1.35	1.39

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	ANP	O1G-PG-N3B	-9.27	98.11	111.77
2	C	401	ANP	O1G-PG-N3B	-8.58	99.14	111.77
2	D	401	ANP	C5-C4-N3	-6.68	117.52	126.72
2	C	401	ANP	O2B-PB-O1B	6.41	123.61	109.87
2	D	401	ANP	N3-C4-N9	5.53	136.57	127.17
2	B	401	ANP	O2B-PB-O1B	5.14	120.89	109.87
2	D	401	ANP	O1G-PG-N3B	-5.11	104.24	111.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	ANP	C5-C4-N3	-5.04	119.78	126.72
2	A	401	ANP	O2G-PG-O3G	4.82	120.54	107.59
2	C	401	ANP	C5-C4-N3	-4.69	120.26	126.72
2	A	401	ANP	O2B-PB-O1B	4.61	119.77	109.87
2	B	401	ANP	O1B-PB-N3B	-4.53	105.11	111.77
2	A	401	ANP	C5-C4-N3	-4.42	120.63	126.72
2	A	401	ANP	O1G-PG-N3B	-4.36	105.36	111.77
2	D	401	ANP	C2-N3-C4	4.24	122.19	111.83
2	B	401	ANP	N3-C4-N9	4.11	134.15	127.17
2	B	401	ANP	N3-C2-N1	-3.87	122.72	128.58
2	D	401	ANP	N3-C2-N1	-3.78	122.86	128.58
2	C	401	ANP	O3G-PG-O1G	3.75	122.86	113.45
2	C	401	ANP	N3-C4-N9	3.68	133.43	127.17
2	D	401	ANP	O1B-PB-N3B	-3.54	106.56	111.77
2	D	401	ANP	O2G-PG-O3G	3.54	117.10	107.59
2	A	401	ANP	C4-C5-N7	-3.51	106.57	110.58
2	A	401	ANP	O2G-PG-O1G	-3.50	104.67	113.45
2	B	401	ANP	C2-N1-C6	3.45	124.40	118.73
2	D	401	ANP	O2B-PB-O1B	3.40	117.16	109.87
2	C	401	ANP	C4-C5-N7	-3.39	106.71	110.58
2	B	401	ANP	C2-N3-C4	3.28	119.83	111.83
2	D	401	ANP	C4-C5-N7	-3.24	106.88	110.58
2	C	401	ANP	N3-C2-N1	-3.21	123.72	128.58
2	A	401	ANP	N3-C2-N1	-3.20	123.74	128.58
2	A	401	ANP	N3-C4-N9	3.06	132.38	127.17
2	A	401	ANP	C2-N3-C4	3.02	119.21	111.83
2	C	401	ANP	C2-N3-C4	2.91	118.94	111.83
2	D	401	ANP	C4-N9-C8	2.86	108.74	105.74
2	B	401	ANP	C2'-C1'-N9	-2.77	106.43	113.30
2	A	401	ANP	C2-N1-C6	2.70	123.16	118.73
2	B	401	ANP	C4-N9-C8	2.66	108.53	105.74
2	C	401	ANP	C4-N9-C8	2.61	108.48	105.74
2	C	401	ANP	O2A-PA-O3A	2.57	114.23	107.27
2	B	401	ANP	C4-C5-N7	-2.56	107.65	110.58
2	C	401	ANP	C2-N1-C6	2.53	122.89	118.73
2	A	401	ANP	O3G-PG-O1G	2.52	119.77	113.45
2	A	401	ANP	C6-C5-N7	2.39	136.69	132.09
2	D	401	ANP	O2A-PA-O1A	2.37	123.48	112.44
2	C	401	ANP	C5-N7-C8	2.35	107.15	103.45
2	C	401	ANP	O3'-C3'-C4'	2.29	117.67	111.08
2	C	401	ANP	O1B-PB-N3B	-2.29	108.41	111.77
2	D	401	ANP	C5-N7-C8	2.28	107.04	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	ANP	C5-N7-C8	2.25	106.99	103.45
2	C	401	ANP	C6-C5-N7	2.18	136.29	132.09
2	A	401	ANP	C4-N9-C8	2.14	107.98	105.74
2	C	401	ANP	O2G-PG-O3G	2.09	113.21	107.59
2	D	401	ANP	C6-C5-N7	2.08	136.10	132.09
2	A	401	ANP	C2'-C3'-C4'	-2.03	98.68	102.61
2	C	401	ANP	C4-N9-C1'	-2.01	121.92	126.63

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	ANP	PB-N3B-PG-O1G
2	A	401	ANP	C5'-O5'-PA-O1A
2	A	401	ANP	C5'-O5'-PA-O3A
2	B	401	ANP	PB-N3B-PG-O1G
2	B	401	ANP	PG-N3B-PB-O1B
2	C	401	ANP	PB-N3B-PG-O1G
2	C	401	ANP	PG-N3B-PB-O1B
2	D	401	ANP	PB-N3B-PG-O1G
2	D	401	ANP	PB-O3A-PA-O1A
2	A	401	ANP	C5'-O5'-PA-O2A
2	B	401	ANP	PB-O3A-PA-O1A
2	C	401	ANP	PB-O3A-PA-O2A
2	A	401	ANP	O4'-C4'-C5'-O5'
2	C	401	ANP	PB-O3A-PA-O1A

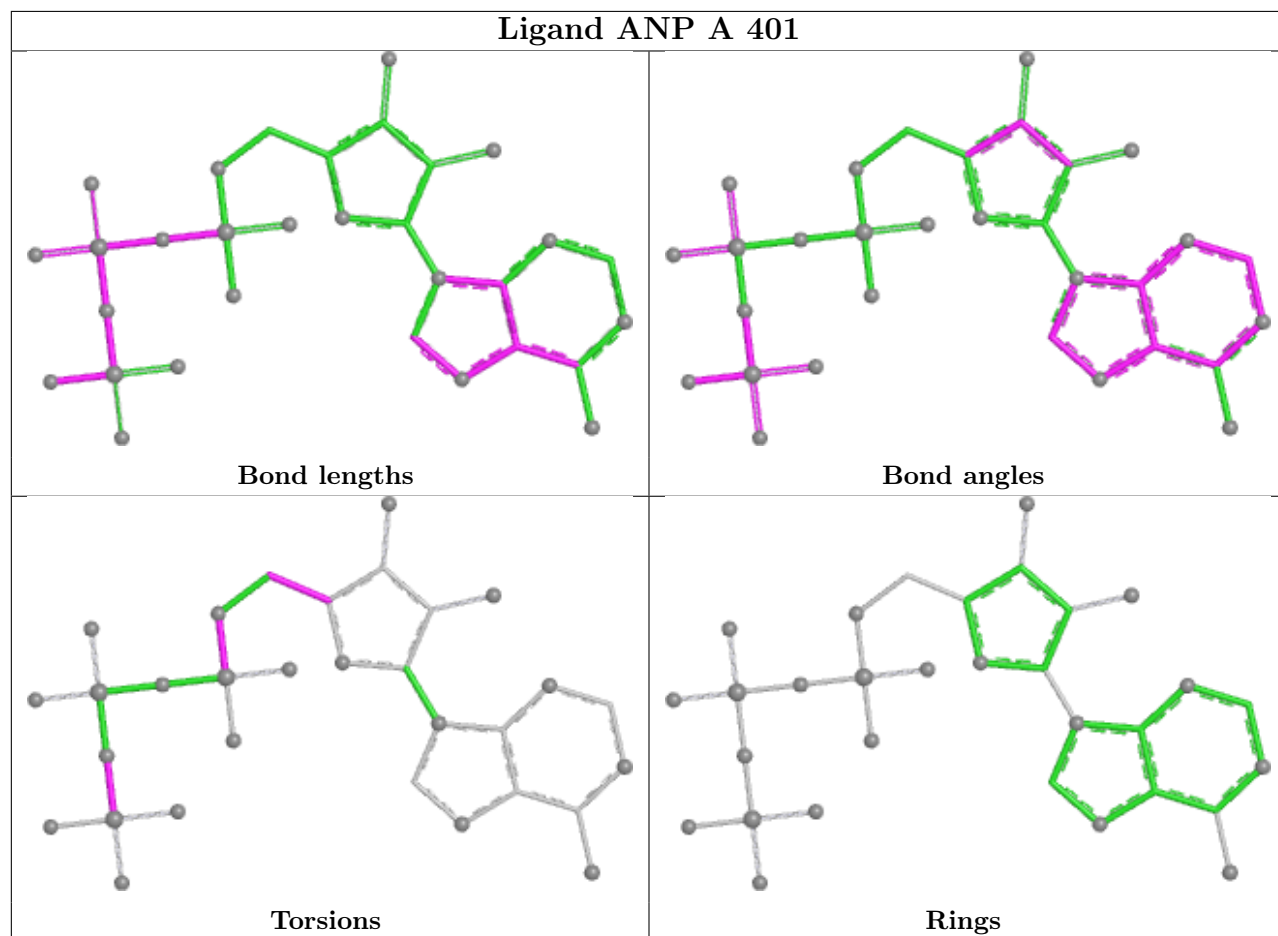
There are no ring outliers.

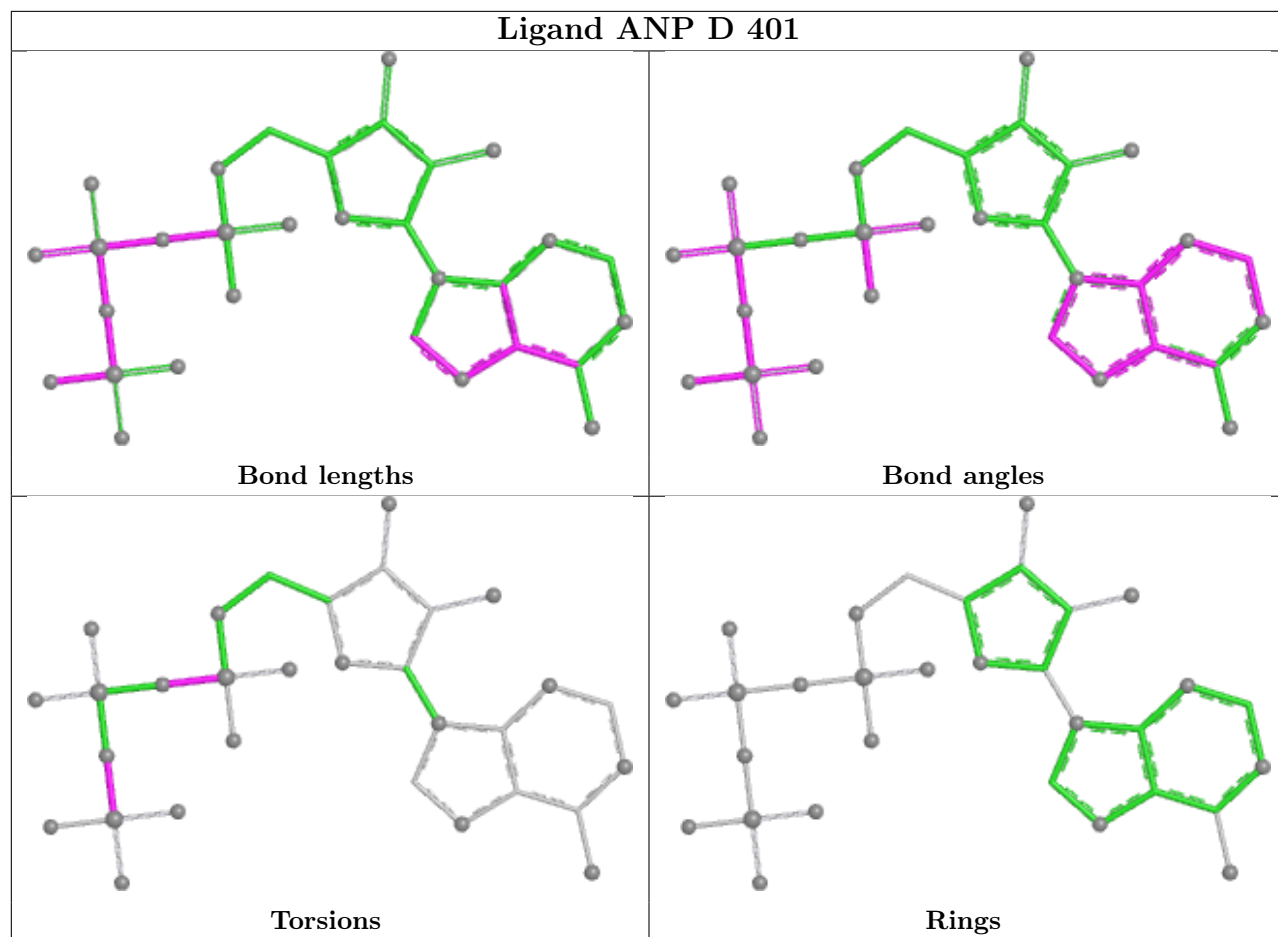
1 monomer is involved in 3 short contacts:

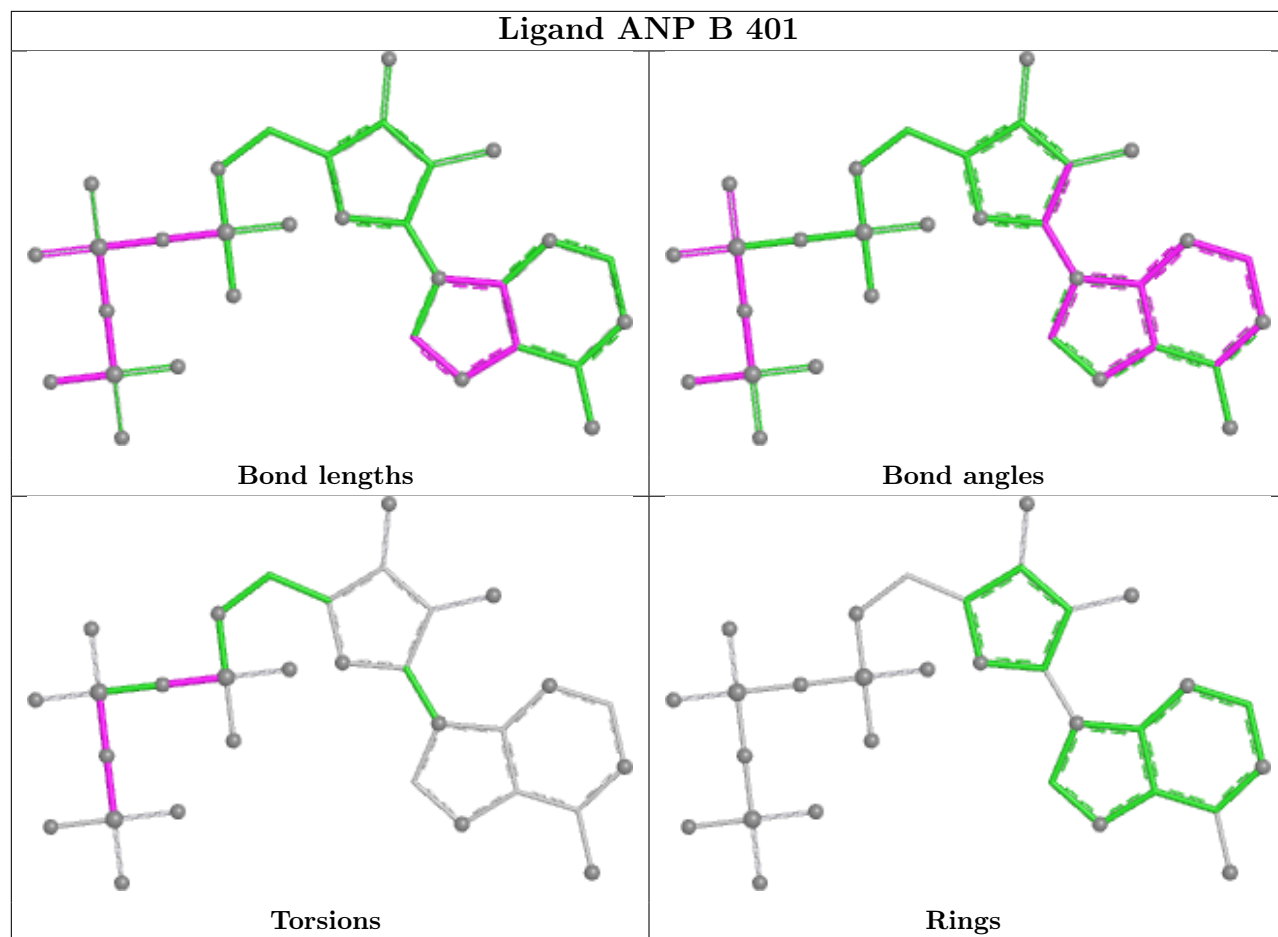
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	ANP	3	0

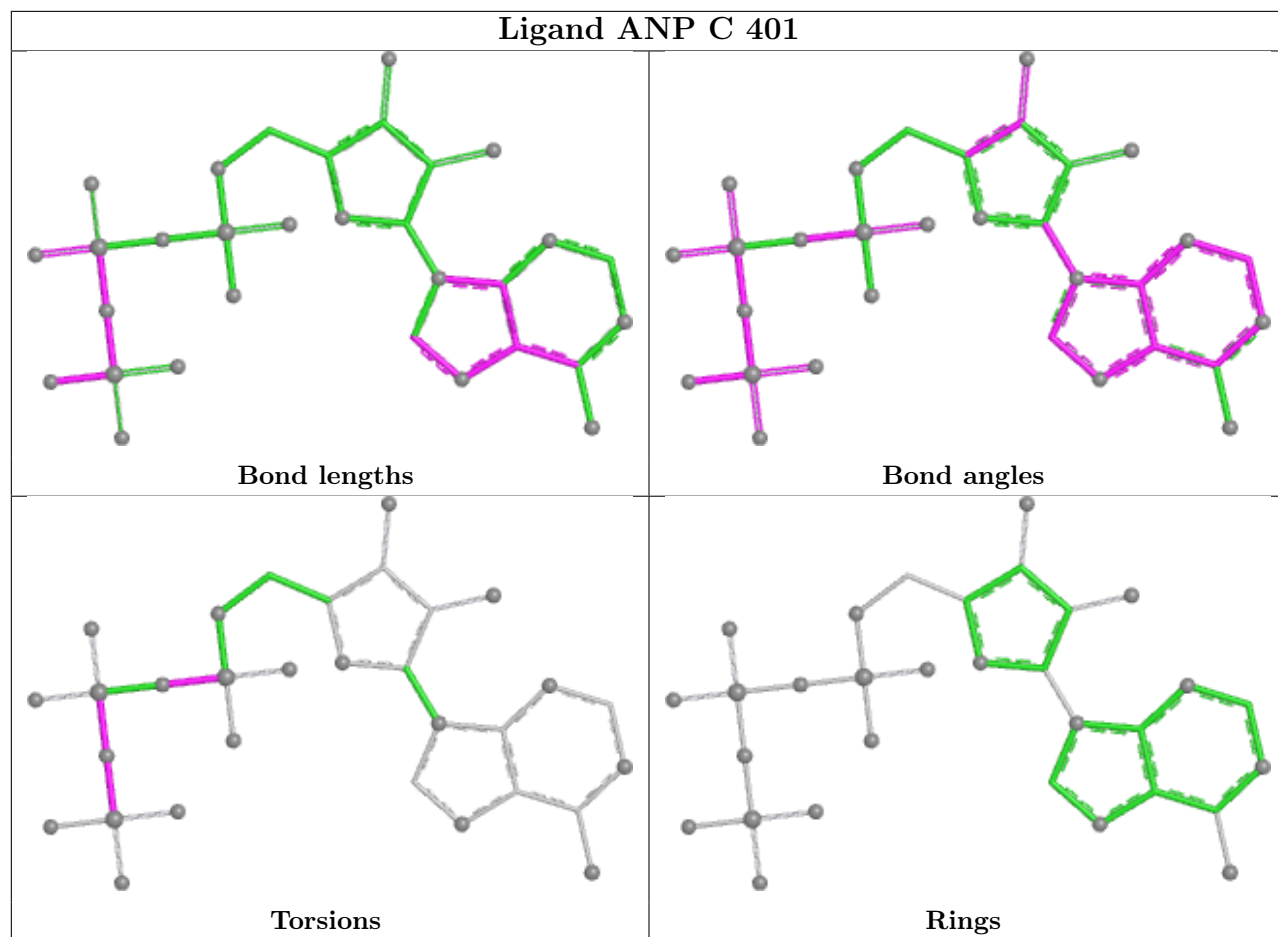
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

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	297/306 (97%)	0.11	13 (4%) 39 34	27, 51, 112, 127	0
1	B	290/306 (94%)	-0.12	6 (2%) 63 59	28, 46, 85, 104	0
1	C	306/306 (100%)	-0.24	3 (0%) 79 76	29, 44, 73, 99	0
1	D	295/306 (96%)	0.22	12 (4%) 41 37	33, 57, 100, 116	0
All	All	1188/1224 (97%)	-0.01	34 (2%) 53 49	27, 50, 95, 127	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	152	VAL	5.9
1	D	208	VAL	4.2
1	C	1	MET	3.7
1	A	212	TYR	3.5
1	A	1	MET	3.5
1	A	208	VAL	3.5
1	D	221	THR	3.1
1	D	170	ALA	3.0
1	A	173	HIS	3.0
1	C	124	THR	2.9
1	B	1	MET	2.9
1	B	152	VAL	2.9
1	B	205	PRO	2.9
1	D	209	PHE	2.8
1	A	170	ALA	2.7
1	A	153	GLY	2.7
1	A	168	VAL	2.6
1	A	147	HIS	2.6
1	D	152	VAL	2.5
1	D	220	LYS	2.5
1	B	2	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	206	PRO	2.4
1	D	226	PRO	2.4
1	D	148	GLU	2.4
1	B	206	PRO	2.4
1	D	1	MET	2.3
1	A	220	LYS	2.3
1	D	225	CYS	2.3
1	A	122	PHE	2.3
1	C	149	GLY	2.2
1	A	157	VAL	2.1
1	D	124	THR	2.1
1	B	225	CYS	2.1
1	D	122	PHE	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.

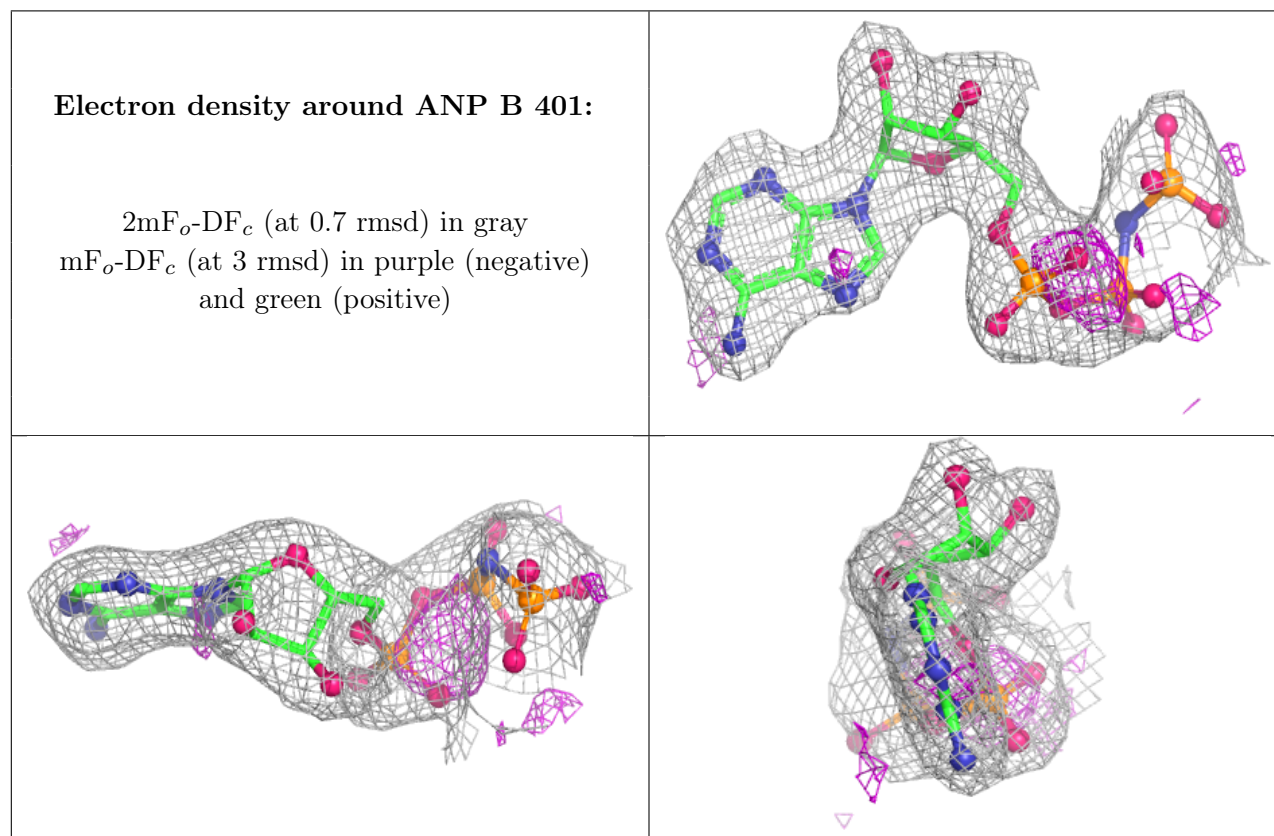
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NA	D	402	1/1	0.80	0.10	47,47,47,47	0
2	ANP	B	401	31/31	0.89	0.09	43,58,97,100	0
2	ANP	D	401	31/31	0.91	0.08	50,64,80,86	0
2	ANP	A	401	31/31	0.91	0.08	48,58,109,115	0
4	MG	C	403	1/1	0.91	0.08	63,63,63,63	0
3	NA	B	404	1/1	0.93	0.24	39,39,39,39	0
4	MG	A	403	1/1	0.95	0.09	77,77,77,77	0
4	MG	D	403	1/1	0.95	0.10	76,76,76,76	0
2	ANP	C	401	31/31	0.96	0.07	39,46,72,78	0
4	MG	B	403	1/1	0.96	0.14	88,88,88,88	0

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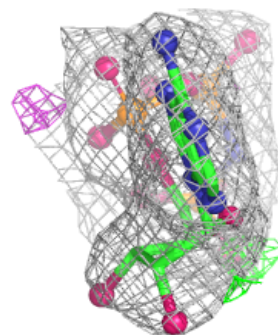
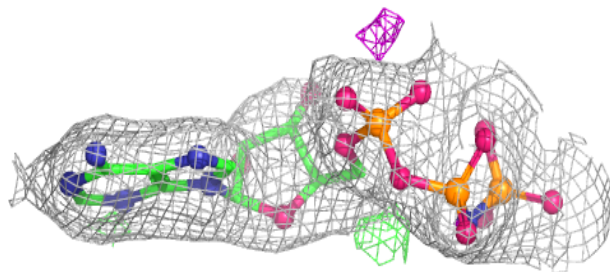
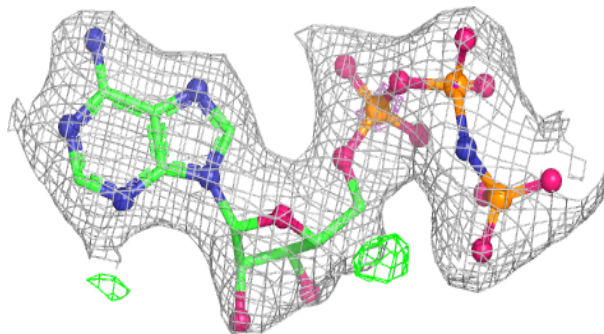
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	A	402	1/1	0.97	0.04	35,35,35,35	0
3	NA	C	402	1/1	0.98	0.08	38,38,38,38	0
3	NA	B	402	1/1	0.98	0.04	36,36,36,36	0

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.

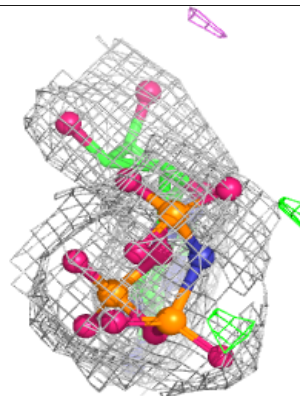
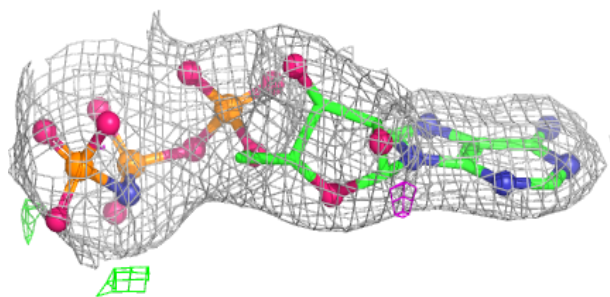
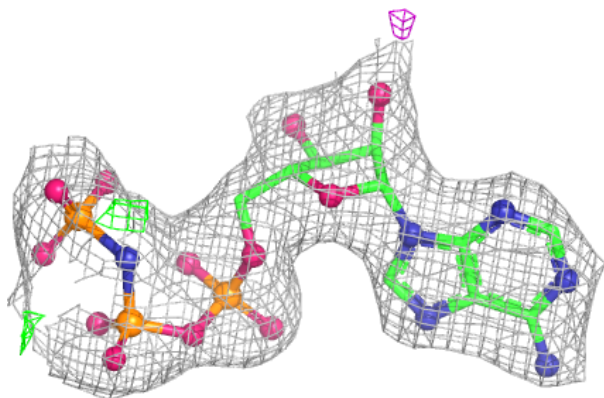


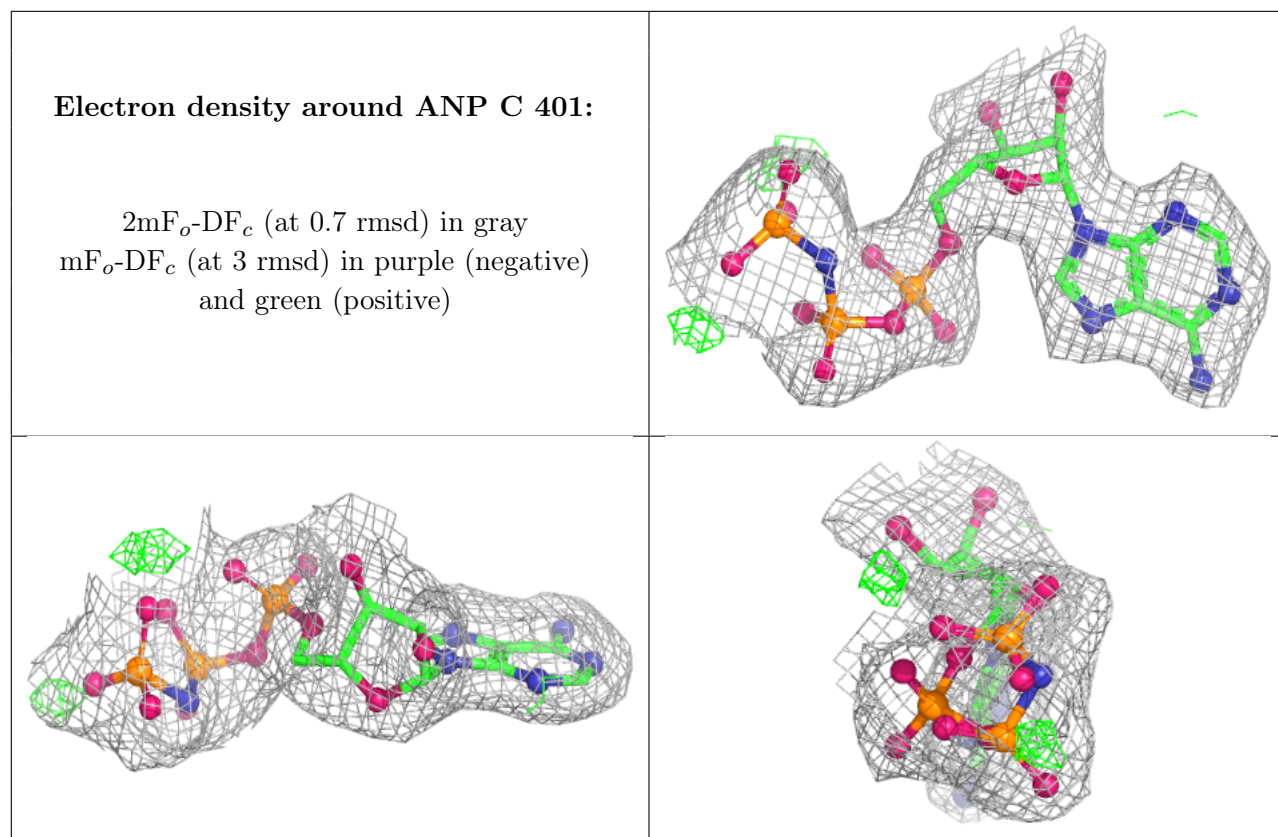
Electron density around ANP D 401:

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

**Electron density around ANP A 401:**

$2mF_o-DF_c$ (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.