



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 11:13 AM UTC

PDB ID : 8RRJ / pdb_00008rrj
Title : The structural basis of aldo-keto reductase 1C3 inhibition by 17alpha-picolyl and 17(E)-picolinylidene androstane derivatives
Authors : Petri, E.T.; Skerlova, J.; Plavska, J.J.; Brynda, J.; Ajdukovic, J.J.; Bekic, S.; Celic, A.S.; Rezacova, P.
Deposited on : 2024-01-22
Resolution : 1.70 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

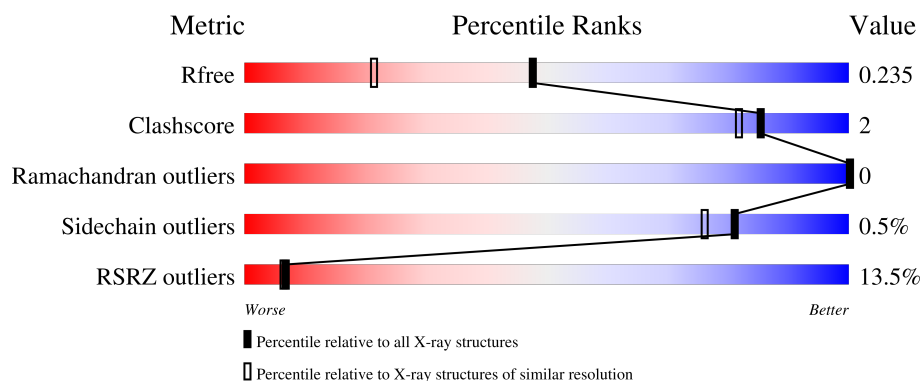
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	326	9% 91% 6%
1	B	326	17% 91% 6%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5905 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.

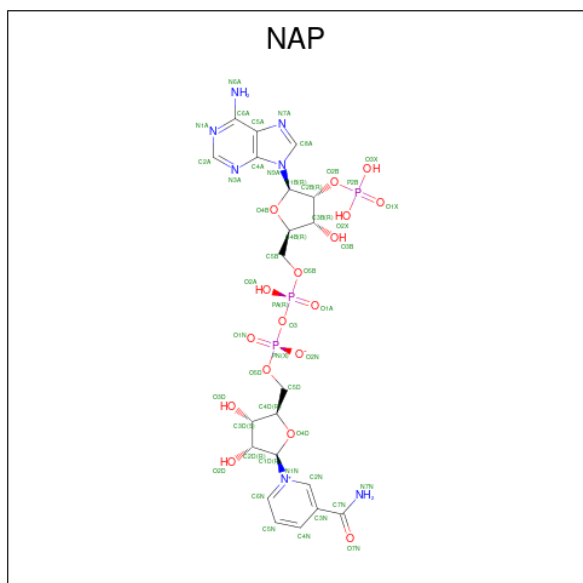
- Molecule 1 is a protein called Aldo-keto reductase family 1 member C3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	5	0
			2578	1643	445	476	14			
1	B	318	Total	C	N	O	S	0	3	0
			2581	1643	450	476	12			

There are 6 discrepancies between the modelled and reference sequences:

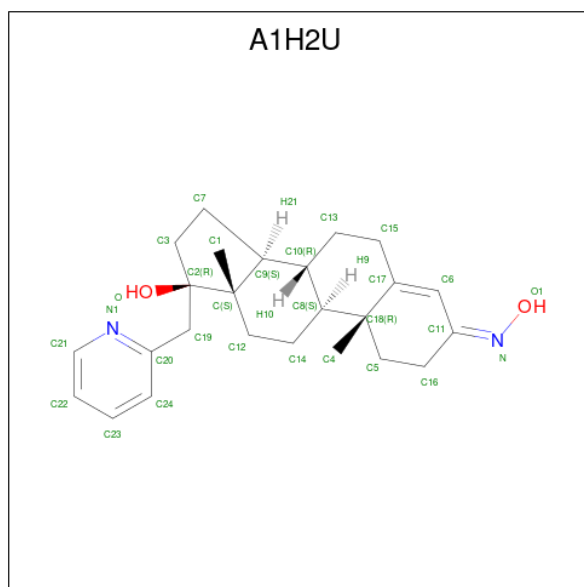
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P42330
A	-1	SER	-	expression tag	UNP P42330
A	0	HIS	-	expression tag	UNP P42330
B	-2	GLY	-	expression tag	UNP P42330
B	-1	SER	-	expression tag	UNP P42330
B	0	HIS	-	expression tag	UNP P42330

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	
			48	21	7	17	3	
2	B	1	Total	C	N	O	P	
			48	21	7	17	3	

- Molecule 3 is (3 {Z},8 {R},9 {S},10 {R},13 {S},14 {S},17 {R})-3-hydroxyimino-10,13-dimethyl-17-(pyridin-2-ylmethyl)-2,6,7,8,9,11,12,14,15,16-decahydro-1 {H}-cyclopenta[a]phenanthren-17-ol (CCD ID: A1H2U) (formula: $C_{25}H_{34}N_2O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O		
			29	25	2	2		
3	B	1	Total	C	N	O		
			29	25	2	2		

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	5	Total	Na		
			5	5	0	0
4	B	6	Total	Na		
			6	6	0	0

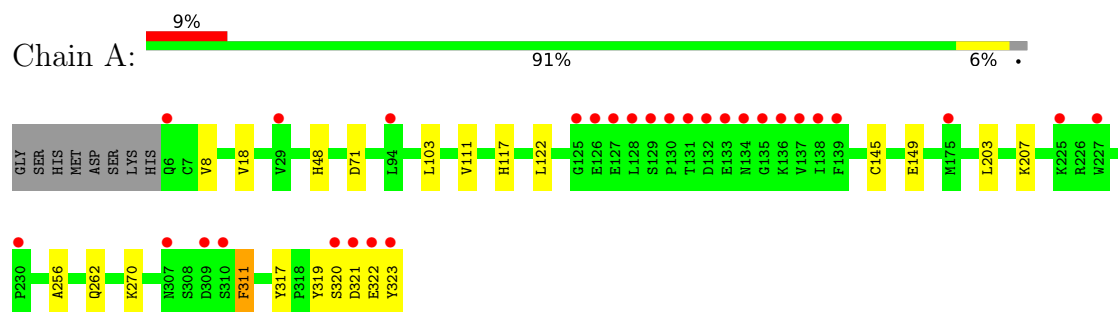
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	328	Total 328	O 328	0	1
5	B	253	Total 253	O 253	0	1

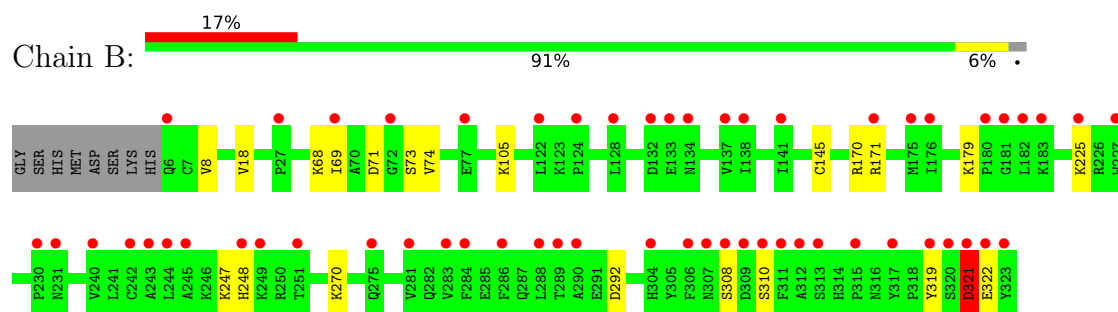
3 Residue-property plots [i](#)

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: Aldo-keto reductase family 1 member C3



- Molecule 1: Aldo-keto reductase family 1 member C3



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	41.17Å 53.52Å 75.62Å 77.71° 85.61° 76.06°	Depositor
Resolution (Å)	46.49 – 1.70 46.49 – 1.70	Depositor EDS
% Data completeness (in resolution range)	83.2 (46.49-1.70) 83.1 (46.49-1.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.186 , 0.221 0.205 , 0.235	Depositor DCC
R_{free} test set	1120 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5905	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAP, NA, A1H2U

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	1.08	2/2649 (0.1%)	1.32	6/3584 (0.2%)
1	B	1.05	0/2640	1.32	5/3571 (0.1%)
All	All	1.07	2/5289 (0.0%)	1.32	11/7155 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	HIS	C-O	5.24	1.30	1.24
1	A	256	ALA	C-O	5.06	1.30	1.24

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	319	TYR	CA-C-N	6.89	130.20	120.28
1	A	319	TYR	C-N-CA	6.89	130.20	120.28
1	A	71	ASP	CA-CB-CG	6.35	118.95	112.60
1	B	71	ASP	CA-CB-CG	6.04	118.64	112.60
1	A	311	PHE	CA-CB-CG	-5.82	107.98	113.80
1	B	292	ASP	CA-CB-CG	5.61	118.21	112.60
1	A	262	GLN	N-CA-C	-5.55	106.46	113.18
1	A	48	HIS	CA-CB-CG	5.23	119.03	113.80
1	B	321	ASP	CA-CB-CG	5.20	117.80	112.60
1	B	105	LYS	CB-CA-C	-5.18	102.75	110.88
1	B	225	LYS	CB-CA-C	-5.16	102.23	110.79

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	2578	0	2579	11	0
1	B	2581	0	2576	13	0
2	A	48	0	25	2	0
2	B	48	0	25	2	0
3	A	29	0	0	0	0
3	B	29	0	0	0	0
4	A	5	0	0	0	0
4	B	6	0	0	0	0
5	A	328	0	0	0	0
5	B	253	0	0	0	0
All	All	5905	0	5205	23	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 2.

All (23) 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:321:ASP:HB2	1:B:321:ASP:HB3	1.71	0.73
1:A:8:VAL:HG13	1:A:18:VAL:HG12	1.78	0.66
1:B:319:TYR:O	1:B:322:GLU:HG2	1.96	0.66
1:A:321:ASP:HB2	1:B:321:ASP:CB	2.28	0.63
1:B:8:VAL:HG13	1:B:18:VAL:CG1	2.32	0.59
1:B:8:VAL:HG13	1:B:18:VAL:HG12	1.84	0.59
1:A:8:VAL:HG13	1:A:18:VAL:CG1	2.33	0.59
1:A:145:CYS:O	1:A:149:GLU:HG3	2.06	0.56
1:B:308:SER:OG	1:B:310:SER:HB2	2.06	0.55
1:B:270:LYS:O	2:B:401:NAP:H8A	2.10	0.52
1:B:145:CYS:HB3	1:B:179:LYS:HD2	1.92	0.52
1:A:323:TYR:OH	1:B:170:ARG:NH2	2.44	0.50
1:A:270:LYS:O	2:A:401:NAP:H8A	2.13	0.49
2:A:401:NAP:H6N	2:A:401:NAP:H3D	1.97	0.47
1:A:122:LEU:HD21	1:A:311:PHE:HE1	1.79	0.46
1:A:103:LEU:HD21	1:A:111:VAL:HG13	1.97	0.46
1:B:68:LYS:HE3	1:B:73:SER:CB	2.45	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:LYS:HG2	1:B:248:HIS:CD2	2.52	0.45
1:B:68:LYS:HE3	1:B:73:SER:HB2	1.99	0.44
1:A:317:TYR:HB3	1:A:320:SER:HB2	1.99	0.44
1:A:203:LEU:HD11	1:A:207:LYS:HE3	2.02	0.41
2:B:401:NAP:H6N	2:B:401:NAP:H3D	2.02	0.41
1:B:69:ILE:HA	1:B:74:VAL:O	2.20	0.41

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	321/326 (98%)	315 (98%)	6 (2%)	0	100	100
1	B	319/326 (98%)	308 (97%)	11 (3%)	0	100	100
All	All	640/652 (98%)	623 (97%)	17 (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	287/289 (99%)	286 (100%)	1 (0%)	86	83
1	B	285/289 (99%)	282 (99%)	3 (1%)	65	54

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	572/578 (99%)	568 (99%)	4 (1%)	81	69

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	322	GLU
1	B	171[A]	ARG
1	B	171[B]	ARG
1	B	321	ASP

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	167	ASN
1	A	275	GLN
1	B	6	GLN
1	B	282	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 15 ligands modelled in this entry, 11 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	NAP	B	401	-	50,52,52	1.32	8 (16%)	71,80,80	1.50	12 (16%)
3	A1H2U	A	402	-	32,33,33	1.59	8 (25%)	48,52,52	1.70	15 (31%)
2	NAP	A	401	-	50,52,52	1.22	5 (10%)	71,80,80	1.46	10 (14%)
3	A1H2U	B	402	-	32,33,33	1.58	7 (21%)	48,52,52	2.01	14 (29%)

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	NAP	B	401	-	-	4/35/67/67	0/5/5/5
3	A1H2U	A	402	-	-	1/7/69/69	0/5/5/5
2	NAP	A	401	-	-	4/35/67/67	0/5/5/5
3	A1H2U	B	402	-	-	2/7/69/69	0/5/5/5

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAP	C5A-C4A	4.15	1.46	1.39
2	A	401	NAP	C5A-C4A	4.05	1.46	1.39
3	B	402	A1H2U	C20-N1	3.82	1.42	1.34
3	A	402	A1H2U	C24-C20	-3.62	1.31	1.38
3	B	402	A1H2U	C24-C20	-3.54	1.31	1.38
3	A	402	A1H2U	C20-N1	3.44	1.41	1.34
3	B	402	A1H2U	C2-C	3.18	1.59	1.56
3	A	402	A1H2U	C14-C8	2.93	1.58	1.53
3	A	402	A1H2U	C21-N1	2.91	1.40	1.34
2	A	401	NAP	C5A-N7A	-2.86	1.33	1.39
2	B	401	NAP	C5A-N7A	-2.86	1.33	1.39
2	A	401	NAP	C4A-N9A	-2.77	1.31	1.37
2	B	401	NAP	PA-O3	2.72	1.62	1.59
2	B	401	NAP	C4A-N9A	-2.71	1.32	1.37
3	A	402	A1H2U	C2-C	2.65	1.59	1.56
3	B	402	A1H2U	C23-C24	-2.60	1.34	1.38
3	B	402	A1H2U	C21-N1	2.56	1.40	1.34
3	A	402	A1H2U	C23-C24	-2.46	1.34	1.38
2	B	401	NAP	PN-O3	2.43	1.62	1.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAP	C8A-N9A	-2.39	1.33	1.37
2	A	401	NAP	C5A-C6A	2.35	1.47	1.41
3	A	402	A1H2U	C6-C11	2.31	1.49	1.44
2	B	401	NAP	C5A-C6A	2.30	1.47	1.41
2	A	401	NAP	C8A-N9A	-2.28	1.33	1.37
3	A	402	A1H2U	C19-C2	2.22	1.57	1.54
3	B	402	A1H2U	C16-C11	-2.16	1.47	1.50
3	B	402	A1H2U	C14-C8	2.14	1.57	1.53
2	B	401	NAP	O7N-C7N	-2.13	1.20	1.24

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	NAP	C5A-C4A-N3A	-5.36	119.34	126.72
2	B	401	NAP	C5A-C4A-N3A	-4.98	119.85	126.72
3	B	402	A1H2U	C4-C18-C8	4.88	117.14	111.66
2	A	401	NAP	N3A-C4A-N9A	4.57	134.93	127.17
2	B	401	NAP	N3A-C4A-N9A	4.31	134.49	127.17
3	B	402	A1H2U	C19-C20-C24	-4.27	115.58	121.73
3	B	402	A1H2U	C19-C20-N1	4.24	122.79	117.08
3	B	402	A1H2U	C16-C11-C6	-4.19	113.29	117.59
2	B	401	NAP	N3A-C2A-N1A	-3.63	123.09	128.58
2	A	401	NAP	C2A-N3A-C4A	3.47	120.30	111.83
3	B	402	A1H2U	C3-C2-C	3.40	105.29	103.06
2	B	401	NAP	C4A-C5A-N7A	-3.40	106.70	110.58
3	A	402	A1H2U	C22-C21-N1	-3.39	118.05	123.42
2	A	401	NAP	C4A-C5A-N7A	-3.38	106.72	110.58
2	B	401	NAP	C2A-N3A-C4A	3.35	120.02	111.83
2	B	401	NAP	C4A-N9A-C8A	3.33	109.24	105.74
3	A	402	A1H2U	C19-C20-N1	3.33	121.57	117.08
3	B	402	A1H2U	C8-C18-C17	-3.32	104.80	109.65
2	A	401	NAP	N3A-C2A-N1A	-3.31	123.57	128.58
2	A	401	NAP	C4A-N9A-C8A	3.28	109.18	105.74
3	A	402	A1H2U	C8-C18-C17	-3.25	104.89	109.65
3	A	402	A1H2U	C22-C23-C24	3.14	124.12	120.24
3	A	402	A1H2U	O1-N-C11	3.13	117.81	112.15
3	B	402	A1H2U	C22-C23-C24	3.01	123.96	120.24
3	B	402	A1H2U	C22-C21-N1	-3.01	118.65	123.42
3	A	402	A1H2U	C19-C20-C24	-2.98	117.44	121.73
3	B	402	A1H2U	C2-C-C9	-2.82	99.02	101.43
3	B	402	A1H2U	C1-C-C12	-2.78	105.89	109.69
3	B	402	A1H2U	C12-C-C2	2.74	122.41	116.97

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	A1H2U	C5-C16-C11	2.62	113.97	110.97
3	A	402	A1H2U	C16-C11-C6	-2.61	114.91	117.59
3	A	402	A1H2U	C1-C-C12	-2.58	106.16	109.69
2	B	401	NAP	C2A-N1A-C6A	2.53	122.89	118.73
2	B	401	NAP	C5A-N7A-C8A	2.53	107.42	103.45
2	A	401	NAP	C5A-N7A-C8A	2.52	107.42	103.45
3	A	402	A1H2U	C12-C-C2	2.42	121.77	116.97
2	B	401	NAP	C6A-C5A-N7A	2.41	136.74	132.09
2	B	401	NAP	C6N-N1N-C2N	-2.33	119.90	121.88
3	B	402	A1H2U	O1-N-C11	2.31	116.33	112.15
2	A	401	NAP	C2A-N1A-C6A	2.23	122.39	118.73
2	A	401	NAP	C6A-C5A-N7A	2.22	136.36	132.09
3	A	402	A1H2U	C12-C14-C8	2.21	116.89	113.14
2	A	401	NAP	N9A-C8A-N7A	-2.10	110.95	113.94
3	A	402	A1H2U	C5-C18-C8	-2.09	105.97	108.74
3	B	402	A1H2U	C9-C10-C8	2.08	111.80	109.09
2	B	401	NAP	N9A-C8A-N7A	-2.08	110.99	113.94
3	A	402	A1H2U	C21-N1-C20	2.06	120.31	117.41
3	B	402	A1H2U	C13-C10-C9	-2.06	108.66	112.08
3	A	402	A1H2U	C4-C18-C17	2.05	111.51	108.38
2	B	401	NAP	O3X-P2B-O2X	2.03	115.42	107.80
3	A	402	A1H2U	C4-C18-C8	2.00	113.91	111.66

There are no chirality outliers.

All (11) torsion outliers are listed below:

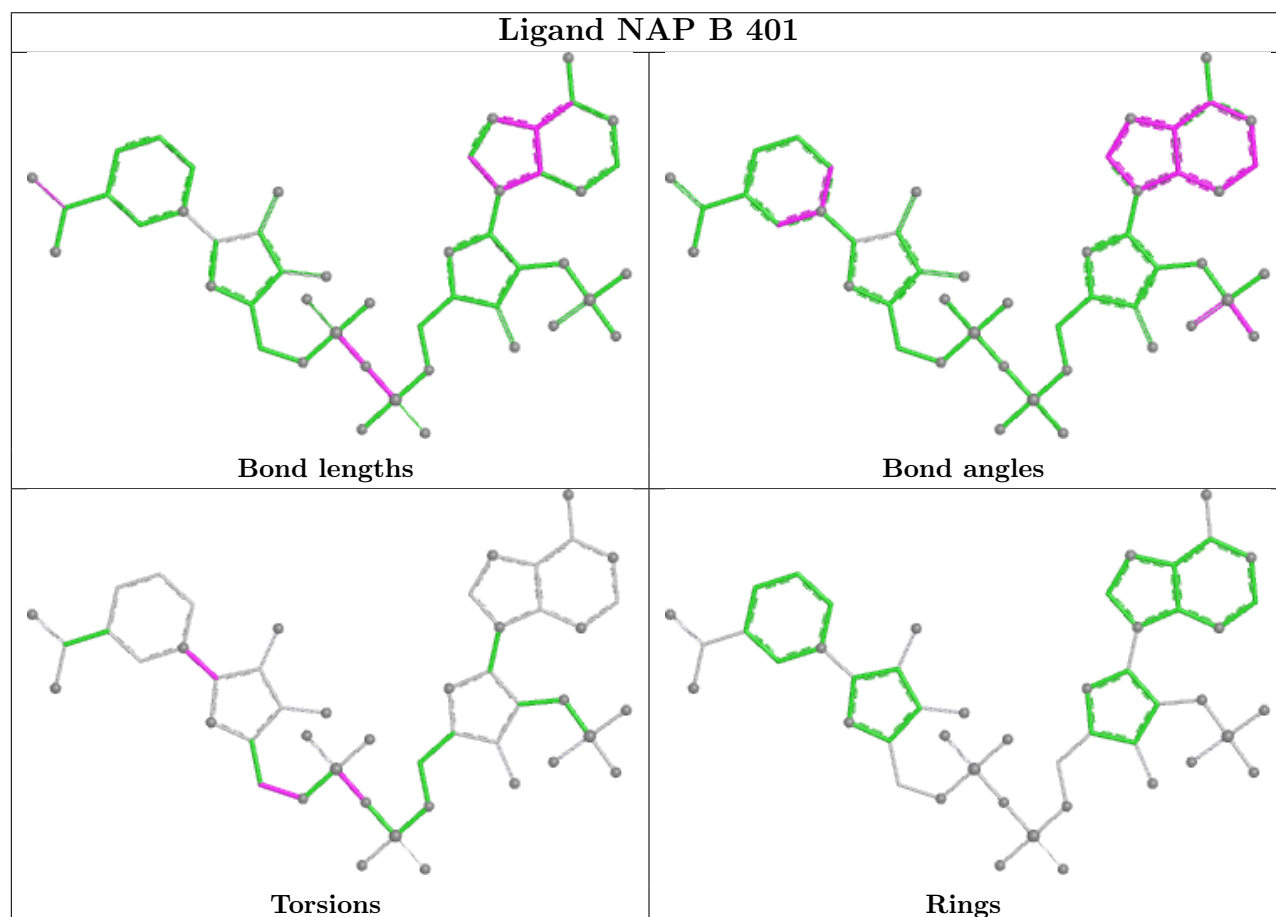
Mol	Chain	Res	Type	Atoms
2	B	401	NAP	O4D-C1D-N1N-C6N
3	B	402	A1H2U	C20-C19-C2-C
2	A	401	NAP	PA-O3-PN-O5D
2	B	401	NAP	PA-O3-PN-O5D
2	A	401	NAP	C4D-C5D-O5D-PN
2	B	401	NAP	C4D-C5D-O5D-PN
2	A	401	NAP	O4D-C1D-N1N-C2N
2	A	401	NAP	O4D-C1D-N1N-C6N
2	B	401	NAP	O4D-C1D-N1N-C2N
3	A	402	A1H2U	C16-C11-N-O1
3	B	402	A1H2U	C16-C11-N-O1

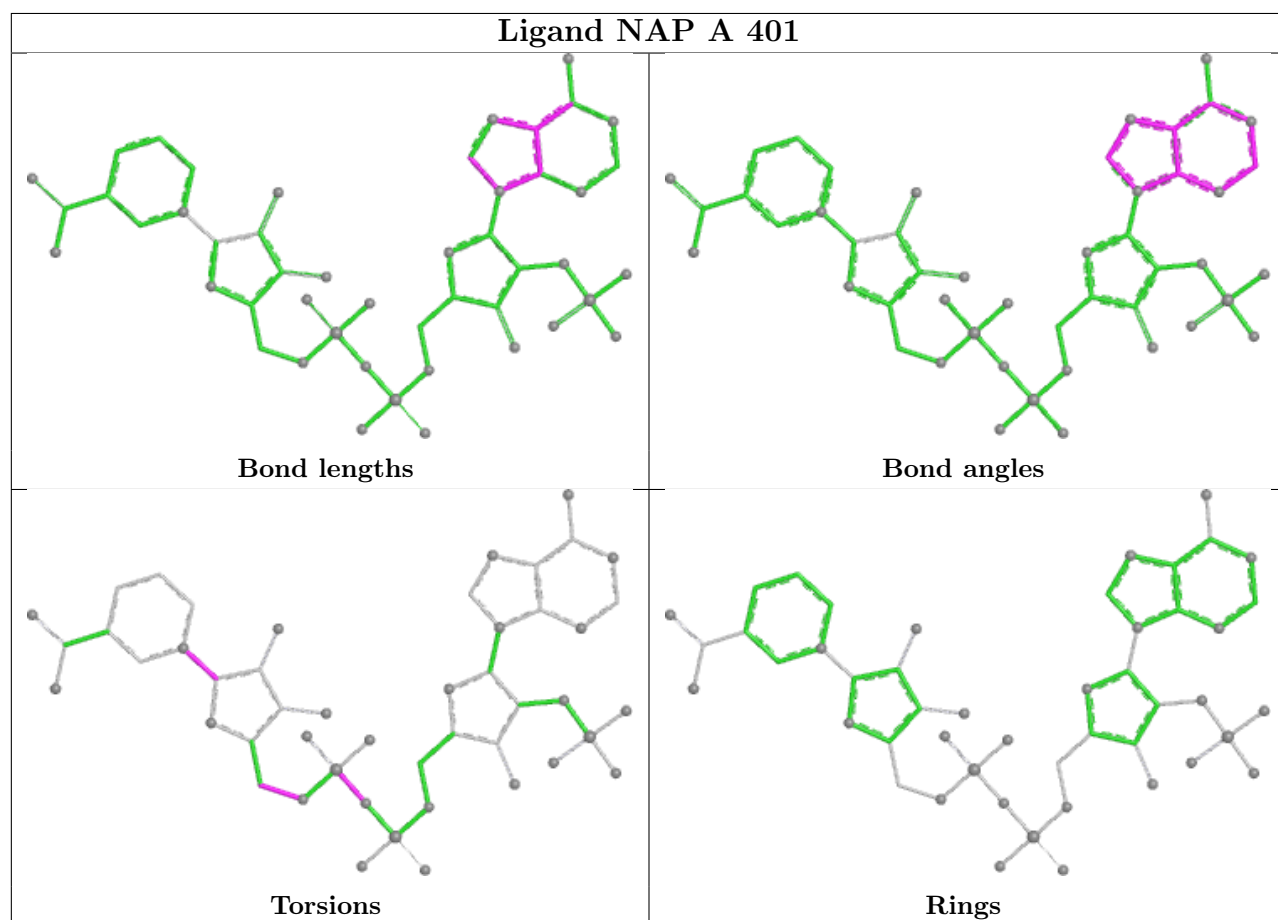
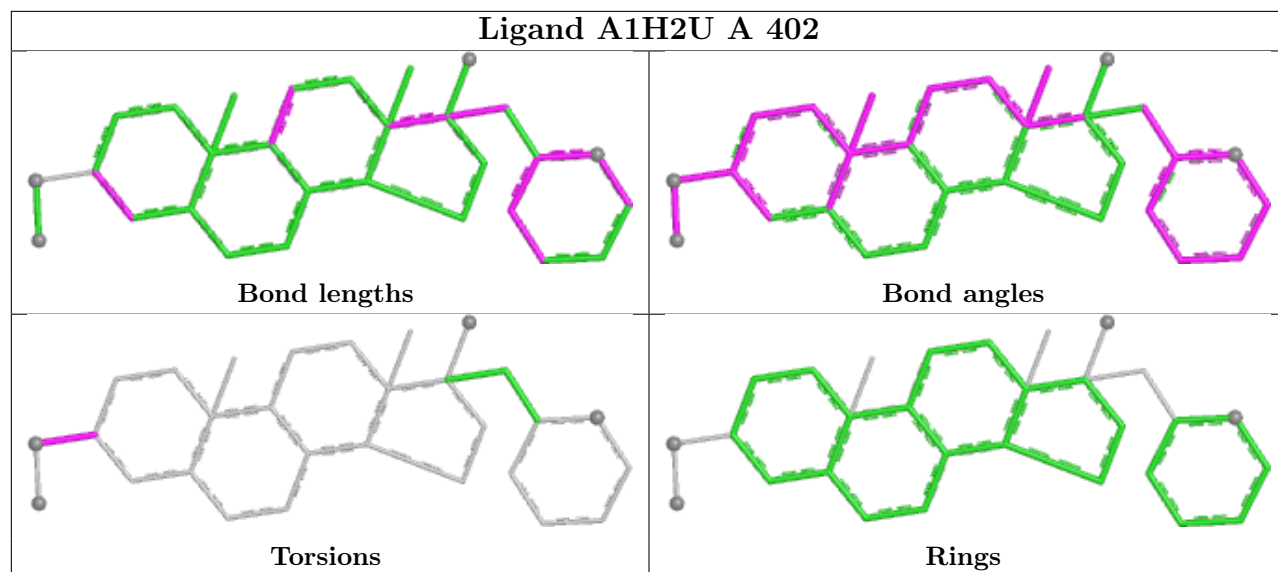
There are no ring outliers.

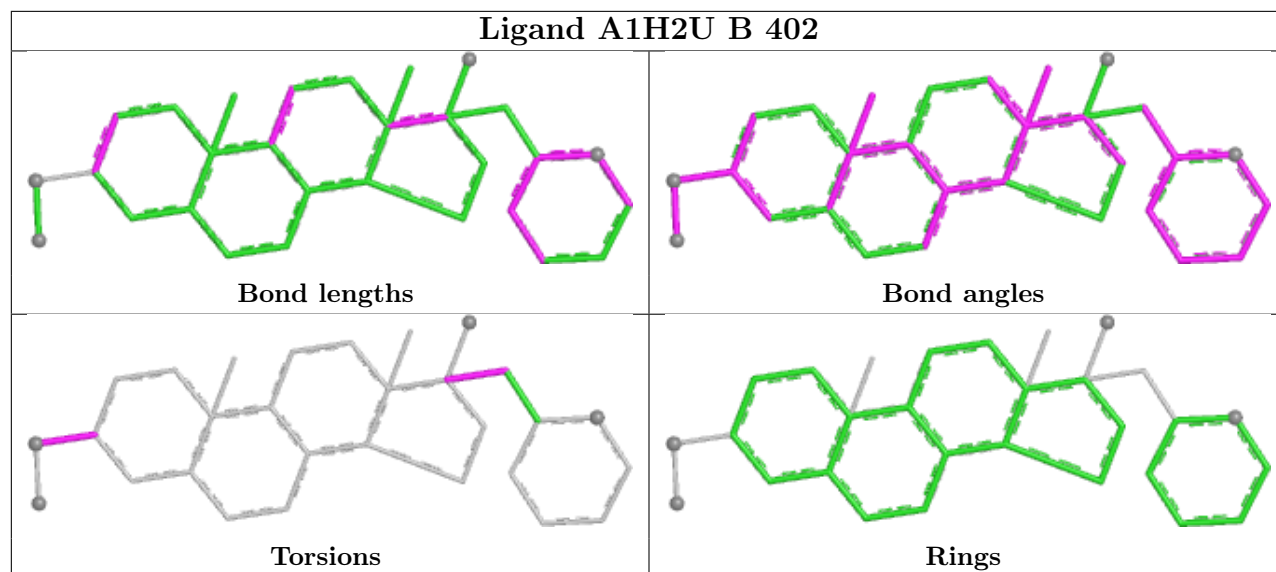
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	NAP	2	0
2	A	401	NAP	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	318/326 (97%)	0.64	29 (9%) 15 15	10, 30, 62, 98	5 (1%)
1	B	318/326 (97%)	1.07	57 (17%) 3 3	12, 37, 62, 95	3 (0%)
All	All	636/652 (97%)	0.85	86 (13%) 7 6	10, 33, 62, 98	8 (1%)

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	323	TYR	6.6
1	A	323	TYR	6.3
1	A	125	GLY	5.0
1	A	128	LEU	4.9
1	A	130	PRO	4.7
1	A	135	GLY	4.3
1	B	307[A]	ASN	4.3
1	B	306	PHE	4.2
1	A	137	VAL	4.1
1	A	322	GLU	4.0
1	B	290	ALA	3.8
1	B	309	ASP	3.6
1	A	132	ASP	3.5
1	B	308	SER	3.5
1	B	311	PHE	3.4
1	B	312	ALA	3.4
1	A	133	GLU	3.4
1	B	249	LYS	3.3
1	A	309	ASP	3.3
1	A	134	ASN	3.3
1	B	243	ALA	3.3
1	B	6	GLN	3.1
1	B	310	SER	3.1
1	A	320	SER	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	313	SER	3.0
1	A	127	GLU	3.0
1	A	131	THR	2.9
1	A	310	SER	2.9
1	B	227	TRP	2.9
1	A	94	LEU	2.8
1	B	320	SER	2.8
1	B	245	ALA	2.8
1	B	242	CYS	2.8
1	A	138	ILE	2.8
1	A	307[A]	ASN	2.8
1	B	134	ASN	2.8
1	B	137	VAL	2.8
1	B	171[A]	ARG	2.7
1	B	322	GLU	2.7
1	B	304	HIS	2.7
1	B	317	TYR	2.7
1	B	275	GLN	2.7
1	B	180	PRO	2.6
1	A	129	SER	2.6
1	B	284	PHE	2.6
1	A	136	LYS	2.6
1	B	321	ASP	2.5
1	B	128	LEU	2.5
1	B	244	LEU	2.4
1	B	315	PRO	2.4
1	A	321	ASP	2.4
1	B	231	ASN	2.4
1	B	124	PRO	2.4
1	B	77	GLU	2.4
1	A	230	PRO	2.3
1	B	248	HIS	2.3
1	B	283	VAL	2.3
1	B	225	LYS	2.3
1	A	175	MET	2.3
1	B	72	GLY	2.3
1	B	141	ILE	2.3
1	B	175	MET	2.3
1	B	288	LEU	2.2
1	B	138	ILE	2.2
1	B	281	VAL	2.2
1	B	176	ILE	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	286	PHE	2.2
1	B	133	GLU	2.1
1	A	225	LYS	2.1
1	B	183	LYS	2.1
1	B	230	PRO	2.1
1	A	29	VAL	2.1
1	B	240	VAL	2.1
1	B	132	ASP	2.1
1	B	122	LEU	2.1
1	B	27	PRO	2.1
1	B	182	LEU	2.0
1	B	251	THR	2.0
1	B	289	THR	2.0
1	A	126	GLU	2.0
1	A	6	GLN	2.0
1	B	181	GLY	2.0
1	A	227	TRP	2.0
1	A	139	PHE	2.0
1	B	69	ILE	2.0
1	B	319	TYR	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
4	NA	A	406	1/1	0.73	0.16	54,54,54,54	0
4	NA	B	404	1/1	0.74	0.27	43,43,43,43	0
4	NA	B	406	1/1	0.78	0.19	46,46,46,46	0

Continued on next page...

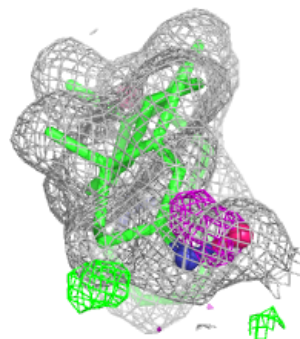
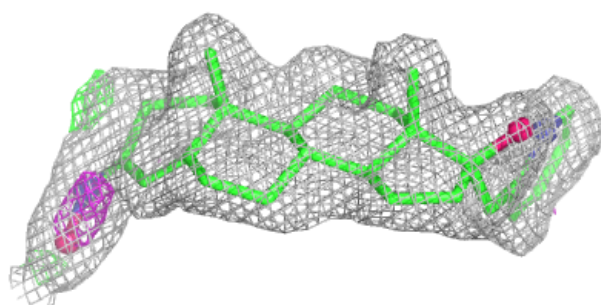
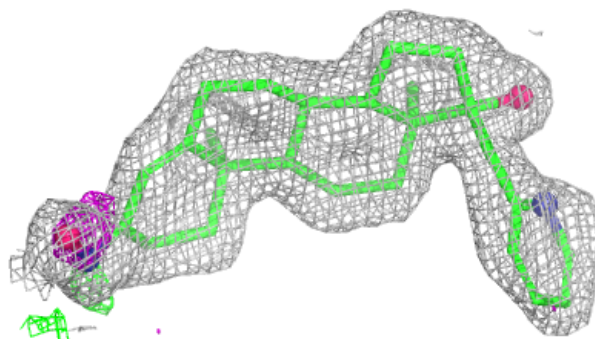
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NA	B	407	1/1	0.79	0.18	49,49,49,49	0
4	NA	A	407	1/1	0.80	0.17	49,49,49,49	0
4	NA	B	408	1/1	0.83	0.12	53,53,53,53	0
4	NA	A	405	1/1	0.84	0.15	52,52,52,52	0
4	NA	B	405	1/1	0.88	0.12	51,51,51,51	0
4	NA	A	404	1/1	0.89	0.11	48,48,48,48	0
3	A1H2U	B	402	29/29	0.91	0.09	20,24,36,38	0
3	A1H2U	A	402	29/29	0.91	0.08	20,22,25,26	0
4	NA	A	403	1/1	0.93	0.25	31,31,31,31	0
2	NAP	A	401	48/48	0.96	0.07	12,15,16,17	0
2	NAP	B	401	48/48	0.96	0.07	13,17,21,22	0
4	NA	B	403	1/1	0.97	0.16	26,26,26,26	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.

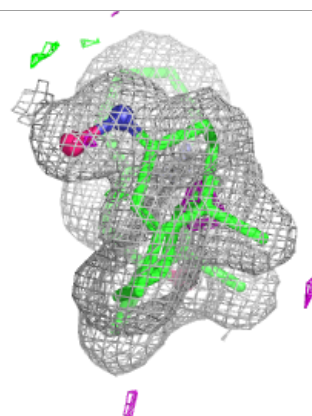
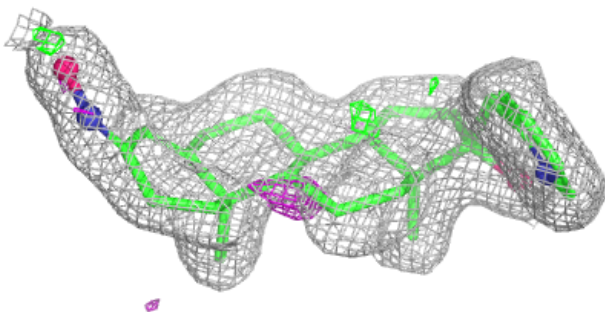
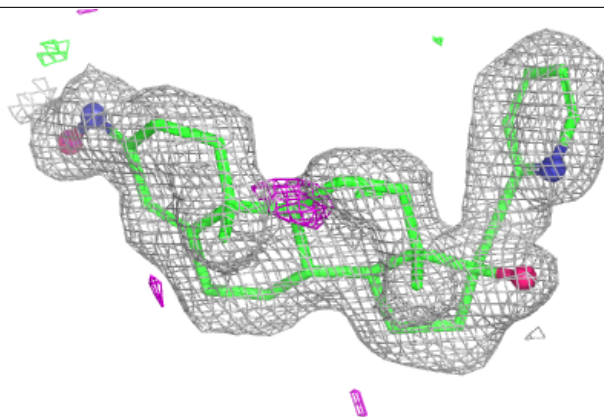
Electron density around A1H2U B 402:

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

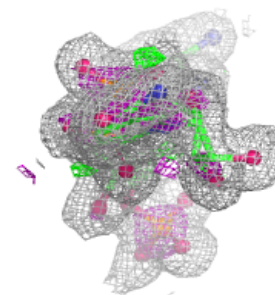
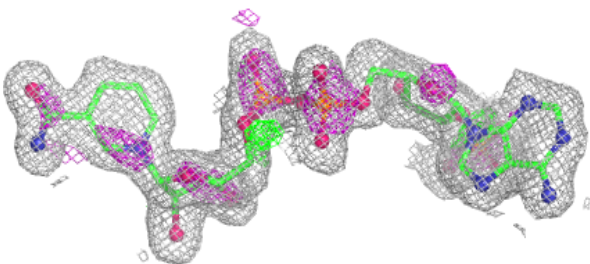
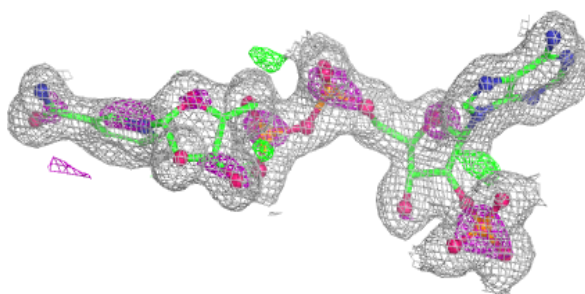


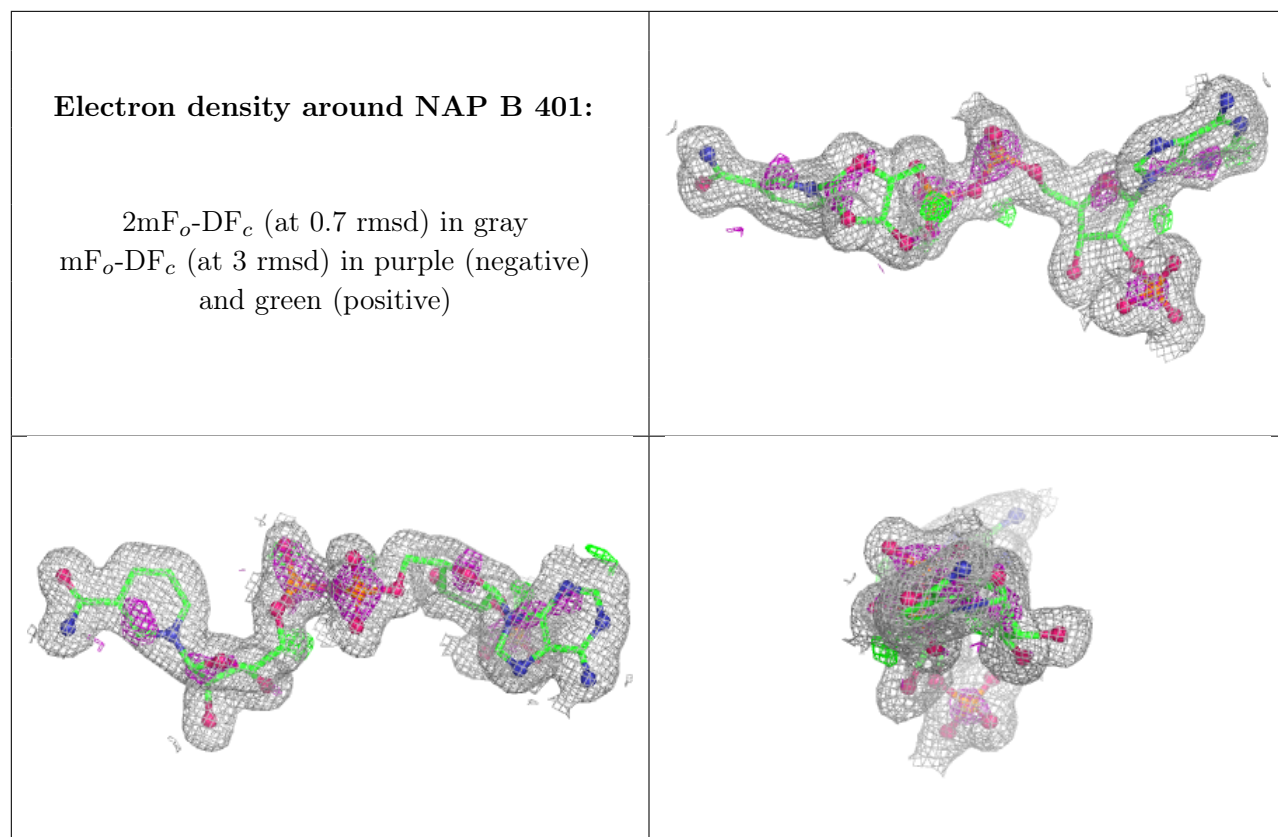
Electron density around A1H2U A 402:

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

**Electron density around NAP A 401:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.