



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 18, 2026 – 11:43 PM UTC

PDB ID : 5H5X / pdb_00005h5x
Title : Crystal structure of NADH bound carbonyl reductase from *Streptomyces coelicolor*
Authors : Kong, X.-D.; Xu, J.-H.; Zhou, J.
Deposited on : 2016-11-10
Resolution : 2.30 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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

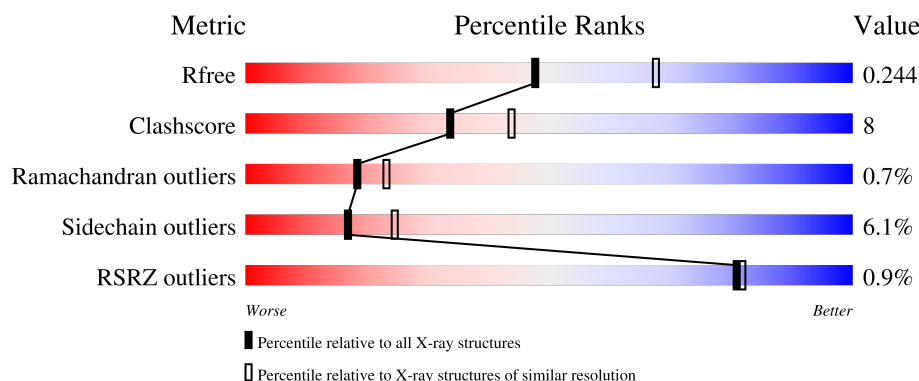
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	263	81% 11% • 6%
1	B	263	81% 13% • •
1	C	263	80% 14% • •
1	D	263	74% 17% • 5%
1	E	263	78% 13% • 6%

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Mol	Chain	Length	Quality of chain
1	F	263	
1	G	263	
1	H	263	
1	I	263	
1	J	263	
1	K	263	
1	L	263	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	IPA	G	302	-	-	X	-
3	IPA	K	303	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 23002 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 Putative oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1758	1106	308	342	2			
1	B	253	Total	C	N	O	S	0	0	0
			1797	1133	314	348	2			
1	C	257	Total	C	N	O	S	0	0	0
			1823	1149	318	354	2			
1	D	250	Total	C	N	O	S	0	0	0
			1774	1116	311	345	2			
1	E	248	Total	C	N	O	S	0	0	0
			1757	1106	308	342	1			
1	F	253	Total	C	N	O	S	0	0	0
			1797	1133	314	348	2			
1	G	257	Total	C	N	O	S	0	0	0
			1823	1149	318	354	2			
1	H	252	Total	C	N	O	S	0	0	0
			1789	1127	313	347	2			
1	I	243	Total	C	N	O	S	0	1	0
			1728	1090	304	333	1			
1	J	253	Total	C	N	O	S	0	0	0
			1797	1133	314	348	2			
1	K	257	Total	C	N	O	S	0	0	0
			1823	1149	318	354	2			
1	L	246	Total	C	N	O	S	0	0	0
			1743	1097	306	339	1			

There are 36 discrepancies between the modelled and reference sequences:

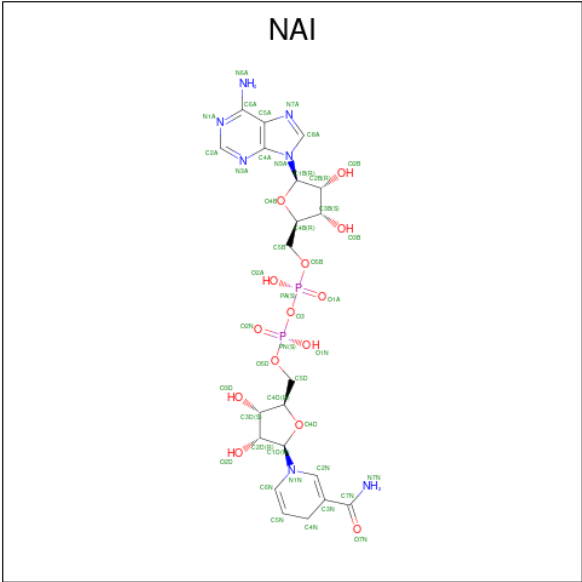
Chain	Residue	Modelled	Actual	Comment	Reference
A	212	GLU	ASP	conflict	UNP Q9KYM4
A	232	ASP	GLU	conflict	UNP Q9KYM4
A	239	VAL	ALA	conflict	UNP Q9KYM4
B	212	GLU	ASP	conflict	UNP Q9KYM4
B	232	ASP	GLU	conflict	UNP Q9KYM4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	239	VAL	ALA	conflict	UNP Q9KYM4
C	212	GLU	ASP	conflict	UNP Q9KYM4
C	232	ASP	GLU	conflict	UNP Q9KYM4
C	239	VAL	ALA	conflict	UNP Q9KYM4
D	212	GLU	ASP	conflict	UNP Q9KYM4
D	232	ASP	GLU	conflict	UNP Q9KYM4
D	239	VAL	ALA	conflict	UNP Q9KYM4
E	212	GLU	ASP	conflict	UNP Q9KYM4
E	232	ASP	GLU	conflict	UNP Q9KYM4
E	239	VAL	ALA	conflict	UNP Q9KYM4
F	212	GLU	ASP	conflict	UNP Q9KYM4
F	232	ASP	GLU	conflict	UNP Q9KYM4
F	239	VAL	ALA	conflict	UNP Q9KYM4
G	212	GLU	ASP	conflict	UNP Q9KYM4
G	232	ASP	GLU	conflict	UNP Q9KYM4
G	239	VAL	ALA	conflict	UNP Q9KYM4
H	212	GLU	ASP	conflict	UNP Q9KYM4
H	232	ASP	GLU	conflict	UNP Q9KYM4
H	239	VAL	ALA	conflict	UNP Q9KYM4
I	212	GLU	ASP	conflict	UNP Q9KYM4
I	232	ASP	GLU	conflict	UNP Q9KYM4
I	239	VAL	ALA	conflict	UNP Q9KYM4
J	212	GLU	ASP	conflict	UNP Q9KYM4
J	232	ASP	GLU	conflict	UNP Q9KYM4
J	239	VAL	ALA	conflict	UNP Q9KYM4
K	212	GLU	ASP	conflict	UNP Q9KYM4
K	232	ASP	GLU	conflict	UNP Q9KYM4
K	239	VAL	ALA	conflict	UNP Q9KYM4
L	212	GLU	ASP	conflict	UNP Q9KYM4
L	232	ASP	GLU	conflict	UNP Q9KYM4
L	239	VAL	ALA	conflict	UNP Q9KYM4

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	H	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	I	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	J	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	K	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	L	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C₃H₈O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	3	1		
3	A	1	Total	C	O	0	0
			4	3	1		
3	B	1	Total	C	O	0	0
			4	3	1		
3	B	1	Total	C	O	0	0
			4	3	1		
3	C	1	Total	C	O	0	0
			4	3	1		
3	C	1	Total	C	O	0	0
			4	3	1		
3	D	1	Total	C	O	0	0
			4	3	1		
3	E	1	Total	C	O	0	0
			4	3	1		
3	E	1	Total	C	O	0	0
			4	3	1		
3	F	1	Total	C	O	0	0
			4	3	1		
3	F	1	Total	C	O	0	0
			4	3	1		
3	G	1	Total	C	O	0	0
			4	3	1		
3	G	1	Total	C	O	0	0
			4	3	1		
3	H	1	Total	C	O	0	0
			4	3	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	H	1	Total C O 4 3 1	0	0
3	I	1	Total C O 4 3 1	0	0
3	I	1	Total C O 4 3 1	0	0
3	J	1	Total C O 4 3 1	0	0
3	K	1	Total C O 4 3 1	0	0
3	K	1	Total C O 4 3 1	0	0
3	L	1	Total C O 4 3 1	0	0
3	L	1	Total C O 4 3 1	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	C	1	Total Mg 1 1	0	0
4	D	1	Total Mg 1 1	0	0
4	E	1	Total Mg 1 1	0	0
4	H	1	Total Mg 1 1	0	0
4	I	1	Total Mg 1 1	0	0
4	L	1	Total Mg 1 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	93	Total O 93 93	0	0
5	B	85	Total O 85 85	0	0
5	C	97	Total O 97 97	0	0

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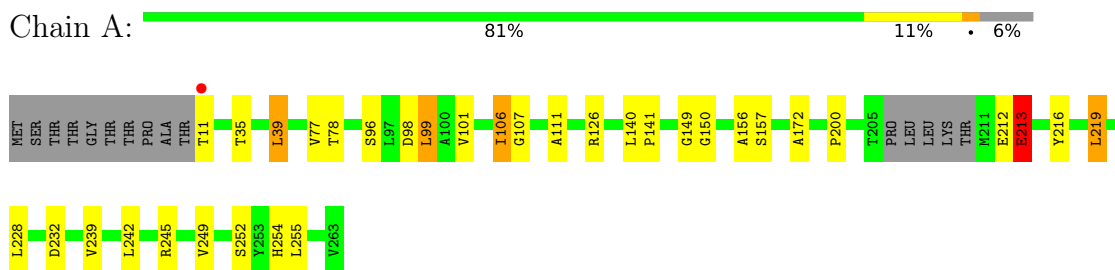
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	76	Total 76	O 76	0	0
5	E	75	Total 75	O 75	0	0
5	F	78	Total 78	O 78	0	0
5	G	80	Total 80	O 80	0	0
5	H	77	Total 77	O 77	0	0
5	I	52	Total 52	O 52	0	0
5	J	93	Total 93	O 93	0	0
5	K	98	Total 98	O 98	0	0
5	L	67	Total 67	O 67	0	0

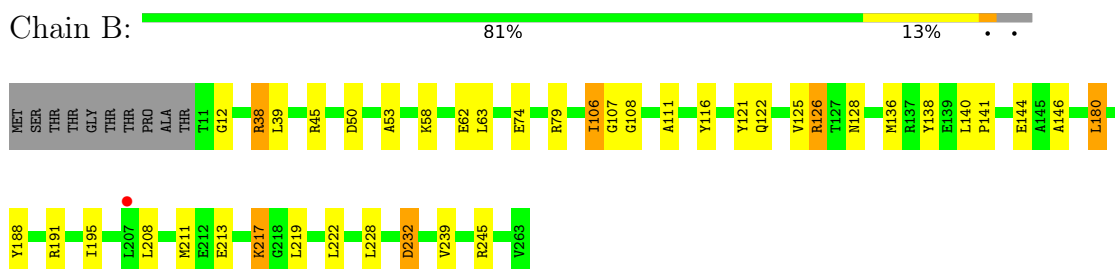
3 Residue-property plots [i](#)

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.

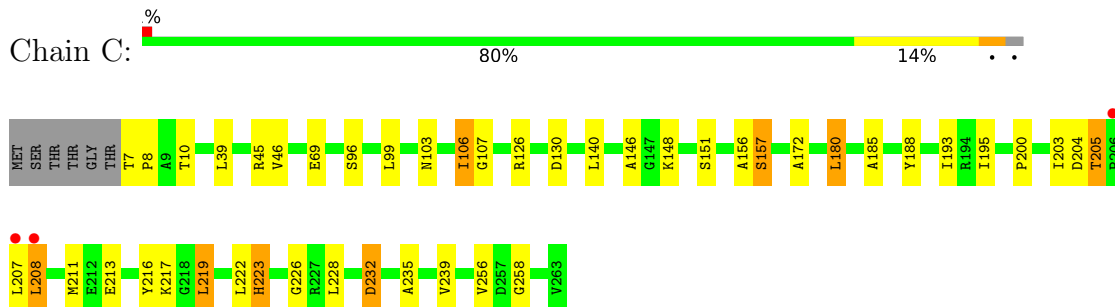
- Molecule 1: Putative oxidoreductase



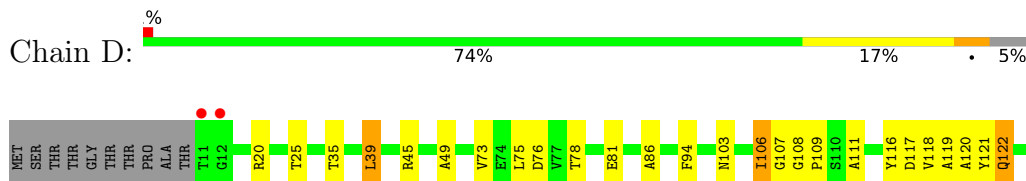
- Molecule 1: Putative oxidoreductase



- Molecule 1: Putative oxidoreductase

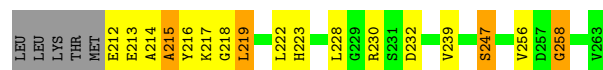
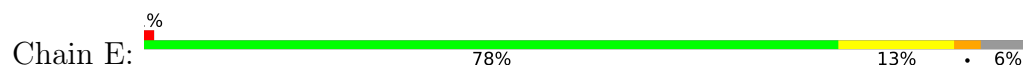


- Molecule 1: Putative oxidoreductase

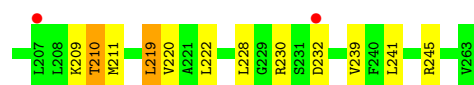
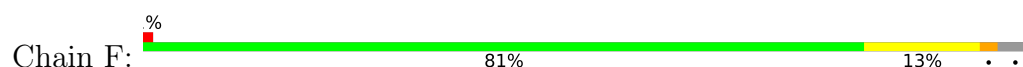




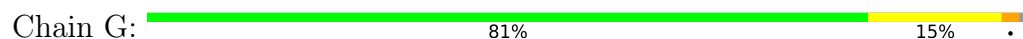
- Molecule 1: Putative oxidoreductase



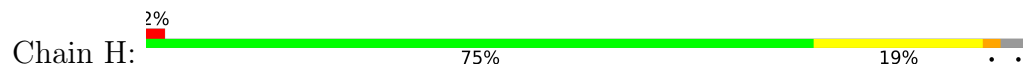
- Molecule 1: Putative oxidoreductase



- Molecule 1: Putative oxidoreductase

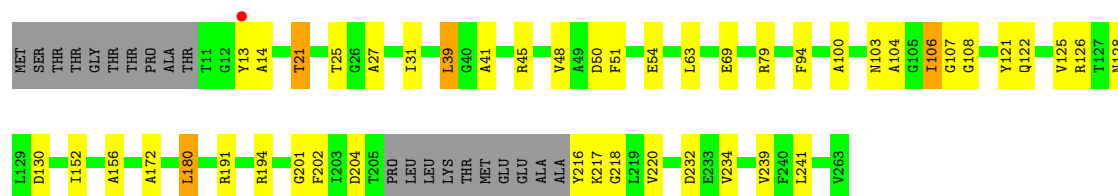


- Molecule 1: Putative oxidoreductase

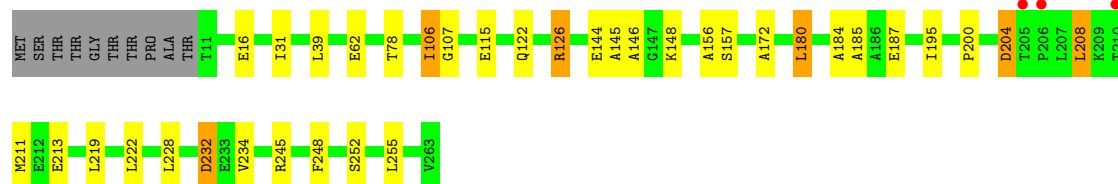
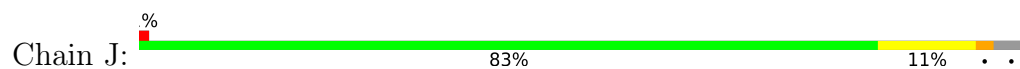


- Molecule 1: Putative oxidoreductase

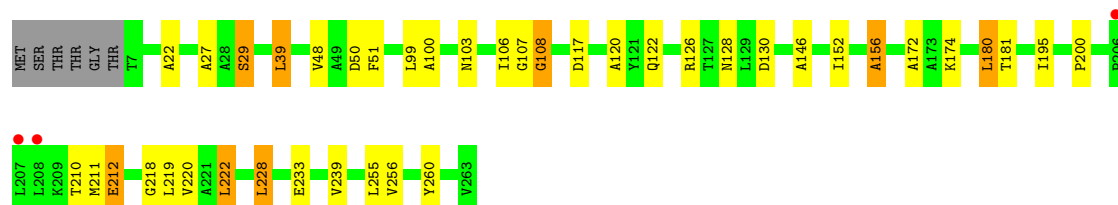
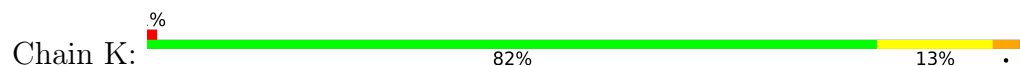




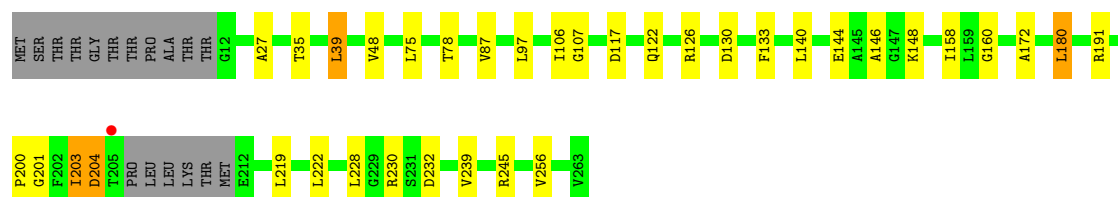
- Molecule 1: Putative oxidoreductase



- Molecule 1: Putative oxidoreductase



- Molecule 1: Putative oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	187.79Å 187.79Å 80.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.50 – 2.30 32.50 – 2.30	Depositor EDS
% Data completeness (in resolution range)	84.0 (32.50-2.30) 83.9 (32.50-2.30)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.179 , 0.243 0.182 , 0.244	Depositor DCC
R_{free} test set	5979 reflections (4.22%)	wwPDB-VP
Wilson B-factor (Å ²)	41.7	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 31.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.058 for -h,-k,l 0.054 for h,-h-k,-l 0.055 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	23002	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 2.05% 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: MG, NAI, IPA

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	0.54	0/1785	0.96	3/2423 (0.1%)
1	B	0.56	1/1826 (0.1%)	0.93	3/2481 (0.1%)
1	C	0.58	1/1853 (0.1%)	0.98	5/2520 (0.2%)
1	D	0.59	1/1801 (0.1%)	0.93	0/2444
1	E	0.55	0/1785	0.89	2/2425 (0.1%)
1	F	0.59	0/1826	0.95	2/2481 (0.1%)
1	G	0.56	0/1853	0.93	2/2520 (0.1%)
1	H	0.55	0/1818	0.89	2/2470 (0.1%)
1	I	0.50	0/1758	0.86	1/2386 (0.0%)
1	J	0.56	1/1826 (0.1%)	0.90	0/2481
1	K	0.56	0/1853	0.92	3/2520 (0.1%)
1	L	0.55	0/1770	0.92	0/2403
All	All	0.56	4/21754 (0.0%)	0.92	23/29554 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1
1	E	0	1
1	G	0	1
1	H	0	1
1	J	0	1
All	All	0	5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	232	ASP	C-O	-5.92	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	232	ASP	C-O	-5.49	1.17	1.24
1	D	232	ASP	C-O	-5.31	1.17	1.24
1	B	232	ASP	C-O	-5.22	1.17	1.24

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	146	ALA	N-CA-C	9.21	122.51	111.82
1	A	106	ILE	N-CA-C	-7.98	97.15	108.80
1	C	146	ALA	N-CA-C	7.49	120.51	111.82
1	G	259	ALA	N-CA-C	-6.58	105.21	113.23
1	C	46	VAL	N-CA-C	6.54	117.27	108.11

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	106	ILE	Peptide
1	E	258	GLY	Peptide
1	G	258	GLY	Peptide
1	H	106	ILE	Peptide
1	J	106	ILE	Peptide

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	1758	0	1737	23	0
1	B	1797	0	1787	22	0
1	C	1823	0	1813	25	0
1	D	1774	0	1757	43	0
1	E	1757	0	1735	35	0
1	F	1797	0	1787	23	0
1	G	1823	0	1813	25	0
1	H	1789	0	1776	31	0
1	I	1728	0	1719	33	0
1	J	1797	0	1787	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	1823	0	1813	25	0
1	L	1743	0	1721	22	0
2	A	44	0	27	3	0
2	B	44	0	27	3	0
2	C	44	0	27	6	0
2	D	44	0	27	3	0
2	E	44	0	27	6	0
2	F	44	0	27	3	0
2	G	44	0	27	3	0
2	H	44	0	27	3	0
2	I	44	0	27	5	0
2	J	44	0	27	3	0
2	K	44	0	27	6	0
2	L	44	0	27	5	0
3	A	8	0	16	2	0
3	B	8	0	16	2	0
3	C	8	0	16	3	0
3	D	4	0	8	2	0
3	E	8	0	16	2	0
3	F	8	0	16	3	0
3	G	8	0	16	4	0
3	H	8	0	16	3	0
3	I	8	0	16	4	0
3	J	4	0	8	2	0
3	K	8	0	16	4	0
3	L	8	0	16	3	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	H	1	0	0	0	0
4	I	1	0	0	0	0
4	L	1	0	0	0	0
5	A	93	0	0	4	0
5	B	85	0	0	4	0
5	C	97	0	0	4	0
5	D	76	0	0	7	0
5	E	75	0	0	3	0
5	F	78	0	0	2	0
5	G	80	0	0	2	0
5	H	77	0	0	2	0
5	I	52	0	0	4	0
5	J	93	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	K	98	0	0	2	0
5	L	67	0	0	2	0
All	All	23002	0	21745	335	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 8.

The worst 5 of 335 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:F:301:NAI:H4N	3:F:302:IPA:H2	1.42	1.00
1:E:205:THR:HG21	2:E:301:NAI:H72N	1.40	0.87
2:H:301:NAI:H4N	3:H:303:IPA:H2	1.60	0.84
2:B:301:NAI:H4N	3:B:302:IPA:H2	1.63	0.81
1:K:172:ALA:HB2	1:L:180:LEU:HD13	1.64	0.79

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	244/263 (93%)	229 (94%)	12 (5%)	3 (1%)	10	12
1	B	251/263 (95%)	239 (95%)	11 (4%)	1 (0%)	30	38
1	C	255/263 (97%)	244 (96%)	9 (4%)	2 (1%)	16	20
1	D	246/263 (94%)	236 (96%)	10 (4%)	0	100	100
1	E	244/263 (93%)	230 (94%)	12 (5%)	2 (1%)	16	20
1	F	251/263 (95%)	238 (95%)	10 (4%)	3 (1%)	10	12
1	G	255/263 (97%)	243 (95%)	11 (4%)	1 (0%)	30	38
1	H	250/263 (95%)	238 (95%)	11 (4%)	1 (0%)	30	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	I	240/263 (91%)	223 (93%)	16 (7%)	1 (0%)	30	38
1	J	251/263 (95%)	237 (94%)	11 (4%)	3 (1%)	10	12
1	K	255/263 (97%)	243 (95%)	9 (4%)	3 (1%)	10	12
1	L	242/263 (92%)	228 (94%)	13 (5%)	1 (0%)	30	38
All	All	2984/3156 (95%)	2828 (95%)	135 (4%)	21 (1%)	18	23

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	213	GLU
1	C	156	ALA
1	C	204	ASP
1	F	210	THR
1	H	211	MET

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	162/175 (93%)	156 (96%)	6 (4%)	30	45
1	B	167/175 (95%)	157 (94%)	10 (6%)	17	25
1	C	170/175 (97%)	157 (92%)	13 (8%)	12	16
1	D	164/175 (94%)	151 (92%)	13 (8%)	11	15
1	E	162/175 (93%)	156 (96%)	6 (4%)	30	45
1	F	167/175 (95%)	159 (95%)	8 (5%)	23	35
1	G	170/175 (97%)	158 (93%)	12 (7%)	13	19
1	H	166/175 (95%)	152 (92%)	14 (8%)	10	14
1	I	160/175 (91%)	152 (95%)	8 (5%)	22	33
1	J	167/175 (95%)	153 (92%)	14 (8%)	10	14
1	K	170/175 (97%)	162 (95%)	8 (5%)	23	35
1	L	160/175 (91%)	150 (94%)	10 (6%)	16	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1985/2100 (94%)	1863 (94%)	122 (6%)	17	24

5 of 122 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	62	GLU
1	K	228	LEU
1	H	118	VAL
1	K	222	LEU
1	L	219	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	175	HIS
1	K	102	ASN
1	L	122	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 40 ligands modelled in this entry, 6 are monoatomic - leaving 34 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	IPA	H	303	-	3,3,3	0.51	0	3,3,3	0.27	0
2	NAI	B	301	-	47,48,48	2.05	18 (38%)	64,73,73	1.75	15 (23%)
2	NAI	J	301	-	47,48,48	2.11	18 (38%)	64,73,73	1.78	12 (18%)
3	IPA	L	303	-	3,3,3	0.49	0	3,3,3	0.37	0
3	IPA	G	302	-	3,3,3	0.40	0	3,3,3	0.23	0
3	IPA	G	303	-	3,3,3	0.50	0	3,3,3	0.38	0
3	IPA	K	302	-	3,3,3	0.58	0	3,3,3	0.12	0
3	IPA	B	302	-	3,3,3	0.52	0	3,3,3	0.19	0
3	IPA	C	302	-	3,3,3	0.45	0	3,3,3	0.27	0
2	NAI	E	301	-	47,48,48	1.94	17 (36%)	64,73,73	1.81	14 (21%)
3	IPA	B	303	-	3,3,3	0.48	0	3,3,3	0.40	0
2	NAI	G	301	-	47,48,48	2.13	19 (40%)	64,73,73	1.72	14 (21%)
2	NAI	F	301	-	47,48,48	2.03	19 (40%)	64,73,73	1.68	10 (15%)
3	IPA	I	304	-	3,3,3	0.47	0	3,3,3	0.28	0
3	IPA	F	303	-	3,3,3	0.55	0	3,3,3	0.35	0
3	IPA	L	304	-	3,3,3	0.44	0	3,3,3	0.30	0
2	NAI	I	301	-	47,48,48	2.02	19 (40%)	64,73,73	1.74	13 (20%)
3	IPA	A	303	-	3,3,3	0.52	0	3,3,3	0.36	0
3	IPA	F	302	-	3,3,3	0.48	0	3,3,3	0.32	0
2	NAI	K	301	-	47,48,48	1.99	15 (31%)	64,73,73	1.71	12 (18%)
3	IPA	I	303	-	3,3,3	0.52	0	3,3,3	0.34	0
2	NAI	D	301	-	47,48,48	1.90	15 (31%)	64,73,73	1.73	12 (18%)
3	IPA	C	304	-	3,3,3	0.52	0	3,3,3	0.35	0
2	NAI	L	301	-	47,48,48	2.01	19 (40%)	64,73,73	1.75	13 (20%)
2	NAI	A	301	-	47,48,48	1.98	17 (36%)	64,73,73	1.68	14 (21%)
2	NAI	H	301	-	47,48,48	1.94	16 (34%)	64,73,73	1.72	13 (20%)
3	IPA	K	303	-	3,3,3	0.56	0	3,3,3	0.54	0
3	IPA	D	302	-	3,3,3	0.58	0	3,3,3	0.34	0
2	NAI	C	301	-	47,48,48	2.08	18 (38%)	64,73,73	1.80	14 (21%)
3	IPA	E	304	-	3,3,3	0.44	0	3,3,3	0.37	0
3	IPA	J	302	-	3,3,3	0.55	0	3,3,3	0.26	0
3	IPA	A	302	-	3,3,3	0.50	0	3,3,3	0.22	0
3	IPA	H	304	-	3,3,3	0.43	0	3,3,3	0.63	0
3	IPA	E	303	-	3,3,3	0.42	0	3,3,3	0.29	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
2	NAI	B	301	-	-	5/29/72/72	0/5/5/5
2	NAI	J	301	-	-	3/29/72/72	0/5/5/5
2	NAI	K	301	-	-	8/29/72/72	0/5/5/5
2	NAI	D	301	-	-	5/29/72/72	0/5/5/5
2	NAI	G	301	-	-	7/29/72/72	0/5/5/5
2	NAI	F	301	-	-	12/29/72/72	0/5/5/5
2	NAI	C	301	-	-	7/29/72/72	0/5/5/5
2	NAI	H	301	-	-	9/29/72/72	0/5/5/5
2	NAI	L	301	-	-	9/29/72/72	0/5/5/5
2	NAI	A	301	-	-	9/29/72/72	0/5/5/5
2	NAI	I	301	-	-	13/29/72/72	0/5/5/5
2	NAI	E	301	-	-	4/29/72/72	0/5/5/5

The worst 5 of 210 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	301	NAI	PN-O3	4.60	1.64	1.59
2	J	301	NAI	PN-O3	4.58	1.64	1.59
2	F	301	NAI	PN-O3	4.36	1.64	1.59
2	C	301	NAI	C4N-C5N	4.23	1.60	1.49
2	B	301	NAI	PN-O3	4.19	1.64	1.59

The worst 5 of 156 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	NAI	C5A-C4A-N3A	-5.50	119.14	126.72
2	L	301	NAI	C5A-C4A-N3A	-5.50	119.14	126.72
2	I	301	NAI	C5A-C4A-N3A	-5.46	119.20	126.72
2	G	301	NAI	C5A-C4A-N3A	-5.42	119.26	126.72
2	H	301	NAI	C5A-C4A-N3A	-5.14	119.64	126.72

There are no chirality outliers.

5 of 91 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	NAI	C5D-O5D-PN-O3
2	A	301	NAI	C5D-O5D-PN-O1N

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Mol	Chain	Res	Type	Atoms
2	A	301	NAI	C5D-O5D-PN-O2N
2	C	301	NAI	C5D-O5D-PN-O3
2	C	301	NAI	C5D-O5D-PN-O1N

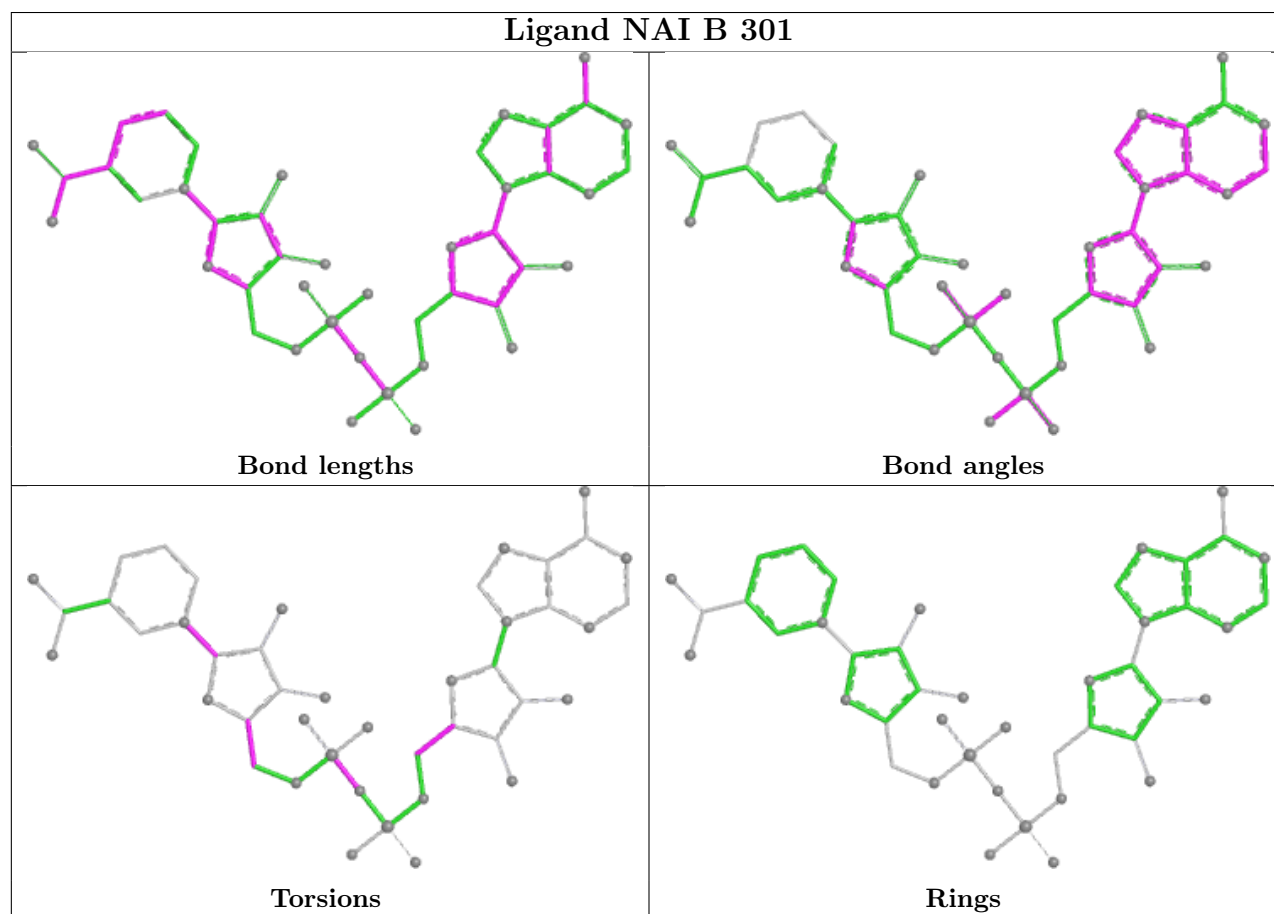
There are no ring outliers.

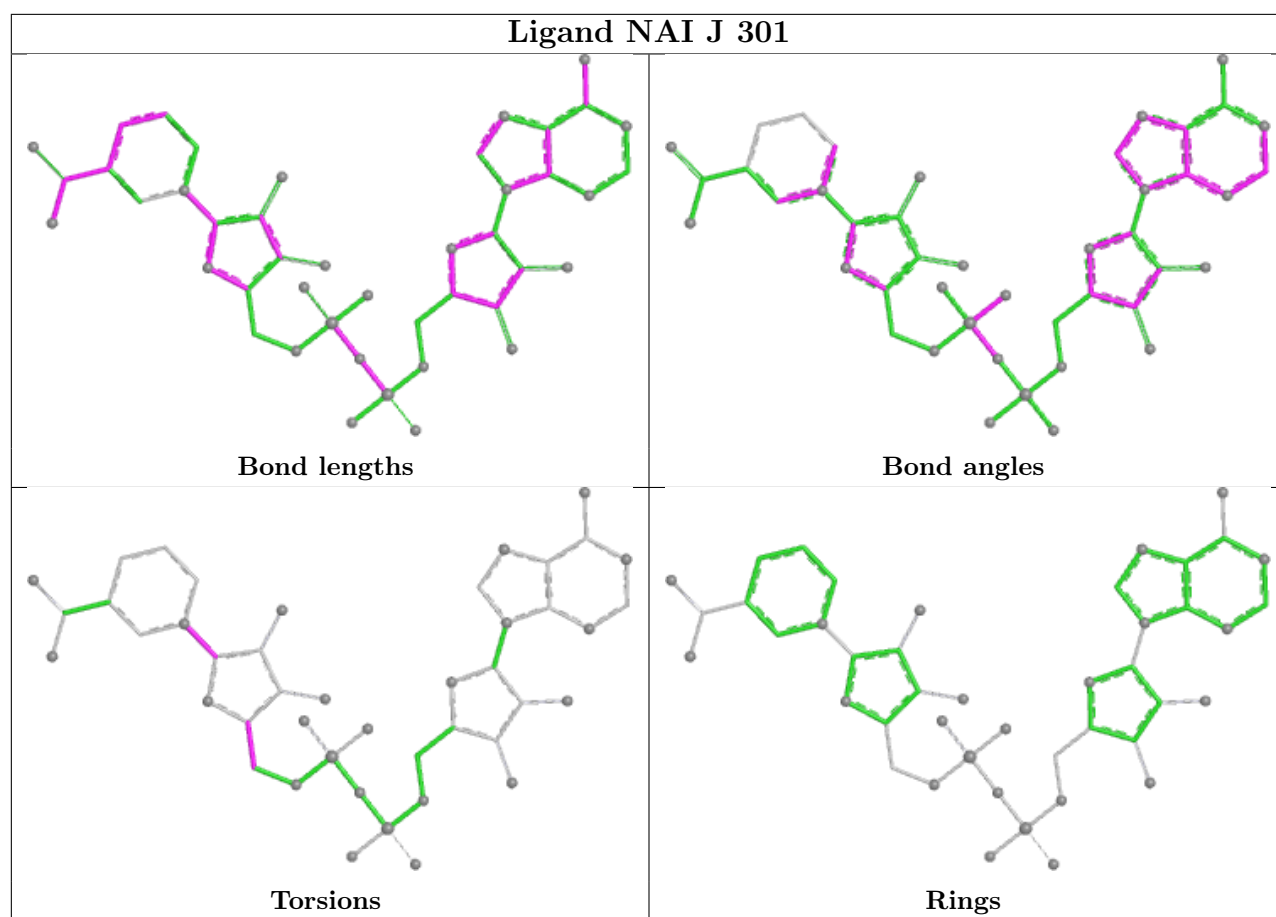
30 monomers are involved in 57 short contacts:

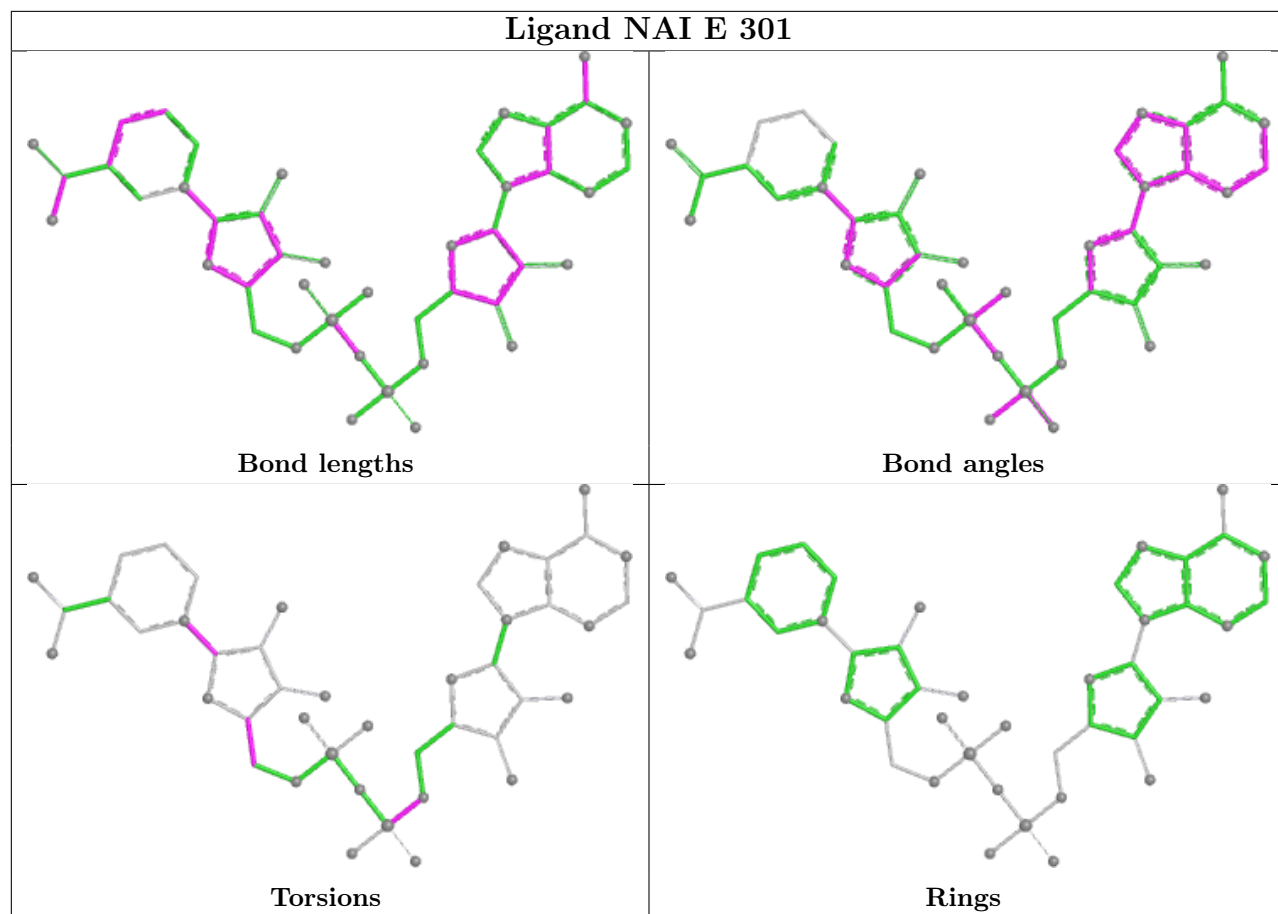
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	303	IPA	3	0
2	B	301	NAI	3	0
2	J	301	NAI	3	0
3	L	303	IPA	1	0
3	G	302	IPA	4	0
3	G	303	IPA	1	0
3	K	302	IPA	1	0
3	B	302	IPA	2	0
3	C	302	IPA	3	0
2	E	301	NAI	6	0
2	G	301	NAI	3	0
2	F	301	NAI	3	0
3	I	304	IPA	3	0
3	F	303	IPA	1	0
3	L	304	IPA	3	0
2	I	301	NAI	5	0
3	F	302	IPA	3	0
2	K	301	NAI	6	0
3	I	303	IPA	1	0
2	D	301	NAI	3	0
2	L	301	NAI	5	0
2	A	301	NAI	3	0
2	H	301	NAI	3	0
3	K	303	IPA	4	0
3	D	302	IPA	2	0
2	C	301	NAI	6	0
3	J	302	IPA	2	0
3	A	302	IPA	2	0
3	H	304	IPA	1	0
3	E	303	IPA	2	0

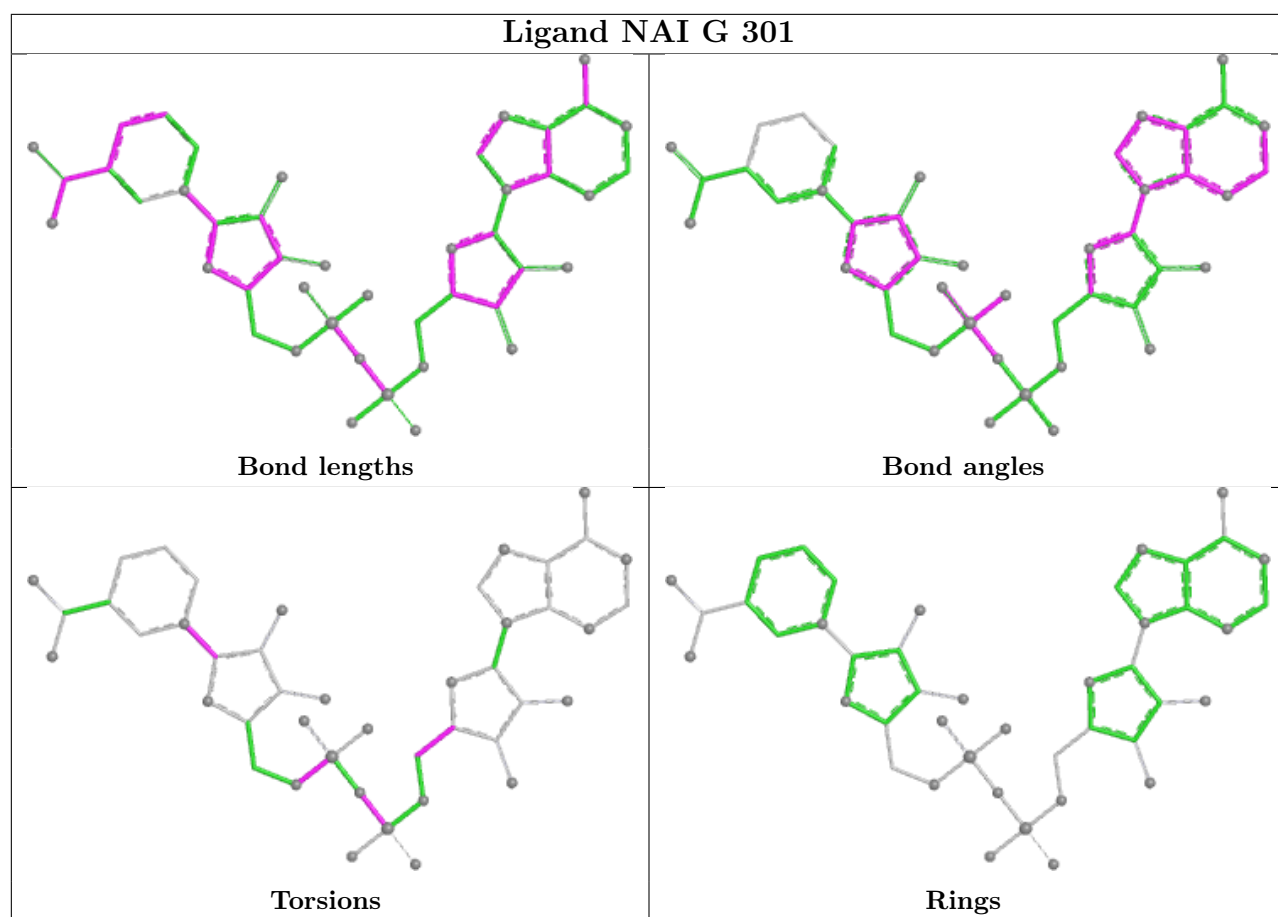
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

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.

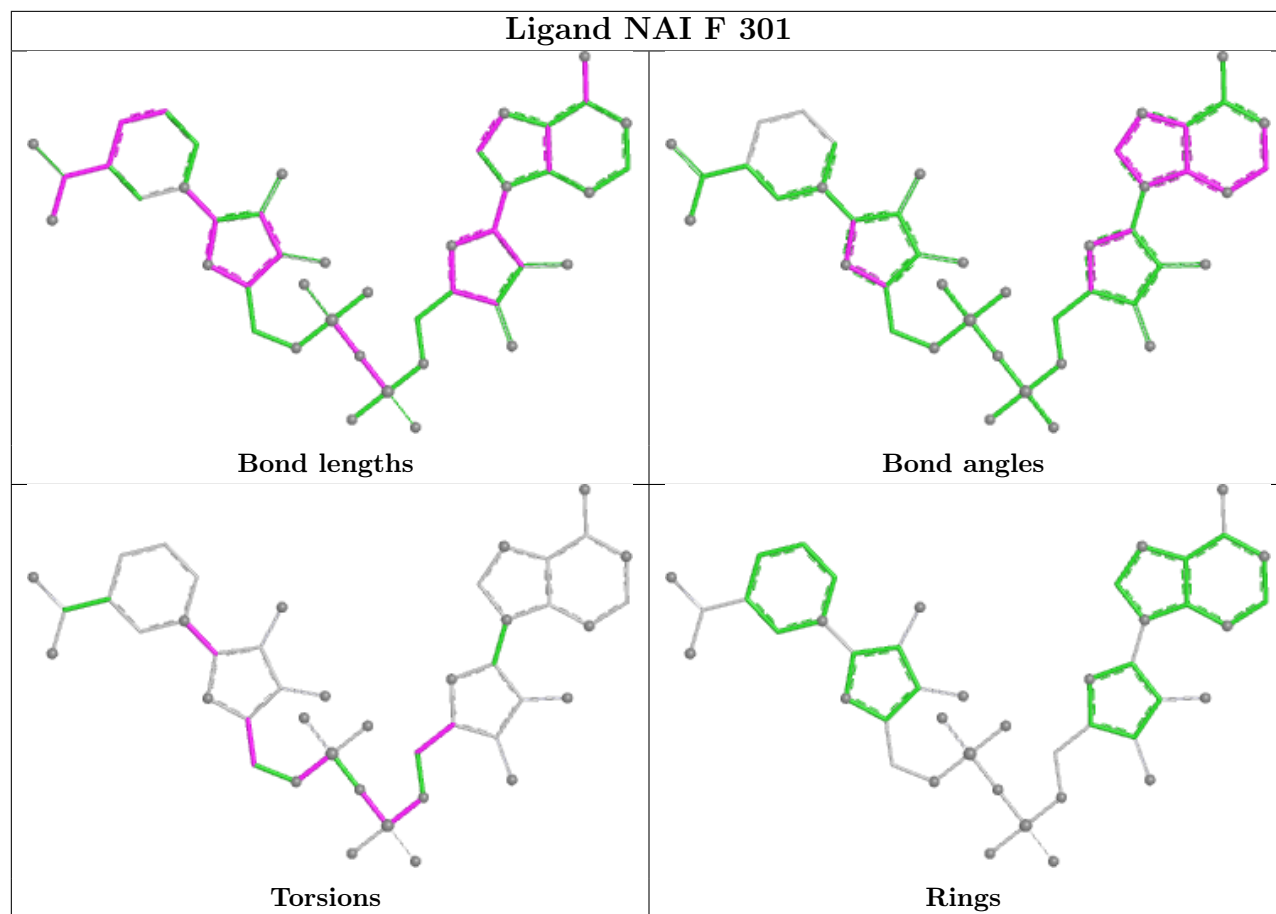




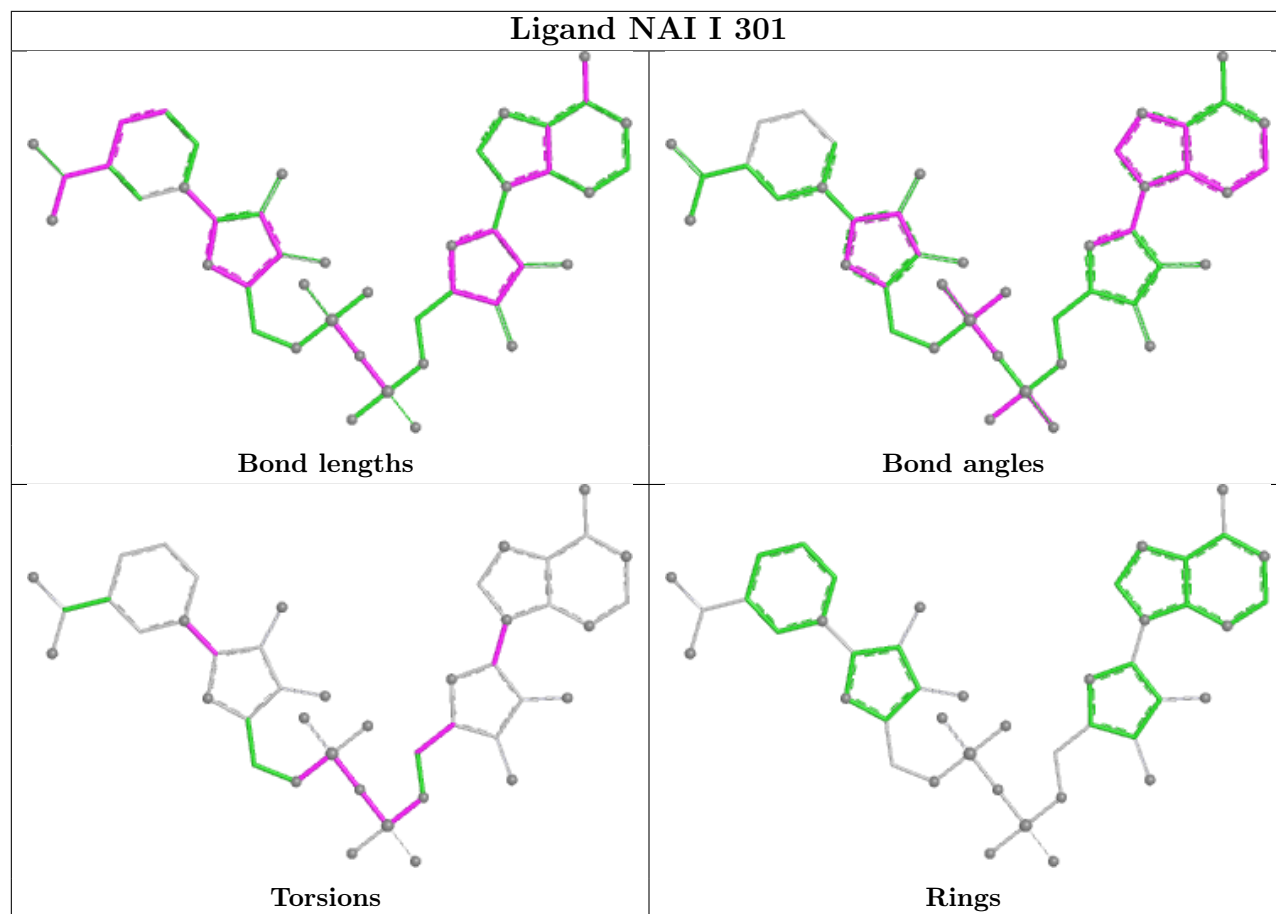


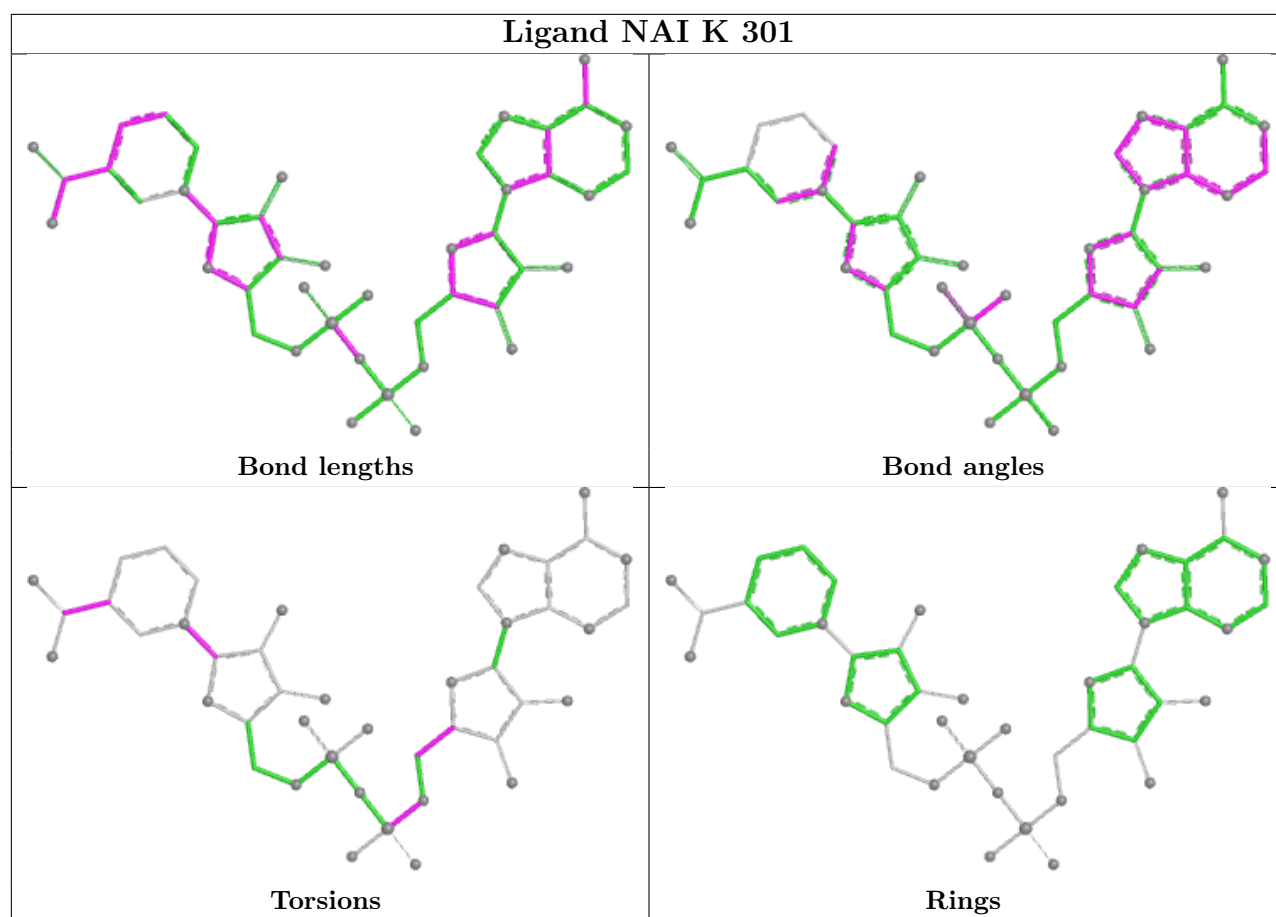


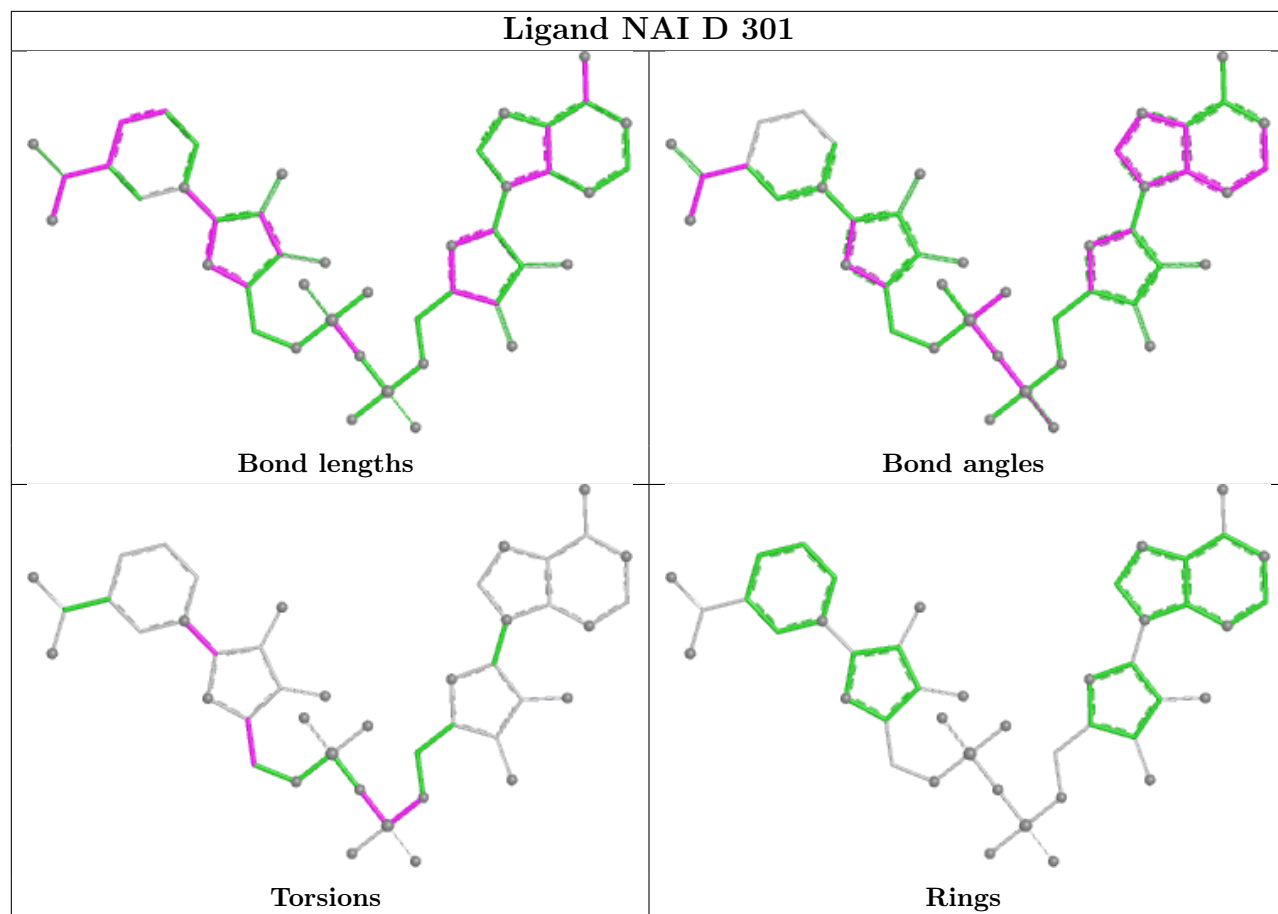
Ligand NAI F 301



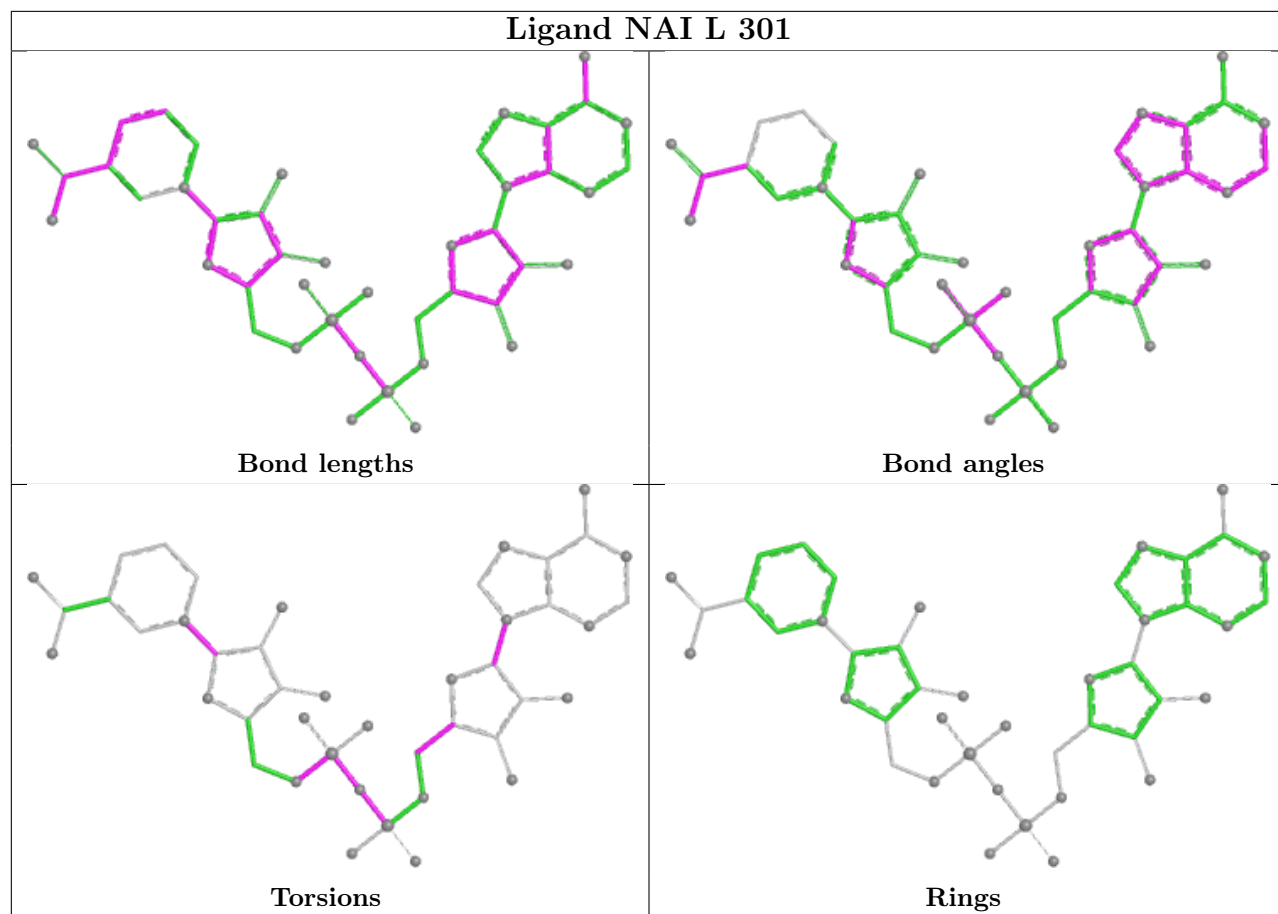
Ligand NAI I 301

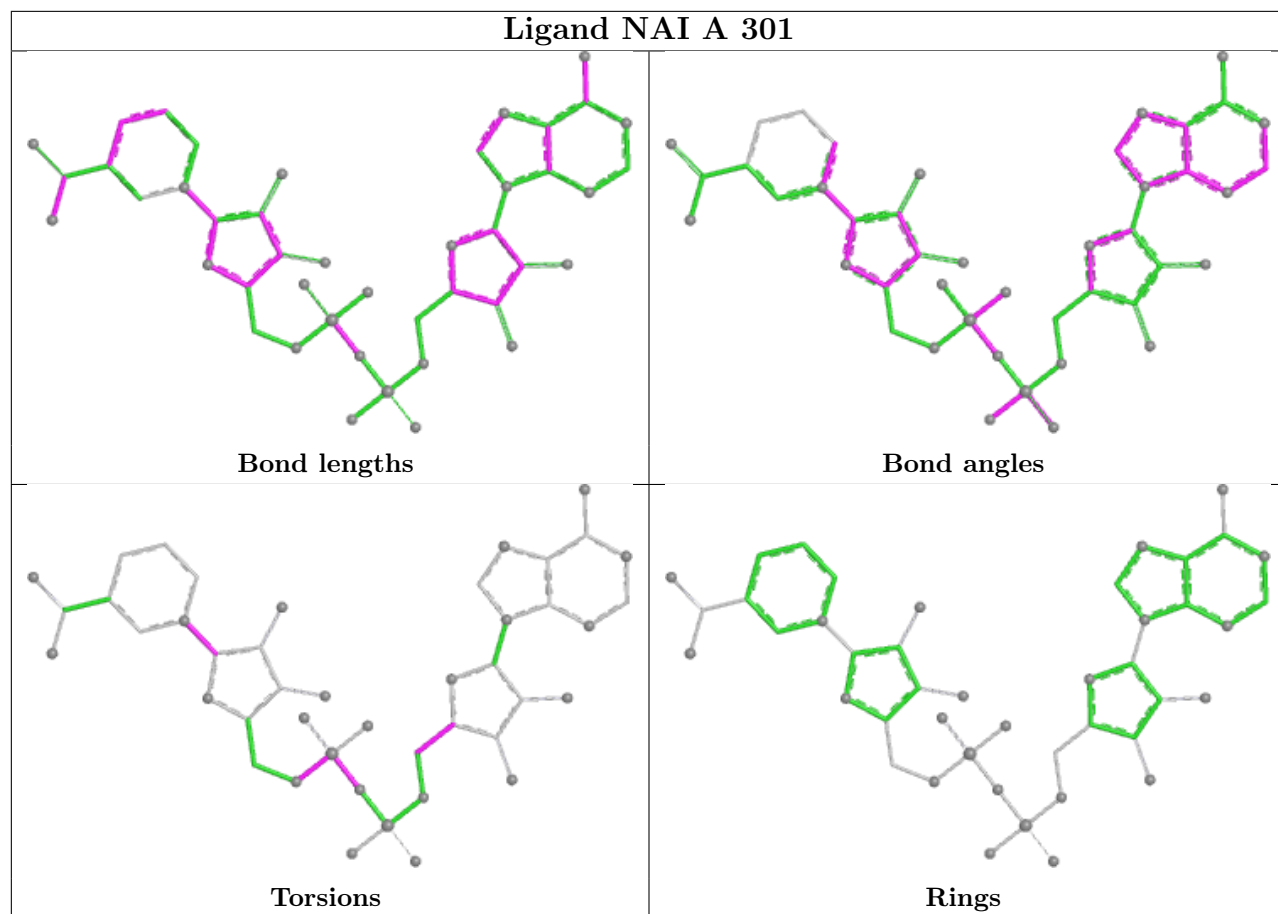




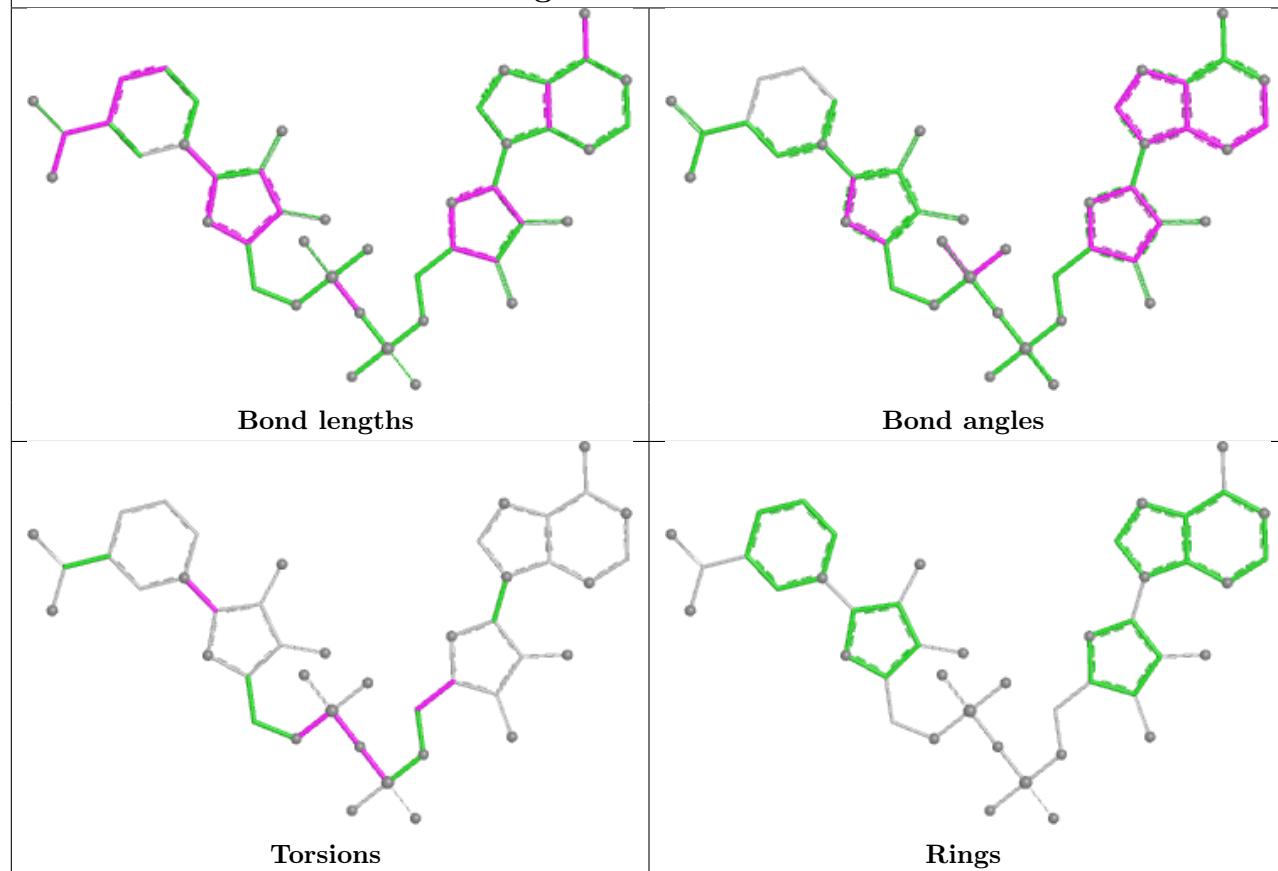


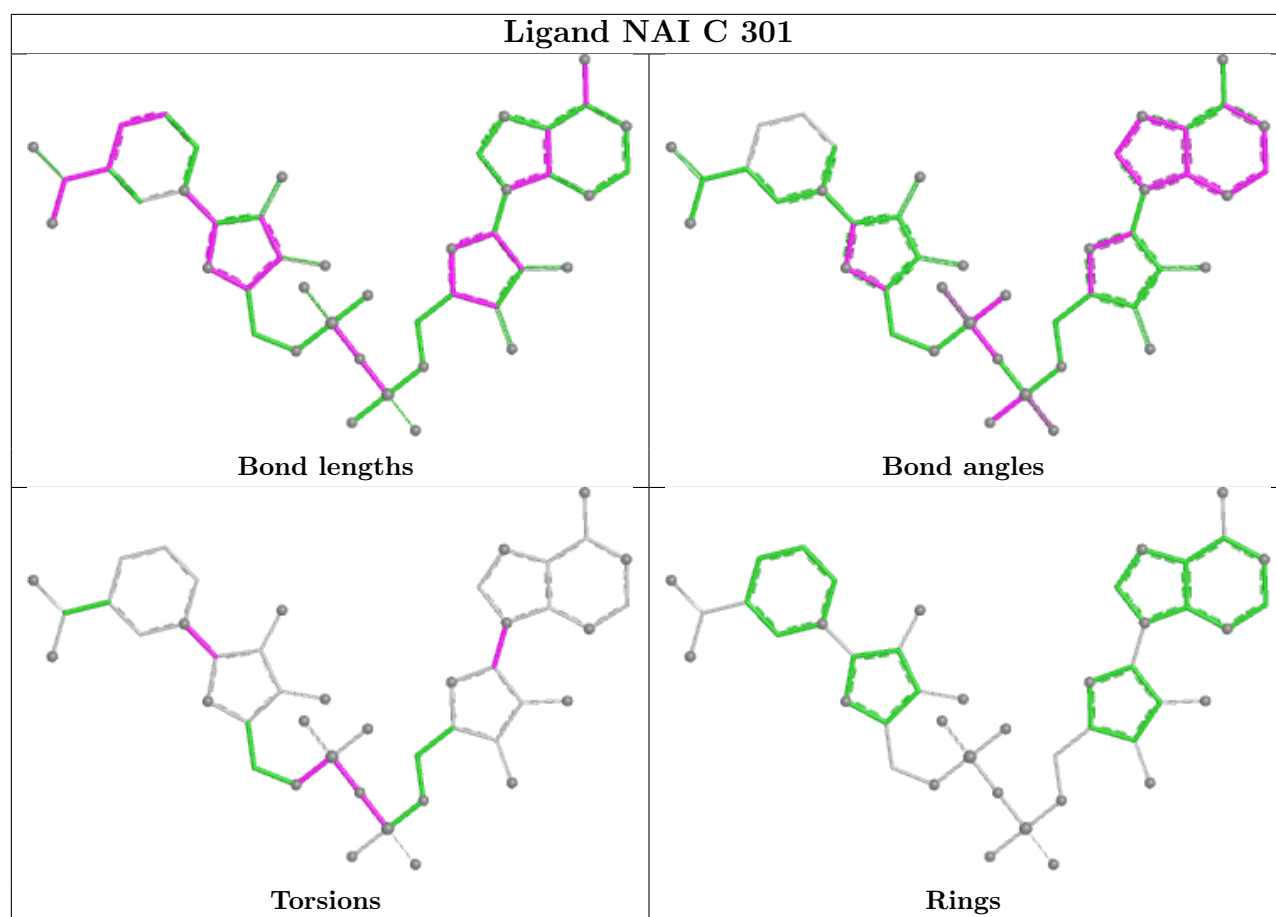
Ligand NAI L 301





Ligand NAI H 301





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	248/263 (94%)	-0.34	1 (0%) 88 89	24, 32, 48, 64	0
1	B	253/263 (96%)	-0.34	1 (0%) 88 89	22, 31, 47, 72	0
1	C	257/263 (97%)	-0.38	3 (1%) 76 77	20, 30, 54, 96	0
1	D	250/263 (95%)	-0.36	2 (0%) 82 83	22, 30, 50, 79	0
1	E	248/263 (94%)	-0.32	3 (1%) 76 77	24, 33, 52, 73	0
1	F	253/263 (96%)	-0.32	3 (1%) 76 77	22, 32, 52, 90	0
1	G	257/263 (97%)	-0.28	1 (0%) 88 89	22, 33, 51, 90	0
1	H	252/263 (95%)	-0.24	4 (1%) 70 72	23, 34, 61, 100	0
1	I	243/263 (92%)	-0.06	1 (0%) 88 89	26, 41, 61, 72	1 (0%)
1	J	253/263 (96%)	-0.31	3 (1%) 76 77	24, 32, 49, 78	0
1	K	257/263 (97%)	-0.32	3 (1%) 76 77	23, 31, 51, 90	0
1	L	246/263 (93%)	-0.30	1 (0%) 88 89	24, 34, 51, 75	0
All	All	3017/3156 (95%)	-0.30	26 (0%) 81 82	20, 32, 55, 100	1 (0%)

The worst 5 of 26 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	207	LEU	3.9
1	H	206	PRO	3.6
1	K	207	LEU	3.6
1	C	207	LEU	3.1
1	G	207	LEU	3.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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(Å ²)	Q<0.9
3	IPA	G	303	4/4	0.86	0.14	36,40,40,48	0
2	NAI	G	301	44/44	0.88	0.09	37,46,53,62	0
3	IPA	K	302	4/4	0.88	0.12	23,37,40,43	0
3	IPA	B	302	4/4	0.90	0.13	29,30,30,31	0
3	IPA	G	302	4/4	0.91	0.11	33,36,39,40	0
3	IPA	B	303	4/4	0.91	0.10	31,36,37,43	0
3	IPA	C	302	4/4	0.91	0.12	31,31,35,35	0
3	IPA	K	303	4/4	0.91	0.10	27,27,28,32	0
3	IPA	H	304	4/4	0.92	0.12	30,31,36,37	0
2	NAI	K	301	44/44	0.92	0.07	31,41,55,61	0
2	NAI	C	301	44/44	0.92	0.07	31,38,44,51	0
3	IPA	H	303	4/4	0.93	0.14	30,31,33,37	0
3	IPA	C	304	4/4	0.93	0.10	31,38,38,40	0
3	IPA	J	302	4/4	0.93	0.10	36,38,41,45	0
2	NAI	F	301	44/44	0.93	0.08	30,37,45,48	0
2	NAI	J	301	44/44	0.93	0.07	29,38,43,46	0
3	IPA	E	304	4/4	0.94	0.10	30,33,33,39	0
2	NAI	I	301	44/44	0.94	0.07	38,48,54,60	0
3	IPA	A	302	4/4	0.94	0.11	31,33,36,37	0
2	NAI	D	301	44/44	0.94	0.07	24,35,39,45	0
3	IPA	L	303	4/4	0.94	0.11	25,29,33,34	0
4	MG	C	303	1/1	0.94	0.04	38,38,38,38	0
2	NAI	L	301	44/44	0.95	0.06	27,35,41,42	0
2	NAI	A	301	44/44	0.95	0.06	26,35,40,48	0
3	IPA	I	303	4/4	0.95	0.08	34,35,39,41	0
3	IPA	D	302	4/4	0.95	0.10	26,28,29,30	0
3	IPA	E	303	4/4	0.95	0.10	28,29,30,31	0
3	IPA	A	303	4/4	0.95	0.09	35,35,36,46	0
2	NAI	B	301	44/44	0.95	0.06	29,39,44,52	0
2	NAI	H	301	44/44	0.95	0.06	28,37,43,45	0
3	IPA	F	302	4/4	0.96	0.09	27,28,29,30	0
3	IPA	I	304	4/4	0.97	0.07	32,35,36,38	0

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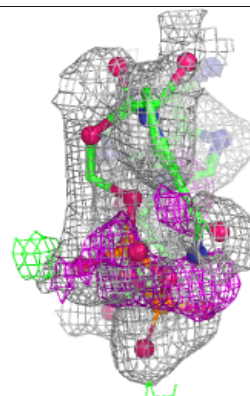
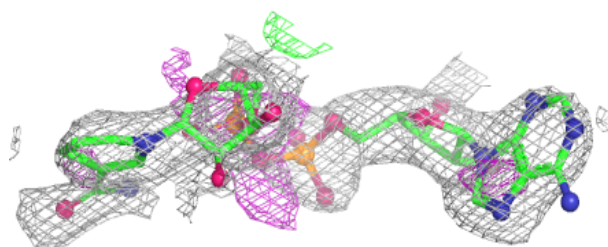
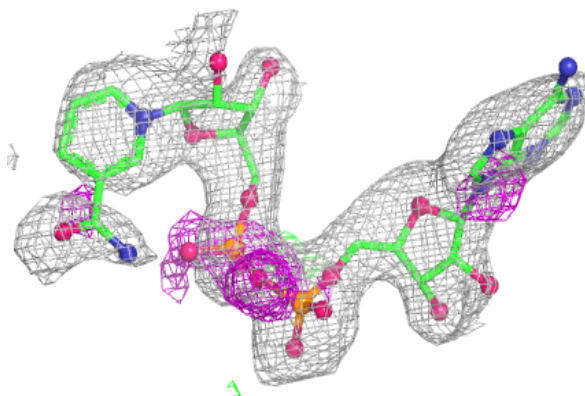
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	IPA	F	303	4/4	0.97	0.07	22,30,31,32	0
3	IPA	L	304	4/4	0.97	0.15	32,32,34,34	0
2	NAI	E	301	44/44	0.97	0.05	29,36,41,48	0
4	MG	I	302	1/1	0.97	0.10	37,37,37,37	0
4	MG	E	302	1/1	0.99	0.03	30,30,30,30	0
4	MG	H	302	1/1	0.99	0.04	28,28,28,28	0
4	MG	D	303	1/1	0.99	0.02	36,36,36,36	0
4	MG	L	302	1/1	0.99	0.02	33,33,33,33	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.

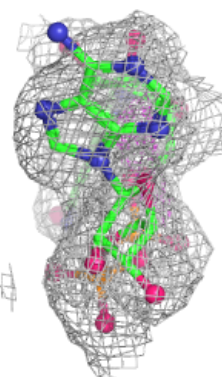
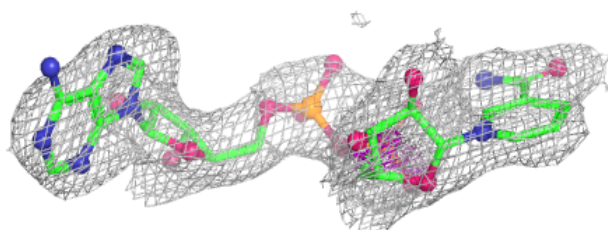
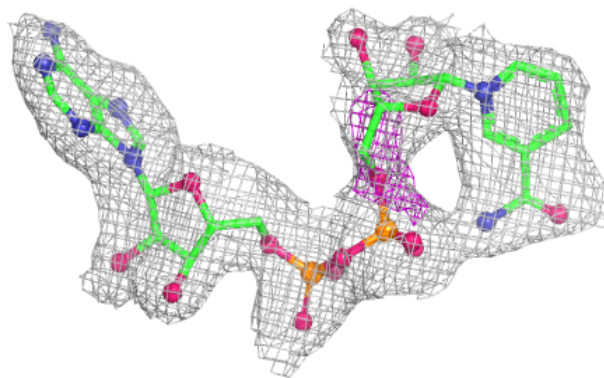
Electron density around NAI G 301:

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

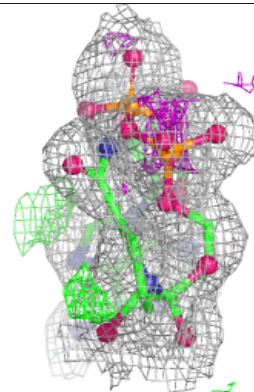
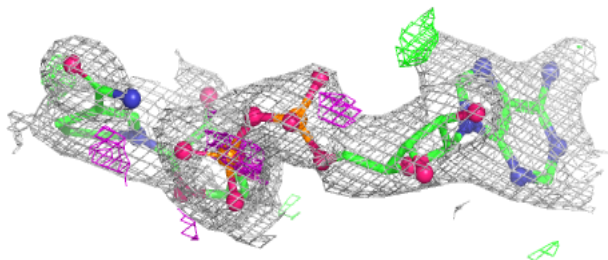
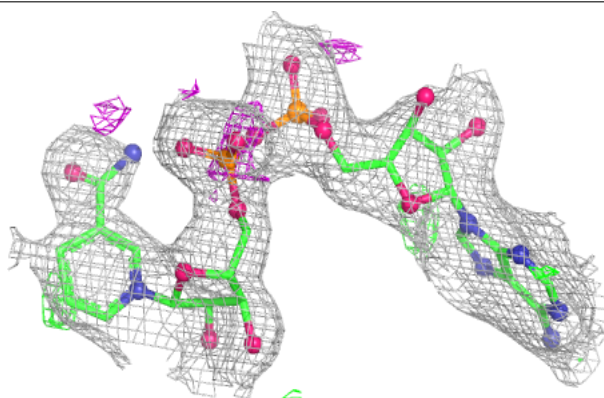


Electron density around NAI K 301:

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

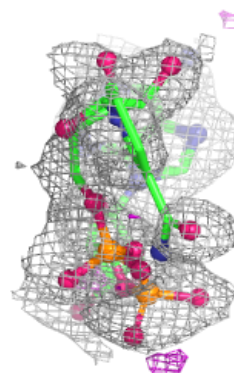
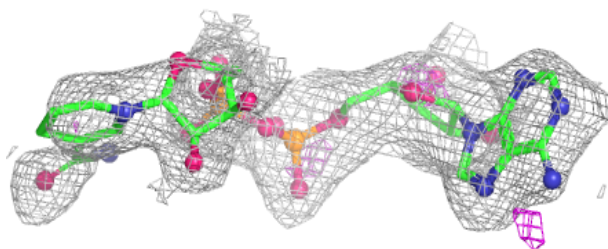
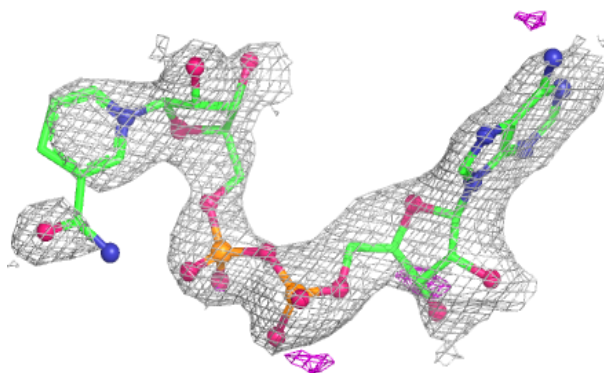
**Electron density around NAI C 301:**

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

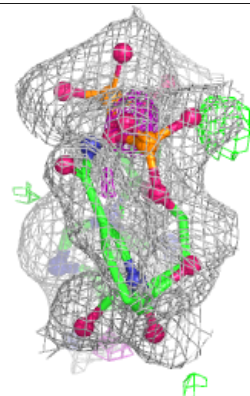
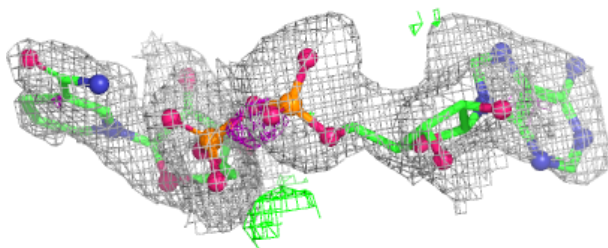
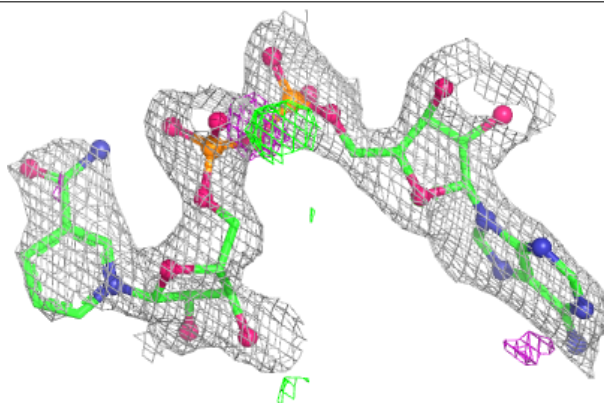


Electron density around NAI F 301:

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

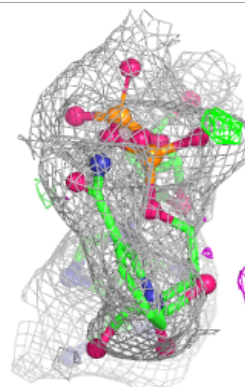
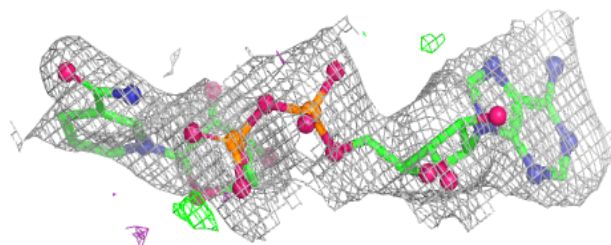
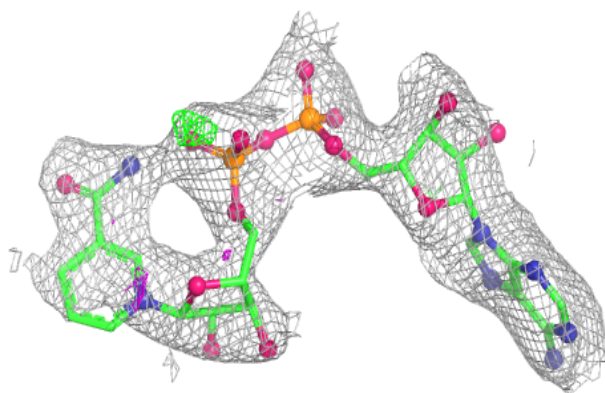
**Electron density around NAI J 301:**

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

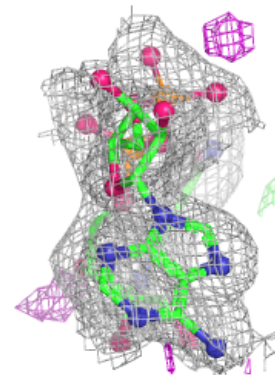
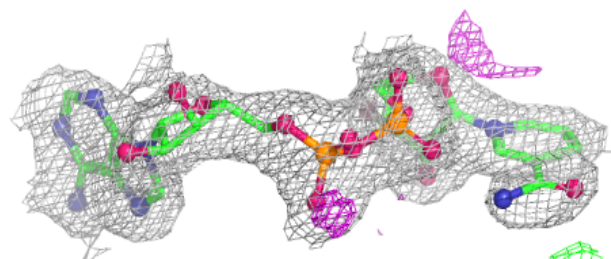
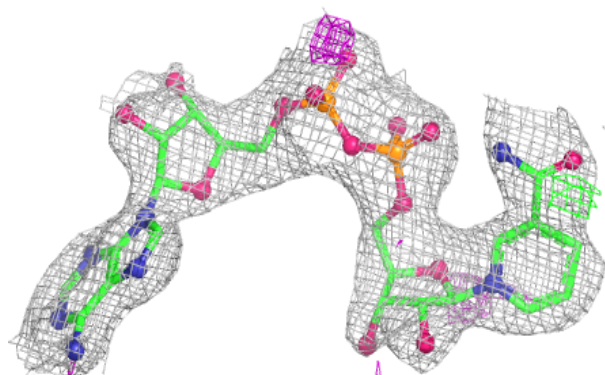


Electron density around NAI I 301:

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

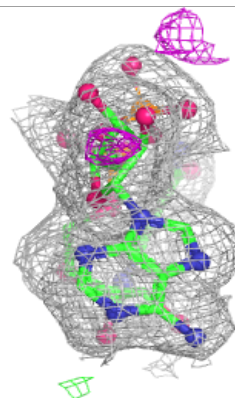
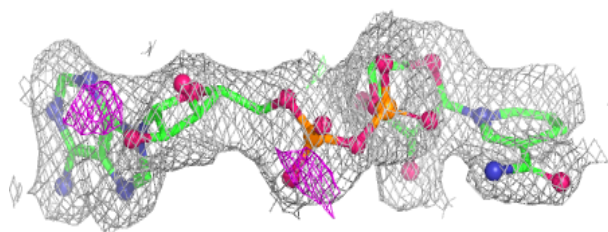
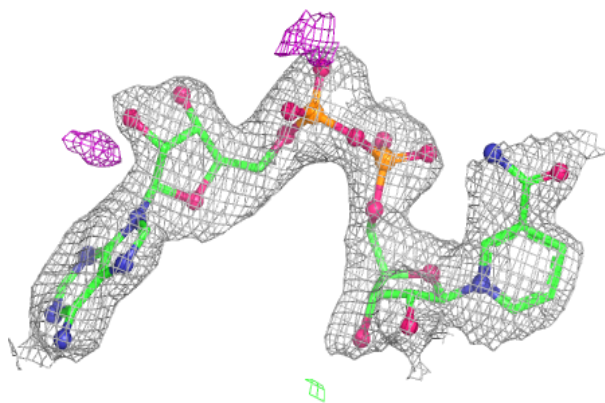
**Electron density around NAI D 301:**

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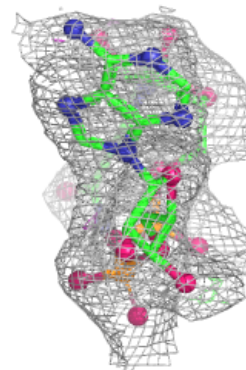
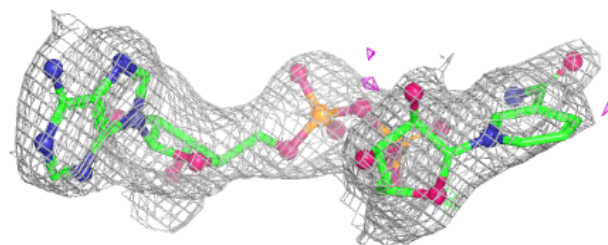
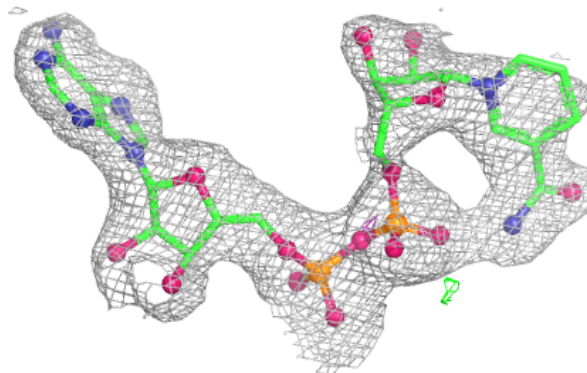


Electron density around NAI L 301:

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and green (positive)

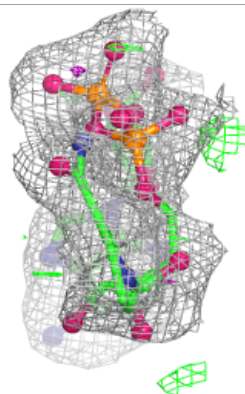
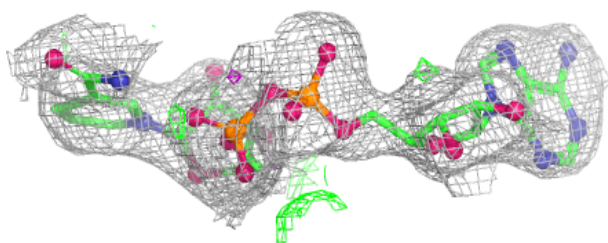
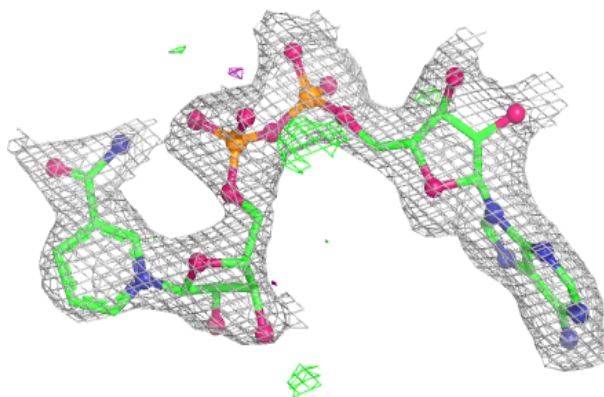
**Electron density around NAI A 301:**

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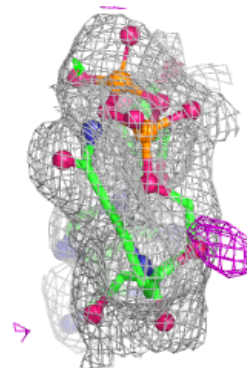
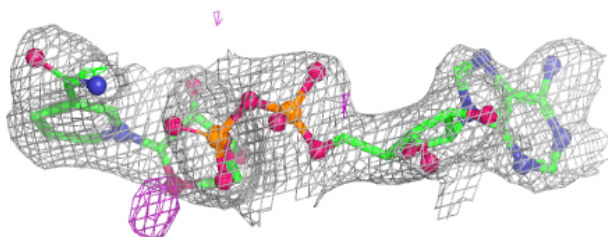
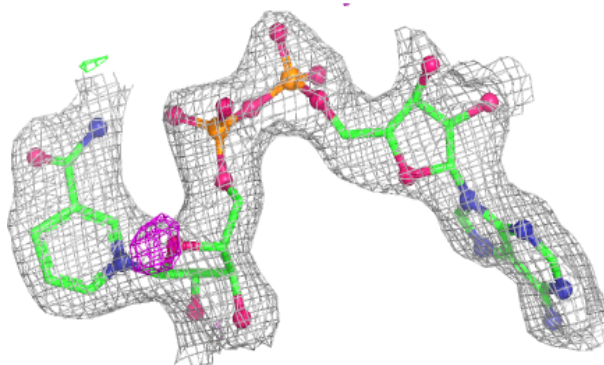


Electron density around NAI B 301:

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and green (positive)

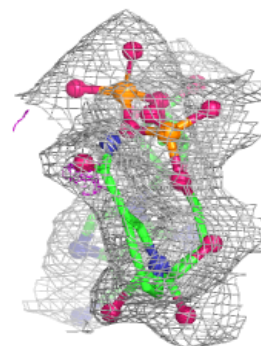
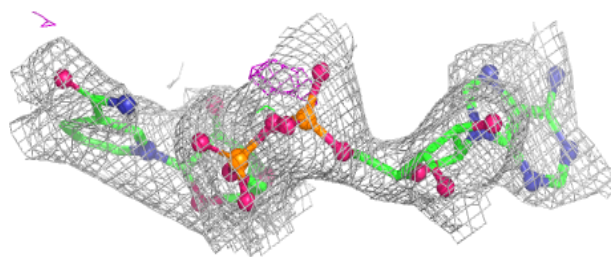
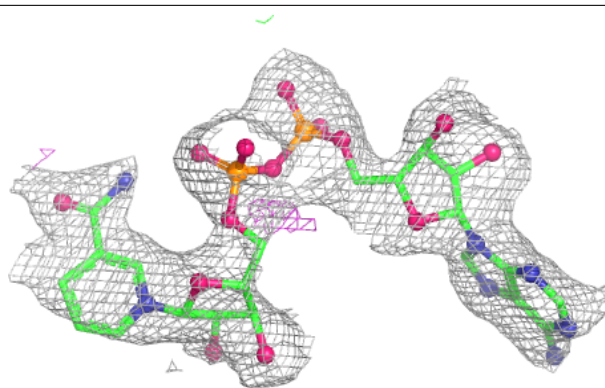
**Electron density around NAI H 301:**

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and green (positive)



Electron density around NAI E 301:

$2mF_o-DF_c$ (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.