



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 08:44 PM JST

PDB ID : 6A88
Title : Crystal Structure of *T. gondii* prolyl tRNA synthetase with Febrifugine and ATP Analog
Authors : Kumari, S.; Mishra, S.; Yogavel, M.; Sharma, A.
Deposited on : 2018-07-06
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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

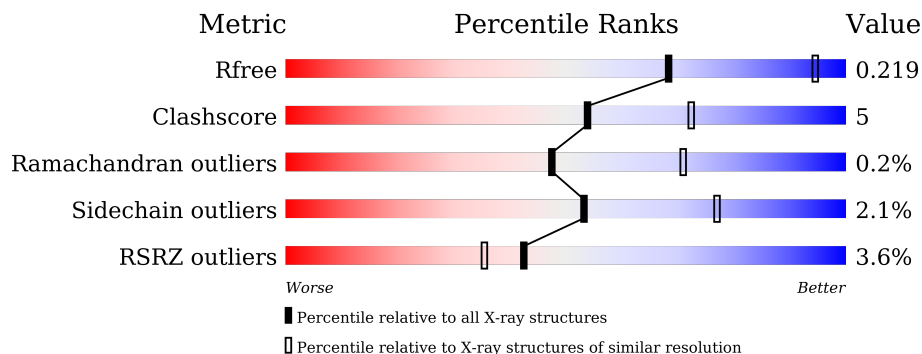
1 Overall quality at a glance

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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (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	500	 2% 82% 14% ••
1	B	500	 4% 82% 14% •
1	C	500	 6% 79% 16% • 5%
1	D	500	 2% 82% 13% • 5%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 16177 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.

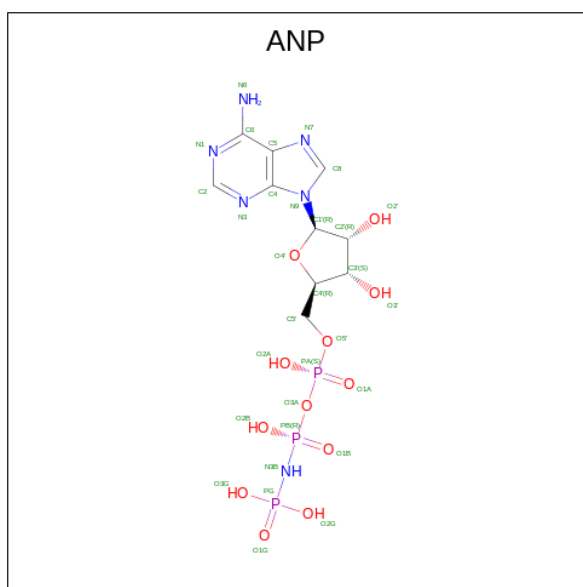
- Molecule 1 is a protein called Prolyl-tRNA synthetase (ProRS).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	482	3959	2543	678	716	22	0	2	0
1	B	481	3938	2529	673	714	22	0	0	0
1	C	477	3915	2515	671	707	22	0	1	0
1	D	475	3893	2503	664	704	22	0	1	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	331	GLY	-	expression tag	UNP S8G8I1
A	332	ALA	-	expression tag	UNP S8G8I1
A	333	MET	-	expression tag	UNP S8G8I1
B	331	GLY	-	expression tag	UNP S8G8I1
B	332	ALA	-	expression tag	UNP S8G8I1
B	333	MET	-	expression tag	UNP S8G8I1
C	331	GLY	-	expression tag	UNP S8G8I1
C	332	ALA	-	expression tag	UNP S8G8I1
C	333	MET	-	expression tag	UNP S8G8I1
D	331	GLY	-	expression tag	UNP S8G8I1
D	332	ALA	-	expression tag	UNP S8G8I1
D	333	MET	-	expression tag	UNP S8G8I1

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

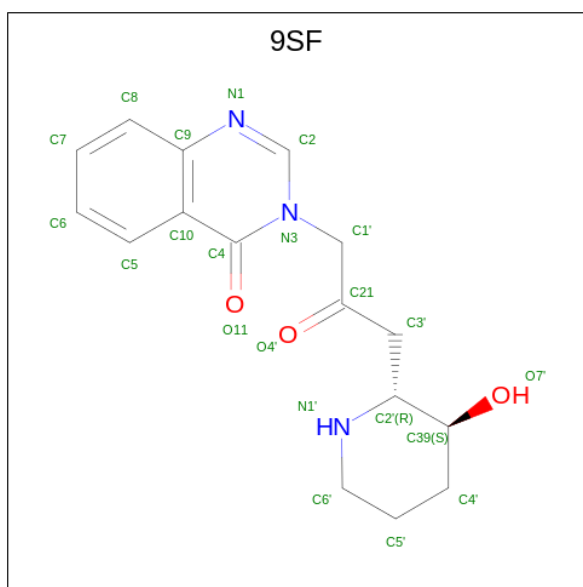


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	31	10	6	12	3	0	0
2	B	1	31	10	6	12	3	0	0
2	C	1	31	10	6	12	3	0	0
2	D	1	31	10	6	12	3	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is 3-{3-[(2R,3S)-3-hydroxypiperidin-2-yl]-2-oxopropyl}quinazolin-4(3H)-one (three-letter code: 9SF) (formula: C₁₆H₁₉N₃O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			22	16	3	3		
4	B	1	Total	C	N	O	0	0
			22	16	3	3		
4	C	1	Total	C	N	O	0	0
			22	16	3	3		
4	D	1	Total	C	N	O	0	0
			22	16	3	3		

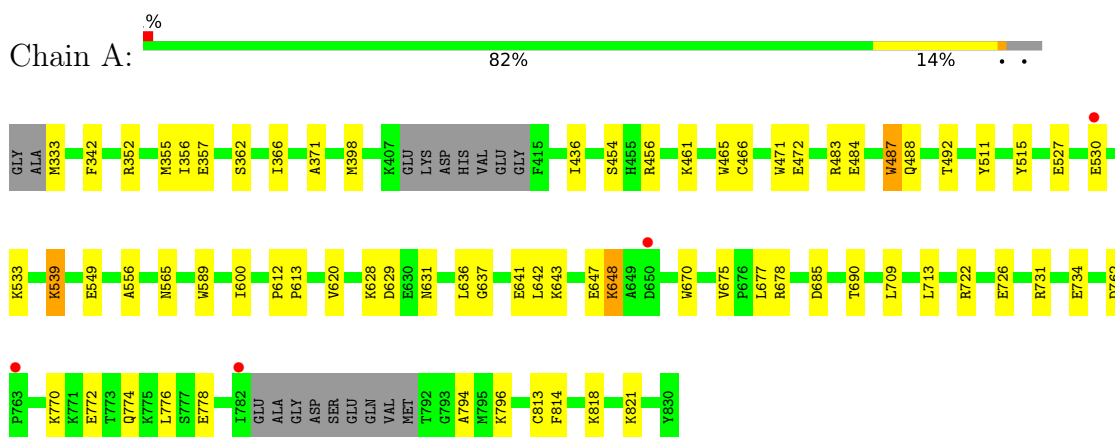
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	104	Total	O	0	0
			104	104		
5	B	53	Total	O	0	0
			53	53		
5	C	36	Total	O	0	0
			36	36		
5	D	63	Total	O	0	0
			63	63		

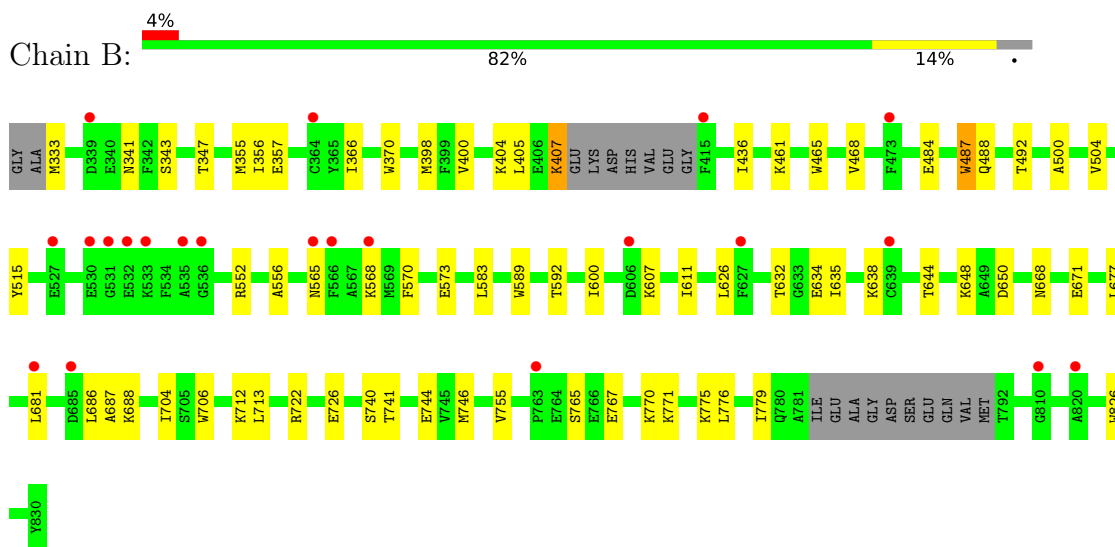
3 Residue-property plots [i](#)

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.

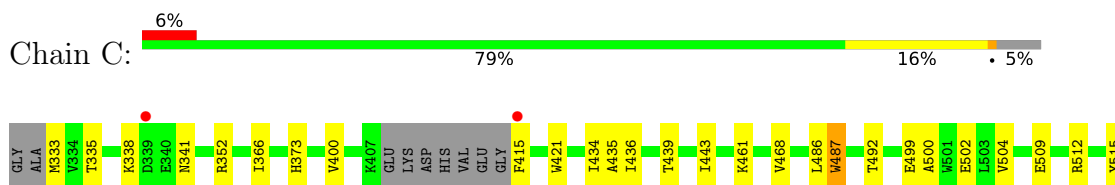
- Molecule 1: Prolyl-tRNA synthetase (ProRS)

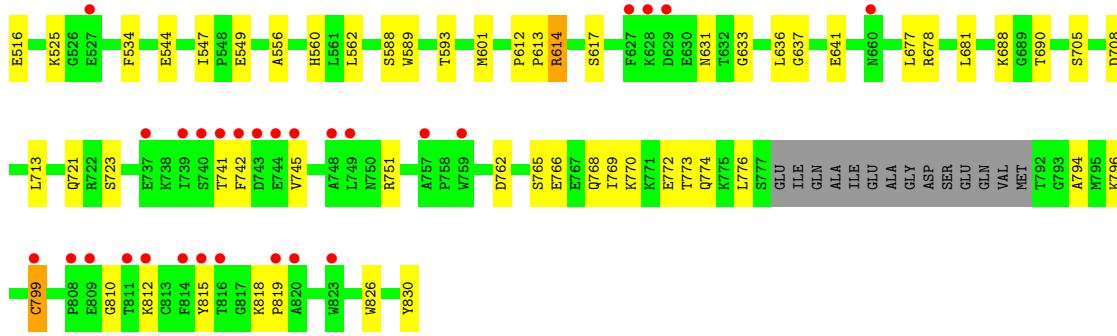


- Molecule 1: Prolyl-tRNA synthetase (ProRS)

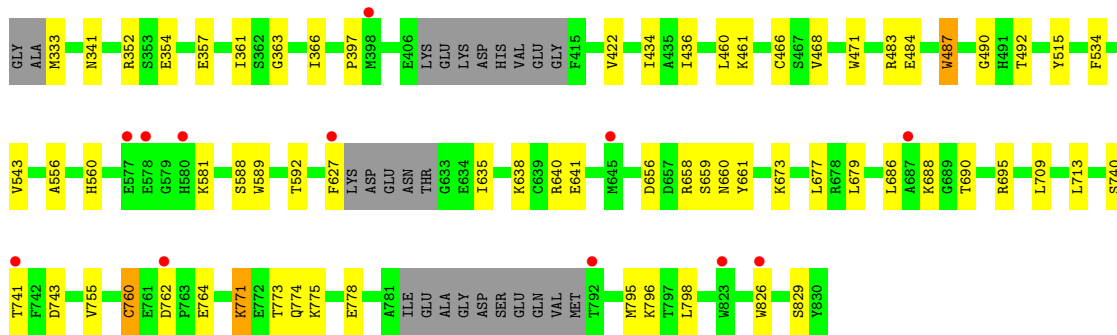
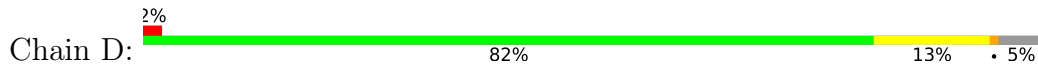


- Molecule 1: Prolyl-tRNA synthetase (ProRS)





• Molecule 1: Prolyl-tRNA synthetase (ProRS)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	76.91Å 90.88Å 93.03Å 89.91° 80.30° 75.58°	Depositor
Resolution (Å)	39.94 – 2.60 42.29 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.3 (39.94-2.60) 98.3 (42.29-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.61Å)	Xtrriage
Refinement program	PHENIX 1.15rc2_3428	Depositor
R, R_{free}	0.172 , 0.220 0.172 , 0.219	Depositor DCC
R_{free} test set	3662 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	41.0	Xtrriage
Anisotropy	0.414	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16177	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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, 9SF

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	0.28	0/4071	0.47	0/5503
1	B	0.27	0/4044	0.46	0/5467
1	C	0.28	0/4024	0.47	0/5439
1	D	0.28	0/4001	0.48	0/5409
All	All	0.28	0/16140	0.47	0/21818

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	3959	0	3931	43	0
1	B	3938	0	3901	41	0
1	C	3915	0	3884	46	0
1	D	3893	0	3855	41	0
2	A	31	0	13	1	0
2	B	31	0	13	0	0
2	C	31	0	13	0	0
2	D	31	0	13	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	22	0	0	0	0
4	B	22	0	0	0	0
4	C	22	0	0	0	0
4	D	22	0	0	0	0
5	A	104	0	0	1	0
5	B	53	0	0	0	0
5	C	36	0	0	2	0
5	D	63	0	0	0	0
All	All	16177	0	15623	166	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 5.

All (166) 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:515:TYR:CZ	1:C:556:ALA:HB1	2.26	0.71
1:B:565:ASN:HA	1:B:568:LYS:HD2	1.73	0.71
1:B:515:TYR:CZ	1:B:556:ALA:HB1	2.27	0.69
1:C:677:LEU:HD13	1:C:713:LEU:HD22	1.74	0.68
1:D:677:LEU:HD13	1:D:713:LEU:HD22	1.75	0.68
1:D:688:LYS:HE2	1:D:690:THR:HG21	1.75	0.68
1:B:333:MET:HG3	1:B:668:ASN:ND2	2.08	0.68
1:D:515:TYR:CZ	1:D:556:ALA:HB1	2.28	0.67
1:A:677:LEU:HD13	1:A:713:LEU:HD22	1.77	0.66
1:A:515:TYR:CZ	1:A:556:ALA:HB1	2.30	0.66
1:B:767:GLU:HA	1:B:770:LYS:HG2	1.77	0.65
1:D:773:THR:HG21	1:D:795:MET:HB3	1.78	0.65
1:C:773:THR:HA	1:C:776:LEU:HD23	1.80	0.63
1:C:774:GLN:HB2	1:C:794:ALA:HB2	1.81	0.61
1:A:731:ARG:HA	1:A:734:GLU:HG2	1.83	0.60
1:B:767:GLU:O	1:B:771:LYS:HG3	2.01	0.60
1:C:688:LYS:HE2	1:C:690:THR:OG1	2.01	0.60
1:C:468:VAL:HG21	1:C:487:TRP:CE2	2.38	0.58
1:C:499:GLU:O	1:C:502:GLU:HG3	2.04	0.58
1:C:500:ALA:O	1:C:504:VAL:HG23	2.04	0.58
1:B:357:GLU:HB3	1:B:366:ILE:HB	1.85	0.58
1:A:643:LYS:O	1:A:647:GLU:HG3	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:772:GLU:O	1:A:776:LEU:HG	2.06	0.56
1:D:762:ASP:HB3	1:D:764:GLU:OE2	2.06	0.56
1:B:704:ILE:HD13	1:B:712:LYS:HG2	1.88	0.56
1:C:768:GLN:NE2	1:C:815:TYR:HE2	2.04	0.55
1:B:634:GLU:HG2	1:B:686:LEU:HD21	1.88	0.55
1:D:638:LYS:HA	1:D:641:GLU:HG3	1.88	0.55
1:C:637:GLY:O	1:C:641:GLU:HG3	2.07	0.55
1:B:638:LYS:HE2	1:B:706:TRP:CE3	2.42	0.55
1:C:818:LYS:HG3	1:C:819:PRO:HD2	1.89	0.54
1:D:490:GLY:O	1:D:588:SER:HA	2.07	0.54
1:A:355:MET:HE1	1:A:600:ILE:HD11	1.88	0.54
1:C:400:VAL:HG22	1:C:443:ILE:HD11	1.89	0.54
1:B:400:VAL:HG11	1:B:405:LEU:HD13	1.90	0.54
1:A:355:MET:CE	1:A:600:ILE:HD11	2.36	0.54
1:A:436:ILE:HD11	1:B:436:ILE:HD11	1.90	0.53
1:C:769:ILE:HD13	1:C:826:TRP:CH2	2.44	0.53
1:B:407:LYS:HE2	1:B:570:PHE:CZ	2.44	0.53
1:B:573:GLU:HG2	1:B:583:LEU:HD23	1.91	0.53
1:D:333:MET:O	1:D:352:ARG:NH1	2.42	0.52
1:B:775:LYS:O	1:B:779:ILE:HG12	2.10	0.52
1:B:746:MET:SD	1:B:776:LEU:HB2	2.49	0.52
1:B:333:MET:HG3	1:B:668:ASN:HD21	1.73	0.52
1:A:530:GLU:HG2	1:A:533:LYS:HE3	1.90	0.52
1:C:631:ASN:ND2	1:C:633:GLY:H	2.08	0.52
1:D:741:THR:OG1	1:D:743:ASP:OD1	2.24	0.52
1:D:679:LEU:HD11	1:D:709:LEU:HD21	1.92	0.52
1:D:466:CYS:HG	1:D:487:TRP:HE1	1.57	0.51
1:A:488:GLN:NE2	1:A:511:TYR:OH	2.40	0.51
1:D:461:LYS:HA	1:D:492:THR:HG22	1.92	0.51
1:D:635:ILE:HG13	1:D:686:LEU:HD12	1.92	0.51
1:D:755:VAL:HG23	1:D:826:TRP:HB2	1.94	0.50
1:C:544:GLU:HG3	1:C:799:CYS:HB3	1.94	0.50
1:A:465:TRP:CE3	1:A:488:GLN:HB3	2.47	0.49
1:B:638:LYS:HE2	1:B:706:TRP:CD2	2.48	0.49
1:A:642:LEU:HD22	1:A:709:LEU:HD12	1.94	0.49
1:A:774:GLN:O	1:A:778:GLU:HG3	2.13	0.49
1:B:468:VAL:HG21	1:B:487:TRP:CE2	2.48	0.49
1:B:607:LYS:HE2	1:B:671:GLU:OE2	2.12	0.49
1:A:731:ARG:O	1:A:734:GLU:HG2	2.13	0.48
1:B:635:ILE:HG23	1:B:681:LEU:HD22	1.94	0.48
1:D:771:LYS:HA	1:D:774:GLN:HG2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:404:LYS:O	1:B:407:LYS:NZ	2.45	0.48
1:C:762:ASP:O	1:C:765:SER:OG	2.24	0.48
1:C:338:LYS:HG2	1:C:549:GLU:HG2	1.95	0.48
1:D:460:LEU:O	1:D:492:THR:HA	2.14	0.48
1:B:687:ALA:HB3	1:B:688:LYS:NZ	2.29	0.48
1:C:436:ILE:HD11	1:D:436:ILE:HD11	1.96	0.48
1:C:772:GLU:HG3	1:C:773:THR:N	2.28	0.48
1:B:487:TRP:HB3	1:B:592:THR:HG22	1.95	0.48
1:D:760:CYS:SG	1:D:762:ASP:HB2	2.54	0.48
1:B:626:LEU:HD22	1:B:632:THR:HA	1.96	0.47
1:A:637:GLY:O	1:A:641:GLU:HG3	2.14	0.47
1:C:810:GLY:O	1:C:812:LYS:HE3	2.14	0.47
1:D:755:VAL:CG2	1:D:826:TRP:HB2	2.45	0.47
1:A:631:ASN:ND2	1:A:636:LEU:HD12	2.30	0.47
1:C:614:ARG:HH21	1:C:721:GLN:HG2	1.80	0.47
1:C:512:ARG:HD3	1:C:525:LYS:HE3	1.95	0.47
1:A:472:GLU:H	1:A:483:ARG:HD3	1.80	0.47
1:D:543:VAL:HG23	1:D:556:ALA:HB3	1.97	0.47
1:B:677:LEU:HD13	1:B:713:LEU:HD22	1.96	0.47
1:C:742:PHE:O	1:C:745:VAL:HG13	2.14	0.47
1:B:573:GLU:HG2	1:B:583:LEU:CD2	2.45	0.46
1:D:357:GLU:HB3	1:D:366:ILE:HB	1.96	0.46
1:A:631:ASN:HD21	1:A:636:LEU:HD12	1.81	0.46
1:A:770:LYS:HD2	1:A:796:LYS:HB3	1.96	0.46
1:C:415:PHE:N	5:C:1103:HOH:O	2.48	0.46
1:B:355:MET:HE1	1:B:600:ILE:CD1	2.45	0.46
1:B:465:TRP:CE3	1:B:488:GLN:HB3	2.50	0.46
1:A:620:VAL:HG21	1:A:713:LEU:HD13	1.97	0.46
1:C:366:ILE:HG12	1:D:397:PRO:HD3	1.97	0.46
1:D:688:LYS:HG2	1:D:690:THR:HG23	1.97	0.46
1:A:355:MET:HE3	1:A:371:ALA:HB2	1.98	0.46
1:C:335:THR:HG23	1:C:352:ARG:HH12	1.81	0.46
1:C:486:LEU:HB2	1:C:593:THR:HG23	1.98	0.45
1:C:562:LEU:HD11	1:C:588:SER:HB3	1.99	0.45
1:A:648:LYS:HD3	1:A:648:LYS:HA	1.54	0.45
1:D:709:LEU:HD12	1:D:709:LEU:HA	1.85	0.45
1:D:656:ASP:OD1	1:D:658:ARG:HB2	2.17	0.45
1:D:774:GLN:O	1:D:778:GLU:HG2	2.17	0.44
1:A:484:GLU:CD	1:B:398:MET:HG2	2.37	0.44
1:A:774:GLN:HG2	1:A:794:ALA:HB2	1.98	0.44
1:C:373:HIS:CE1	1:C:617:SER:HA	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:MET:HE1	1:B:600:ILE:HD11	1.99	0.44
1:D:534:PHE:HB2	1:D:560:HIS:CE1	2.52	0.44
1:B:461:LYS:HA	1:B:492:THR:HG22	1.98	0.44
1:C:547:ILE:HD12	1:C:830:TYR:CE2	2.53	0.44
1:C:705:SER:OG	1:C:708:ASP:OD2	2.36	0.44
1:D:774:GLN:HG3	1:D:775:LYS:N	2.32	0.44
1:D:659:SER:O	1:D:661:TYR:N	2.51	0.43
1:A:722:ARG:HG2	1:A:726:GLU:OE2	2.19	0.43
1:D:677:LEU:HD23	1:D:695:ARG:HA	1.99	0.43
1:A:762:ASP:OD2	1:A:818:LYS:NZ	2.50	0.43
1:D:762:ASP:HB3	1:D:764:GLU:CD	2.38	0.43
1:A:527:GLU:CD	1:A:539:LYS:HE3	2.39	0.43
1:A:355:MET:HE3	1:A:355:MET:HB3	1.40	0.43
1:C:512:ARG:NH1	1:C:516:GLU:OE1	2.50	0.43
1:A:454:SER:OG	1:A:456:ARG:HG3	2.19	0.43
1:C:612:PRO:HA	1:C:613:PRO:HD3	1.93	0.43
1:D:363:GLY:HA3	1:D:483:ARG:HG3	2.00	0.43
1:B:487:TRP:CB	1:B:592:THR:HG22	2.48	0.43
1:B:722:ARG:O	1:B:726:GLU:HG2	2.18	0.43
1:A:678:ARG:HD3	5:A:1138:HOH:O	2.19	0.42
1:C:636:LEU:HD12	1:C:636:LEU:HA	1.87	0.42
1:C:678:ARG:HD3	5:C:1120:HOH:O	2.20	0.42
1:A:342:PHE:CD1	1:A:549:GLU:HG2	2.55	0.42
1:A:356:ILE:HA	1:A:366:ILE:O	2.19	0.42
1:C:434:ILE:HD13	1:D:361:ILE:HD13	2.00	0.42
1:C:765:SER:O	1:C:768:GLN:HB2	2.19	0.42
1:C:766:GLU:HG2	1:C:770:LYS:HD3	2.00	0.42
1:A:398:MET:HG2	1:B:484:GLU:CD	2.39	0.42
1:A:466:CYS:HG	1:A:487:TRP:HE1	1.67	0.42
1:C:421:TRP:CZ3	1:C:435:ALA:HB2	2.54	0.42
1:D:581:LYS:HB3	1:D:581:LYS:HE3	1.71	0.42
1:D:771:LYS:O	1:D:774:GLN:HG2	2.19	0.42
1:D:798:LEU:HD21	1:D:829:SER:HB3	2.01	0.42
1:B:741:THR:HG22	1:B:744:GLU:OE1	2.20	0.42
1:C:509:GLU:OE1	1:C:525:LYS:NZ	2.40	0.41
1:C:333:MET:O	1:C:352:ARG:NH1	2.53	0.41
1:A:461:LYS:HA	1:A:492:THR:HG22	2.02	0.41
1:A:670:TRP:CE3	1:A:675:VAL:HG21	2.56	0.41
1:B:644:THR:HG23	1:B:648:LYS:HE3	2.02	0.41
1:C:461:LYS:HA	1:C:492:THR:HG22	2.01	0.41
1:C:547:ILE:HD13	1:C:601:MET:SD	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:422:VAL:HB	1:D:434:ILE:HG12	2.03	0.41
1:D:487:TRP:CB	1:D:592:THR:HG22	2.51	0.41
2:A:1001:ANP:O3G	2:A:1001:ANP:O2B	2.38	0.41
1:C:534:PHE:HB2	1:C:560:HIS:CE1	2.56	0.41
1:A:549:GLU:N	1:A:549:GLU:OE2	2.53	0.41
1:A:515:TYR:CE2	1:A:556:ALA:HB1	2.56	0.41
1:A:612:PRO:HA	1:A:613:PRO:HD3	1.94	0.41
1:A:685:ASP:HB3	1:A:690:THR:O	2.21	0.41
1:B:343:SER:O	1:B:347:THR:HG23	2.21	0.41
1:B:755:VAL:CG2	1:B:826:TRP:HB2	2.50	0.41
1:C:768:GLN:HE21	1:C:815:TYR:HE2	1.67	0.41
1:A:821:LYS:HE2	1:A:821:LYS:HB3	1.80	0.41
1:B:370:TRP:HE1	1:B:611:ILE:HD11	1.86	0.41
1:B:500:ALA:O	1:B:504:VAL:HG23	2.21	0.41
1:A:333:MET:O	1:A:352:ARG:NH1	2.53	0.40
1:D:468:VAL:O	1:D:484:GLU:HA	2.21	0.40
1:A:357:GLU:HB3	1:A:366:ILE:HB	2.03	0.40
1:D:354:GLU:OE1	1:D:673:LYS:NZ	2.54	0.40
1:D:796:LYS:HD3	1:D:829:SER:OG	2.22	0.40
1:C:741:THR:OG1	1:C:742:PHE:N	2.53	0.40
1:B:356:ILE:HD13	1:B:356:ILE:HA	1.90	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	478/500 (96%)	464 (97%)	13 (3%)	1 (0%)	47 71
1	B	475/500 (95%)	460 (97%)	15 (3%)	0	100 100
1	C	472/500 (94%)	455 (96%)	16 (3%)	1 (0%)	47 71
1	D	468/500 (94%)	455 (97%)	12 (3%)	1 (0%)	47 71

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1893/2000 (95%)	1834 (97%)	56 (3%)	3 (0%)	47 71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	629	ASP
1	C	439	THR
1	D	660	ASN

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	425/436 (98%)	415 (98%)	10 (2%)	49 74
1	B	422/436 (97%)	414 (98%)	8 (2%)	57 79
1	C	420/436 (96%)	411 (98%)	9 (2%)	53 77
1	D	417/436 (96%)	408 (98%)	9 (2%)	52 76
All	All	1684/1744 (97%)	1648 (98%)	36 (2%)	53 77

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	362	SER
1	A	471	TRP
1	A	487	TRP
1	A	539	LYS
1	A	565	ASN
1	A	589	TRP
1	A	628	LYS
1	A	648	LYS
1	A	813	CYS
1	A	814	PHE
1	B	341	ASN
1	B	407	LYS

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Mol	Chain	Res	Type
1	B	487	TRP
1	B	552	ARG
1	B	589	TRP
1	B	650	ASP
1	B	740	SER
1	B	765	SER
1	C	341	ASN
1	C	487	TRP
1	C	589	TRP
1	C	614	ARG
1	C	681	LEU
1	C	723	SER
1	C	751	ARG
1	C	796	LYS
1	C	799	CYS
1	D	341	ASN
1	D	471	TRP
1	D	487	TRP
1	D	589	TRP
1	D	627	PHE
1	D	640	ARG
1	D	740	SER
1	D	760	CYS
1	D	771	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	488	GLN
1	A	585	HIS
1	A	631	ASN
1	A	768	GLN
1	B	341	ASN
1	B	488	GLN
1	B	669	HIS
1	B	750	ASN
1	C	488	GLN
1	C	631	ASN
1	C	750	ASN
1	C	768	GLN
1	D	341	ASN
1	D	488	GLN

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Mol	Chain	Res	Type
1	D	750	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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
4	9SF	D	1003	-	23,24,24	0.85	0	25,33,33	1.61	1 (4%)
4	9SF	C	1003	-	23,24,24	0.80	0	25,33,33	1.09	1 (4%)
2	ANP	C	1001	3	29,33,33	1.10	5 (17%)	31,52,52	1.26	4 (12%)
4	9SF	B	1003	-	23,24,24	0.88	0	25,33,33	0.52	0
2	ANP	B	1001	3	29,33,33	1.08	3 (10%)	31,52,52	1.04	2 (6%)
4	9SF	A	1003	-	23,24,24	0.17	0	25,33,33	0.49	0
2	ANP	D	1001	3	29,33,33	1.09	4 (13%)	31,52,52	1.08	3 (9%)
2	ANP	A	1001	3	29,33,33	1.09	4 (13%)	31,52,52	1.12	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
4	9SF	D	1003	-	-	0/8/19/19	0/3/3/3
4	9SF	C	1003	-	-	0/8/19/19	0/3/3/3
2	ANP	C	1001	3	-	6/14/38/38	0/3/3/3
4	9SF	B	1003	-	-	0/8/19/19	0/3/3/3
2	ANP	B	1001	3	-	5/14/38/38	0/3/3/3
4	9SF	A	1003	-	-	1/8/19/19	0/3/3/3
2	ANP	D	1001	3	-	7/14/38/38	0/3/3/3
2	ANP	A	1001	3	-	7/14/38/38	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	ANP	PG-N3B	2.68	1.70	1.63
2	C	1001	ANP	PG-N3B	2.66	1.70	1.63
2	D	1001	ANP	PG-N3B	2.57	1.70	1.63
2	B	1001	ANP	PG-N3B	2.55	1.70	1.63
2	C	1001	ANP	PG-O1G	2.49	1.50	1.46
2	D	1001	ANP	PG-O1G	2.47	1.50	1.46
2	B	1001	ANP	PG-O1G	2.46	1.50	1.46
2	A	1001	ANP	PG-O1G	2.42	1.50	1.46
2	C	1001	ANP	PB-O1B	2.40	1.50	1.46
2	B	1001	ANP	PB-O1B	2.36	1.49	1.46
2	D	1001	ANP	PB-O1B	2.23	1.49	1.46
2	D	1001	ANP	PB-O3A	-2.14	1.56	1.59
2	A	1001	ANP	PB-O1B	2.13	1.49	1.46
2	A	1001	ANP	PB-N3B	2.10	1.68	1.63
2	C	1001	ANP	PB-N3B	2.04	1.68	1.63
2	C	1001	ANP	PB-O3A	-2.02	1.56	1.59

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1003	9SF	C6'-N1'-C2'	-7.56	106.90	111.62
4	C	1003	9SF	C6'-N1'-C2'	-5.13	108.42	111.62
2	C	1001	ANP	PB-O3A-PA	-3.25	121.17	132.62
2	B	1001	ANP	PB-O3A-PA	-2.97	122.15	132.62
2	C	1001	ANP	C3'-C2'-C1'	2.88	105.31	100.98
2	A	1001	ANP	PB-O3A-PA	-2.55	123.62	132.62
2	D	1001	ANP	PB-O3A-PA	-2.49	123.84	132.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1001	ANP	O1G-PG-N3B	-2.30	108.38	111.77
2	C	1001	ANP	C5-C6-N6	2.29	123.83	120.35
2	A	1001	ANP	O1G-PG-N3B	-2.29	108.40	111.77
2	D	1001	ANP	C5-C6-N6	2.21	123.72	120.35
2	B	1001	ANP	C5-C6-N6	2.21	123.72	120.35
2	A	1001	ANP	C5-C6-N6	2.11	123.55	120.35
2	A	1001	ANP	C4-C5-N7	2.07	111.56	109.40
2	C	1001	ANP	O1B-PB-N3B	-2.04	108.76	111.77
2	A	1001	ANP	O1B-PB-N3B	-2.01	108.81	111.77

There are no chirality outliers.

All (26) torsion outliers are listed below:

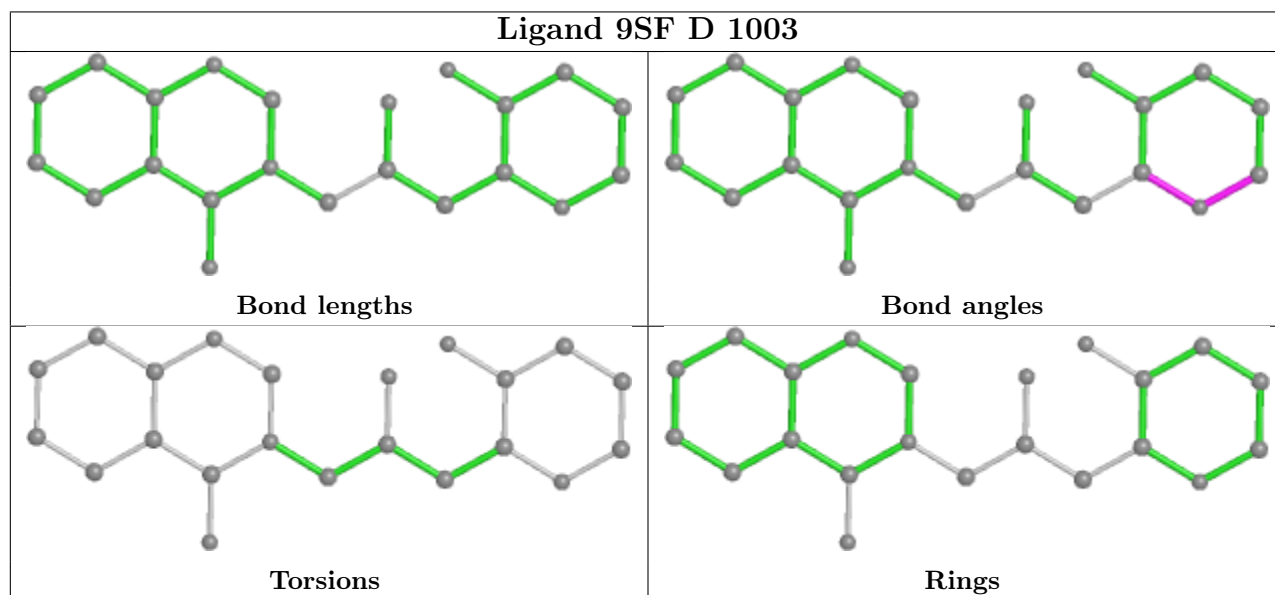
Mol	Chain	Res	Type	Atoms
2	A	1001	ANP	PB-N3B-PG-O1G
2	A	1001	ANP	PG-N3B-PB-O1B
2	B	1001	ANP	PB-N3B-PG-O1G
2	B	1001	ANP	C5'-O5'-PA-O1A
2	C	1001	ANP	PB-N3B-PG-O1G
2	C	1001	ANP	PG-N3B-PB-O1B
2	C	1001	ANP	C5'-O5'-PA-O1A
2	D	1001	ANP	PB-N3B-PG-O1G
2	D	1001	ANP	PG-N3B-PB-O1B
2	D	1001	ANP	PA-O3A-PB-O1B
2	D	1001	ANP	C5'-O5'-PA-O1A
2	A	1001	ANP	C5'-O5'-PA-O3A
2	B	1001	ANP	C5'-O5'-PA-O3A
2	C	1001	ANP	C5'-O5'-PA-O3A
2	D	1001	ANP	C5'-O5'-PA-O3A
2	A	1001	ANP	C5'-O5'-PA-O1A
2	A	1001	ANP	C5'-O5'-PA-O2A
2	B	1001	ANP	C5'-O5'-PA-O2A
2	C	1001	ANP	C5'-O5'-PA-O2A
2	D	1001	ANP	C5'-O5'-PA-O2A
4	A	1003	9SF	N3-C1'-C21-O4'
2	D	1001	ANP	C4'-C5'-O5'-PA
2	A	1001	ANP	C4'-C5'-O5'-PA
2	B	1001	ANP	C4'-C5'-O5'-PA
2	A	1001	ANP	O4'-C4'-C5'-O5'
2	C	1001	ANP	O4'-C4'-C5'-O5'

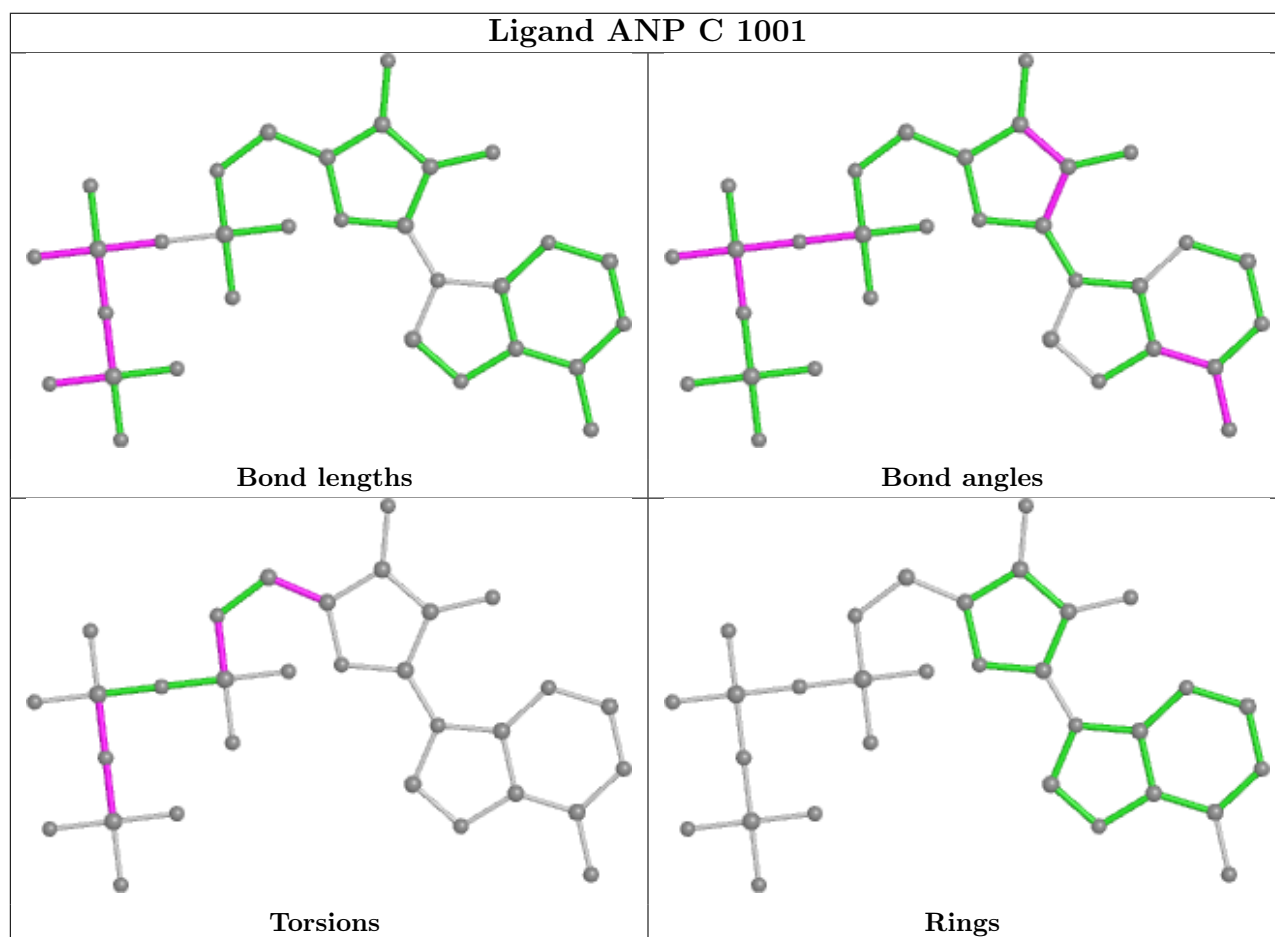
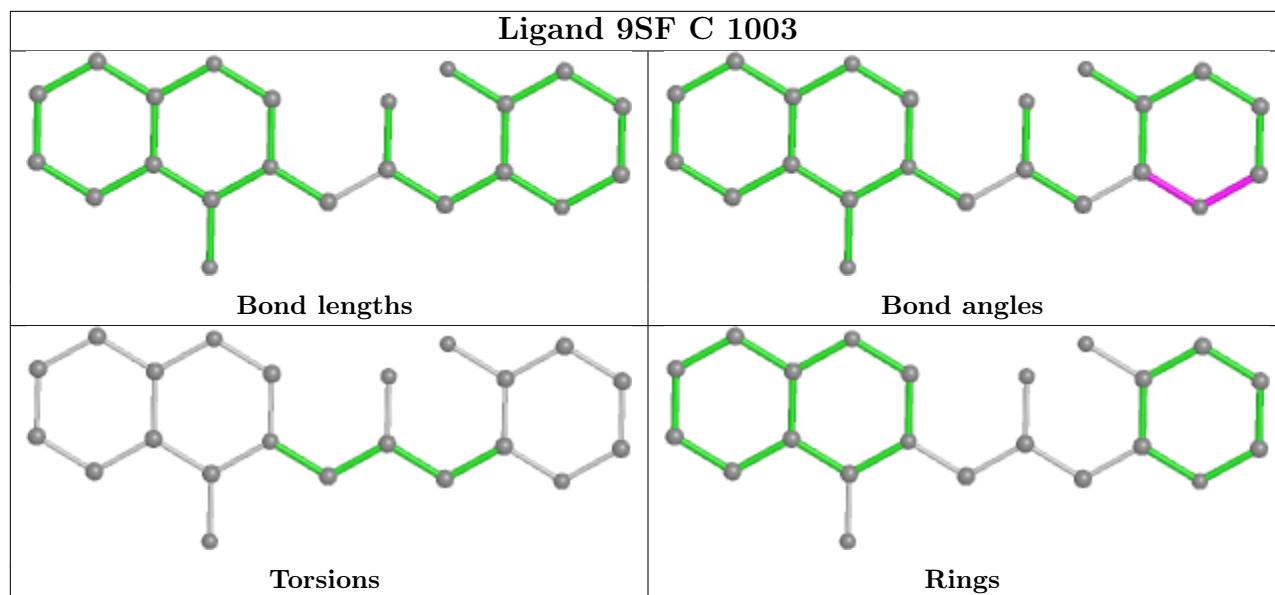
There are no ring outliers.

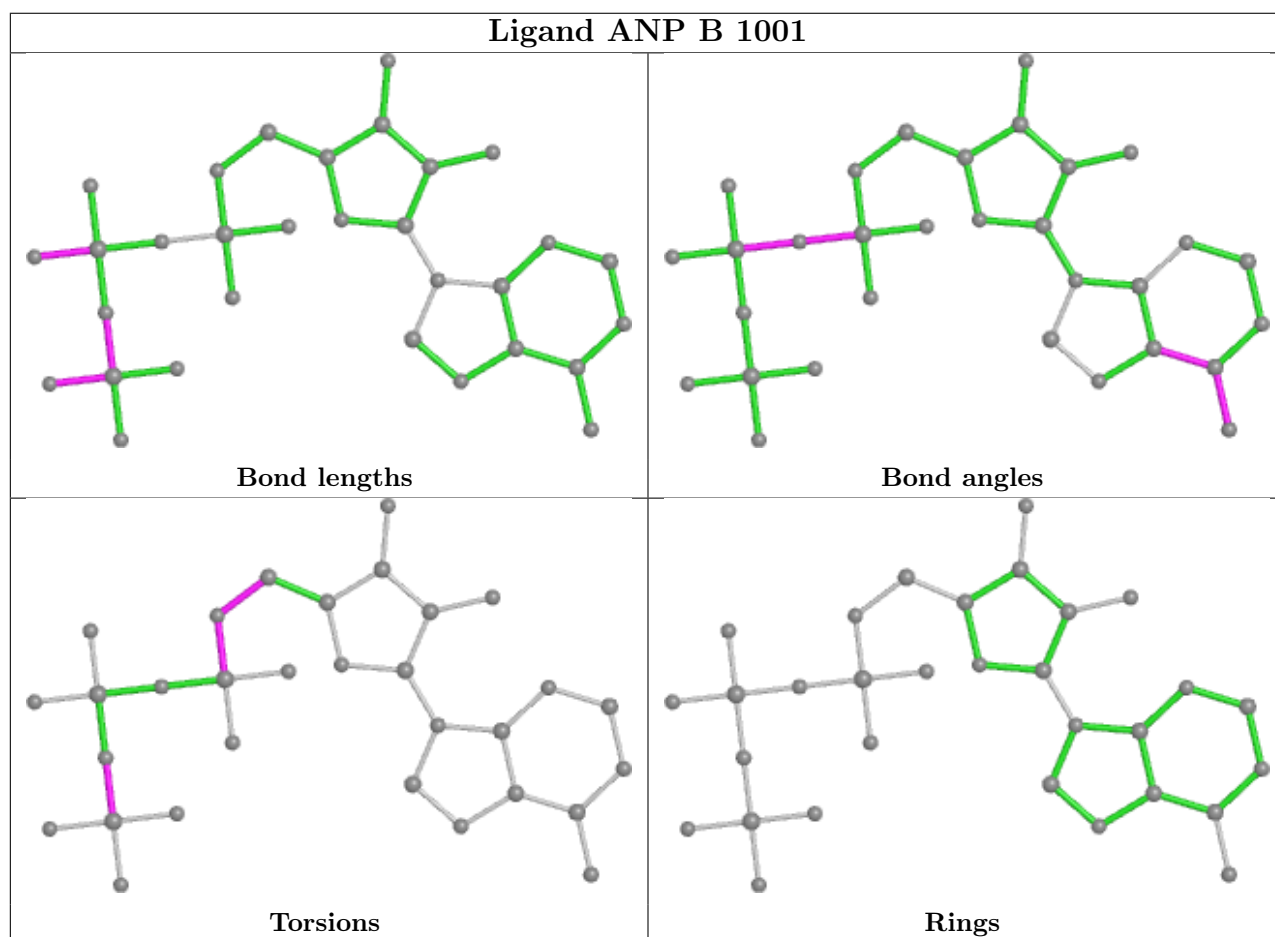
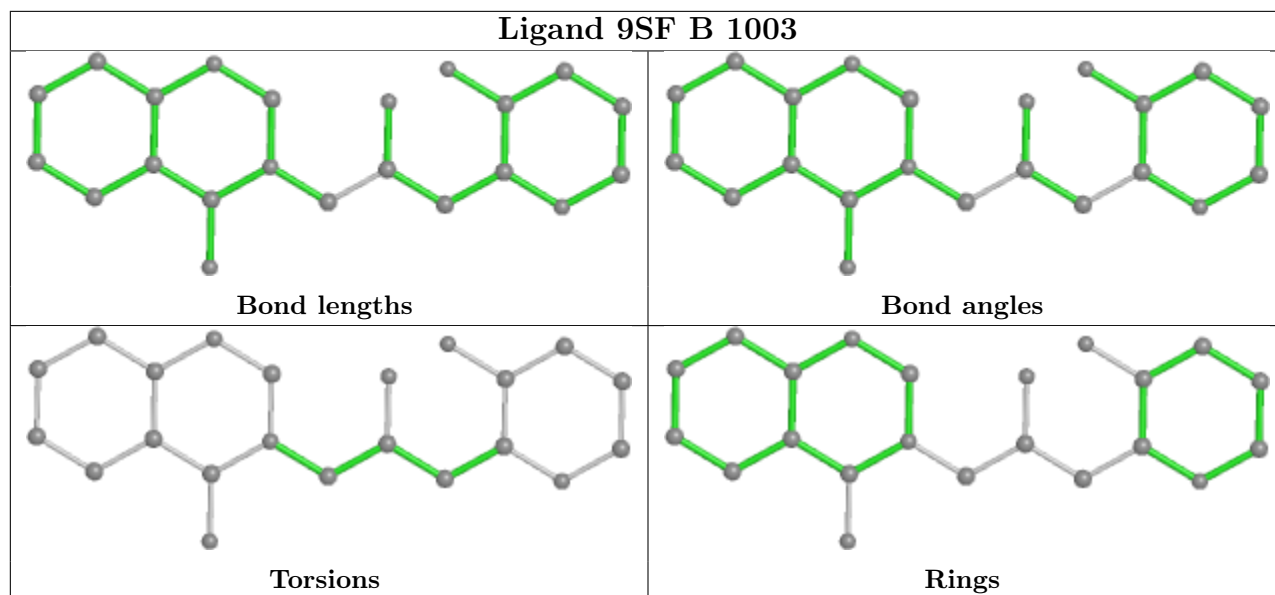
1 monomer is involved in 1 short contact:

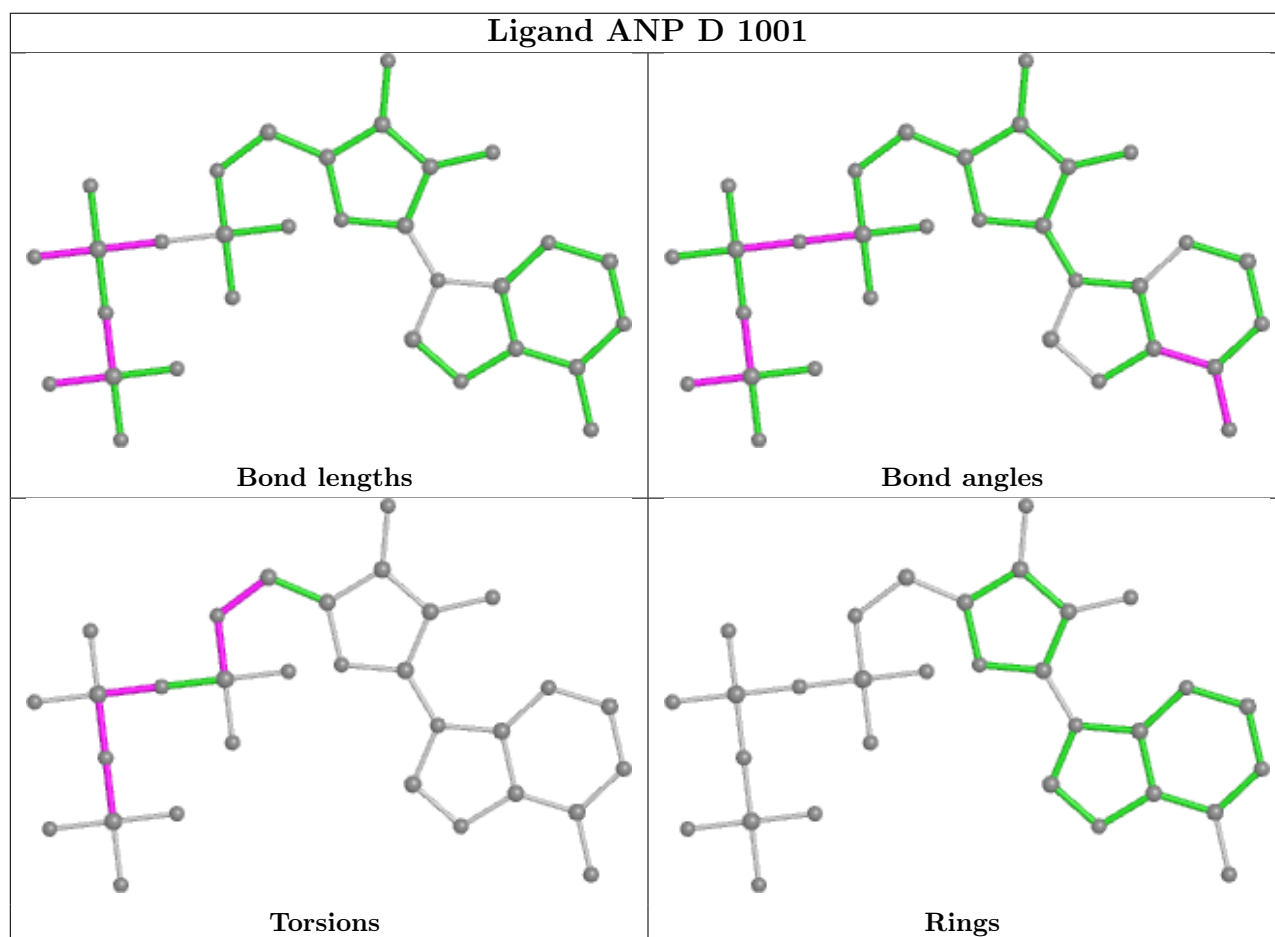
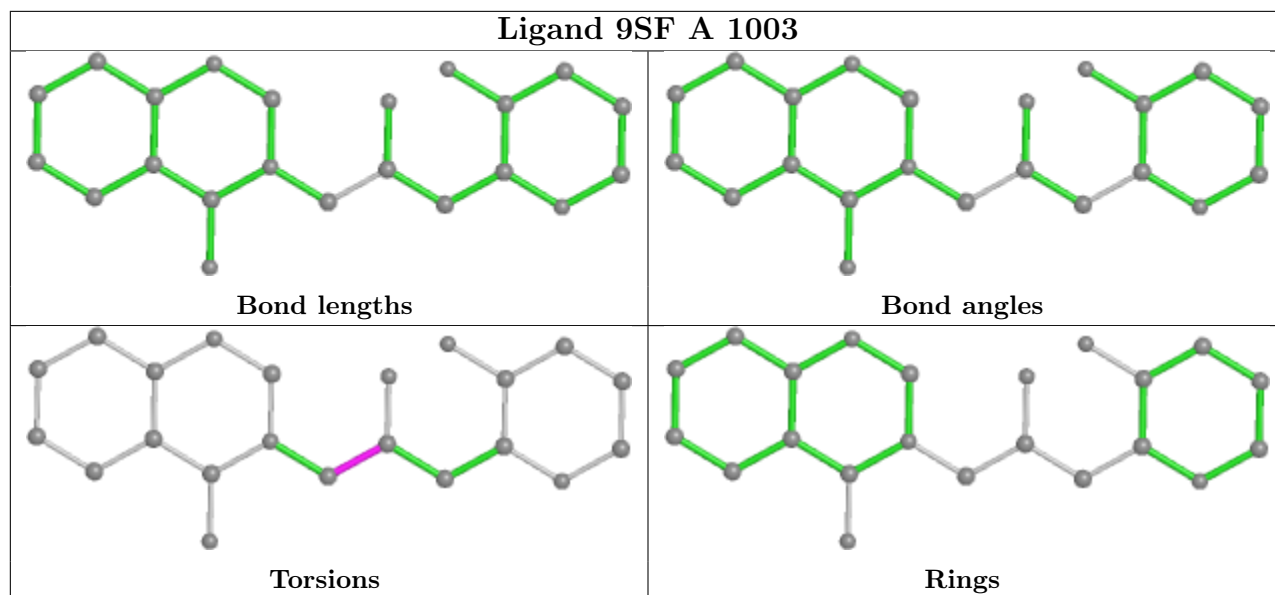
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	ANP	1	0

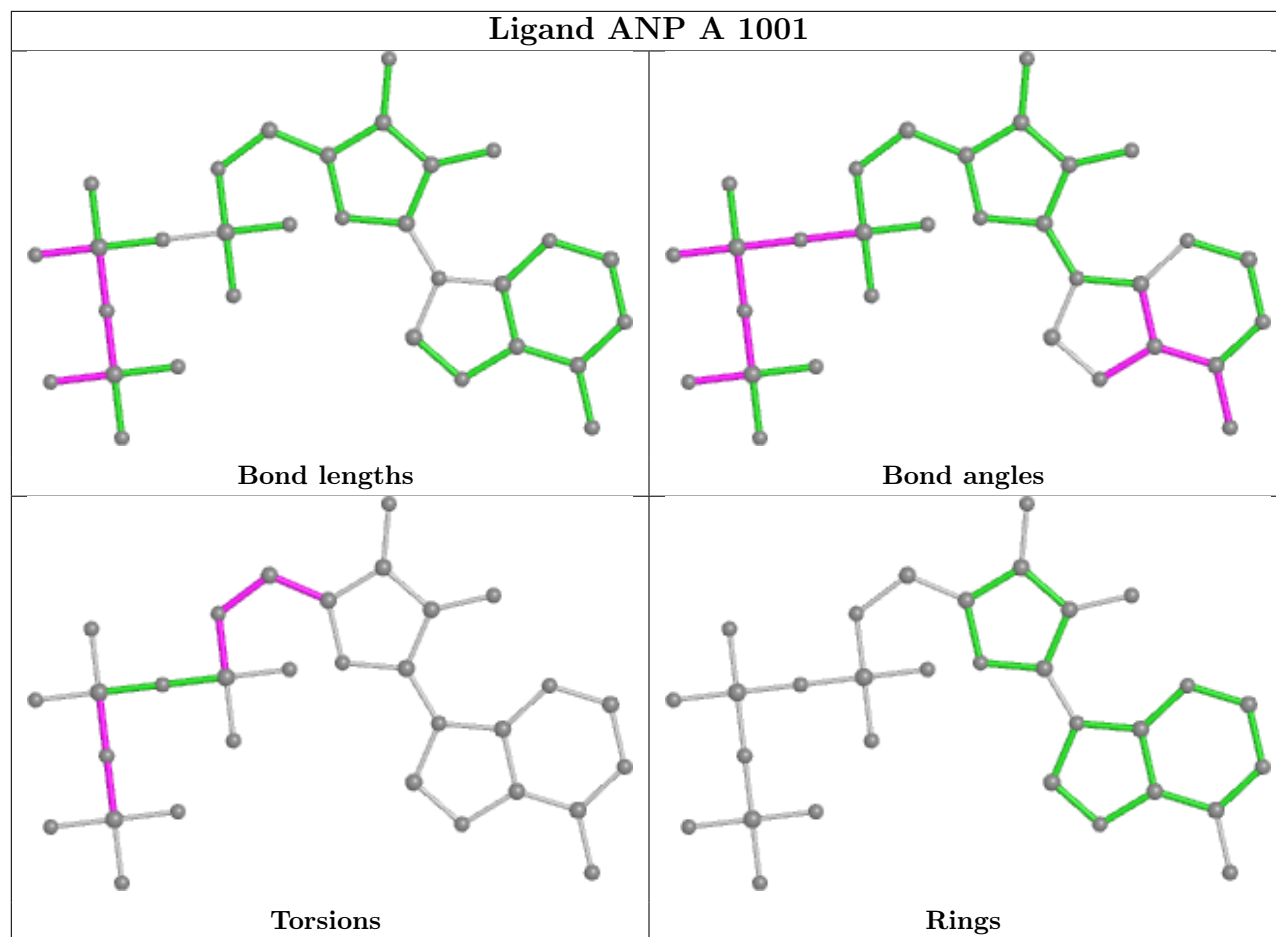
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	482/500 (96%)	-0.00	4 (0%) 86 84	17, 36, 65, 110	1 (0%)
1	B	481/500 (96%)	0.29	22 (4%) 32 26	23, 46, 79, 97	0
1	C	477/500 (95%)	0.28	30 (6%) 20 15	22, 47, 89, 106	0
1	D	475/500 (95%)	0.20	12 (2%) 57 51	20, 43, 73, 96	2 (0%)
All	All	1915/2000 (95%)	0.19	68 (3%) 42 35	17, 43, 79, 110	3 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	814	PHE	4.8
1	C	749	LEU	4.6
1	D	687	ALA	4.5
1	C	627	PHE	4.1
1	B	627	PHE	4.0
1	A	782	ILE	3.9
1	C	629	ASP	3.8
1	C	739	ILE	3.7
1	C	743	ASP	3.7
1	C	812	LYS	3.6
1	D	627	PHE	3.5
1	B	820	ALA	3.5
1	B	763	PRO	3.4
1	B	530	GLU	3.3
1	B	531	GLY	3.2
1	C	742	PHE	3.1
1	C	759	TRP	3.1
1	C	741	THR	3.1
1	B	810	GLY	3.1
1	C	740	SER	3.0
1	C	815	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	650	ASP	2.8
1	B	535	ALA	2.8
1	C	628	LYS	2.8
1	C	745	VAL	2.7
1	C	527	GLU	2.7
1	C	748	ALA	2.7
1	C	415	PHE	2.7
1	B	639	CYS	2.7
1	C	819	PRO	2.7
1	D	577	GLU	2.7
1	C	823	TRP	2.7
1	B	339	ASP	2.7
1	B	533	LYS	2.6
1	B	565	ASN	2.6
1	B	606	ASP	2.5
1	C	660	ASN	2.5
1	C	757	ALA	2.5
1	C	811	THR	2.5
1	B	566	PHE	2.4
1	D	823	TRP	2.4
1	C	816	THR	2.4
1	C	820	ALA	2.4
1	B	681	LEU	2.4
1	D	398	MET	2.3
1	C	339	ASP	2.3
1	B	685	ASP	2.3
1	C	744	GLU	2.3
1	C	808	PRO	2.2
1	B	536	GLY	2.2
1	B	527	GLU	2.2
1	B	532	GLU	2.2
1	C	737	GLU	2.2
1	C	809	GLU	2.2
1	D	762	ASP	2.2
1	D	792	THR	2.2
1	B	415	PHE	2.2
1	A	763	PRO	2.1
1	D	826	TRP	2.1
1	B	473	PHE	2.1
1	D	580	HIS	2.1
1	A	530	GLU	2.1
1	B	568	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	578	GLU	2.1
1	B	364	CYS	2.1
1	C	799	CYS	2.0
1	D	741	THR	2.0
1	D	645	MET	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](#)

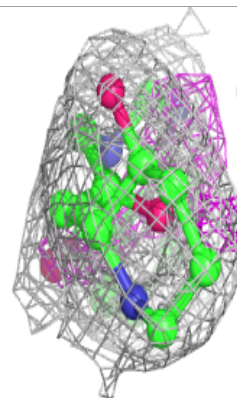
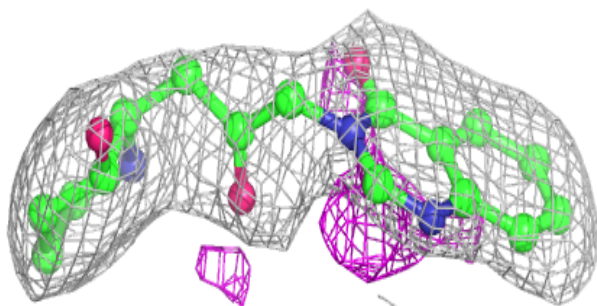
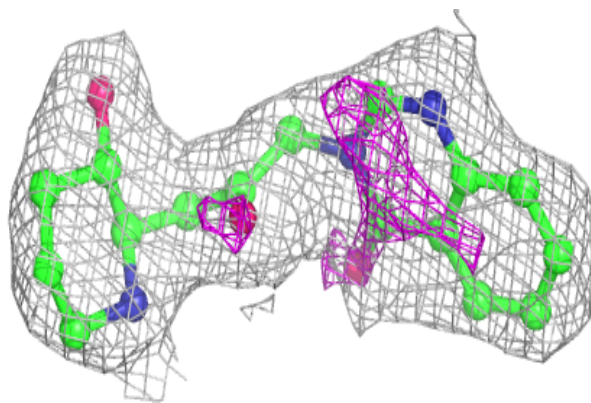
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MG	C	1002	1/1	0.93	0.14	28,28,28,28	0
4	9SF	A	1003	22/22	0.93	0.17	22,33,36,38	0
3	MG	A	1002	1/1	0.95	0.17	29,29,29,29	0
3	MG	B	1002	1/1	0.95	0.15	32,32,32,32	0
4	9SF	B	1003	22/22	0.96	0.15	25,34,41,42	0
4	9SF	C	1003	22/22	0.96	0.15	30,35,40,43	0
4	9SF	D	1003	22/22	0.97	0.14	22,30,35,39	0
2	ANP	D	1001	31/31	0.98	0.15	27,34,42,50	0
2	ANP	A	1001	31/31	0.98	0.14	19,30,41,48	0
2	ANP	B	1001	31/31	0.98	0.15	26,34,49,56	0
2	ANP	C	1001	31/31	0.98	0.15	27,37,50,53	0
3	MG	D	1002	1/1	0.99	0.18	25,25,25,25	0

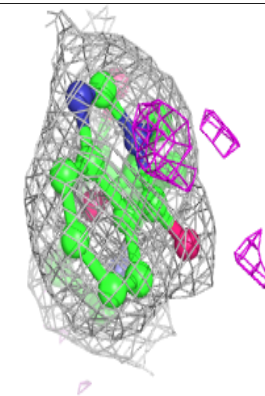
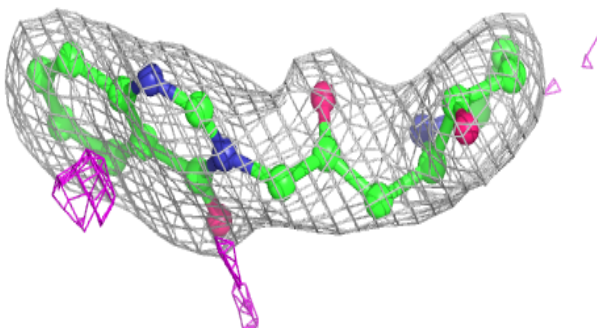
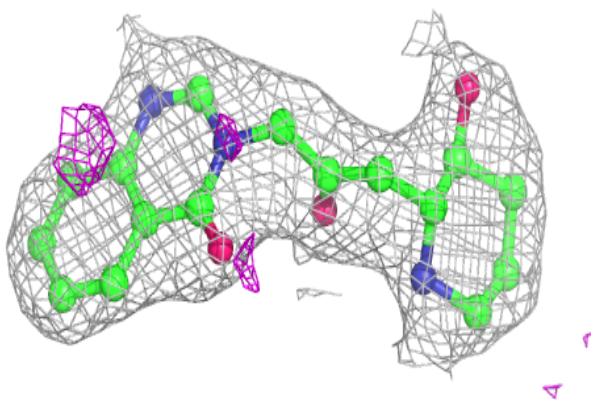
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

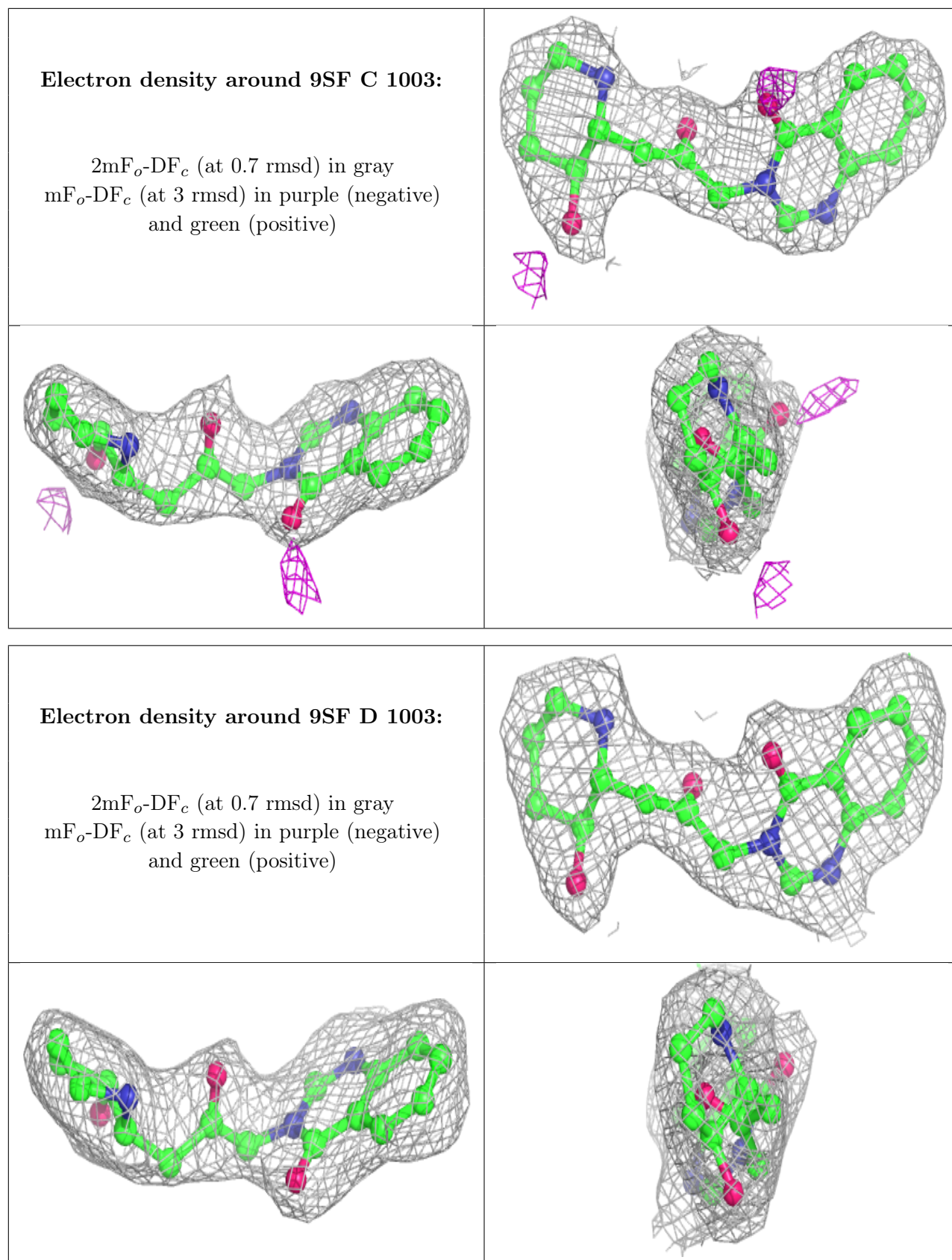
Electron density around 9SF A 1003:

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

**Electron density around 9SF B 1003:**

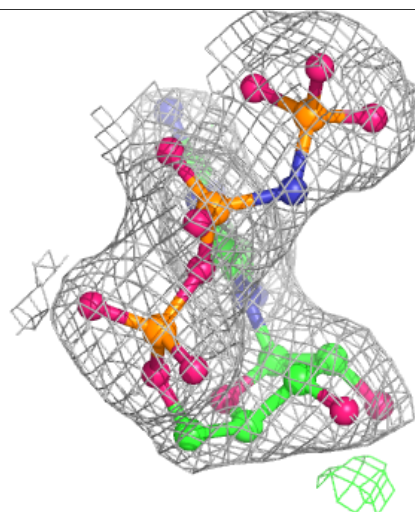
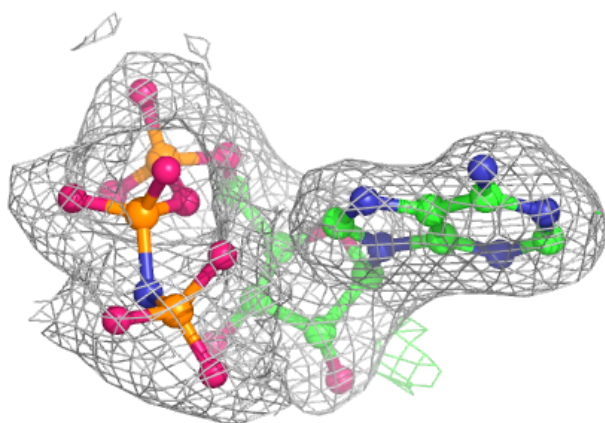
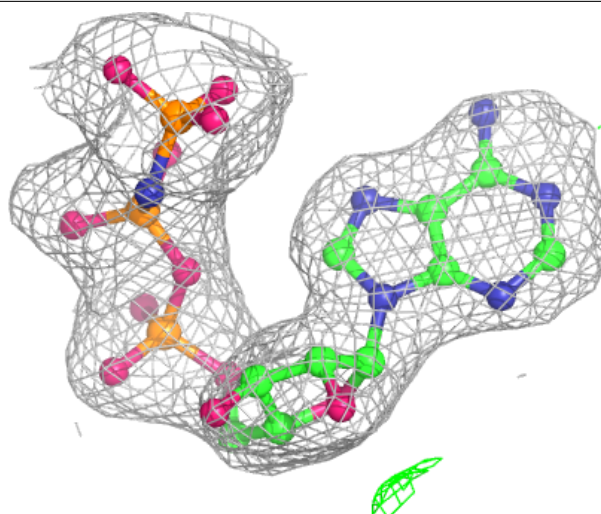
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





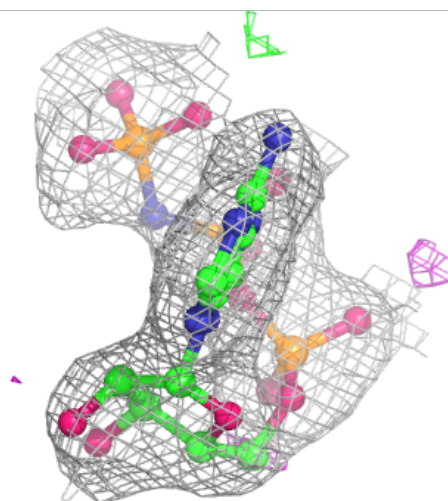
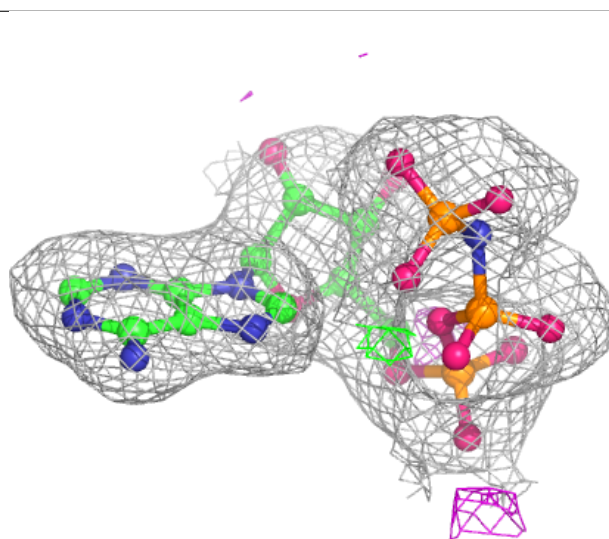
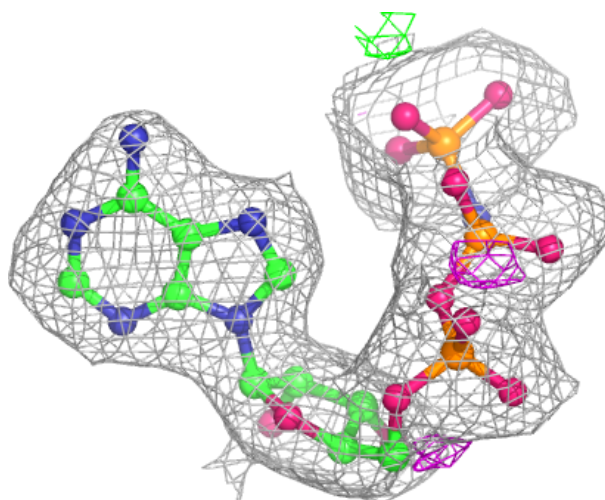
Electron density around ANP D 1001:

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



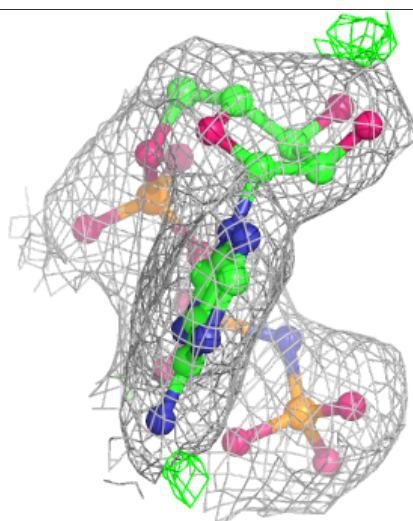
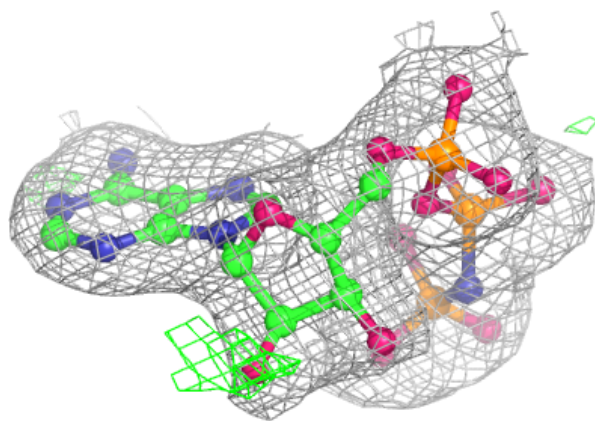
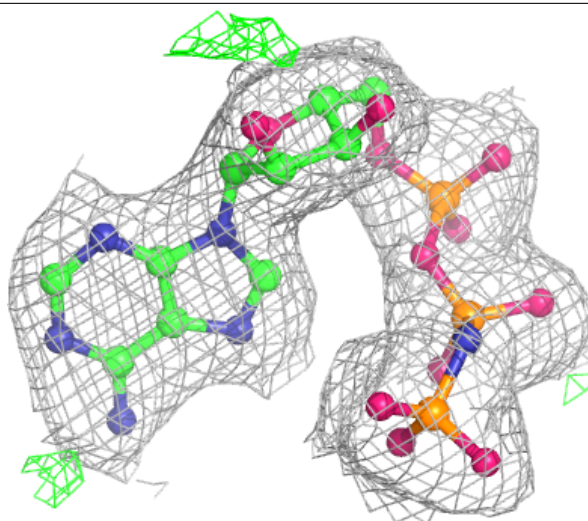
Electron density around ANP A 1001:

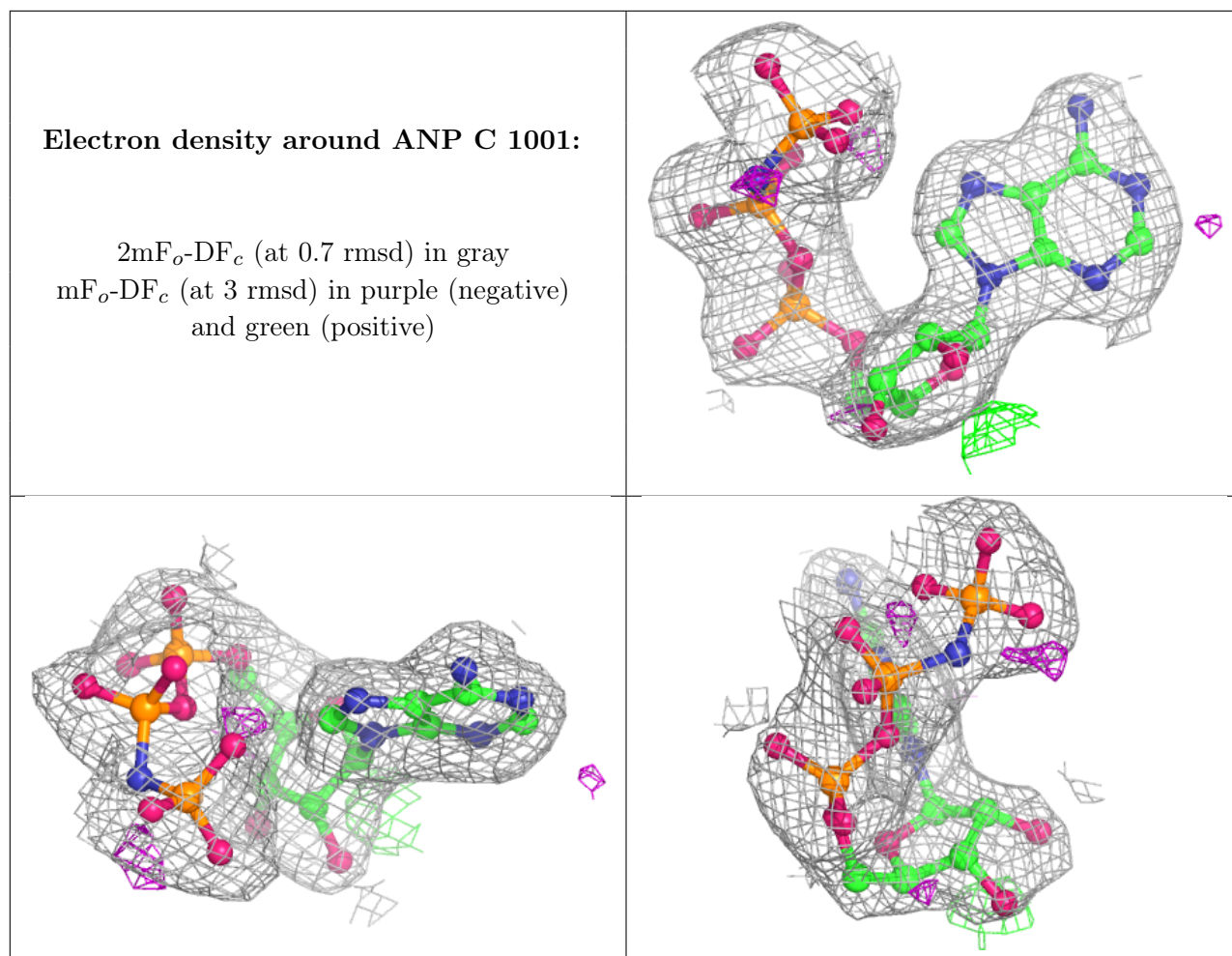
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ANP B 1001:

$2mF_o-DF_c$ (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.