



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:44 PM UTC

PDB ID : 5XIO / pdb_00005xio
Title : Crystal Structure of Cryptosporidium parvum Prolyl-tRNA Synthetase (Cp-PRS) in complex with Halofuginone
Authors : Jain, V.; Manickam, Y.; Sharma, A.
Deposited on : 2017-04-26
Resolution : 2.46 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

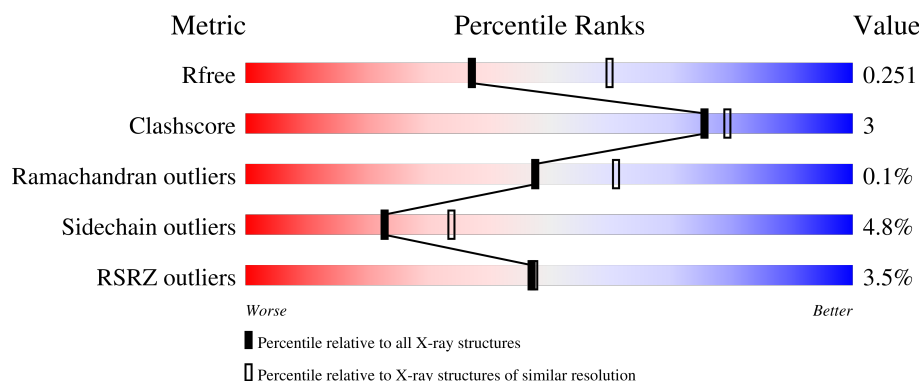
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	498	 3% 82% 12% • 5%
1	B	498	 4% 82% 11% • 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7892 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.

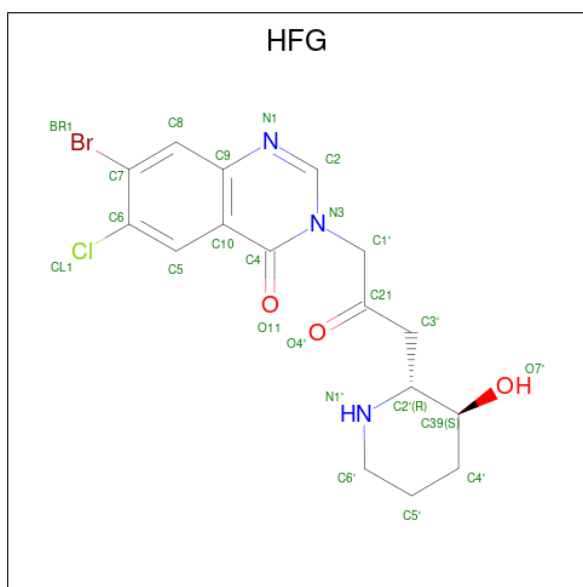
- Molecule 1 is a protein called Proline-tRNA synthetase class II aaRS (Ybak RNA binding domain plus tRNA synthetase).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	474	Total	C	N	O	S	0	2	0
			3793	2446	639	687	21			
1	B	471	Total	C	N	O	S	0	2	0
			3788	2443	642	682	21			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	GLY	-	expression tag	UNP Q5CWN3
A	223	ALA	-	expression tag	UNP Q5CWN3
A	224	MET	-	expression tag	UNP Q5CWN3
B	222	GLY	-	expression tag	UNP Q5CWN3
B	223	ALA	-	expression tag	UNP Q5CWN3
B	224	MET	-	expression tag	UNP Q5CWN3

- Molecule 2 is 7-bromo-6-chloro-3-{3-[(2R,3S)-3-hydroxypiperidin-2-yl]-2-oxopropyl}quinazolin-4(3H)-one (CCD ID: HFG) (formula: C₁₆H₁₇BrClN₃O₃).

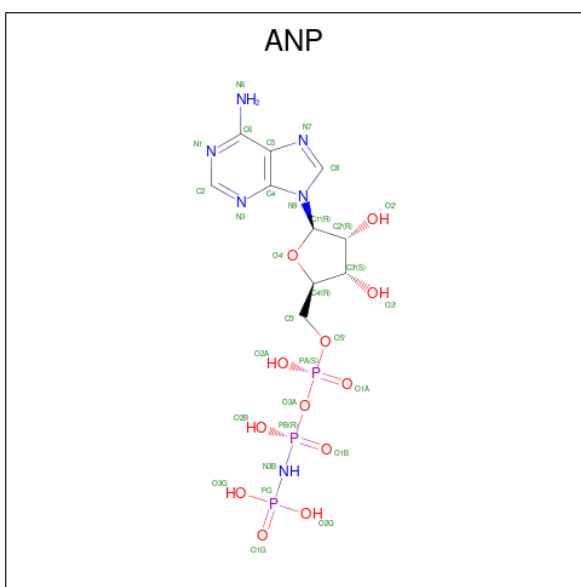


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	Br	C	Cl	N	O	0	0
			24	1	16	1	3	3		
2	B	1	Total	Br	C	Cl	N	O	0	0
			24	1	16	1	3	3		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is PHOSPHOAMINOPHOSPHONIC ACID-ADENYLATE ESTER (CCD ID: ANP) (formula: C₁₀H₁₇N₆O₁₂P₃).



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

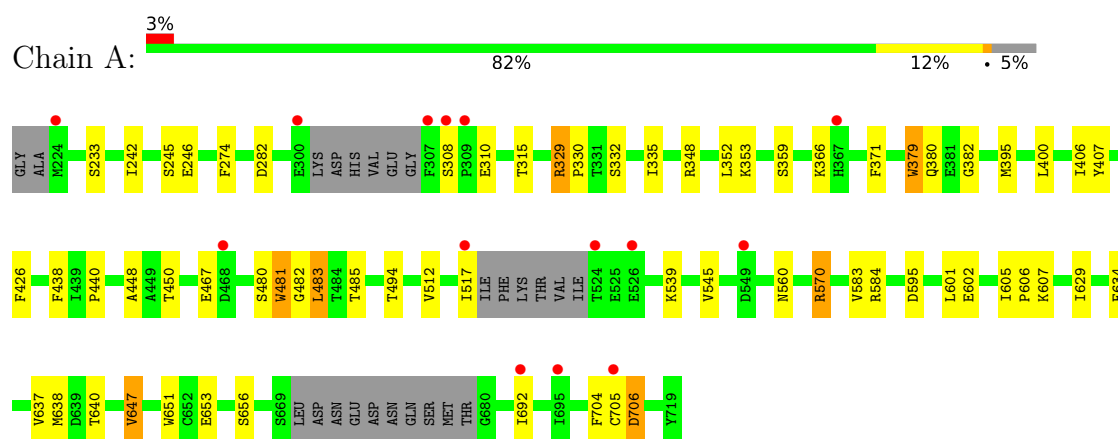
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	96	Total O 96 96	0	0
5	B	102	Total O 103 103	0	1

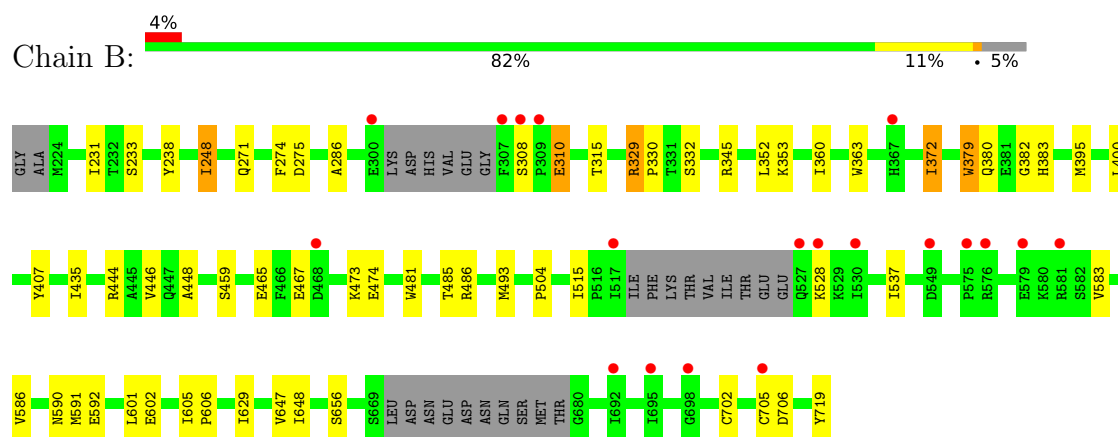
3 Residue-property plots [i](#)

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: Proline-tRNA synthetase class II aaRS (Ybak RNA binding domain plus tRNA synthetase)



- Molecule 1: Proline-tRNA synthetase class II aaRS (Ybak RNA binding domain plus tRNA synthetase)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.77Å 112.64Å 140.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.31 – 2.46 29.31 – 2.46	Depositor EDS
% Data completeness (in resolution range)	98.5 (29.31-2.46) 98.6 (29.31-2.46)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.06 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.196 , 0.252 0.206 , 0.251	Depositor DCC
R_{free} test set	2428 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	37.9	Xtriage
Anisotropy	0.494	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7892	wwPDB-VP
Average B, all atoms (Å ²)	41.0	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 52.53 % 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 4.7972e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: ZN, HFG, 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	1.23	12/3888 (0.3%)	1.17	10/5259 (0.2%)
1	B	1.20	10/3883 (0.3%)	1.18	9/5248 (0.2%)
All	All	1.21	22/7771 (0.3%)	1.18	19/10507 (0.2%)

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	704	PHE	CA-C	6.95	1.59	1.52
1	B	400	LEU	N-CA	-6.48	1.38	1.46
1	B	504	PRO	C-O	-6.25	1.17	1.24
1	A	482	GLY	N-CA	5.92	1.51	1.45
1	B	383	HIS	CA-C	5.72	1.58	1.52
1	B	332	SER	C-O	-5.63	1.16	1.24
1	A	330	PRO	C-O	-5.56	1.16	1.24
1	A	494	THR	CA-C	5.48	1.59	1.52
1	B	382	GLY	CA-C	5.46	1.56	1.52
1	A	382	GLY	CA-C	5.45	1.56	1.52
1	B	352	LEU	CA-C	5.37	1.59	1.52
1	A	308	SER	CA-C	5.27	1.59	1.52
1	A	647	VAL	N-CA	-5.24	1.40	1.46
1	B	286	ALA	CA-C	-5.21	1.46	1.52
1	A	483	LEU	N-CA	5.21	1.52	1.46
1	A	400	LEU	N-CA	-5.19	1.40	1.46
1	B	330	PRO	C-O	-5.16	1.17	1.24
1	A	359	SER	N-CA	-5.14	1.39	1.45
1	B	274	PHE	C-O	-5.14	1.18	1.24
1	B	310	GLU	CA-C	5.12	1.59	1.52
1	A	352	LEU	N-CA	-5.12	1.40	1.46
1	A	560	ASN	CA-C	5.08	1.59	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	332	SER	N-CA-C	9.22	124.26	112.92
1	A	332	SER	N-CA-C	9.04	124.04	112.92
1	B	231	ILE	N-CA-C	7.25	117.38	110.42
1	A	637	VAL	N-CA-C	6.78	117.28	110.23
1	B	329	ARG	N-CA-C	6.69	119.21	109.71
1	A	634	PHE	N-CA-C	6.69	120.54	112.38
1	A	335	ILE	N-CA-C	-6.44	104.76	111.58
1	A	308	SER	N-CA-C	6.40	122.41	113.57
1	B	515	ILE	N-CA-C	6.06	114.09	107.60
1	B	372	ILE	N-CA-C	6.01	116.79	110.72
1	B	308	SER	CA-C-N	5.96	127.29	119.84
1	B	308	SER	C-N-CA	5.96	127.29	119.84
1	A	329	ARG	N-CA-C	5.87	118.05	109.71
1	A	426	PHE	CA-C-N	-5.87	113.63	119.56
1	A	426	PHE	C-N-CA	-5.87	113.63	119.56
1	B	586	VAL	N-CA-C	5.46	115.73	107.75
1	B	238	TYR	N-CA-C	-5.43	105.44	111.36
1	A	607	LYS	N-CA-C	5.23	117.89	111.82
1	A	282	ASP	N-CA-C	5.03	118.68	112.24

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	3793	0	3731	21	0
1	B	3788	0	3757	25	0
2	A	24	0	17	1	0
2	B	24	0	17	3	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	31	0	13	0	0
4	B	31	0	13	2	0
5	A	96	0	0	2	0
5	B	103	0	0	2	0
All	All	7892	0	7548	46	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (46) 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)
2:B:802:HFG:O7'	4:B:803:ANP:O1A	1.99	0.79
1:B:329:ARG:HD3	5:B:904:HOH:O	1.85	0.76
1:B:467:GLU:HG3	1:B:473:LYS:HE3	1.74	0.68
1:B:372:ILE:HD12	1:B:493:MET:CE	2.27	0.65
1:A:407:TYR:CZ	1:A:448:ALA:HB1	2.36	0.61
1:B:310:GLU:HB3	2:B:802:HFG:BR1	2.61	0.56
1:A:570:ARG:O	1:A:570:ARG:HG2	2.07	0.54
1:A:651:TRP:CZ3	1:A:653:GLU:HA	2.44	0.53
1:A:517:ILE:HG22	1:A:517:ILE:O	2.10	0.51
1:A:512:VAL:O	1:A:545:VAL:HA	2.10	0.51
1:B:310:GLU:HG3	1:B:363:TRP:O	2.12	0.50
1:B:702:CYS:HB3	1:B:705:CYS:O	2.13	0.49
1:A:570:ARG:HD3	5:A:976:HOH:O	2.12	0.49
1:B:601:LEU:O	1:B:602:GLU:C	2.56	0.48
1:A:483:LEU:HD23	1:A:483:LEU:C	2.39	0.48
1:A:438:PHE:O	1:A:440:PRO:HD3	2.15	0.47
1:A:274:PHE:HB2	1:A:406:ILE:HD11	1.96	0.47
1:A:651:TRP:CH2	1:A:653:GLU:HA	2.50	0.47
1:B:467:GLU:HG3	1:B:473:LYS:CE	2.44	0.47
1:B:372:ILE:HD12	1:B:493:MET:HE3	1.97	0.46
1:A:315:THR:OG1	1:B:315:THR:OG1	2.15	0.46
1:B:605:ILE:O	1:B:606:PRO:C	2.56	0.46
1:B:407:TYR:CZ	1:B:448:ALA:HB1	2.52	0.45
1:B:435:ILE:HG13	1:B:448:ALA:HB3	1.98	0.44
1:A:353:LYS:HE2	1:A:395:MET:CE	2.46	0.44
1:A:705:CYS:SG	1:A:706:ASP:N	2.91	0.44
1:B:248:ILE:HA	1:B:248:ILE:HD13	1.77	0.44
1:B:353:LYS:HE2	1:B:395:MET:CE	2.48	0.44
1:B:271:GLN:O	1:B:275:ASP:HB2	2.18	0.43
1:B:444:ARG:HD3	1:B:719:TYR:HA	2.00	0.42
1:B:705:CYS:SG	5:B:917:HOH:O	2.56	0.42
1:B:446:VAL:HG22	1:B:719:TYR:HB2	2.00	0.42
1:A:638:MET:HG2	5:A:987:HOH:O	2.18	0.42
1:A:379:TRP:CH2	2:A:801:HFG:H8	2.55	0.42
1:A:245:SER:O	1:A:246:GLU:HB2	2.19	0.42
1:B:379:TRP:CH2	2:B:802:HFG:H8	2.55	0.42
1:A:605:ILE:O	1:A:606:PRO:C	2.61	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:480:SER:C	1:A:481:TRP:CD1	2.98	0.41
1:B:329:ARG:HG2	1:B:360:ILE:HD13	2.02	0.41
1:B:486:ARG:HD2	4:B:803:ANP:O2'	2.19	0.41
1:B:590:ASN:OD1	1:B:592:GLU:HB3	2.21	0.41
1:A:584:ARG:NH1	1:A:595:ASP:OD2	2.54	0.41
1:B:590:ASN:O	1:B:591:MET:HB2	2.20	0.41
1:A:601:LEU:O	1:A:602:GLU:C	2.63	0.40
1:A:651:TRP:CE3	1:A:653:GLU:HA	2.57	0.40
1:B:372:ILE:HD12	1:B:493:MET:HE1	2.01	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	468/498 (94%)	452 (97%)	15 (3%)	1 (0%)	43	54
1	B	465/498 (93%)	448 (96%)	17 (4%)	0	100	100
All	All	933/996 (94%)	900 (96%)	32 (3%)	1 (0%)	48	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	371	PHE

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	403/449 (90%)	382 (95%)	21 (5%)	21	30
1	B	406/449 (90%)	388 (96%)	18 (4%)	25	37
All	All	809/898 (90%)	770 (95%)	39 (5%)	23	34

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	233	SER
1	A	242	ILE
1	A	310	GLU
1	A	329	ARG
1	A	348	ARG
1	A	366	LYS
1	A	379	TRP
1	A	380	GLN
1	A	450	THR
1	A	467	GLU
1	A	481	TRP
1	A	485	THR
1	A	539	LYS
1	A	570	ARG
1	A	583	VAL
1	A	629	ILE
1	A	640	THR
1	A	647	VAL
1	A	656	SER
1	A	692	ILE
1	A	706	ASP
1	B	233	SER
1	B	248	ILE
1	B	345	ARG
1	B	379	TRP
1	B	380	GLN
1	B	459	SER
1	B	465	GLU
1	B	474	GLU
1	B	481	TRP
1	B	485	THR
1	B	528	LYS
1	B	537	ILE
1	B	583	VAL

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Mol	Chain	Res	Type
1	B	629	ILE
1	B	647	VAL
1	B	648	ILE
1	B	656	SER
1	B	706	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	280	GLN
1	A	281	HIS
1	A	380	GLN
1	A	423	ASN
1	A	511	GLN
1	A	532	ASN
1	A	561	HIS
1	A	625	ASN
1	B	271	GLN
1	B	380	GLN
1	B	511	GLN
1	B	532	ASN
1	B	622	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	HFG	A	801	-	26,26,26	0.87	1 (3%)	29,37,37	1.11	1 (3%)
2	HFG	B	802	-	26,26,26	0.95	1 (3%)	29,37,37	1.28	2 (6%)
4	ANP	A	803	-	33,33,33	1.47	4 (12%)	45,52,52	1.04	3 (6%)
4	ANP	B	803	-	33,33,33	2.10	5 (15%)	45,52,52	0.99	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
2	HFG	A	801	-	-	0/8/19/19	0/3/3/3
2	HFG	B	802	-	-	0/8/19/19	0/3/3/3
4	ANP	A	803	-	-	7/18/38/38	0/3/3/3
4	ANP	B	803	-	-	7/18/38/38	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	803	ANP	PG-O1G	6.71	1.56	1.46
4	B	803	ANP	PG-N3B	6.04	1.79	1.63
4	A	803	ANP	PB-O1B	5.15	1.54	1.46
4	A	803	ANP	PB-O2B	-4.87	1.43	1.56
4	B	803	ANP	PB-N3B	4.75	1.75	1.63
4	B	803	ANP	PG-O3G	-3.94	1.46	1.56
4	A	803	ANP	PA-O2A	-2.20	1.45	1.55
2	B	802	HFG	C39-C2'	-2.12	1.50	1.52
4	B	803	ANP	C5'-C4'	2.07	1.57	1.51
4	A	803	ANP	PA-O1A	-2.05	1.43	1.50
2	A	801	HFG	C2-N3	-2.03	1.34	1.36

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	802	HFG	C4'-C39-C2'	-4.43	108.00	110.89
2	A	801	HFG	C4'-C39-C2'	-3.76	108.43	110.89
4	A	803	ANP	O1G-PG-N3B	-3.18	107.09	111.77
4	B	803	ANP	O2G-PG-O1G	-3.14	105.58	113.45
4	A	803	ANP	C5'-C4'-C3'	-2.89	104.83	115.21
4	A	803	ANP	O1B-PB-N3B	2.52	115.48	111.77
2	B	802	HFG	C6'-N1'-C2'	-2.15	110.28	111.62

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	803	ANP	PB-N3B-PG-O1G
4	A	803	ANP	PA-O3A-PB-O2B
4	A	803	ANP	C5'-O5'-PA-O2A
4	B	803	ANP	PB-N3B-PG-O1G
4	B	803	ANP	PG-N3B-PB-O1B
4	B	803	ANP	PG-N3B-PB-O3A
4	B	803	ANP	C5'-O5'-PA-O2A
4	B	803	ANP	C5'-O5'-PA-O3A
4	A	803	ANP	O4'-C4'-C5'-O5'
4	A	803	ANP	C3'-C4'-C5'-O5'
4	A	803	ANP	C5'-O5'-PA-O1A
4	A	803	ANP	C5'-O5'-PA-O3A
4	B	803	ANP	PA-O3A-PB-O2B
4	B	803	ANP	PA-O3A-PB-O1B

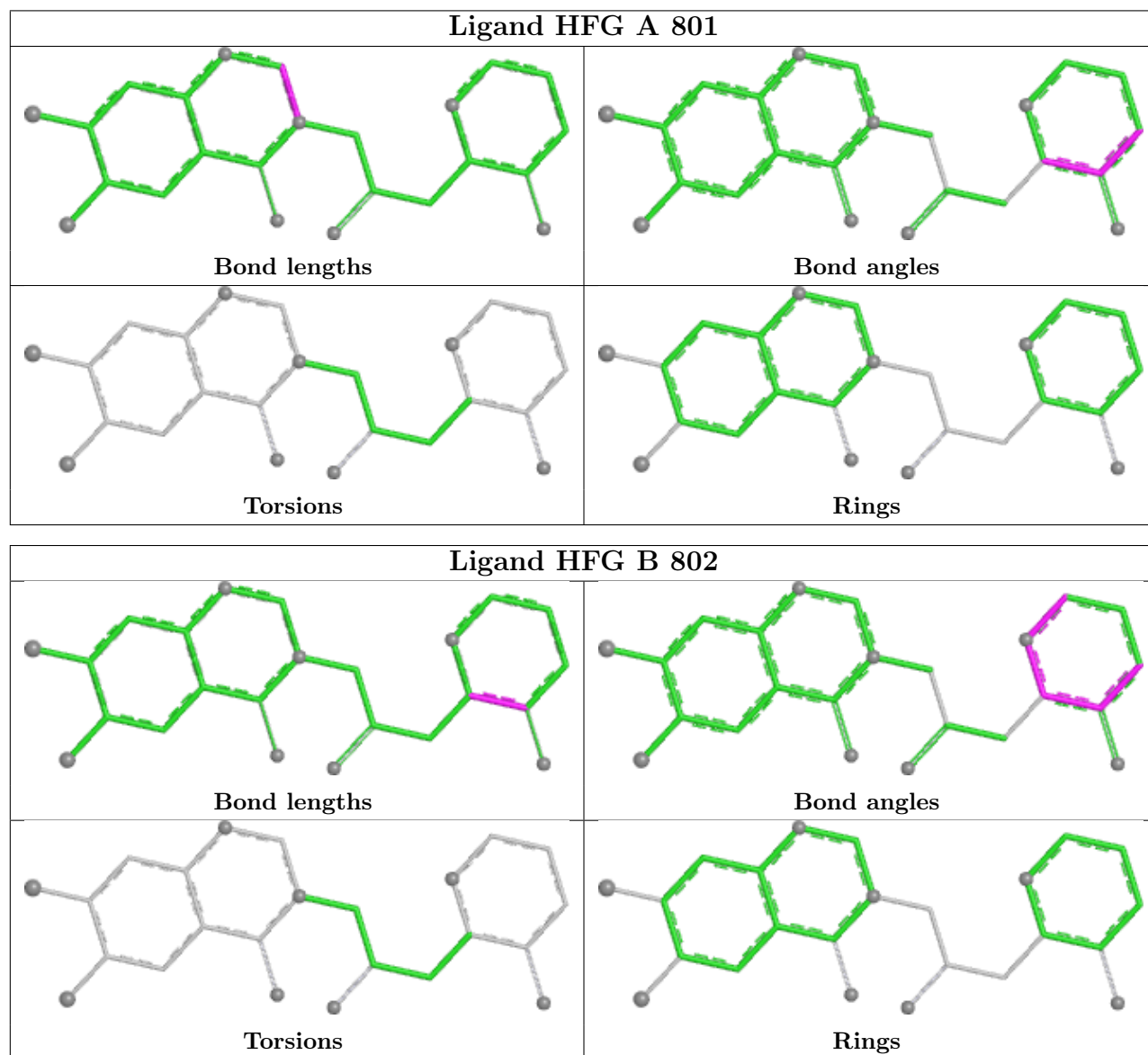
There are no ring outliers.

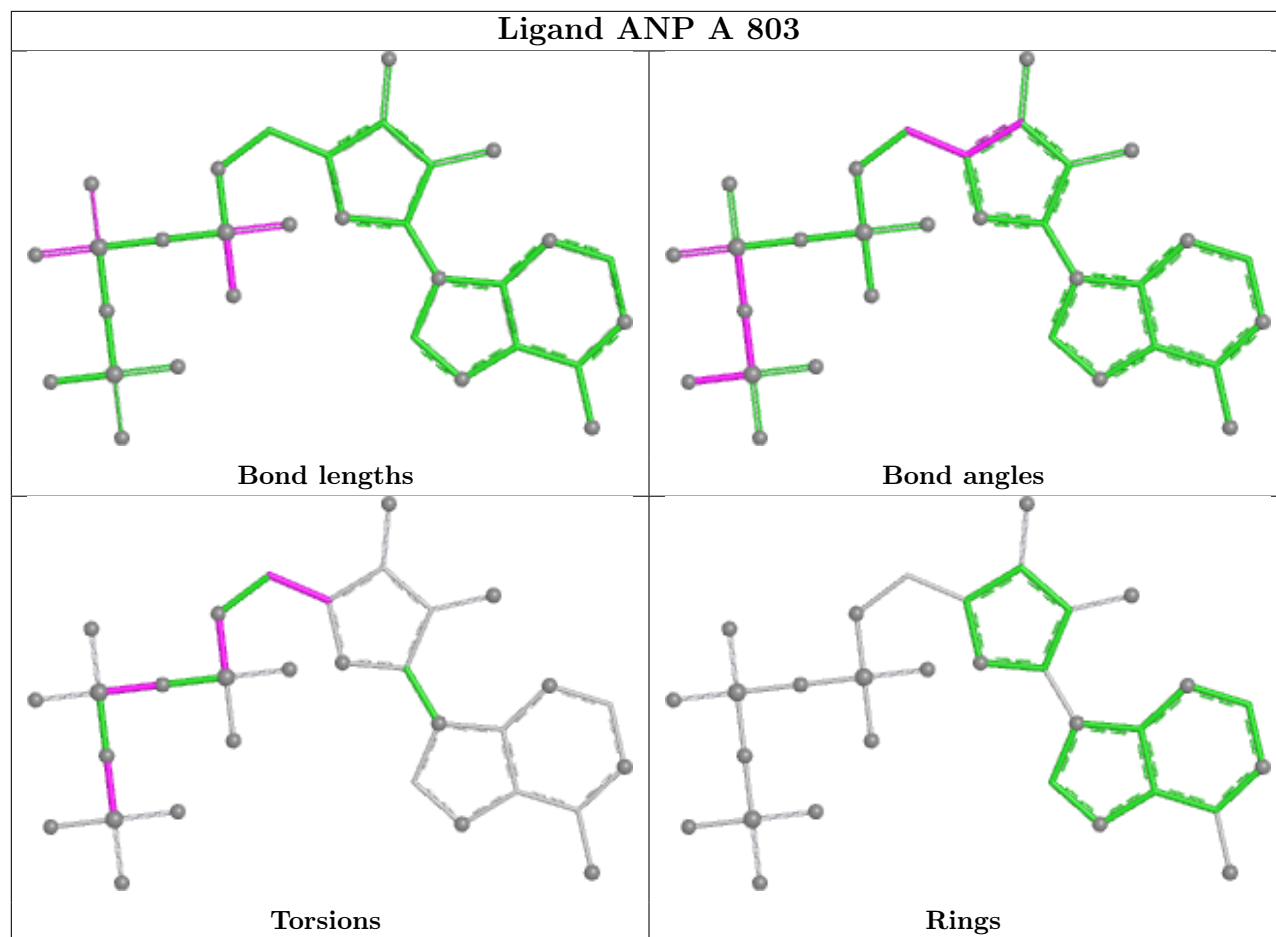
3 monomers are involved in 5 short contacts:

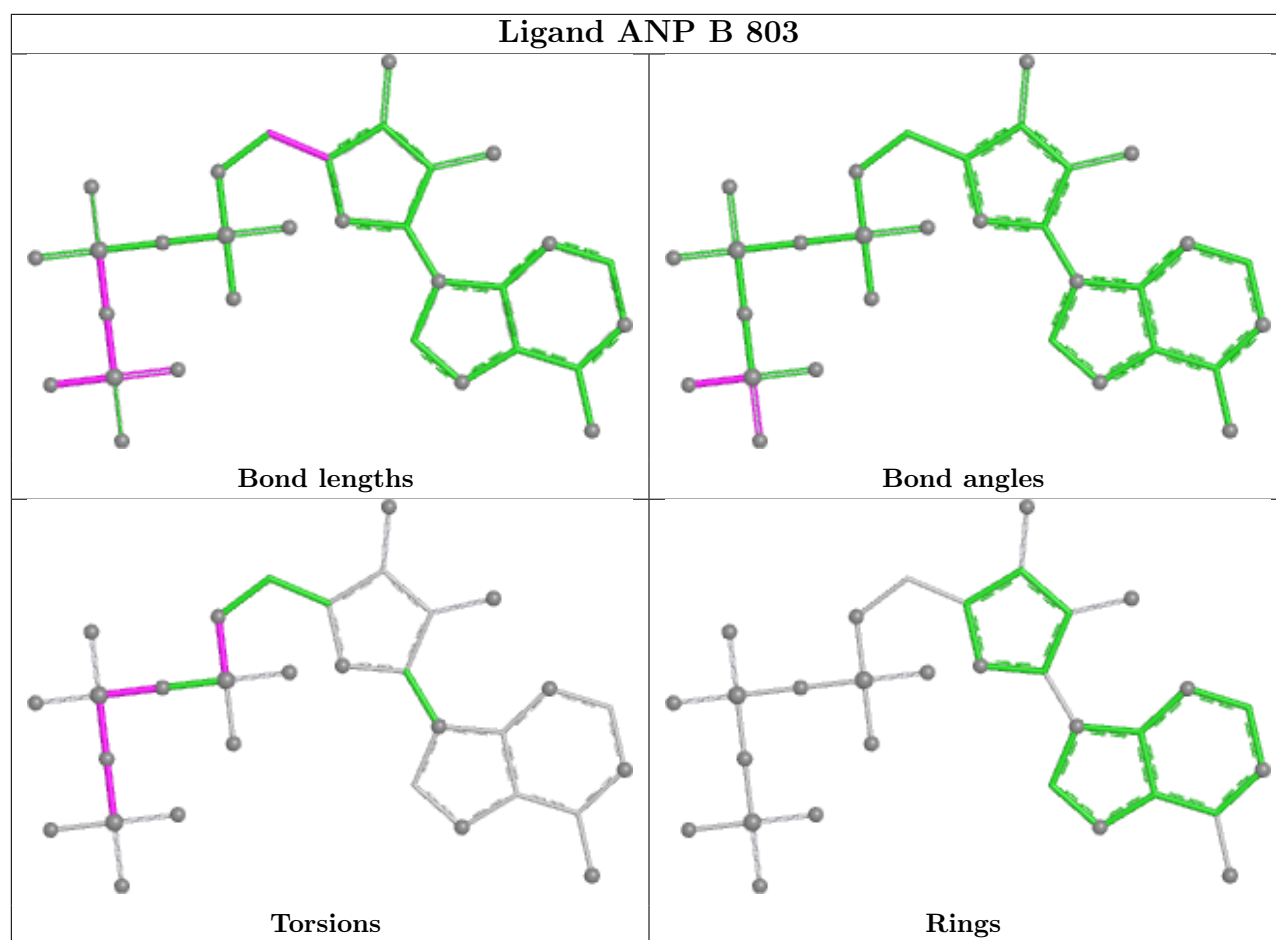
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	HFG	1	0
2	B	802	HFG	3	0
4	B	803	ANP	2	0

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 [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	474/498 (95%)	0.03	14 (2%) 52 53	16, 39, 73, 107	3 (0%)
1	B	471/498 (94%)	-0.01	19 (4%) 42 41	16, 38, 71, 100	2 (0%)
All	All	945/996 (94%)	0.01	33 (3%) 47 47	16, 38, 72, 107	5 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	705	CYS	5.7
1	B	309	PRO	5.2
1	B	705	CYS	4.9
1	A	308	SER	4.1
1	A	300	GLU	4.0
1	B	308	SER	3.9
1	A	307	PHE	3.8
1	B	307	PHE	3.6
1	A	695	ILE	3.5
1	B	695	ILE	3.5
1	A	549	ASP	3.4
1	A	309	PRO	3.2
1	B	698	GLY	3.1
1	B	692	ILE	3.0
1	A	524	THR	2.9
1	B	468	ASP	2.8
1	A	517	ILE	2.8
1	B	549	ASP	2.8
1	B	517	ILE	2.7
1	B	576	ARG	2.6
1	A	526	GLU	2.6
1	A	468	ASP	2.6
1	B	300	GLU	2.5
1	B	528	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	575	PRO	2.5
1	A	692	ILE	2.4
1	B	530	ILE	2.4
1	A	224	MET	2.4
1	B	581[A]	ARG	2.3
1	B	367	HIS	2.3
1	B	579	GLU	2.3
1	A	367	HIS	2.2
1	B	527	GLN	2.1

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](#)

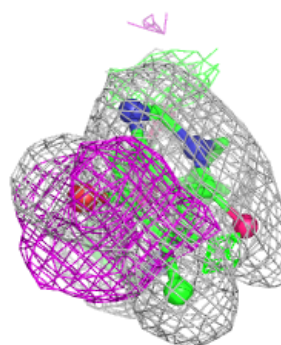
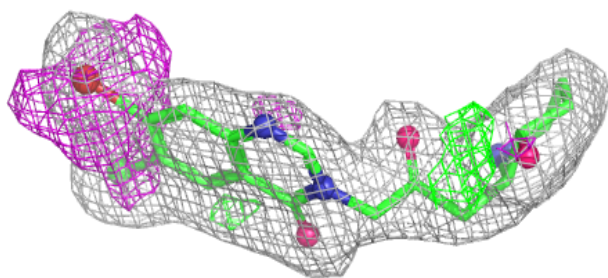
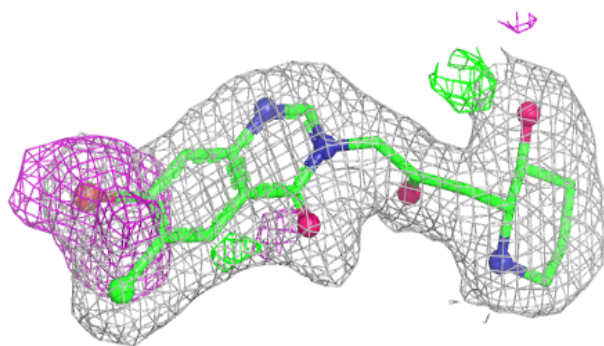
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	HFG	A	801	24/24	0.86	0.10	19,28,44,90	0
2	HFG	B	802	24/24	0.86	0.10	21,28,41,91	0
4	ANP	A	803	31/31	0.90	0.11	26,35,60,72	0
4	ANP	B	803	31/31	0.91	0.11	25,35,71,72	0
3	ZN	A	802	1/1	0.95	0.18	43,43,43,43	1
3	ZN	B	801	1/1	0.97	0.17	43,43,43,43	1

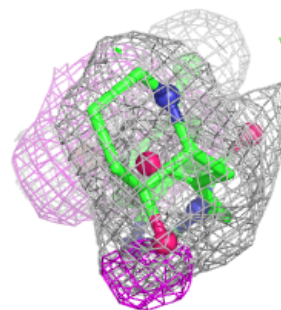
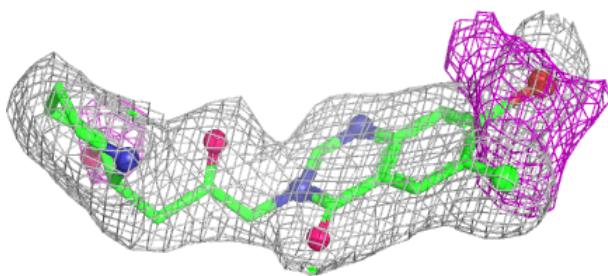
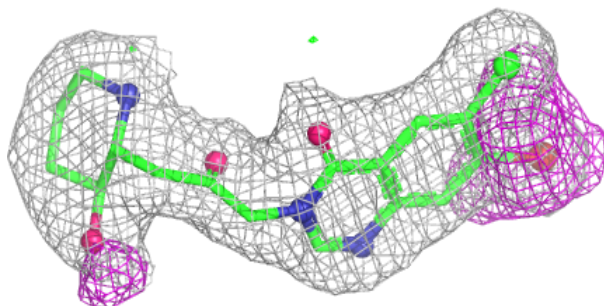
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HFG 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)

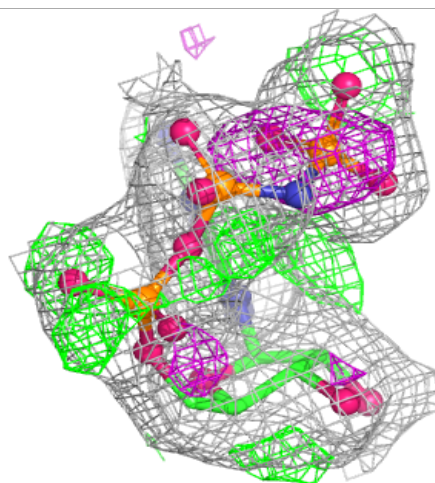
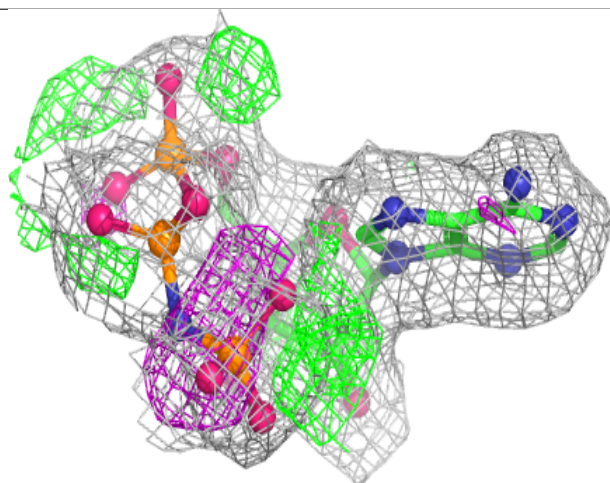
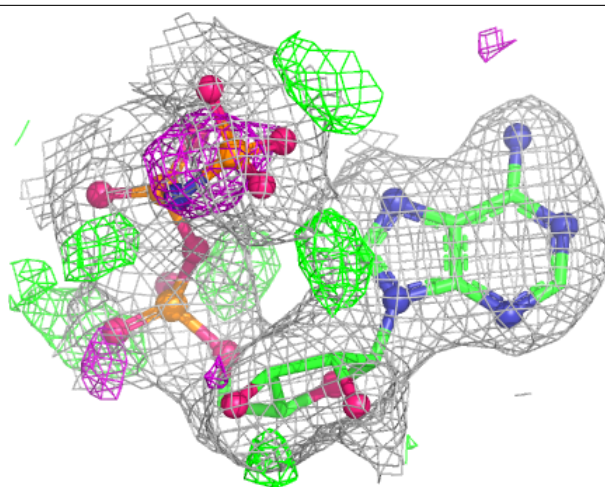
**Electron density around HFG B 802:**

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



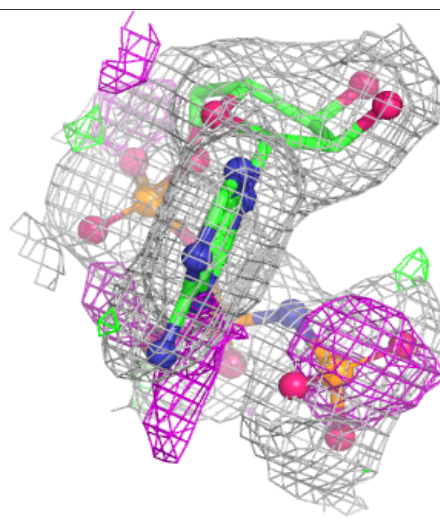
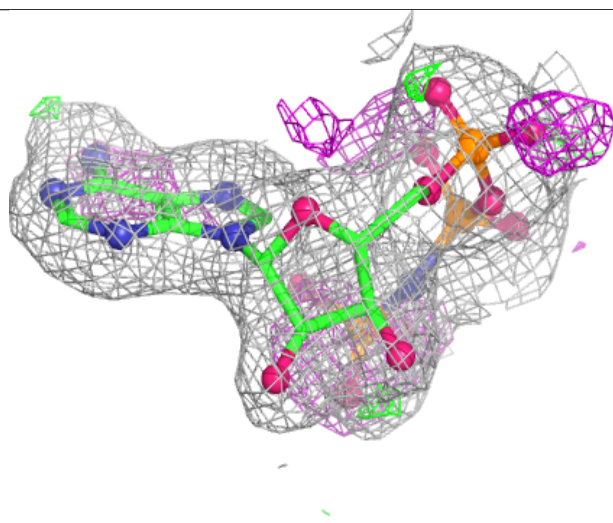
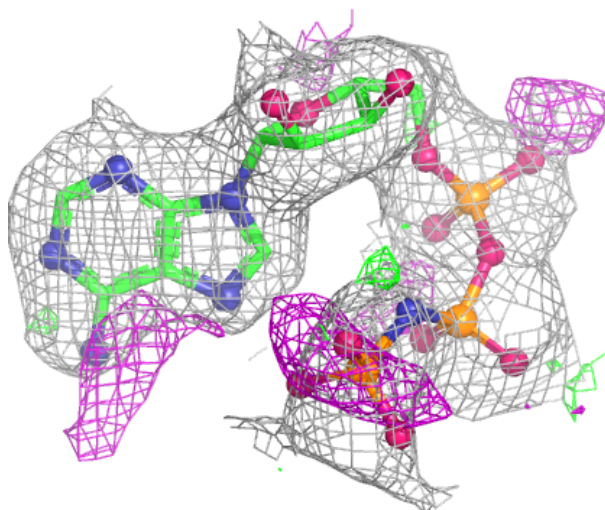
Electron density around ANP A 803:

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



Electron density around ANP 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)



6.5 Other polymers ⓘ

There are no such residues in this entry.