



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 07:57 AM UTC

PDB ID : 6PP9 / pdb_00006pp9
Title : Crystal structure of BRAF:MEK1 complex
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Deposited on : 2019-07-05
Resolution : 2.59 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

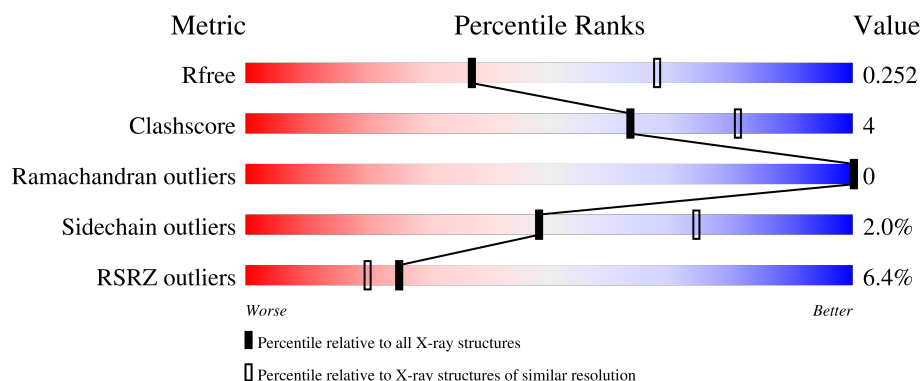
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	283	
2	B	397	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 4858 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 Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	1	0
			2213	1418	386	396	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	441	GLY	-	expression tag	UNP P15056
A	442	GLY	-	expression tag	UNP P15056
A	443	GLY	-	expression tag	UNP P15056
A	444	ARG	-	expression tag	UNP P15056

- Molecule 2 is a protein called Dual specificity mitogen-activated protein kinase kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	314	Total	C	N	O	S	0	2	0
			2465	1574	419	455	17			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP Q02750
B	-2	GLY	-	expression tag	UNP Q02750
B	-1	GLY	-	expression tag	UNP Q02750
B	0	ARG	-	expression tag	UNP Q02750
B	218	ALA	SER	engineered mutation	UNP Q02750
B	222	ALA	SER	engineered mutation	UNP Q02750

- Molecule 3 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃) (labeled as "Ligand of Interest" by depositor).



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

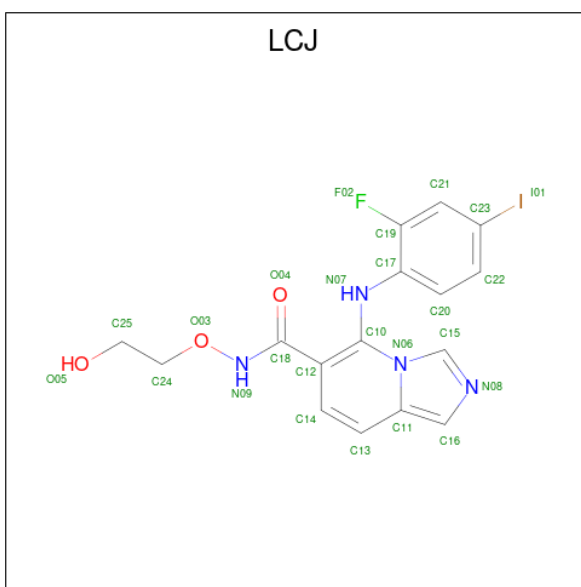
- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Cl 1 1	0	0
5	B	1	Total Cl 1 1	0	0

- Molecule 6 is 5-[(2-fluoro-4-iodophenyl)amino]-N-(2-hydroxyethoxy)imidazo[1,5-a]pyridine-6-carboxamide (CCD ID: LCJ) (formula: C₁₆H₁₄FIN₄O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
6	B	1	Total	C	F	I	N	O	0	0
			25	16	1	1	4	3		

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	O S	0	0
			5	4 1		

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			6	3	3		
8	B	1	Total	C	O	0	0
			6	3	3		

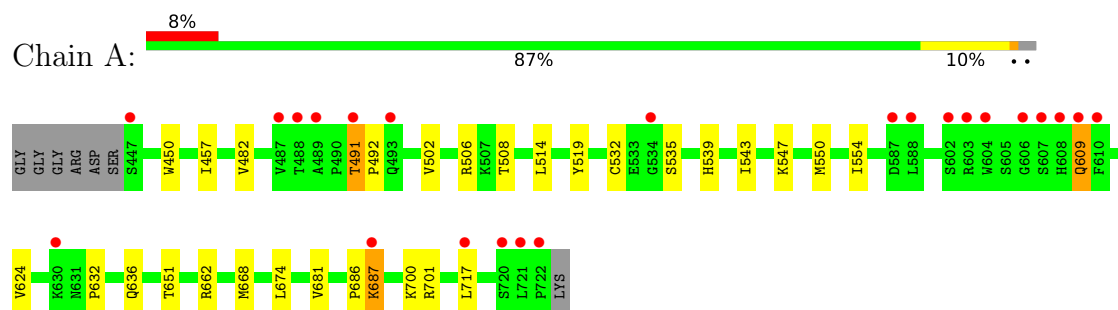
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	37	Total	O	0	0
			37	37		
9	B	35	Total	O	0	0
			35	35		

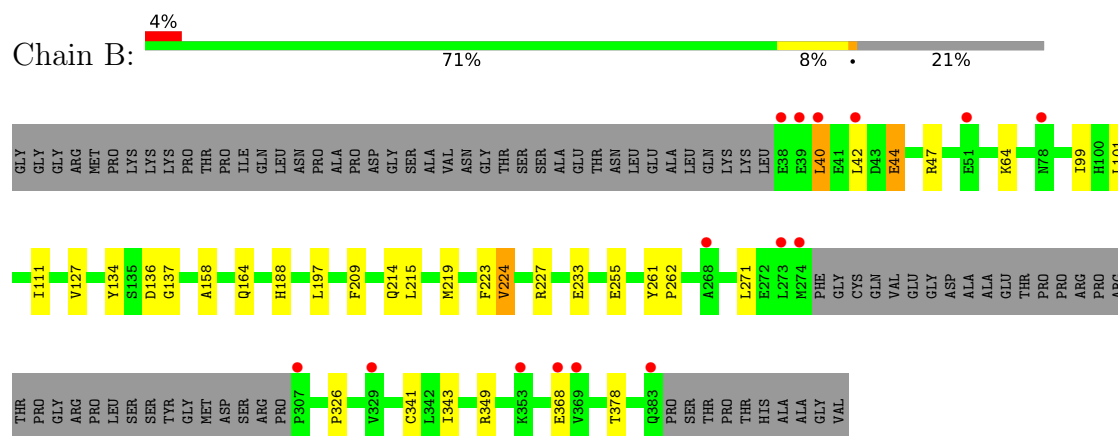
3 Residue-property plots

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: Serine/threonine-protein kinase B-raf



- Molecule 2: Dual specificity mitogen-activated protein kinase kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	116.70Å 116.70Å 130.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.65 – 2.59 43.65 – 2.59	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.65-2.59) 100.0 (43.65-2.59)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.221 , 0.245 0.229 , 0.252	Depositor DCC
R_{free} test set	1543 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	51.8	Xtriage
Anisotropy	0.203	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4858	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 19.81 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.9846e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: SO4, MG, GOL, LCJ, ANP, CL

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	0.26	0/2266	0.32	0/3060
2	B	0.10	0/2518	0.31	0/3391
All	All	0.20	0/4784	0.31	0/6451

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

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	2213	0	2247	20	0
2	B	2465	0	2496	19	0
3	A	31	0	13	0	0
3	B	31	0	13	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	B	25	0	0	0	0
7	B	5	0	0	0	0
8	B	12	0	16	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	37	0	0	0	0
9	B	35	0	0	1	0
All	All	4858	0	4785	38	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 4.

All (38) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:686:PRO:HG2	1:A:717:LEU:HD11	1.85	0.57
1:A:651:THR:HG22	1:A:681:VAL:HA	1.88	0.55
1:A:668:MET:HE3	1:A:674:LEU:HB2	1.88	0.55
1:A:535:SER:HB3	1:A:539:HIS:HB2	1.89	0.55
1:A:491:THR:HG22	1:A:492:PRO:HD2	1.89	0.53
2:B:44:GLU:HG3	2:B:47:ARG:HH12	1.75	0.51
2:B:42:LEU:HD23	2:B:47:ARG:HA	1.92	0.51
2:B:127:VAL:HG21	2:B:197:LEU:HD12	1.94	0.50
1:A:662:ARG:HD2	2:B:224:VAL:HG12	1.94	0.49
2:B:255:GLU:HB2	2:B:262:PRO:HD3	1.93	0.49
2:B:215:LEU:HG	2:B:219:MET:HE2	1.95	0.49
1:A:550:MET:O	1:A:554:ILE:HG12	2.12	0.49
2:B:271:LEU:HD11	2:B:326:PRO:HG3	1.95	0.49
2:B:40:LEU:HB2	2:B:42:LEU:HD13	1.96	0.48
1:A:514[A]:LEU:HD21	1:A:532:CYS:SG	2.54	0.48
1:A:502:VAL:HG11	1:A:519:TYR:CD1	2.49	0.47
1:A:508:THR:O	1:A:508:THR:OG1	2.32	0.47
2:B:134:TYR:OH	2:B:137:GLY:HA2	2.14	0.47
2:B:158:ALA:HB2	2:B:378:THR:HG21	1.96	0.47
2:B:188:HIS:CD2	2:B:209:PHE:HB3	2.51	0.46
2:B:343:ILE:O	2:B:349:ARG:HD3	2.16	0.45
1:A:687:LYS:H	1:A:687:LYS:CD	2.30	0.45
1:A:609:GLN:HE21	1:A:609:GLN:N	2.15	0.45
2:B:219:MET:HG2	2:B:223:PHE:HE1	1.81	0.44
2:B:233:GLU:CD	2:B:349:ARG:HH22	2.25	0.44
1:A:547:LYS:HA	1:A:547:LYS:HD3	1.73	0.44
2:B:99:ILE:HG22	2:B:101:LEU:HG	2.00	0.43
2:B:111:ILE:HG12	2:B:214:GLN:HB3	2.00	0.43
1:A:514[A]:LEU:HD23	1:A:514[A]:LEU:HA	1.83	0.43
1:A:624:VAL:HA	1:A:632:PRO:HB2	2.00	0.42
1:A:636:GLN:HG2	1:A:701:ARG:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:164:GLN:NE2	9:B:907:HOH:O	2.52	0.42
1:A:609:GLN:N	1:A:609:GLN:NE2	2.68	0.42
2:B:341:CYS:O	2:B:349:ARG:HD2	2.19	0.41
1:A:539:HIS:CD2	1:A:543:ILE:HD12	2.56	0.41
1:A:450:TRP:CD2	1:A:506:ARG:HD2	2.56	0.41
1:A:457:ILE:HG21	1:A:482:VAL:HG21	2.03	0.40
2:B:227:ARG:HD2	2:B:261:TYR:CG	2.56	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	275/283 (97%)	266 (97%)	9 (3%)	0	100	100
2	B	312/397 (79%)	303 (97%)	9 (3%)	0	100	100
All	All	587/680 (86%)	569 (97%)	18 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	244/247 (99%)	240 (98%)	4 (2%)	55	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	B	271/337 (80%)	265 (98%)	6 (2%)	45 72
All	All	515/584 (88%)	505 (98%)	10 (2%)	48 75

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	491	THR
1	A	609	GLN
1	A	687	LYS
1	A	700	LYS
2	B	40	LEU
2	B	44	GLU
2	B	64	LYS
2	B	136	ASP
2	B	224	VAL
2	B	368	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	493	GLN
1	A	609	GLN
2	B	164	GLN
2	B	214	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 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 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	801	4	33,33,33	0.92	4 (12%)	45,52,52	0.58	0
8	GOL	B	806	-	5,5,5	0.93	0	5,5,5	1.07	0
3	ANP	B	801	4	33,33,33	0.94	4 (12%)	45,52,52	0.53	0
6	LCJ	B	803	-	26,27,27	2.43	7 (26%)	26,37,37	1.48	5 (19%)
7	SO4	B	804	-	4,4,4	0.83	0	6,6,6	0.08	0
8	GOL	B	807	-	5,5,5	0.92	0	5,5,5	1.07	0

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	801	4	-	4/18/38/38	0/3/3/3
8	GOL	B	806	-	-	4/4/4/4	-
3	ANP	B	801	4	-	3/18/38/38	0/3/3/3
6	LCJ	B	803	-	-	6/13/13/13	0/3/3/3
8	GOL	B	807	-	-	0/4/4/4	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	803	LCJ	C18-N09	7.60	1.45	1.34
6	B	803	LCJ	C11-N06	-6.78	1.31	1.40
6	B	803	LCJ	C10-N07	3.98	1.46	1.37
6	B	803	LCJ	C12-C10	-2.73	1.39	1.43
3	B	801	ANP	PG-N3B	2.58	1.70	1.63
3	A	801	ANP	PG-O1G	2.54	1.50	1.46
3	B	801	ANP	PG-O1G	2.53	1.50	1.46
6	B	803	LCJ	O04-C18	-2.45	1.18	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	ANP	PG-N3B	2.45	1.69	1.63
3	B	801	ANP	PB-O1B	2.43	1.49	1.46
3	A	801	ANP	PB-O1B	2.41	1.49	1.46
6	B	803	LCJ	C17-N07	2.39	1.46	1.41
6	B	803	LCJ	C16-N08	-2.24	1.33	1.37
3	B	801	ANP	PB-O3A	-2.15	1.56	1.59
3	A	801	ANP	PB-O3A	-2.10	1.56	1.59

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	803	LCJ	C16-C11-N06	3.47	108.03	104.58
6	B	803	LCJ	C21-C19-C17	-2.78	120.59	123.43
6	B	803	LCJ	C13-C11-C16	-2.51	131.86	137.90
6	B	803	LCJ	C15-N06-C11	2.31	108.28	106.56
6	B	803	LCJ	C16-N08-C15	2.24	108.17	105.24

There are no chirality outliers.

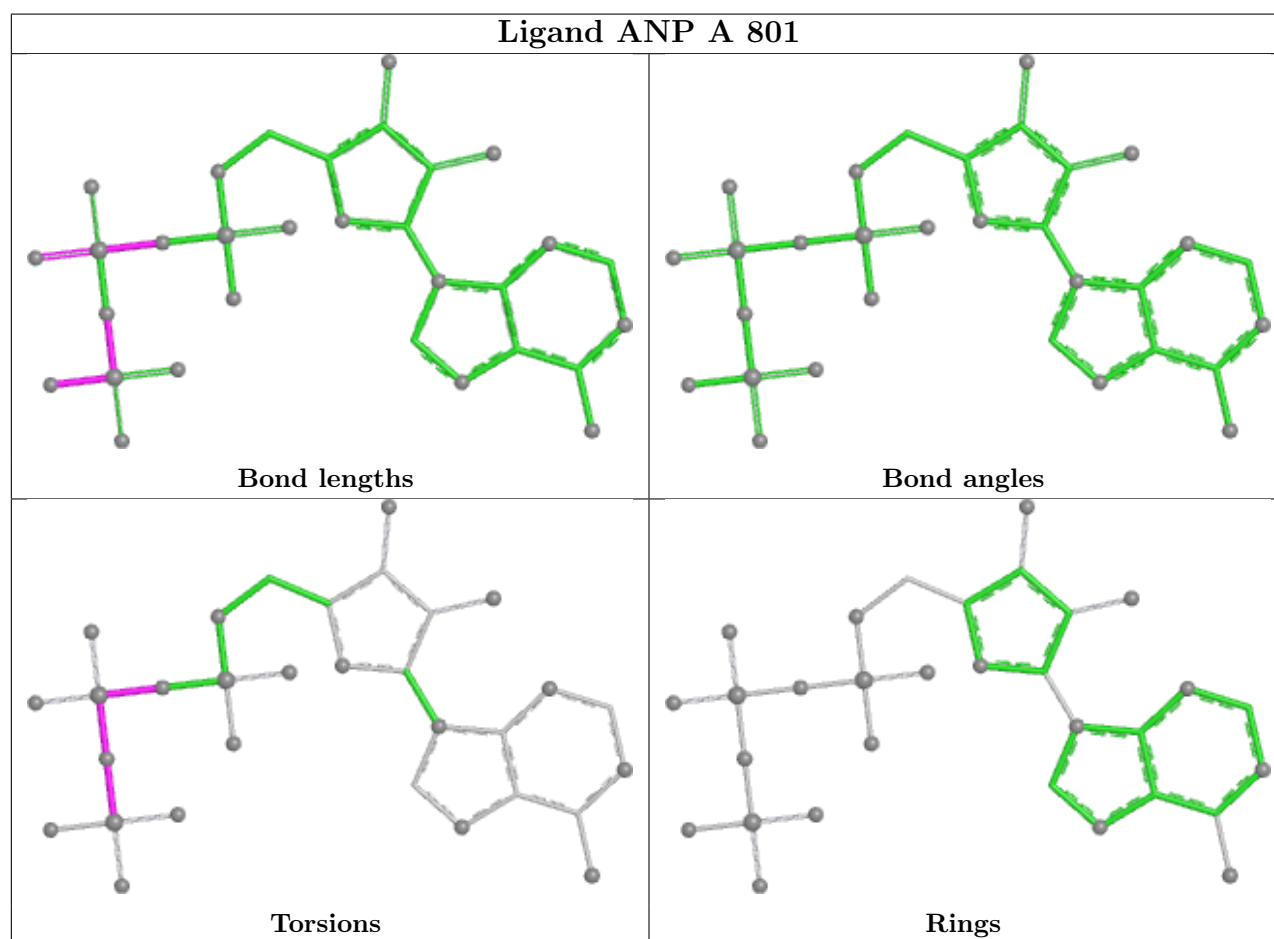
All (17) torsion outliers are listed below:

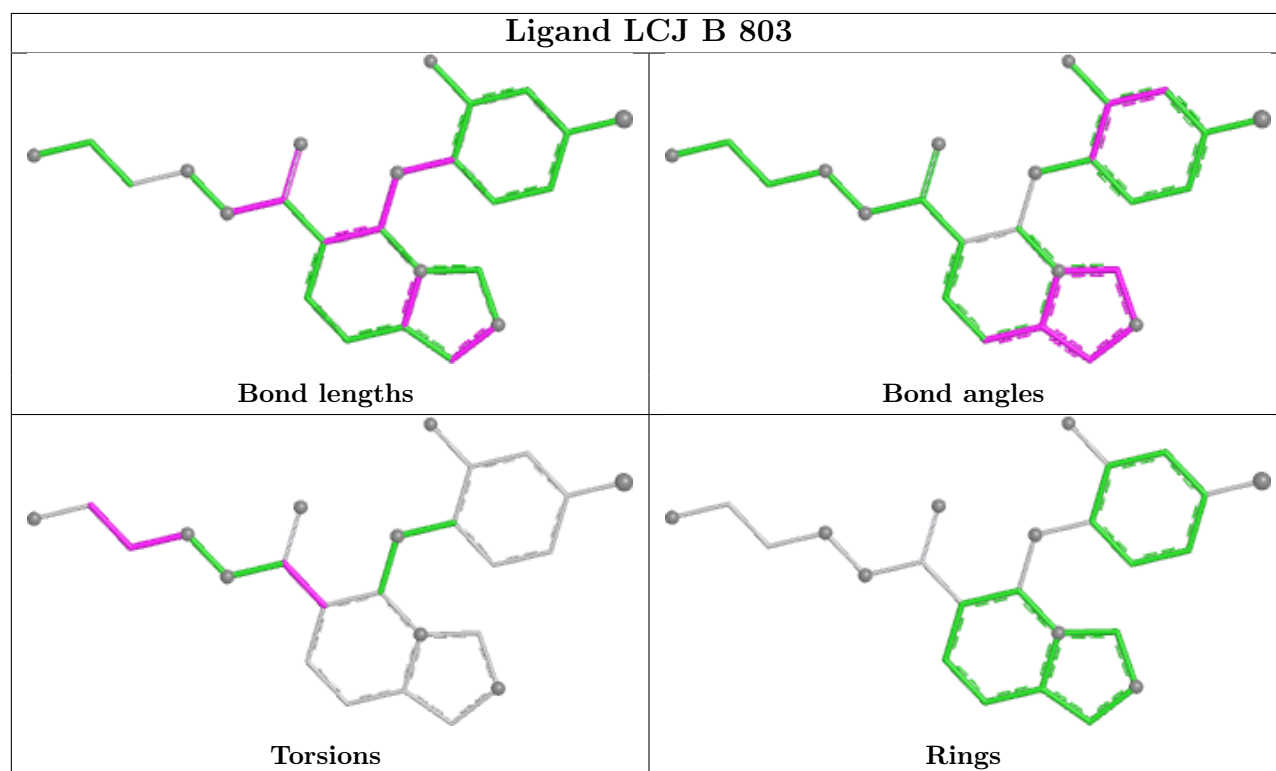
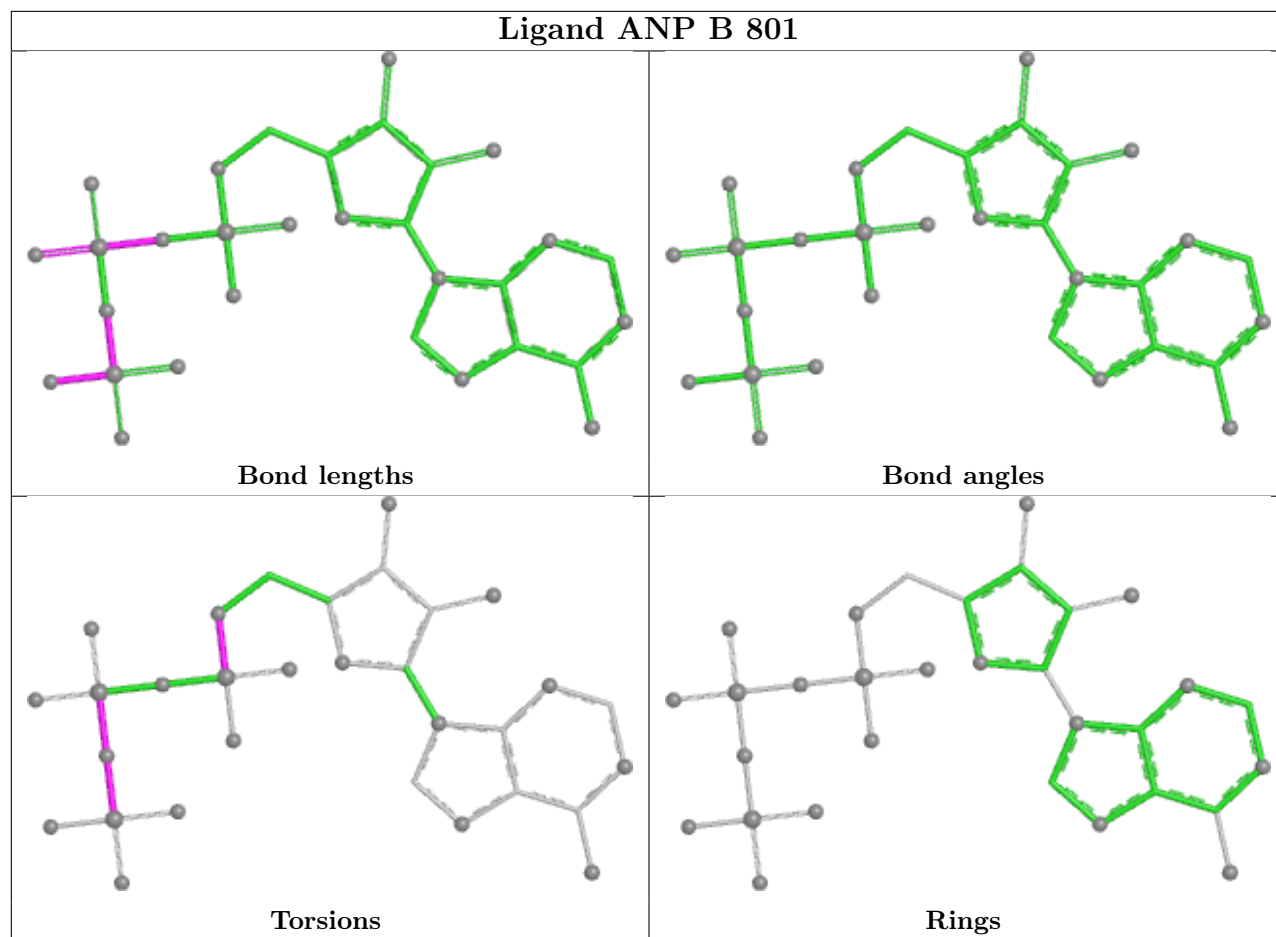
Mol	Chain	Res	Type	Atoms
3	A	801	ANP	PB-N3B-PG-O1G
3	A	801	ANP	PG-N3B-PB-O1B
3	A	801	ANP	PA-O3A-PB-O2B
3	B	801	ANP	PB-N3B-PG-O1G
3	B	801	ANP	PG-N3B-PB-O1B
6	B	803	LCJ	C14-C12-C18-O04
8	B	806	GOL	O1-C1-C2-C3
8	B	806	GOL	C1-C2-C3-O3
8	B	806	GOL	O2-C2-C3-O3
6	B	803	LCJ	C14-C12-C18-N09
8	B	806	GOL	O1-C1-C2-O2
6	B	803	LCJ	C25-C24-O03-N09
6	B	803	LCJ	C10-C12-C18-N09
6	B	803	LCJ	C10-C12-C18-O04
6	B	803	LCJ	O03-C24-C25-O05
3	B	801	ANP	C5'-O5'-PA-O1A
3	A	801	ANP	PA-O3A-PB-O1B

There are no ring outliers.

No monomer is involved in short contacts.

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





5.7 Other polymers

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	276/283 (97%)	0.41	23 (8%) 17 13	28, 51, 84, 117	1 (0%)
2	B	314/397 (79%)	0.40	15 (4%) 35 30	27, 47, 83, 111	2 (0%)
All	All	590/680 (86%)	0.40	38 (6%) 25 20	27, 49, 84, 117	3 (0%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	38	GLU	8.1
2	B	307	PRO	5.4
2	B	273	LEU	4.6
1	A	722	PRO	4.5
1	A	610	PHE	4.0
1	A	608	HIS	3.8
1	A	447	SER	3.6
2	B	268	ALA	3.6
1	A	588	LEU	3.5
1	A	491	THR	3.4
1	A	717	LEU	3.1
1	A	606	GLY	3.1
1	A	604	TRP	3.0
1	A	534	GLY	3.0
1	A	603	ARG	2.9
2	B	40	LEU	2.8
2	B	78	ASN	2.6
1	A	602	SER	2.6
1	A	493	GLN	2.5
2	B	274	MET	2.5
2	B	368	GLU	2.5
1	A	487	VAL	2.5
1	A	630	LYS	2.5
2	B	369	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	720	SER	2.4
1	A	721	LEU	2.4
2	B	329	VAL	2.4
1	A	687	LYS	2.3
2	B	51	GLU	2.3
2	B	383	GLN	2.3
2	B	42	LEU	2.3
1	A	607	SER	2.2
1	A	488	THR	2.2
1	A	489	ALA	2.1
2	B	353	LYS	2.1
1	A	609	GLN	2.1
2	B	39	GLU	2.0
1	A	587	ASP	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.

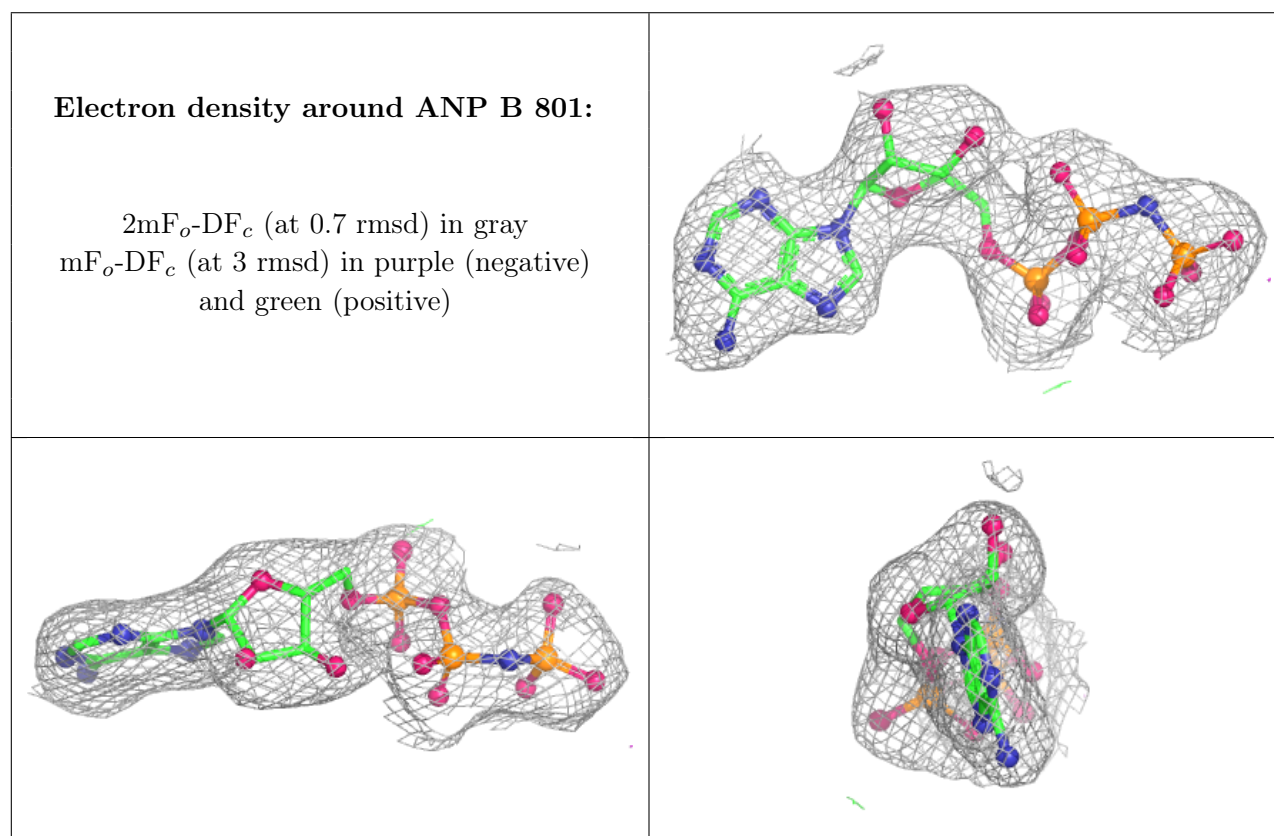
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CL	B	805	1/1	0.86	0.15	79,79,79,79	0
8	GOL	B	807	6/6	0.88	0.12	51,56,60,74	0
5	CL	A	803	1/1	0.92	0.11	79,79,79,79	0
8	GOL	B	806	6/6	0.93	0.12	49,59,63,69	0
7	SO4	B	804	5/5	0.93	0.12	65,67,71,73	0
3	ANP	B	801	31/31	0.97	0.07	36,43,51,58	0
6	LCJ	B	803	25/25	0.98	0.10	38,44,64,75	0
3	ANP	A	801	31/31	0.98	0.05	33,41,48,49	0
4	MG	B	802	1/1	0.99	0.03	51,51,51,51	0

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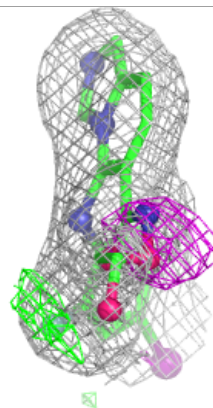
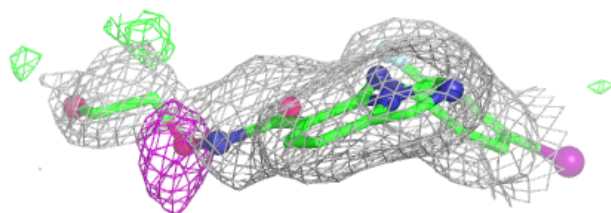
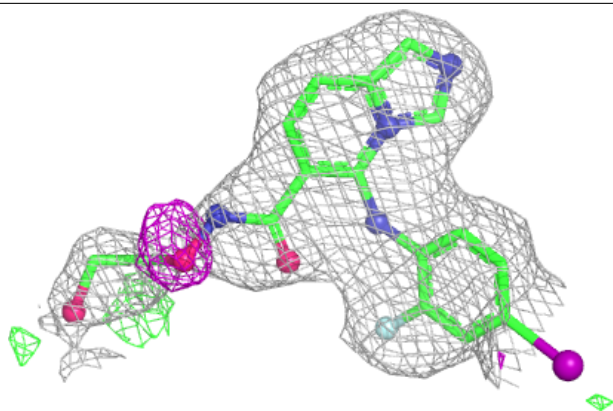
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	802	1/1	1.00	0.03	41,41,41,41	0

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.

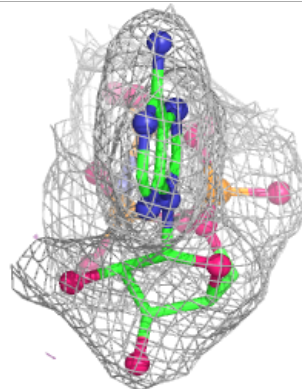
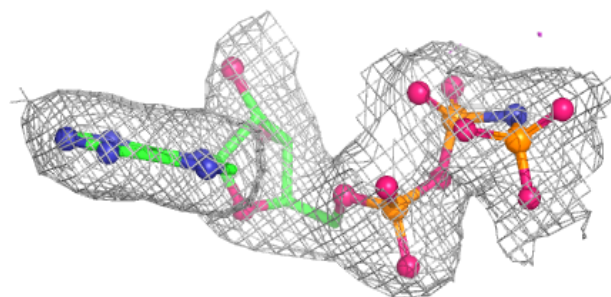
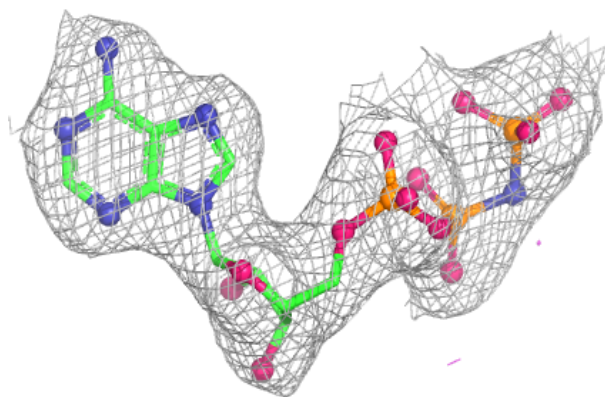


Electron density around LCJ B 803:

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

**Electron density around ANP A 801:**

$2mF_o-DF_c$ (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.