



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 6, 2026 – 03:56 pm BST

PDB ID : 9RTR / pdb_00009rtr
Title : Crystal structure of BRAF:MEK1 complex with asymmetric dimer interface bound to AMPPNP
Authors : Kondo, Y.; Notbohm, J.; Camacho, I.N.; Nagy-Davidescu, G.; Mason, T.; Muhle, J.; Standfuss, J.; Perica, T.
Deposited on : 2025-07-03
Resolution : 2.60 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	:	1.8.4, CSD as541be (2020)
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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

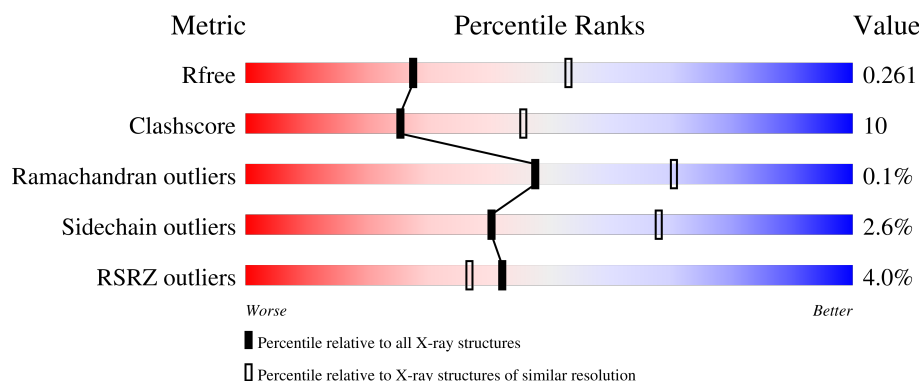
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	358	 3% 68% 18% 13%
1	D	358	 4% 70% 17% 12%
2	B	282	 4% 65% 29% 6%
2	C	282	 4% 71% 24% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9350 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 Dual specificity mitogen-activated protein kinase kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	D	315	Total	C	N	O	S	0	0	0
			2476	1580	421	458	17			
1	A	310	Total	C	N	O	S	0	0	0
			2436	1557	415	448	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	36	SER	-	expression tag	UNP Q02750
A	36	SER	-	expression tag	UNP Q02750

- Molecule 2 is a protein called Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	268	Total	C	N	O	S	0	0	0
			2151	1366	379	393	13			
2	B	266	Total	C	N	O	S	0	0	0
			2131	1352	377	389	13			

There are 36 discrepancies between the modelled and reference sequences:

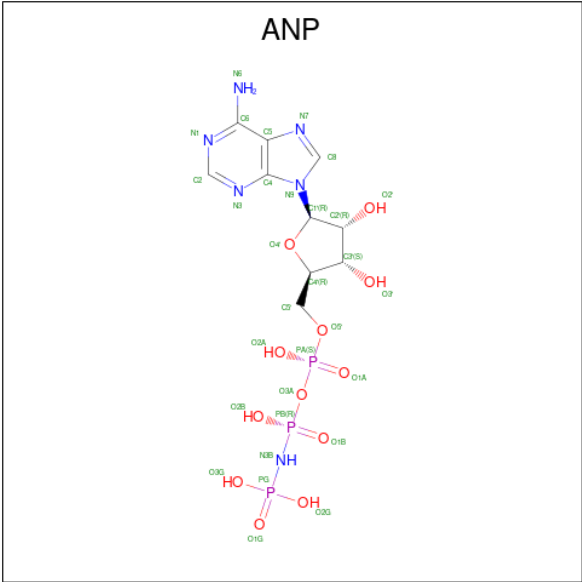
Chain	Residue	Modelled	Actual	Comment	Reference
C	442	SER	-	expression tag	UNP P15056
C	443	ASN	-	expression tag	UNP P15056
C	444	ALA	-	expression tag	UNP P15056
C	543	ALA	ILE	engineered mutation	UNP P15056
C	544	SER	ILE	engineered mutation	UNP P15056
C	551	LYS	ILE	engineered mutation	UNP P15056
C	562	ARG	GLN	engineered mutation	UNP P15056
C	588	ASN	LEU	engineered mutation	UNP P15056
C	600	GLU	VAL	engineered mutation	UNP P15056
C	630	SER	LYS	engineered mutation	UNP P15056

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Chain	Residue	Modelled	Actual	Comment	Reference
C	688	ARG	ALA	engineered mutation	UNP P15056
C	706	SER	LEU	engineered mutation	UNP P15056
C	709	ARG	GLN	engineered mutation	UNP P15056
C	713	GLU	SER	engineered mutation	UNP P15056
C	716	GLU	LEU	engineered mutation	UNP P15056
C	720	GLU	SER	engineered mutation	UNP P15056
C	722	SER	-	expression tag	UNP P15056
C	723	GLY	-	expression tag	UNP P15056
B	442	SER	-	expression tag	UNP P15056
B	443	ASN	-	expression tag	UNP P15056
B	444	ALA	-	expression tag	UNP P15056
B	543	ALA	ILE	engineered mutation	UNP P15056
B	544	SER	ILE	engineered mutation	UNP P15056
B	551	LYS	ILE	engineered mutation	UNP P15056
B	562	ARG	GLN	engineered mutation	UNP P15056
B	588	ASN	LEU	engineered mutation	UNP P15056
B	600	GLU	VAL	engineered mutation	UNP P15056
B	630	SER	LYS	engineered mutation	UNP P15056
B	688	ARG	ALA	engineered mutation	UNP P15056
B	706	SER	LEU	engineered mutation	UNP P15056
B	709	ARG	GLN	engineered mutation	UNP P15056
B	713	GLU	SER	engineered mutation	UNP P15056
B	716	GLU	LEU	engineered mutation	UNP P15056
B	720	GLU	SER	engineered mutation	UNP P15056
B	722	SER	-	expression tag	UNP P15056
B	723	GLY	-	expression tag	UNP P15056

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: $C_{10}H_{17}N_6O_{12}P_3$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	11	Total	O	0	0
			11	11		
5	C	22	Total	O	0	0
			22	22		
5	A	17	Total	O	0	0
			17	17		

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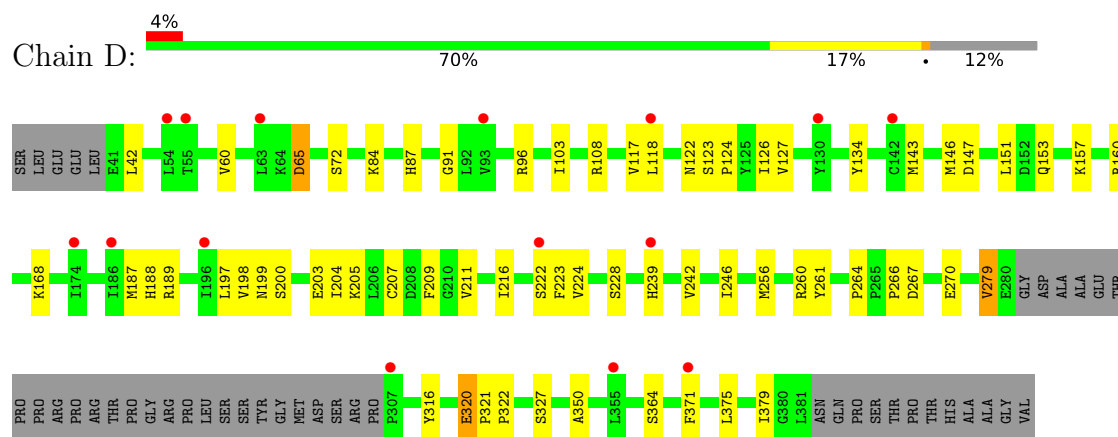
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	9	Total	O	0	0
			9	9		

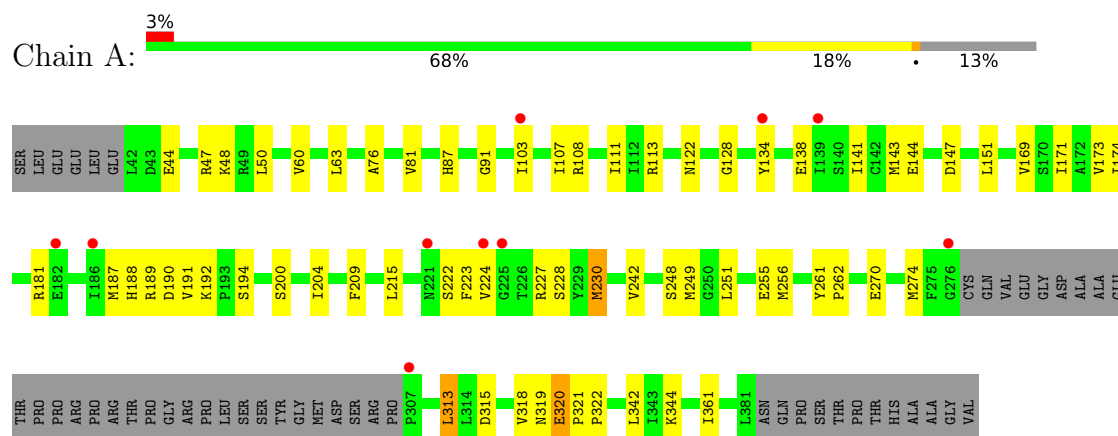
3 Residue-property plots

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.

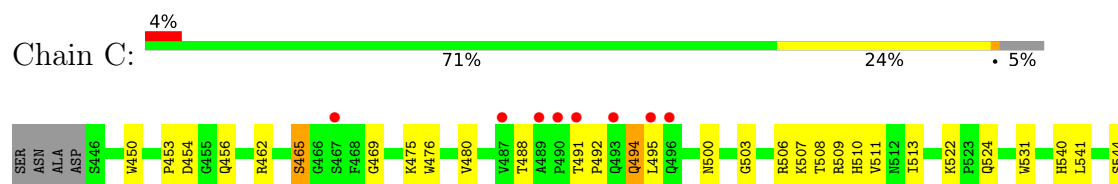
- Molecule 1: Dual specificity mitogen-activated protein kinase kinase 1

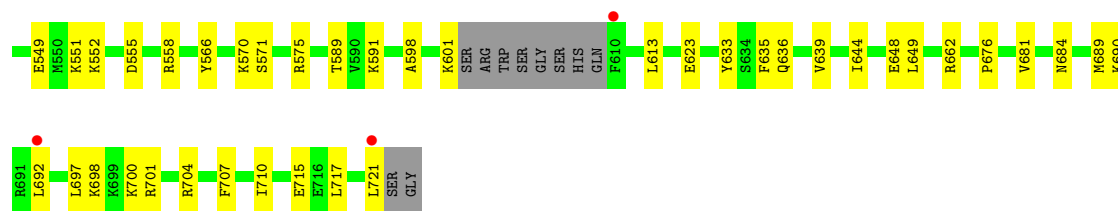


- Molecule 1: Dual specificity mitogen-activated protein kinase kinase 1

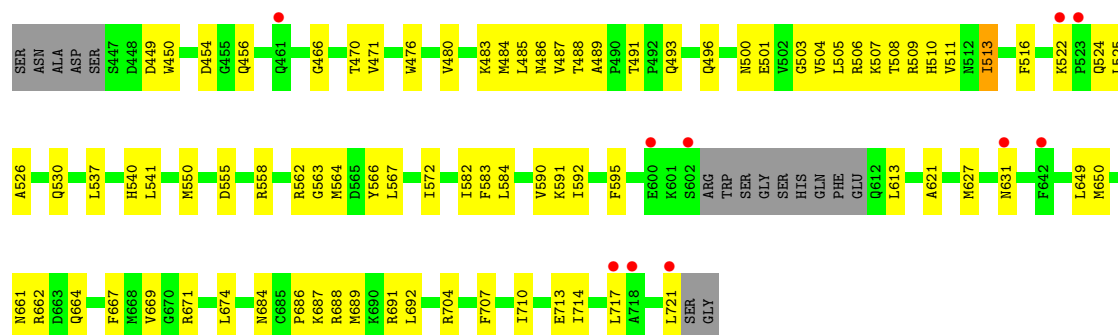


- Molecule 2: Serine/threonine-protein kinase B-raf





● Molecule 2: Serine/threonine-protein kinase B-raf



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.72Å 67.91Å 292.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.70 – 2.60 60.70 – 2.60	Depositor EDS
% Data completeness (in resolution range)	56.3 (60.70-2.60) 56.3 (60.70-2.60)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.28 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.208 , 0.262 0.208 , 0.261	Depositor DCC
R_{free} test set	1178 reflections (2.81%)	wwPDB-VP
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for k,h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9350	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% 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: MG, 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	0.10	0/2484	0.26	0/3343
1	D	0.10	0/2524	0.27	0/3397
2	B	0.11	0/2175	0.33	0/2930
2	C	0.12	0/2196	0.32	0/2958
All	All	0.11	0/9379	0.30	0/12628

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	2436	0	2473	43	0
1	D	2476	0	2507	38	0
2	B	2131	0	2145	56	0
2	C	2151	0	2160	49	0
3	A	31	0	13	4	0
3	B	31	0	13	2	0
3	D	31	0	13	5	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1	0	0	0	0
5	A	17	0	0	2	0
5	B	9	0	0	2	0
5	C	22	0	0	1	0
5	D	11	0	0	0	0
All	All	9350	0	9324	179	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 10.

All (179) 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:D:143:MET:HE2	3:D:901:ANP:HN61	1.47	0.79
3:D:901:ANP:H8	3:D:901:ANP:H5'2	1.66	0.78
1:A:181:ARG:HD2	1:A:242:VAL:HG11	1.69	0.75
2:B:650:MET:HE2	2:B:689:MET:HG2	1.68	0.74
2:B:613:LEU:HG	2:B:627:MET:HE3	1.69	0.74
2:B:503:GLY:O	2:B:507:LYS:NZ	2.24	0.70
2:C:698:LYS:O	2:C:704:ARG:NH1	2.27	0.68
2:C:508:THR:HB	2:C:513:ILE:HG21	1.76	0.68
1:D:146:MET:HE1	1:D:205:LYS:HD2	1.75	0.67
3:A:901:ANP:N3B	3:A:901:ANP:O1A	2.27	0.66
1:A:322:PRO:HG3	1:A:344:LYS:HG3	1.76	0.66
3:A:901:ANP:H5'1	3:A:901:ANP:H8	1.76	0.65
1:A:230:MET:SD	2:B:661:ASN:ND2	2.68	0.65
2:C:541:LEU:HD11	2:C:649:LEU:HD23	1.76	0.65
2:C:500:ASN:ND2	2:C:598:ALA:O	2.25	0.65
1:D:42:LEU:HD11	1:D:124:PRO:HG2	1.80	0.64
3:D:901:ANP:O2G	3:D:901:ANP:O1B	2.15	0.63
1:D:151:LEU:HD13	1:D:256:MET:HE1	1.81	0.63
2:B:496:GLN:NE2	2:B:500:ASN:OD1	2.31	0.62
2:B:508:THR:HB	2:B:513:ILE:HG21	1.82	0.62
1:A:113:ARG:NH2	5:A:1002:HOH:O	2.35	0.59
2:B:686:PRO:HG3	2:B:721:LEU:HD11	1.84	0.58
2:B:504:VAL:HG13	2:B:572:ILE:HD13	1.85	0.58
1:D:108:ARG:HD3	1:D:134:TYR:CD2	2.38	0.58
1:A:50:LEU:HD11	1:A:122:ASN:HB3	1.85	0.58
1:A:187:MET:HE2	1:A:189:ARG:HG2	1.85	0.57
1:A:224:VAL:HG21	2:B:613:LEU:HD22	1.86	0.57
1:A:315:ASP:OD1	1:A:319:ASN:ND2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:510:HIS:CD2	2:B:456:GLN:HE22	2.22	0.57
1:D:143:MET:HE1	1:D:207:CYS:HB2	1.87	0.56
2:B:501:GLU:HG3	2:B:595:PHE:HB2	1.87	0.56
1:D:188:HIS:CD2	1:D:209:PHE:HB3	2.41	0.56
1:A:194:SER:OG	3:A:901:ANP:O1B	2.20	0.56
2:C:462:ARG:NH1	2:C:465:SER:OG	2.38	0.56
3:A:901:ANP:H8	3:A:901:ANP:C5'	2.35	0.55
2:C:549:GLU:HG2	2:C:552:LYS:H	1.71	0.55
1:D:127:VAL:HG21	1:D:143:MET:HE3	1.88	0.55
1:D:123:SER:HB3	1:D:126:ILE:HG12	1.89	0.55
2:C:492:PRO:CD	2:C:495:LEU:HD13	2.37	0.54
1:D:168:LYS:NZ	1:D:364:SER:O	2.38	0.54
1:A:108:ARG:HD2	1:A:134:TYR:CD1	2.43	0.54
2:C:571:SER:HA	2:C:635:PHE:HZ	1.72	0.53
1:D:147:ASP:OD1	1:D:200:SER:N	2.41	0.53
2:C:488:THR:HA	2:C:524:GLN:HE22	1.74	0.53
2:C:692:LEU:HD21	2:C:710:ILE:HG23	1.89	0.53
2:B:484:MET:HE3	2:B:526:ALA:HB2	1.90	0.53
1:A:261:TYR:HE2	1:A:313:LEU:HD21	1.74	0.53
2:B:563:GLY:HA2	2:B:566:TYR:HB3	1.90	0.52
1:D:267:ASP:OD1	1:D:270:GLU:N	2.38	0.52
2:C:454:ASP:HB2	2:C:522:LYS:HB2	1.90	0.52
1:D:157:LYS:HD3	1:D:379:ILE:HA	1.91	0.52
2:C:511:VAL:O	2:C:591:LYS:NZ	2.43	0.51
2:B:541:LEU:HD21	2:B:649:LEU:HD23	1.93	0.51
2:B:583:PHE:HB3	2:B:591:LYS:HB2	1.92	0.51
1:D:187:MET:HE2	1:D:189:ARG:HG2	1.92	0.51
1:D:153:GLN:OE1	3:D:901:ANP:O2'	2.28	0.51
2:B:454:ASP:HB2	2:B:522:LYS:HD2	1.92	0.51
2:B:530:GLN:O	3:B:901:ANP:N6	2.44	0.50
2:C:503:GLY:O	2:C:507:LYS:NZ	2.45	0.50
2:C:698:LYS:HB2	2:C:704:ARG:HG2	1.93	0.50
1:A:318:VAL:HG21	2:B:667:PHE:CZ	2.46	0.50
2:B:558:ARG:NH2	2:B:562:ARG:HH12	2.10	0.50
1:D:127:VAL:HG11	1:D:143:MET:HE3	1.93	0.50
1:D:228:SER:O	1:D:261:TYR:OH	2.25	0.49
2:C:551:LYS:HE3	2:C:721:LEU:HB3	1.94	0.49
1:A:222:SER:HB3	2:B:662:ARG:HH11	1.77	0.49
2:C:509:ARG:HD2	2:B:476:TRP:HH2	1.75	0.49
1:D:65:ASP:OD1	1:D:96:ARG:NH2	2.45	0.49
1:A:103:ILE:HD12	1:A:107:ILE:HG22	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:ILE:HG13	1:A:249:MET:HE1	1.95	0.48
1:A:188:HIS:CD2	1:A:209:PHE:HB3	2.49	0.48
1:A:151:LEU:HD13	1:A:256:MET:HE1	1.95	0.48
1:A:228:SER:O	1:A:261:TYR:OH	2.30	0.48
1:A:76:ALA:HB2	1:A:81:VAL:HG13	1.95	0.48
2:B:450:TRP:CD1	2:B:506:ARG:HG2	2.48	0.48
2:C:707:PHE:HA	2:C:710:ILE:HB	1.96	0.48
2:B:484:MET:HE2	2:B:524:GLN:HB3	1.95	0.48
2:C:509:ARG:HD2	2:B:476:TRP:CH2	2.49	0.47
1:D:316:TYR:CE2	1:D:320:GLU:HG2	2.50	0.47
1:A:320:GLU:HG3	1:A:321:PRO:HD2	1.96	0.47
2:C:453:PRO:HD2	2:C:456:GLN:NE2	2.30	0.47
1:A:270:GLU:O	1:A:274:MET:HG3	2.13	0.47
1:A:181:ARG:NH1	5:A:1003:HOH:O	2.48	0.47
2:B:537:LEU:HD21	2:B:649:LEU:HD21	1.95	0.47
1:A:169:VAL:HA	1:A:204:ILE:HD13	1.97	0.47
2:C:456:GLN:NE2	2:C:476:TRP:HE1	2.13	0.47
1:A:87:HIS:O	1:A:91:GLY:N	2.48	0.46
1:A:318:VAL:HG21	2:B:667:PHE:CE1	2.51	0.46
2:B:471:VAL:HA	2:B:483:LYS:HA	1.98	0.46
1:D:260:ARG:NH2	1:D:264:PRO:O	2.48	0.46
2:B:511:VAL:O	2:B:591:LYS:NZ	2.44	0.46
2:C:566:TYR:CZ	2:C:570:LYS:HD2	2.51	0.46
2:B:485:LEU:HD23	2:B:525:LEU:HB2	1.97	0.46
1:A:319:ASN:ND2	2:B:671:ARG:HH21	2.14	0.46
2:B:669:VAL:HG23	2:B:674:LEU:HD23	1.96	0.46
2:B:631:ASN:HB3	5:B:1006:HOH:O	2.16	0.46
1:A:261:TYR:CE2	1:A:313:LEU:HD21	2.51	0.46
2:B:564:MET:HE2	2:B:567:LEU:HD12	1.98	0.45
2:C:506:ARG:O	2:B:449:ASP:HB3	2.16	0.45
2:C:701:ARG:HG2	2:C:704:ARG:NH2	2.30	0.45
1:D:87:HIS:O	1:D:91:GLY:N	2.48	0.45
2:B:470:THR:OG1	2:B:486:ASN:HB2	2.17	0.45
1:D:187:MET:HG3	1:D:189:ARG:HG3	1.96	0.45
1:D:222:SER:HB3	2:C:662:ARG:HD3	1.98	0.45
2:C:492:PRO:HB2	2:C:494:GLN:NE2	2.31	0.45
2:B:555:ASP:HA	2:B:558:ARG:HG2	1.99	0.45
2:B:550:MET:SD	2:B:650:MET:HE1	2.56	0.45
1:D:375:LEU:O	1:D:379:ILE:HG12	2.17	0.45
1:A:111:ILE:HG23	1:A:215:LEU:HB2	1.97	0.45
2:B:558:ARG:HB3	2:B:714:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:582:ILE:HG23	2:B:590:VAL:HG13	1.99	0.44
2:B:691:ARG:HD3	2:B:691:ARG:HA	1.75	0.44
2:B:505:LEU:HB3	2:B:516:PHE:CG	2.52	0.44
1:A:255:GLU:HB2	1:A:262:PRO:HD3	2.00	0.44
2:C:636:GLN:O	2:C:639:VAL:HG22	2.18	0.44
2:B:504:VAL:O	2:B:508:THR:HG23	2.17	0.44
2:B:688:ARG:HB2	2:B:717:LEU:HD13	1.99	0.44
2:B:684:ASN:OD1	2:B:684:ASN:N	2.46	0.43
2:C:601:LYS:HB3	2:C:601:LYS:HE3	1.67	0.43
1:D:204:ILE:HG12	1:D:371:PHE:HE2	1.83	0.43
2:C:644:ILE:O	2:C:648:GLU:HG3	2.18	0.43
1:A:251:LEU:N	1:A:342:LEU:HD21	2.33	0.43
2:C:681:VAL:HG21	2:C:690:LYS:HD2	2.00	0.43
1:D:65:ASP:CG	1:D:96:ARG:HH22	2.26	0.43
1:D:260:ARG:HD3	1:D:266:PRO:HG3	2.01	0.43
2:C:623:GLU:HG3	2:C:701:ARG:HD3	2.00	0.43
2:C:676:PRO:HG2	2:C:697:LEU:HD13	2.01	0.43
1:A:223:PHE:HB3	5:B:1001:HOH:O	2.17	0.43
1:A:173:VAL:HB	1:A:249:MET:HE3	2.01	0.43
2:B:488:THR:OG1	2:B:489:ALA:N	2.51	0.43
2:B:540:HIS:CD2	2:B:584:LEU:HD12	2.53	0.43
3:D:901:ANP:O2B	3:D:901:ANP:O1A	2.36	0.43
1:A:147:ASP:OD1	1:A:200:SER:N	2.52	0.43
1:A:222:SER:OG	1:A:223:PHE:N	2.52	0.43
2:C:456:GLN:HE21	2:C:476:TRP:HE1	1.67	0.42
1:D:108:ARG:HD3	1:D:134:TYR:CG	2.54	0.42
1:D:222:SER:OG	1:D:223:PHE:N	2.51	0.42
2:C:450:TRP:CD1	2:C:506:ARG:HG2	2.55	0.42
1:D:246:ILE:HD13	1:D:350:ALA:O	2.19	0.42
1:A:44:GLU:O	1:A:48:LYS:HG2	2.20	0.42
2:C:491:THR:N	2:C:492:PRO:HD3	2.34	0.42
1:A:171:ILE:HG12	1:A:361:ILE:HG23	2.02	0.42
2:C:689:MET:HE2	2:C:717:LEU:HD23	2.01	0.42
1:D:224:VAL:HG21	2:C:613:LEU:HB2	2.02	0.42
2:C:488:THR:HG22	2:C:524:GLN:NE2	2.35	0.42
2:C:469:GLY:HA2	5:C:812:HOH:O	2.20	0.42
1:A:151:LEU:HB3	1:A:256:MET:HE3	2.01	0.42
2:B:661:ASN:OD1	2:B:664:GLN:N	2.47	0.42
1:D:279:VAL:HA	1:D:327:SER:O	2.20	0.41
2:B:621:ALA:HB1	2:B:704:ARG:HH22	1.85	0.41
1:D:72:SER:O	1:D:84:LYS:N	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:555:ASP:HA	2:C:558:ARG:HG2	2.03	0.41
2:C:589:THR:OG1	2:C:591:LYS:HE2	2.21	0.41
2:B:466:GLY:HA3	3:B:901:ANP:N3B	2.35	0.41
2:C:700:LYS:HD3	2:C:700:LYS:HA	1.93	0.41
1:A:128:GLY:H	1:A:144:GLU:HG2	1.85	0.41
1:D:199:ASN:ND2	1:D:203:GLU:HB2	2.35	0.41
2:C:480:VAL:HA	2:C:531:TRP:H	1.85	0.41
1:D:146:MET:HB3	1:D:198:VAL:O	2.20	0.41
1:A:190:ASP:OD1	1:A:192:LYS:HE3	2.21	0.41
2:B:661:ASN:HB3	2:B:664:GLN:HB3	2.03	0.41
2:B:692:LEU:HD21	2:B:710:ILE:HG23	2.02	0.41
1:D:146:MET:HG3	1:D:197:LEU:HB3	2.02	0.41
2:C:575:ARG:HD2	2:C:633:TYR:CD2	2.55	0.41
1:A:141:ILE:HG22	1:A:143:MET:HE2	2.02	0.41
2:B:688:ARG:HD3	2:B:717:LEU:HD13	2.02	0.41
2:C:540:HIS:HA	2:C:544:SER:OG	2.21	0.41
2:B:564:MET:HE3	2:B:592:ILE:HD13	2.02	0.41
2:C:456:GLN:HE22	2:B:510:HIS:HA	1.86	0.40
2:C:684:ASN:OD1	2:C:684:ASN:N	2.45	0.40
2:B:491:THR:C	2:B:493:GLN:H	2.29	0.40
1:D:118:LEU:HG	1:D:211:VAL:HG21	2.04	0.40
1:D:321:PRO:HA	1:D:322:PRO:HD3	1.91	0.40
2:C:509:ARG:HD3	2:C:509:ARG:HA	1.85	0.40
2:C:558:ARG:HE	2:C:715:GLU:HG2	1.85	0.40
1:A:44:GLU:HG2	1:A:47:ARG:HH21	1.86	0.40
2:B:707:PHE:HA	2:B:710:ILE:HB	2.02	0.40
1:A:191:VAL:HB	1:A:248:SER:HB2	2.03	0.40
1:A:321:PRO:HA	1:A:322:PRO:HD3	1.94	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	306/358 (86%)	296 (97%)	10 (3%)	0	100	100
1	D	311/358 (87%)	302 (97%)	9 (3%)	0	100	100
2	B	262/282 (93%)	250 (95%)	11 (4%)	1 (0%)	30	51
2	C	264/282 (94%)	249 (94%)	15 (6%)	0	100	100
All	All	1143/1280 (89%)	1097 (96%)	45 (4%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	487	VAL

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	269/309 (87%)	262 (97%)	7 (3%)	40	68
1	D	274/309 (89%)	263 (96%)	11 (4%)	28	55
2	B	234/247 (95%)	229 (98%)	5 (2%)	47	73
2	C	236/247 (96%)	233 (99%)	3 (1%)	61	82
All	All	1013/1112 (91%)	987 (97%)	26 (3%)	40	68

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	60	VAL
1	D	65	ASP
1	D	103	ILE
1	D	117	VAL
1	D	122	ASN
1	D	160	ARG
1	D	216	ILE
1	D	239	HIS
1	D	242	VAL
1	D	279	VAL

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Mol	Chain	Res	Type
1	D	320	GLU
2	C	465	SER
2	C	475	LYS
2	C	494	GLN
1	A	60	VAL
1	A	63	LEU
1	A	138	GLU
1	A	227	ARG
1	A	230	MET
1	A	313	LEU
1	A	320	GLU
2	B	480	VAL
2	B	509	ARG
2	B	513	ILE
2	B	687	LYS
2	B	713	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	100	HIS
1	D	319	ASN
2	C	456	GLN
2	C	494	GLN
2	C	524	GLN
2	C	530	GLN
2	C	539	HIS
1	A	56	GLN
1	A	87	HIS
1	A	319	ASN
2	B	542	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	901	4	32,33,33	0.97	4 (12%)	44,52,52	0.91	2 (4%)
3	ANP	D	901	4	32,33,33	0.94	4 (12%)	44,52,52	0.88	1 (2%)
3	ANP	B	901	4	32,33,33	0.94	4 (12%)	44,52,52	0.83	1 (2%)

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	901	4	-	6/18/38/38	0/3/3/3
3	ANP	D	901	4	-	3/18/38/38	0/3/3/3
3	ANP	B	901	4	-	7/18/38/38	0/3/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	ANP	PB-O3A	-2.76	1.55	1.59
3	D	901	ANP	PG-O1G	2.53	1.50	1.46
3	B	901	ANP	PG-N3B	2.50	1.69	1.63
3	B	901	ANP	PG-O1G	2.48	1.50	1.46
3	D	901	ANP	PB-O3A	-2.47	1.56	1.59
3	A	901	ANP	PG-N3B	2.41	1.69	1.63
3	A	901	ANP	PG-O1G	2.37	1.49	1.46
3	B	901	ANP	PB-O1B	2.36	1.49	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	901	ANP	PG-N3B	2.35	1.69	1.63
3	A	901	ANP	PB-O1B	2.31	1.49	1.46
3	D	901	ANP	PB-O1B	2.26	1.49	1.46
3	B	901	ANP	PB-O3A	-2.19	1.56	1.59

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	901	ANP	PB-O3A-PA	-4.09	118.20	132.62
3	A	901	ANP	PB-O3A-PA	-4.05	118.34	132.62
3	B	901	ANP	PB-O3A-PA	-3.33	120.90	132.62
3	A	901	ANP	O1G-PG-N3B	-2.14	108.62	111.77

There are no chirality outliers.

All (16) torsion outliers are listed below:

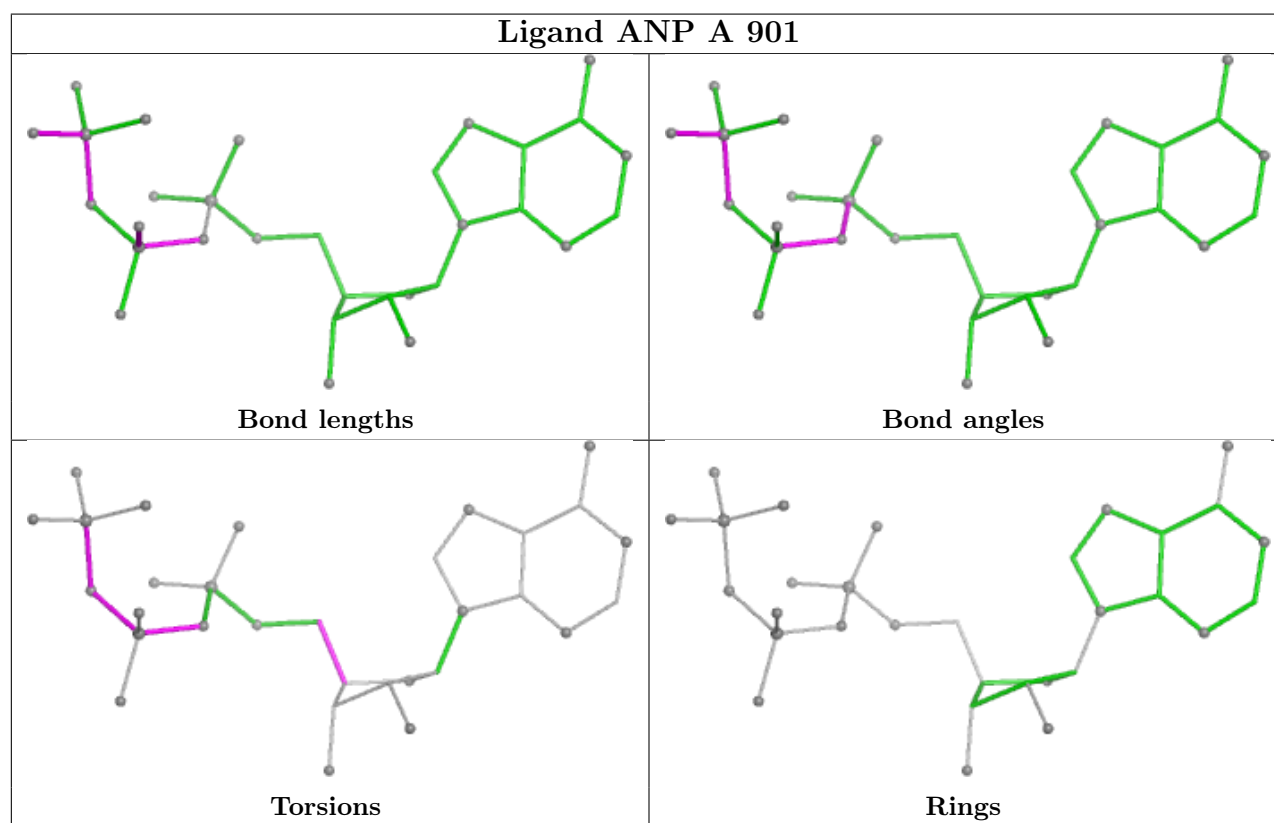
Mol	Chain	Res	Type	Atoms
3	D	901	ANP	PB-N3B-PG-O1G
3	D	901	ANP	PG-N3B-PB-O3A
3	A	901	ANP	PB-N3B-PG-O1G
3	A	901	ANP	PA-O3A-PB-O1B
3	A	901	ANP	PA-O3A-PB-O2B
3	B	901	ANP	PG-N3B-PB-O3A
3	B	901	ANP	C5'-O5'-PA-O1A
3	B	901	ANP	C5'-O5'-PA-O2A
3	B	901	ANP	O4'-C4'-C5'-O5'
3	B	901	ANP	C3'-C4'-C5'-O5'
3	A	901	ANP	O4'-C4'-C5'-O5'
3	A	901	ANP	C3'-C4'-C5'-O5'
3	A	901	ANP	PG-N3B-PB-O3A
3	B	901	ANP	PG-N3B-PB-O1B
3	B	901	ANP	C5'-O5'-PA-O3A
3	D	901	ANP	C5'-O5'-PA-O1A

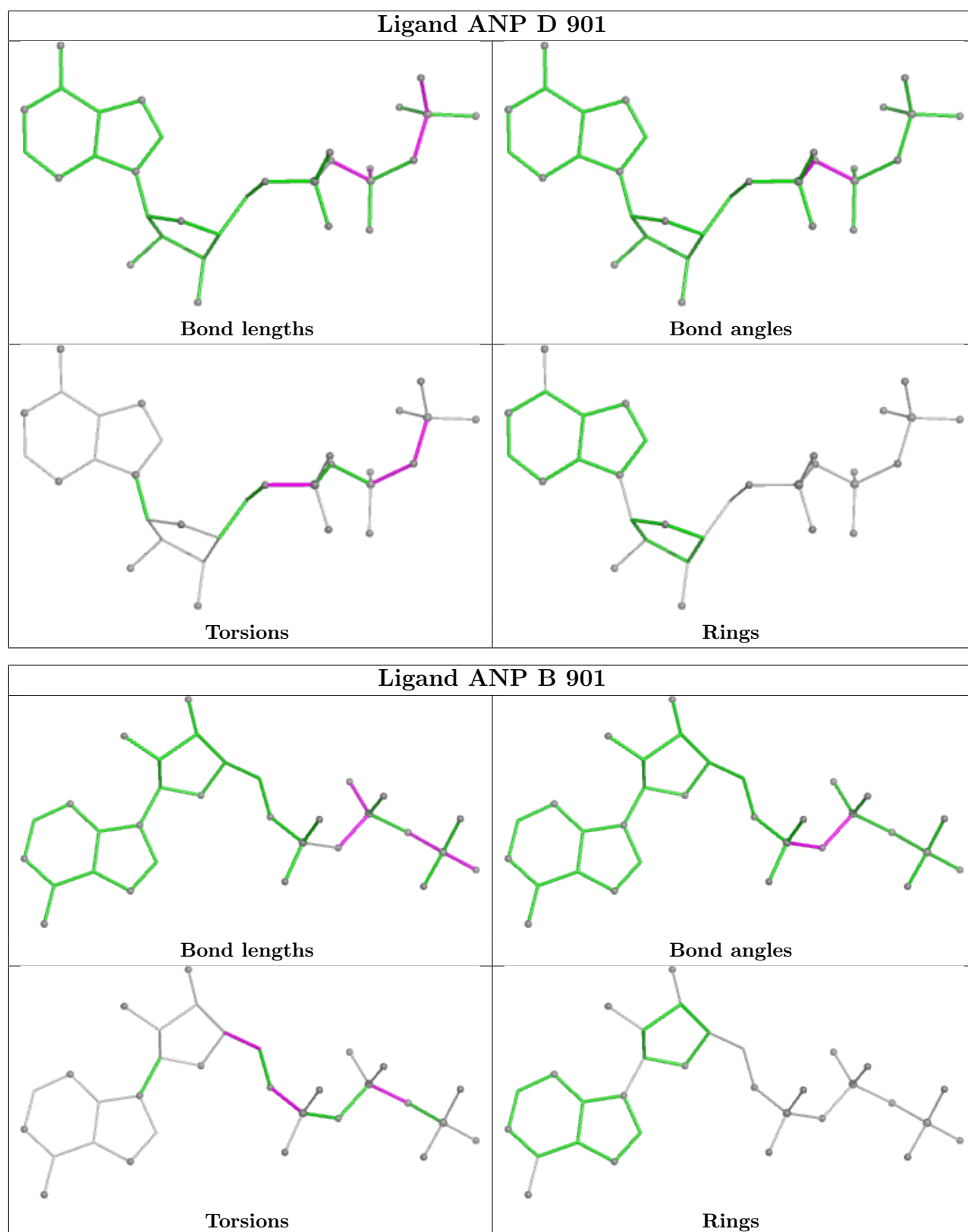
There are no ring outliers.

3 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	ANP	4	0
3	D	901	ANP	5	0
3	B	901	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

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	310/358 (86%)	0.14	10 (3%) 50 44	20, 40, 83, 119	0
1	D	315/358 (87%)	0.53	15 (4%) 35 30	27, 65, 100, 137	0
2	B	266/282 (94%)	0.21	10 (3%) 44 38	17, 38, 75, 110	0
2	C	268/282 (95%)	0.25	11 (4%) 41 36	21, 41, 94, 162	0
All	All	1159/1280 (90%)	0.29	46 (3%) 42 37	17, 46, 94, 162	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	491	THR	6.3
2	C	490	PRO	5.6
1	D	93	VAL	3.9
2	B	717	LEU	3.8
2	B	602	SER	3.7
2	C	489	ALA	3.6
1	A	224	VAL	3.4
2	B	721	LEU	3.2
2	B	718	ALA	3.2
2	C	495	LEU	3.1
2	B	522	LYS	3.0
1	D	186	ILE	2.9
2	B	600	GLU	2.7
1	D	130	TYR	2.7
1	D	174	ILE	2.6
1	D	371	PHE	2.6
1	D	196	ILE	2.6
1	A	276	GLY	2.6
2	C	493	GLN	2.5
1	A	182	GLU	2.5
1	D	142	CYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	63	LEU	2.5
1	D	222	SER	2.5
1	D	355	LEU	2.4
2	C	721	LEU	2.4
1	A	225	GLY	2.4
2	B	523	PRO	2.4
2	C	487	VAL	2.4
1	D	55	THR	2.4
1	D	239	HIS	2.2
1	A	134	TYR	2.2
1	D	307	PRO	2.2
1	A	139	ILE	2.2
1	A	221	ASN	2.2
1	A	307	PRO	2.2
2	C	496	GLN	2.2
1	A	186	ILE	2.2
2	B	461	GLN	2.1
1	D	54	LEU	2.1
2	C	692	LEU	2.1
2	C	467	SER	2.1
1	A	103	ILE	2.1
2	C	610	PHE	2.1
2	B	642	PHE	2.1
2	B	631	ASN	2.0
1	D	118	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

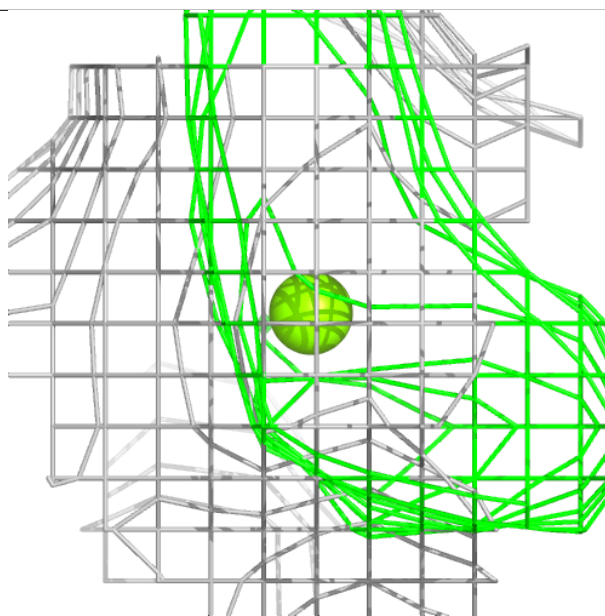
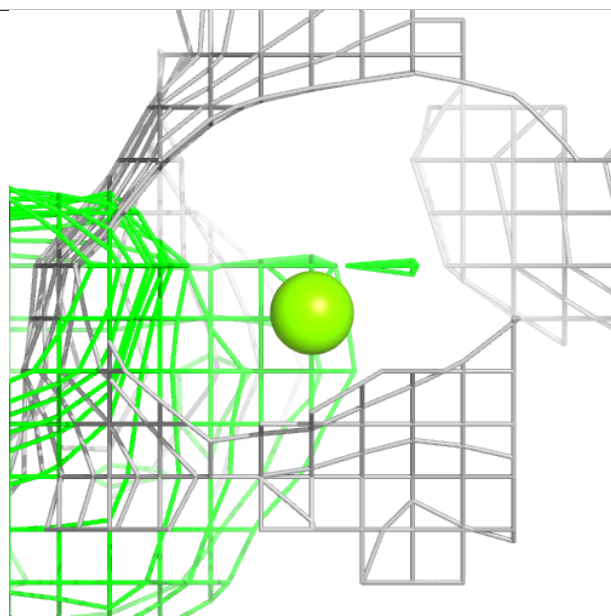
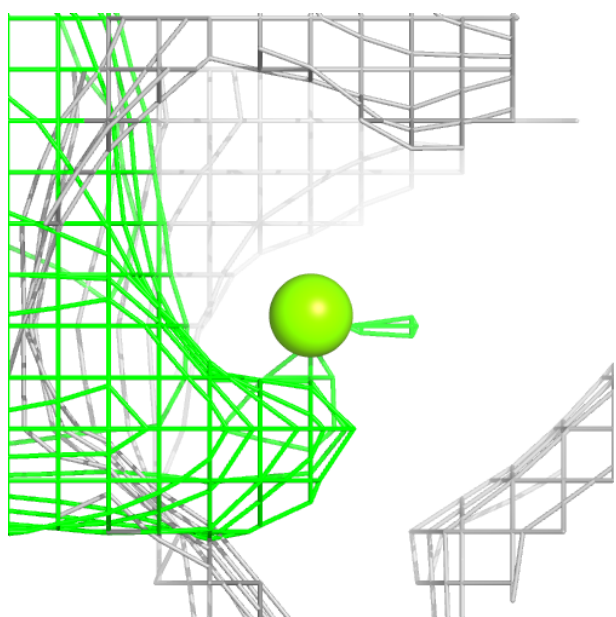
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	902	1/1	0.91	0.15	27,27,27,27	0
3	ANP	B	901	31/31	0.92	0.08	31,44,57,61	0
3	ANP	D	901	31/31	0.94	0.08	32,48,60,71	0
4	MG	B	903	1/1	0.94	0.12	31,31,31,31	0
3	ANP	A	901	31/31	0.95	0.08	24,42,56,59	0
4	MG	B	902	1/1	0.96	0.08	22,22,22,22	0
4	MG	D	902	1/1	0.96	0.07	48,48,48,48	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.

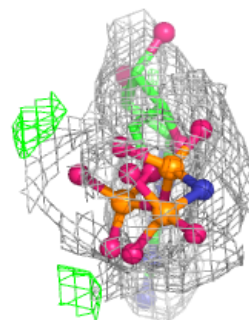
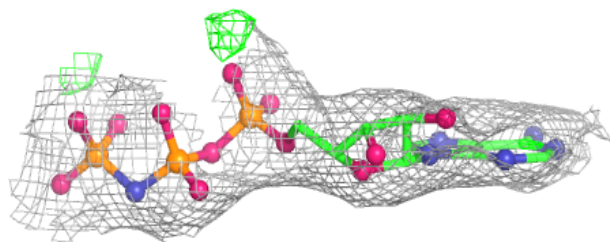
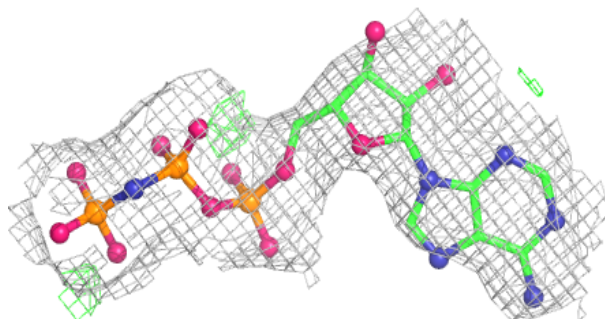
Electron density around MG A 902:

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

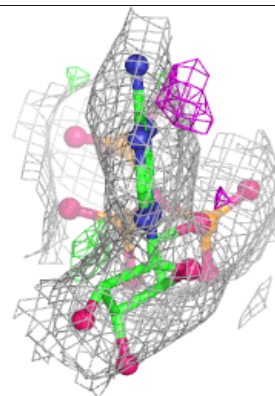
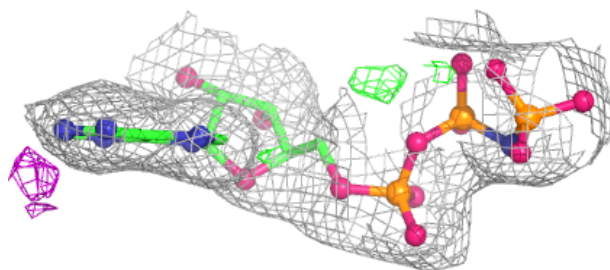
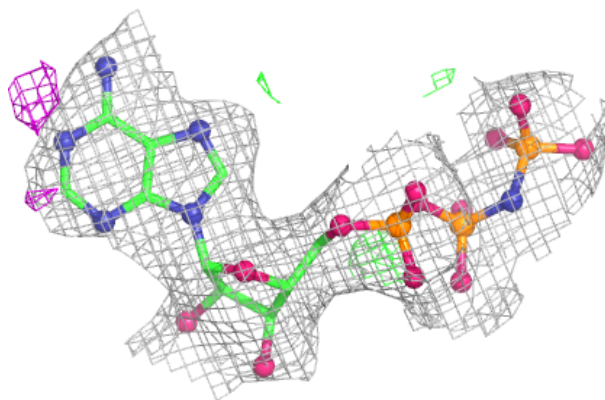


Electron density around ANP B 901:

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

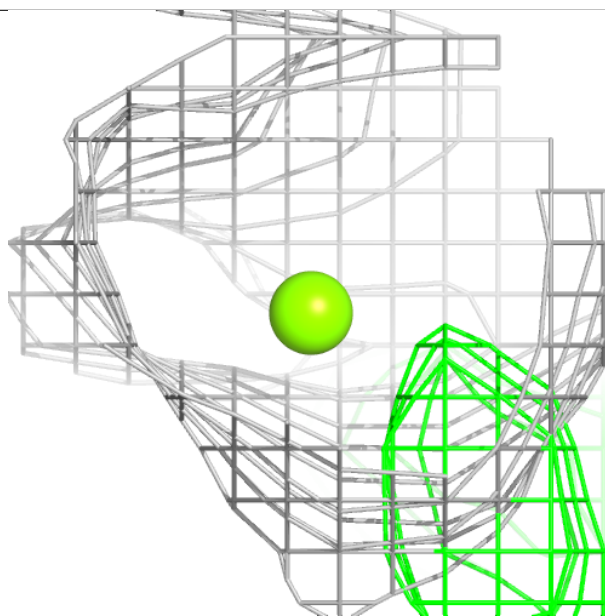
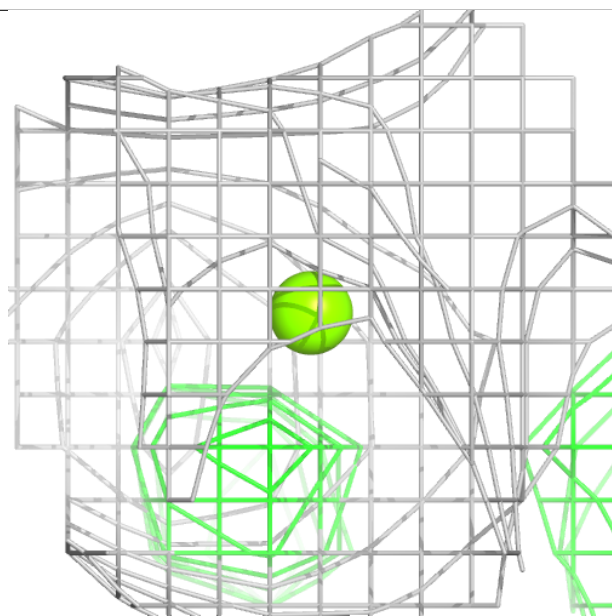
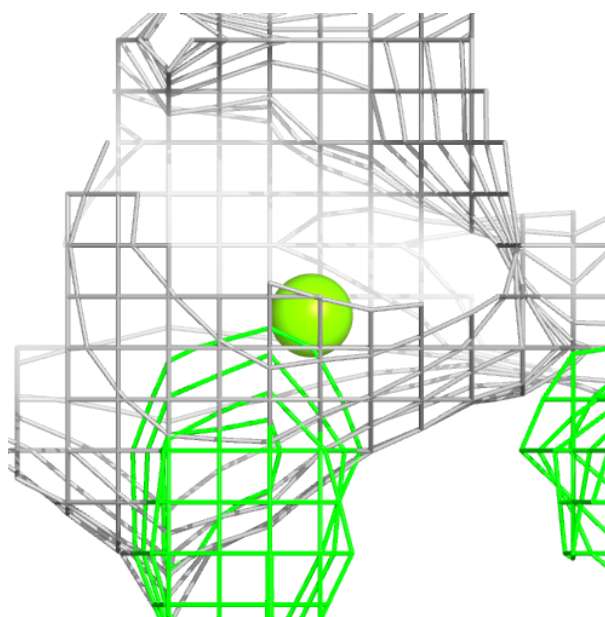
**Electron density around ANP D 901:**

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



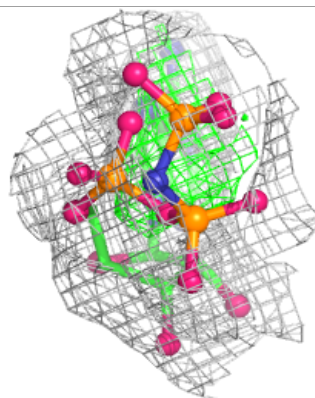
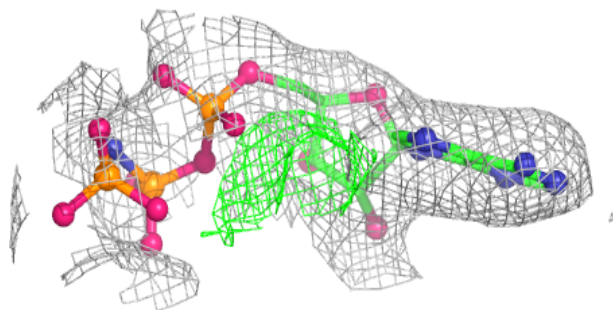
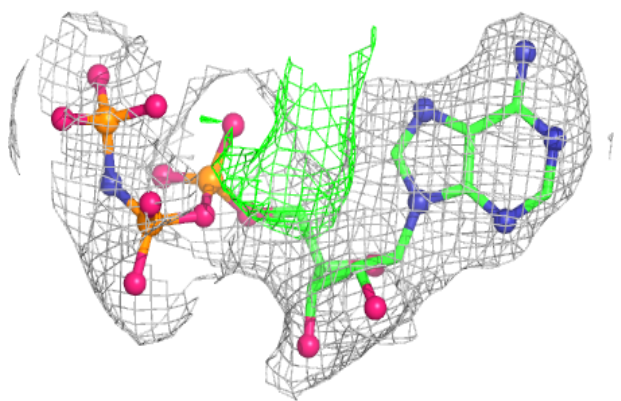
Electron density around MG B 903:

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



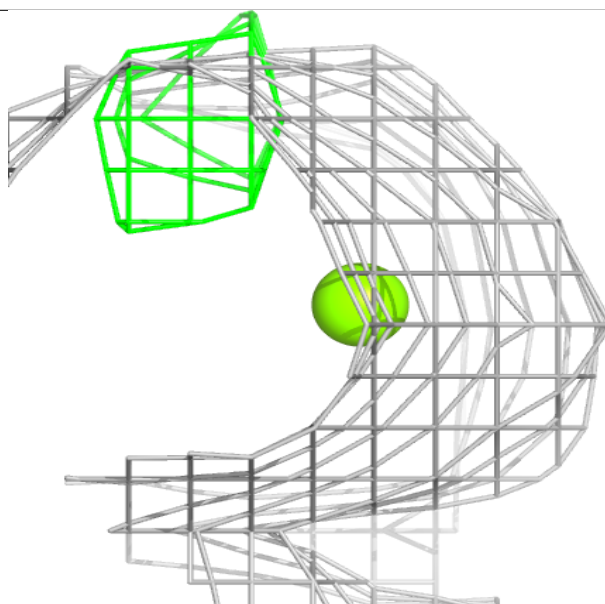
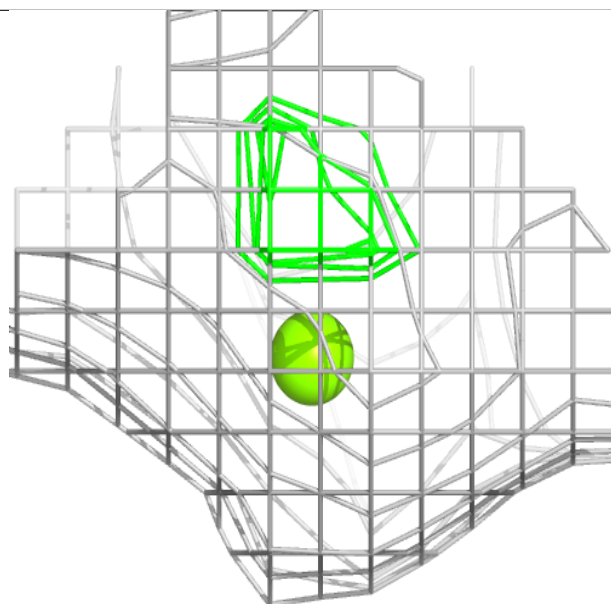
Electron density around ANP A 901:

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



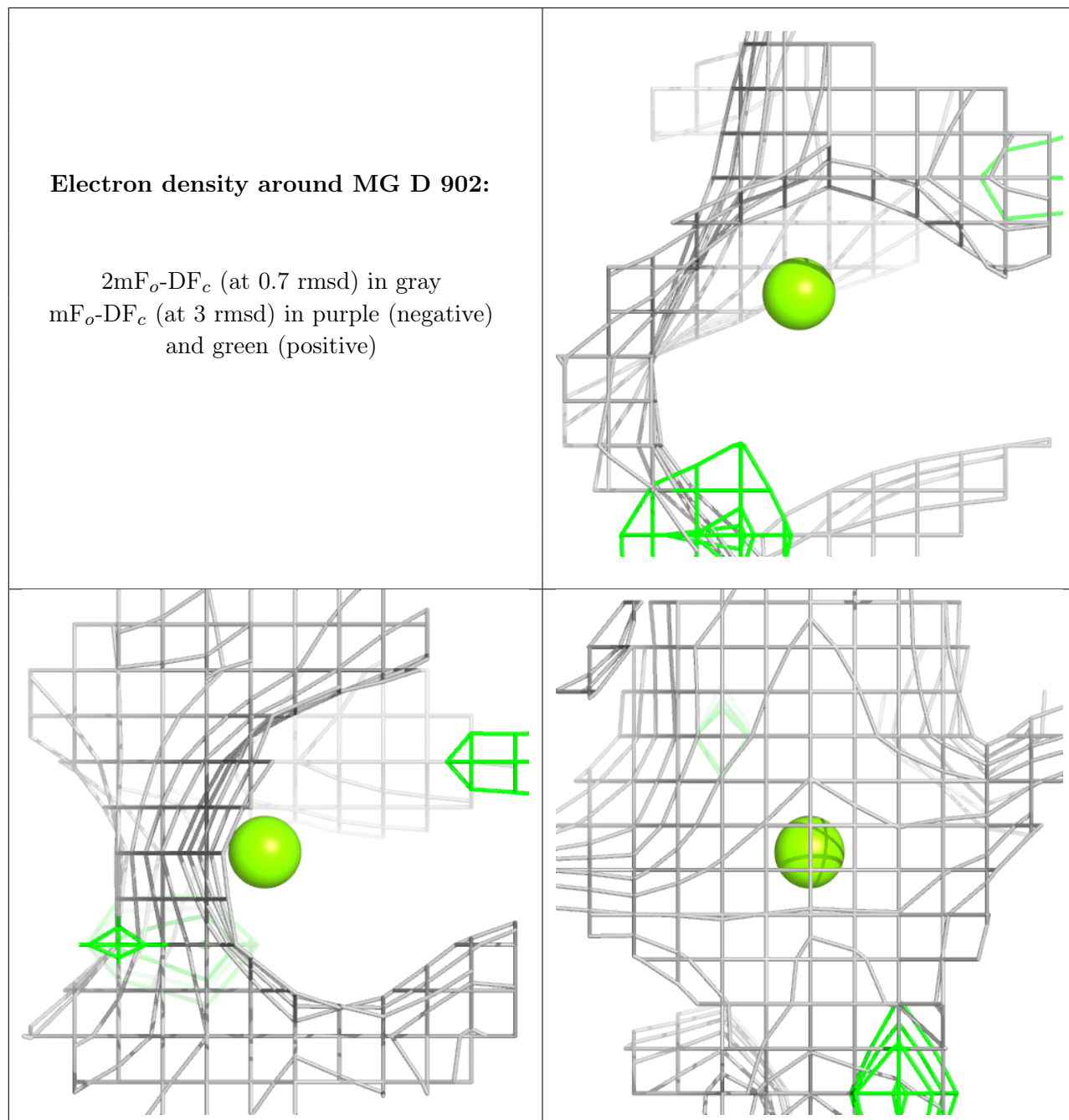
Electron density around MG B 902:

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



Electron density around MG D 902:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.