



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:44 PM UTC

PDB ID : 4HFR / pdb_00004hfr
Title : Human 11beta-Hydroxysteroid Dehydrogenase Type 1 in complex with an orally bioavailable acidic inhibitor AZD4017.
Authors : Ogg, D.J.; Gerhardt, S.; Hargreaves, D.
Deposited on : 2012-10-05
Resolution : 2.73 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

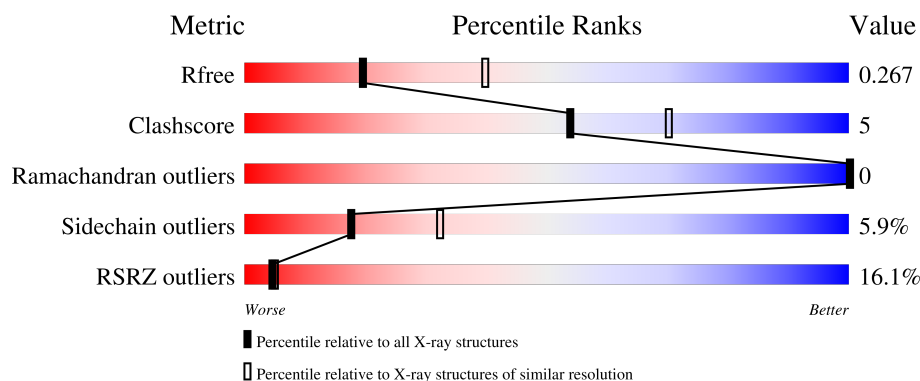
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	272	19% 77% 14% • 6%
1	B	272	11% 78% 12% • 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4053 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.

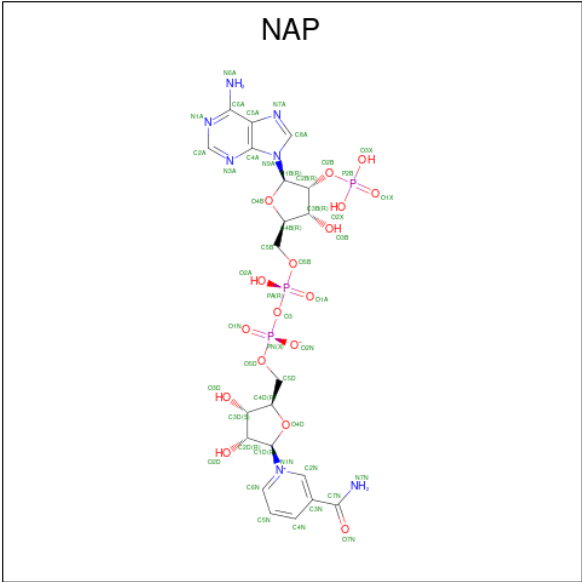
- Molecule 1 is a protein called Corticosteroid 11-beta-dehydrogenase isozyme 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			1956	1246	332	364	14			
1	B	253	Total	C	N	O	S	0	0	0
			1939	1234	330	361	14			

There are 16 discrepancies between the modelled and reference sequences:

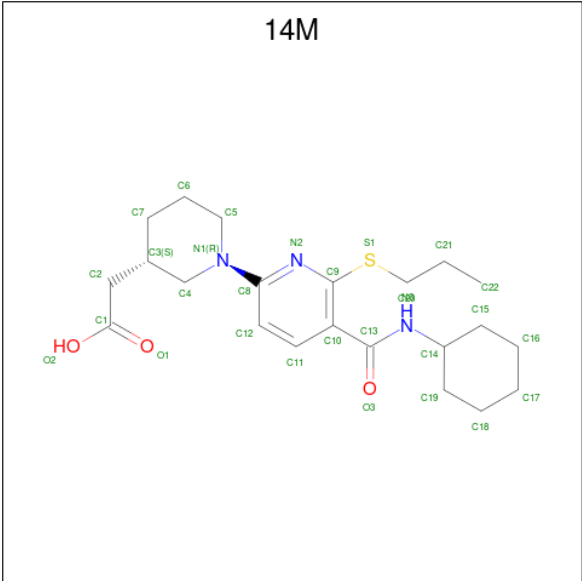
Chain	Residue	Modelled	Actual	Comment	Reference
A	21	GLY	-	expression tag	UNP P28845
A	22	GLY	-	expression tag	UNP P28845
A	23	GLY	-	expression tag	UNP P28845
A	179	LEU	MET	engineered mutation	UNP P28845
A	262	ARG	LEU	engineered mutation	UNP P28845
A	272	SER	CYS	engineered mutation	UNP P28845
A	278	GLU	PHE	engineered mutation	UNP P28845
A	286	TRP	MET	engineered mutation	UNP P28845
B	21	GLY	-	expression tag	UNP P28845
B	22	GLY	-	expression tag	UNP P28845
B	23	GLY	-	expression tag	UNP P28845
B	179	LEU	MET	engineered mutation	UNP P28845
B	262	ARG	LEU	engineered mutation	UNP P28845
B	272	SER	CYS	engineered mutation	UNP P28845
B	278	GLU	PHE	engineered mutation	UNP P28845
B	286	TRP	MET	engineered mutation	UNP P28845

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is {(3*S*)-1-[5-(cyclohexylcarbamoyl)-6-(propylsulfanyl)pyridin-2-yl]piperidin-3-yl}acetic acid (CCD ID: 14M) (formula: C₂₂H₃₃N₃O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			29	22	3	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	S	0	0
			29	22	3	3	1		

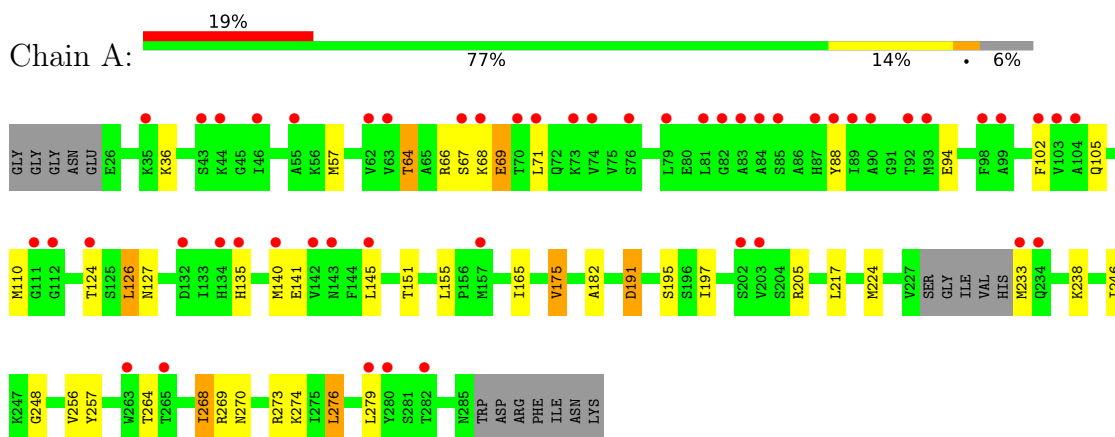
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		
4	B	2	Total	O	0	0
			2	2		

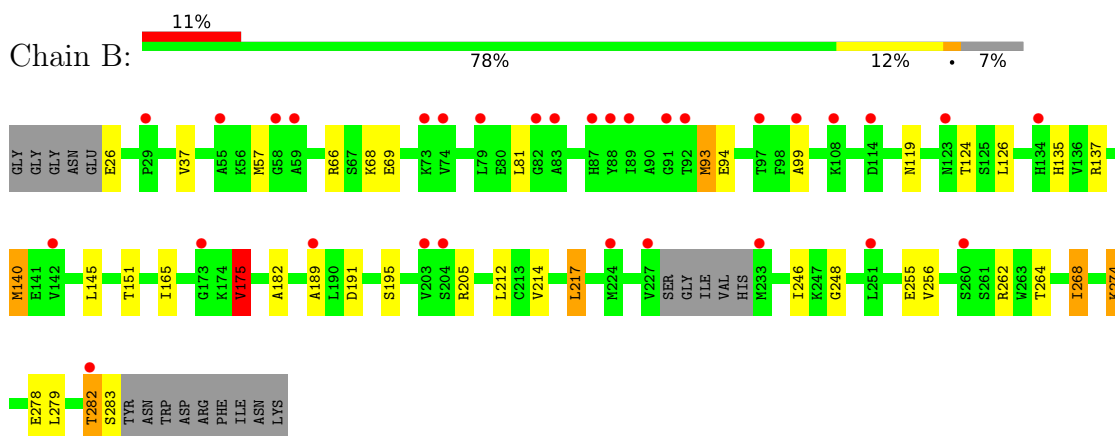
3 Residue-property plots [i](#)

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: Corticosteroid 11-beta-dehydrogenase isozyme 1



- Molecule 1: Corticosteroid 11-beta-dehydrogenase isozyme 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	108.19Å 108.19Å 135.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	21.32 – 2.73 21.32 – 2.73	Depositor EDS
% Data completeness (in resolution range)	96.8 (21.32-2.73) 96.6 (21.32-2.73)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 2.75Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.199 , 0.227 0.252 , 0.267	Depositor DCC
R_{free} test set	1226 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	83.2	Xtriage
Anisotropy	0.354	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 83.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4053	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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, 14M

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	0.87	0/1987	1.44	9/2681 (0.3%)
1	B	0.89	2/1969 (0.1%)	1.39	7/2656 (0.3%)
All	All	0.88	2/3956 (0.1%)	1.42	16/5337 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	MET	SD-CE	-9.32	1.56	1.79
1	B	140	MET	SD-CE	-6.09	1.64	1.79

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	GLU	N-CA-C	-6.37	104.20	112.23
1	B	94	GLU	N-CA-C	-5.94	104.75	112.23
1	A	182	ALA	CA-C-N	5.56	127.67	120.44
1	A	182	ALA	C-N-CA	5.56	127.67	120.44
1	B	182	ALA	CA-C-N	5.52	127.61	120.44
1	B	182	ALA	C-N-CA	5.52	127.61	120.44
1	A	268	ILE	N-CA-CB	5.38	116.85	110.55
1	B	262	ARG	N-CA-C	-5.22	105.67	111.36
1	A	270	ASN	CA-CB-CG	5.19	117.79	112.60
1	B	37	VAL	N-CA-C	5.18	115.58	108.12
1	A	238	LYS	CG-CD-CE	5.16	123.17	111.30
1	A	175	VAL	N-CA-CB	-5.13	103.83	112.44
1	A	191	ASP	CA-CB-CG	5.12	117.72	112.60
1	B	175	VAL	N-CA-CB	-5.12	103.84	112.44
1	A	197	ILE	N-CA-CB	5.04	119.27	110.65
1	B	68	LYS	N-CA-C	5.01	116.55	111.14

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	1956	0	2001	20	0
1	B	1939	0	1990	23	0
2	A	48	0	25	1	0
2	B	48	0	25	2	0
3	A	29	0	32	2	0
3	B	29	0	32	1	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
All	All	4053	0	4105	40	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 (40) 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:B:282:THR:HA	1:B:283:SER:C	2.04	0.82
1:B:66:ARG:HB2	2:B:301:NAP:O3X	2.06	0.56
1:B:274:LYS:O	1:B:278:GLU:HG3	2.06	0.56
1:A:141:GLU:CD	1:B:137:ARG:HH22	2.16	0.52
3:A:302:14M:S1	3:A:302:14M:N3	2.82	0.52
1:B:214:VAL:HG11	1:B:268:ILE:HD13	1.92	0.52
1:A:68:LYS:HG3	1:A:88:TYR:HE2	1.75	0.52
1:A:257:TYR:HE2	1:A:269:ARG:HG3	1.75	0.51
1:A:273:ARG:HG3	1:B:175:VAL:HG22	1.93	0.51
1:A:64:THR:HG21	1:A:102:PHE:CZ	2.45	0.51
1:B:140:MET:HE1	1:B:189:ALA:CB	2.40	0.51
1:A:191:ASP:O	1:A:195:SER:HB2	2.12	0.50
1:A:57:MET:HE1	1:A:246:ILE:HG21	1.93	0.50
1:B:57:MET:HE1	1:B:246:ILE:HG21	1.93	0.49
1:A:64:THR:CG2	1:A:102:PHE:CZ	2.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:MET:HE1	1:B:99:ALA:CB	2.42	0.49
1:B:140:MET:HE1	1:B:189:ALA:HB2	1.94	0.48
1:A:257:TYR:CE2	1:A:269:ARG:HG3	2.48	0.48
1:A:126:LEU:HD11	3:A:302:14M:H17	1.96	0.48
1:B:126:LEU:HD21	3:B:302:14M:H17	1.95	0.48
1:B:248:GLY:HA3	1:B:256:VAL:HG21	1.98	0.46
1:B:119:ASN:ND2	2:B:301:NAP:H4D	2.31	0.45
1:A:248:GLY:HA3	1:A:256:VAL:HG21	1.98	0.45
1:A:151:THR:HG23	1:A:165:ILE:HD13	1.99	0.45
1:A:124:THR:HG22	1:A:135:HIS:CE1	2.52	0.44
1:B:279:LEU:HD23	1:B:279:LEU:HA	1.86	0.44
1:A:276:LEU:HD11	1:B:264:THR:HG23	2.00	0.44
1:A:69:GLU:H	1:A:69:GLU:HG2	1.57	0.43
1:B:151:THR:HG23	1:B:165:ILE:HD13	2.00	0.43
1:B:191:ASP:O	1:B:195:SER:HB2	2.19	0.43
1:B:124:THR:HG22	1:B:135:HIS:CE1	2.54	0.42
1:A:67:SER:O	1:A:71:LEU:HG	2.20	0.42
1:A:66:ARG:HB2	2:A:301:NAP:O1X	2.20	0.42
1:A:140:MET:SD	1:B:140:MET:HE2	2.59	0.42
1:B:93:MET:HE1	1:B:99:ALA:HB2	2.02	0.41
1:B:212:LEU:O	1:B:255:GLU:HA	2.21	0.41
1:B:282:THR:CA	1:B:283:SER:C	2.85	0.41
1:B:217:LEU:HD12	1:B:217:LEU:HA	1.85	0.41
1:A:36:LYS:HG2	1:A:110:MET:HB3	2.04	0.40
1:A:264:THR:O	1:A:268:ILE:HG13	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	251/272 (92%)	243 (97%)	8 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	249/272 (92%)	240 (96%)	9 (4%)	0	100	100
All	All	500/544 (92%)	483 (97%)	17 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	212/226 (94%)	197 (93%)	15 (7%)	13	24
1	B	211/226 (93%)	201 (95%)	10 (5%)	23	43
All	All	423/452 (94%)	398 (94%)	25 (6%)	18	32

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	THR
1	A	69	GLU
1	A	105	GLN
1	A	126	LEU
1	A	127	ASN
1	A	145	LEU
1	A	155	LEU
1	A	175	VAL
1	A	205	ARG
1	A	217	LEU
1	A	224	MET
1	A	233	MET
1	A	274	LYS
1	A	276	LEU
1	A	279	LEU
1	B	26	GLU
1	B	69	GLU
1	B	81	LEU

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Mol	Chain	Res	Type
1	B	145	LEU
1	B	175	VAL
1	B	205	ARG
1	B	217	LEU
1	B	268	ILE
1	B	274	LYS
1	B	282	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	135	HIS
1	A	253	GLN
1	A	270	ASN
1	B	105	GLN
1	B	119	ASN
1	B	135	HIS
1	B	253	GLN
1	B	270	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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	14M	A	302	-	31,31,31	0.87	0	38,41,41	1.87	9 (23%)
2	NAP	A	301	-	50,52,52	1.12	1 (2%)	71,80,80	0.99	6 (8%)
2	NAP	B	301	-	50,52,52	1.13	1 (2%)	71,80,80	0.87	3 (4%)
3	14M	B	302	-	31,31,31	0.81	0	38,41,41	1.67	7 (18%)

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	14M	A	302	-	-	4/20/38/38	0/3/3/3
2	NAP	A	301	-	-	8/35/67/67	0/5/5/5
2	NAP	B	301	-	-	5/35/67/67	0/5/5/5
3	14M	B	302	-	-	4/20/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	NAP	C2N-N1N	7.46	1.43	1.35
2	B	301	NAP	C2N-N1N	7.31	1.43	1.35

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	302	14M	C3-C4-N1	-5.80	104.07	112.35
3	A	302	14M	C3-C4-N1	-5.41	104.62	112.35
3	A	302	14M	C7-C3-C4	4.27	113.65	108.59
2	A	301	NAP	O2B-P2B-O1X	3.71	122.56	109.33
3	A	302	14M	S1-C9-N2	-3.58	115.19	119.19
3	A	302	14M	C11-C10-C9	3.44	120.72	117.29
2	B	301	NAP	O2X-P2B-O2B	-3.42	92.52	105.85
2	A	301	NAP	O3X-P2B-O2B	-3.12	93.68	105.85
3	A	302	14M	C5-N1-C4	3.08	121.27	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	NAP	C2B-C1B-N9A	3.05	118.77	113.75
3	B	302	14M	C11-C10-C9	2.96	120.24	117.29
2	B	301	NAP	O2B-P2B-O1X	2.77	119.22	109.33
3	A	302	14M	C10-C9-S1	2.72	123.94	121.11
3	B	302	14M	C5-N1-C4	2.60	119.89	112.53
3	B	302	14M	S1-C9-N2	-2.53	116.37	119.19
3	B	302	14M	C7-C3-C4	2.53	111.59	108.59
3	B	302	14M	C6-C5-N1	-2.48	106.37	111.06
3	A	302	14M	O3-C13-N3	2.45	127.12	122.47
3	A	302	14M	C10-C13-N3	-2.45	111.36	116.67
3	B	302	14M	O3-C13-N3	2.36	126.96	122.47
2	B	301	NAP	O3-PN-O1N	-2.35	103.65	110.70
3	A	302	14M	C19-C14-C15	2.28	114.74	110.80
2	A	301	NAP	C6N-N1N-C2N	-2.27	119.94	121.88
2	A	301	NAP	O3-PN-O1N	-2.16	104.20	110.70
2	A	301	NAP	O2N-PN-O3	2.10	112.95	107.27

There are no chirality outliers.

All (21) torsion outliers are listed below:

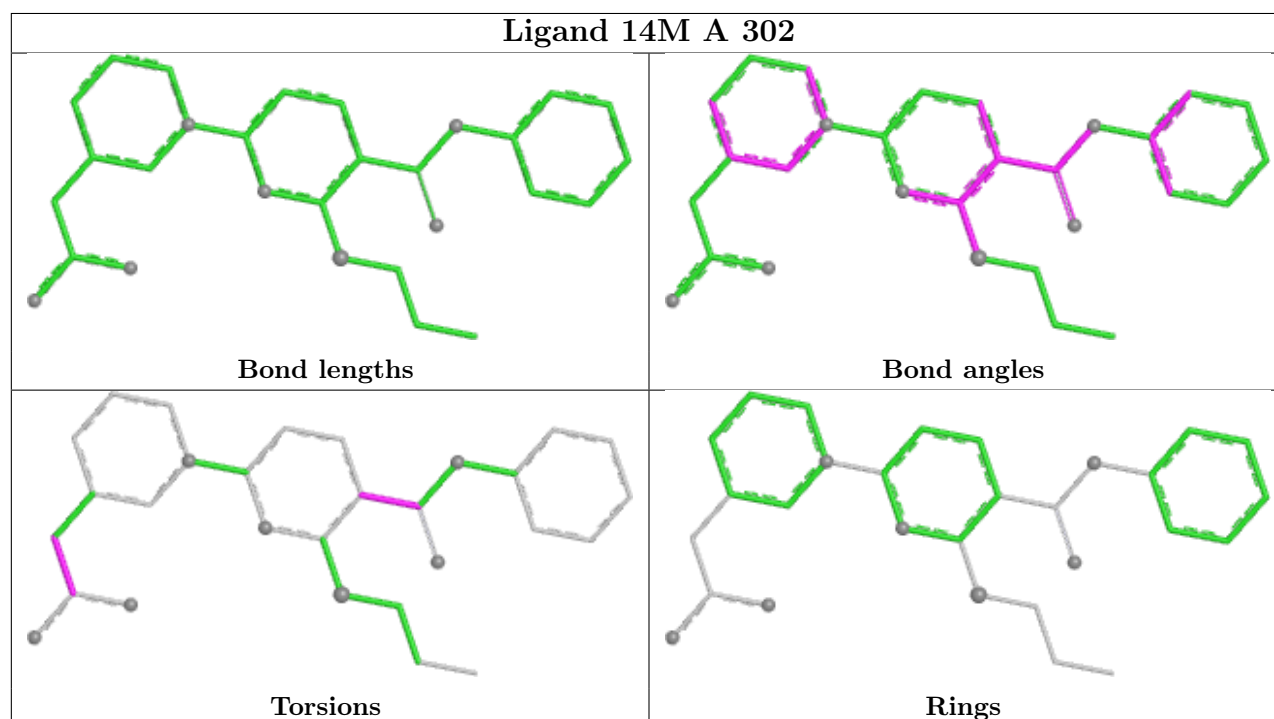
Mol	Chain	Res	Type	Atoms
2	A	301	NAP	C5D-O5D-PN-O3
2	A	301	NAP	C5D-O5D-PN-O1N
2	A	301	NAP	C5D-O5D-PN-O2N
2	B	301	NAP	C5D-O5D-PN-O3
2	B	301	NAP	C5D-O5D-PN-O2N
2	A	301	NAP	C1B-C2B-O2B-P2B
2	A	301	NAP	C3B-C2B-O2B-P2B
3	B	302	14M	O2-C1-C2-C3
3	A	302	14M	C9-C10-C13-O3
3	B	302	14M	C9-C10-C13-O3
3	A	302	14M	C9-C10-C13-N3
3	B	302	14M	C9-C10-C13-N3
3	B	302	14M	O1-C1-C2-C3
3	A	302	14M	O1-C1-C2-C3
3	A	302	14M	O2-C1-C2-C3
2	B	301	NAP	O4B-C4B-C5B-O5B
2	A	301	NAP	PA-O3-PN-O1N
2	A	301	NAP	PA-O3-PN-O2N
2	B	301	NAP	PN-O3-PA-O1A
2	A	301	NAP	C2B-O2B-P2B-O2X
2	B	301	NAP	PA-O3-PN-O1N

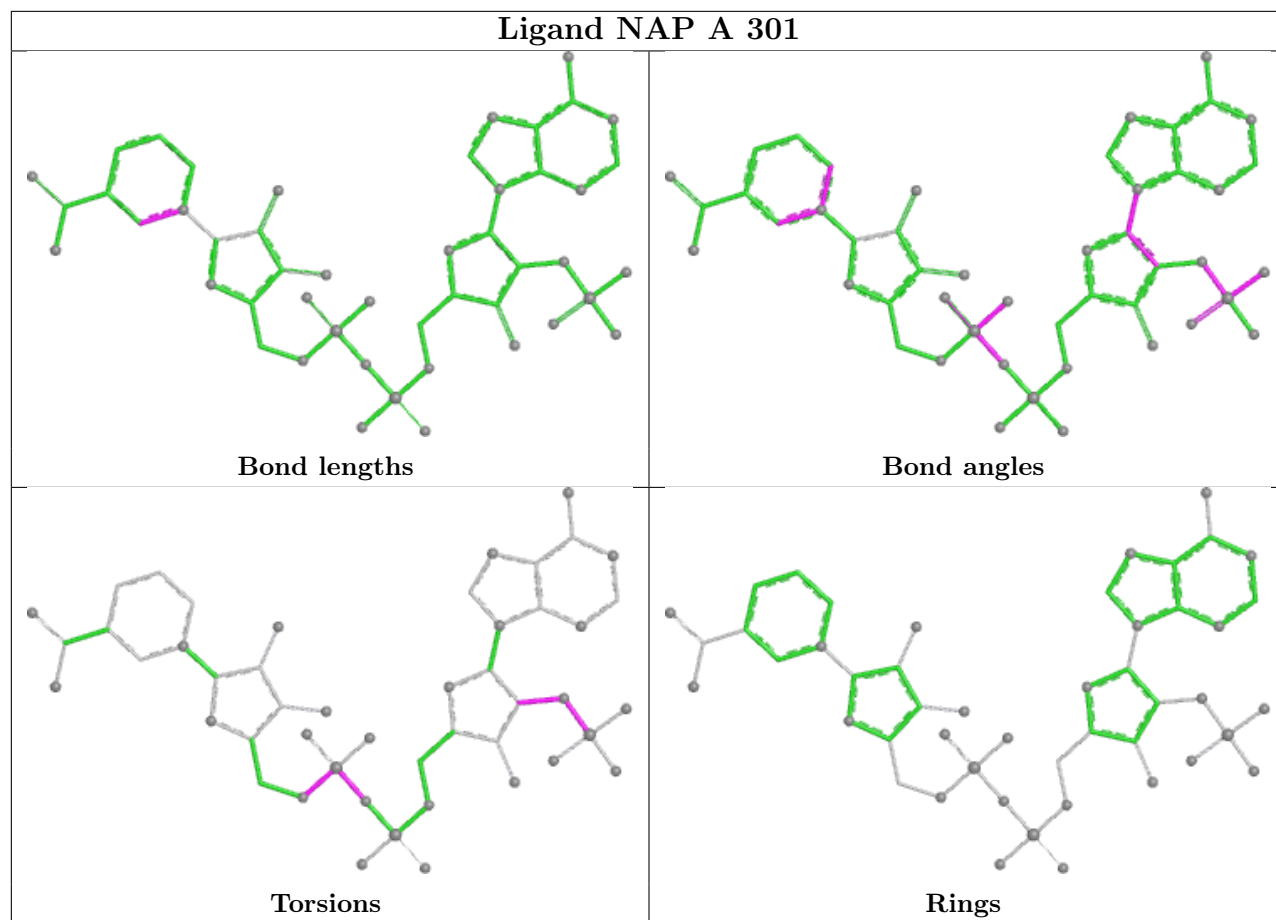
There are no ring outliers.

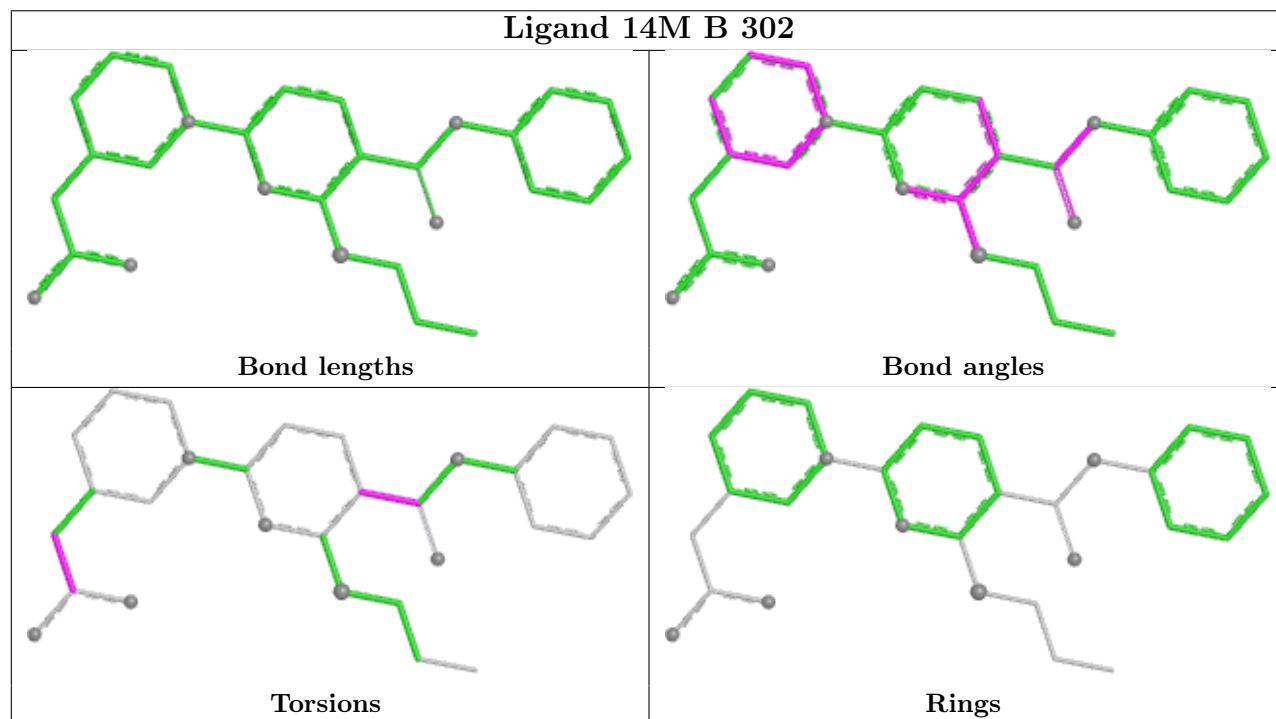
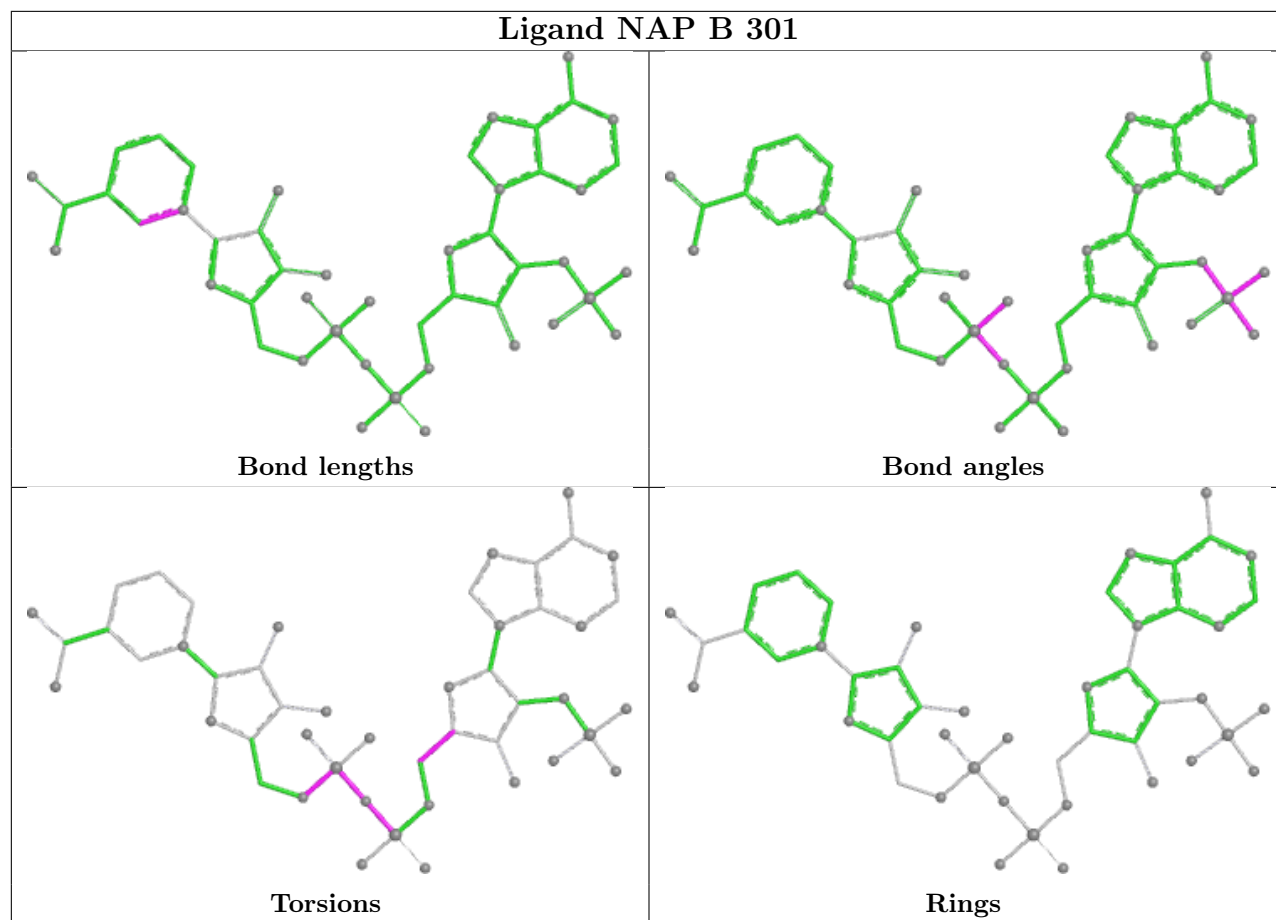
4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	302	14M	2	0
2	A	301	NAP	1	0
2	B	301	NAP	2	0
3	B	302	14M	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	255/272 (93%)	1.23	51 (20%) 3 3	62, 102, 143, 159	0
1	B	253/272 (93%)	0.95	31 (12%) 8 8	60, 96, 134, 158	0
All	All	508/544 (93%)	1.09	82 (16%) 4 5	60, 98, 141, 159	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	ALA	4.9
1	B	55	ALA	4.6
1	A	143	ASN	4.6
1	B	114	ASP	4.5
1	A	142	VAL	4.4
1	A	81	LEU	4.4
1	A	263	TRP	4.2
1	A	282	THR	4.0
1	B	233	MET	4.0
1	A	79	LEU	3.7
1	A	43	SER	3.6
1	B	82	GLY	3.4
1	A	279	LEU	3.3
1	A	99	ALA	3.3
1	A	89	ILE	3.3
1	B	58	GLY	3.3
1	A	82	GLY	3.2
1	B	59	ALA	3.2
1	B	73	LYS	3.1
1	B	97	THR	3.1
1	B	99	ALA	3.1
1	A	87	HIS	3.0
1	B	108	LYS	3.0
1	A	93	MET	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	85	SER	3.0
1	A	44	LYS	3.0
1	A	104	ALA	2.9
1	B	227	VAL	2.9
1	A	76	SER	2.9
1	A	112	GLY	2.9
1	B	83	ALA	2.9
1	A	233	MET	2.8
1	A	203	VAL	2.8
1	A	67	SER	2.8
1	A	132	ASP	2.8
1	A	157	MET	2.8
1	B	142	VAL	2.8
1	A	90	ALA	2.7
1	A	62	VAL	2.7
1	B	203	VAL	2.7
1	A	74	VAL	2.6
1	A	202	SER	2.6
1	A	88	TYR	2.6
1	B	89	ILE	2.6
1	B	79	LEU	2.6
1	A	280	TYR	2.5
1	B	74	VAL	2.5
1	A	102	PHE	2.5
1	A	145	LEU	2.5
1	A	234	GLN	2.5
1	B	91	GLY	2.5
1	A	92	THR	2.4
1	A	135	HIS	2.4
1	A	103	VAL	2.4
1	A	134	HIS	2.3
1	A	111	GLY	2.3
1	B	87	HIS	2.3
1	A	71	LEU	2.3
1	A	35	LYS	2.3
1	B	282	THR	2.3
1	A	63	VAL	2.3
1	A	46	ILE	2.2
1	A	83	ALA	2.2
1	A	68	LYS	2.2
1	B	123	ASN	2.2
1	B	173	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	84	ALA	2.2
1	A	98	PHE	2.2
1	A	124	THR	2.1
1	B	92	THR	2.1
1	A	73	LYS	2.1
1	B	224	MET	2.1
1	B	260	SER	2.1
1	B	29	PRO	2.1
1	B	189	ALA	2.1
1	B	204	SER	2.0
1	A	70	THR	2.0
1	B	251	LEU	2.0
1	B	88	TYR	2.0
1	A	140	MET	2.0
1	A	265	THR	2.0
1	B	134	HIS	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.

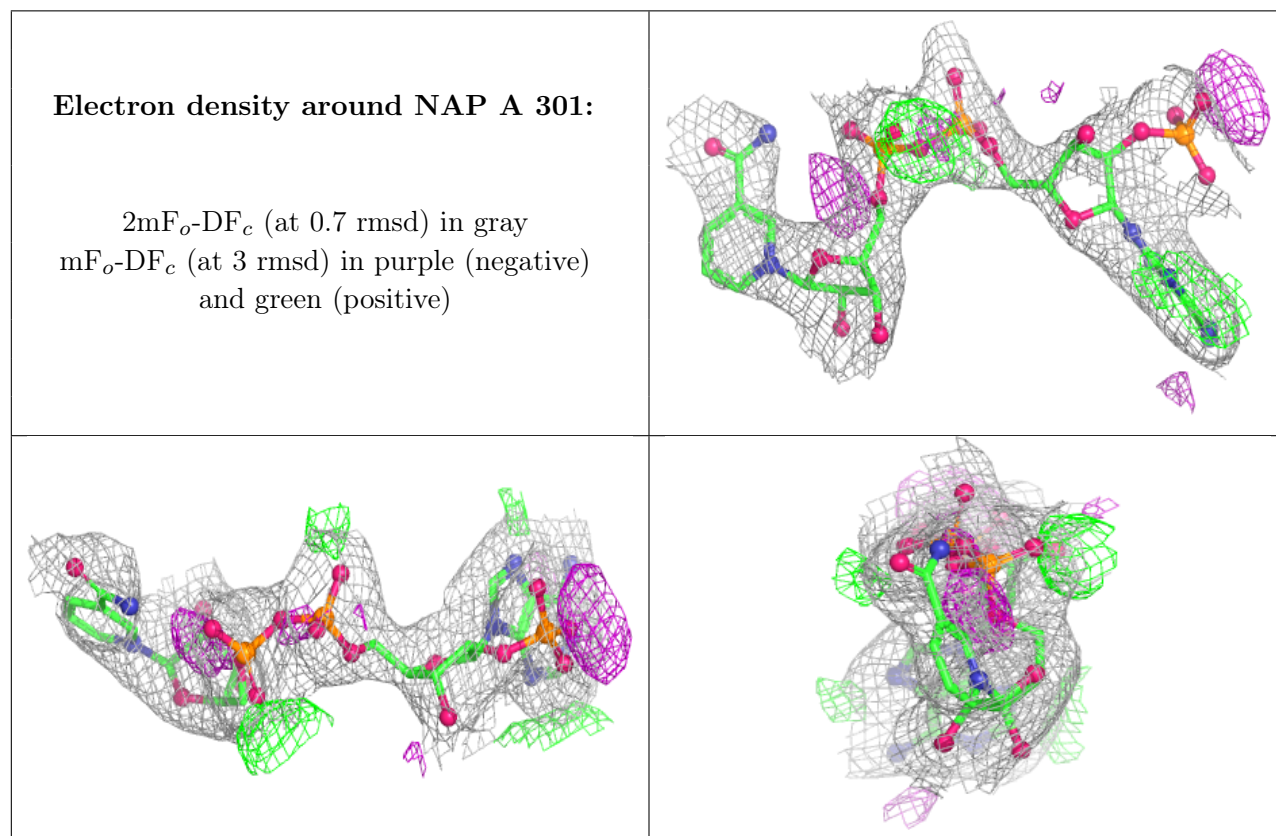
6.4 Ligands [i](#)

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
2	NAP	A	301	48/48	0.87	0.13	77,88,98,100	0
3	14M	A	302	29/29	0.92	0.10	64,77,91,93	0
2	NAP	B	301	48/48	0.93	0.09	70,81,90,91	0
3	14M	B	302	29/29	0.95	0.09	68,77,84,84	0

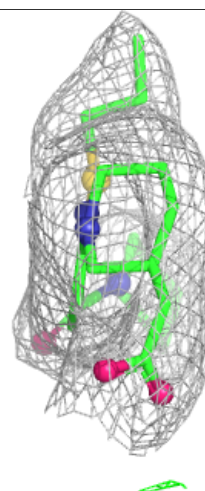
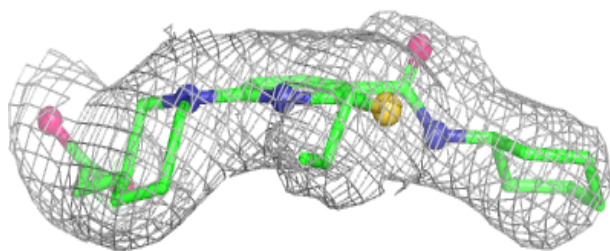
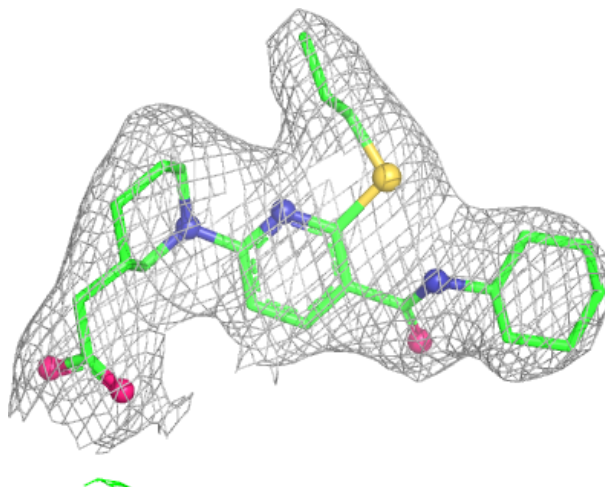
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



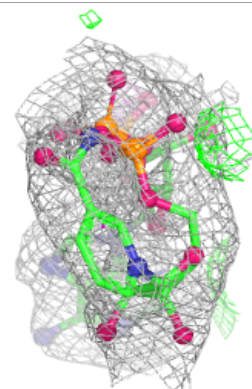
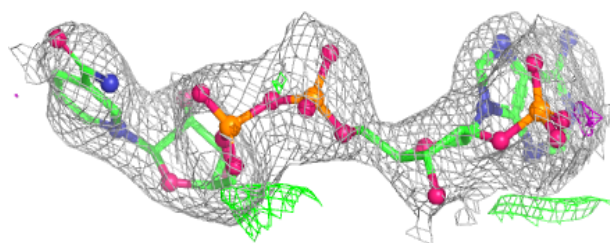
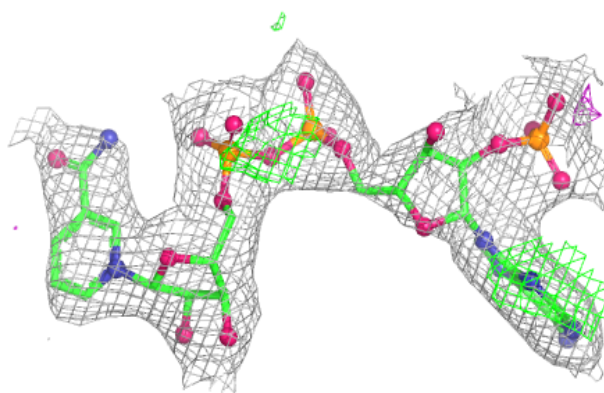
Electron density around 14M A 302:

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

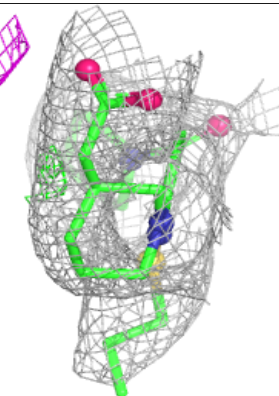
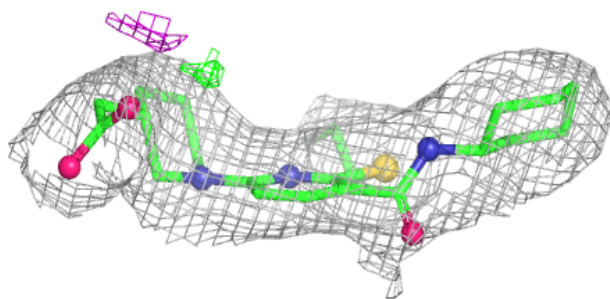
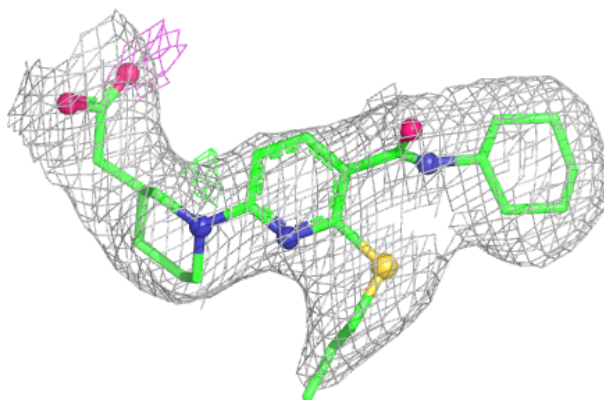


Electron density around NAP B 301:

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

**Electron density around 14M B 302:**

$2mF_o-DF_c$ (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.