



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 08:44 PM UTC

PDB ID : 4BNN / pdb_00004bnn
Title : Crystal structure of S. aureus FabI in complex with NADP and 2-(2-cyanophenoxy)-5-hexylphenol
Authors : Schiebel, J.; Chang, A.; Bommineni, G.R.; Tonge, P.J.; Kisker, C.
Deposited on : 2013-05-15
Resolution : 2.25 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

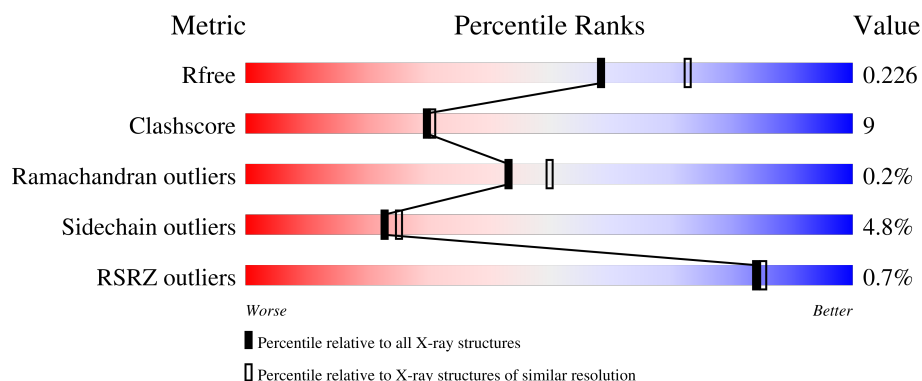
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

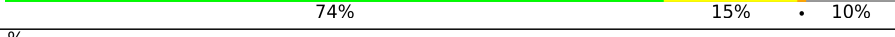
The reported resolution of this entry is 2.25 Å.

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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	282	
1	B	282	
1	C	282	
1	D	282	
1	E	282	

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Mol	Chain	Length	Quality of chain
1	F	282	 % 73% 16% • 10%
1	G	282	 % 77% 12% • 10%
1	H	282	 % 74% 15% • 10%

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
2	GLU	A	1001	-	-	X	-
2	GLU	B	1001	-	-	X	-
2	GLU	E	1001	-	-	X	-
2	GLU	F	1001	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17807 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 ENOYL-[ACYL-CARRIER-PROTEIN] REDUCTASE [NADPH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	254	Total	C	N	O	S	0	13	0
			2052	1289	359	399	5			
1	B	255	Total	C	N	O	S	0	12	0
			2051	1289	357	400	5			
1	C	255	Total	C	N	O	S	0	6	0
			2000	1260	347	389	4			
1	D	254	Total	C	N	O	S	0	3	0
			1971	1241	340	386	4			
1	E	255	Total	C	N	O	S	0	13	0
			2059	1294	360	400	5			
1	F	255	Total	C	N	O	S	0	11	0
			2045	1285	357	398	5			
1	G	255	Total	C	N	O	S	0	5	0
			1992	1255	344	389	4			
1	H	254	Total	C	N	O	S	0	4	0
			1979	1246	343	386	4			

There are 216 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-25	MET	-	expression tag	UNP Q7A6D8
A	-24	LYS	-	expression tag	UNP Q7A6D8
A	-23	HIS	-	expression tag	UNP Q7A6D8
A	-22	HIS	-	expression tag	UNP Q7A6D8
A	-21	HIS	-	expression tag	UNP Q7A6D8
A	-20	HIS	-	expression tag	UNP Q7A6D8
A	-19	HIS	-	expression tag	UNP Q7A6D8
A	-18	HIS	-	expression tag	UNP Q7A6D8
A	-17	PRO	-	expression tag	UNP Q7A6D8
A	-16	MET	-	expression tag	UNP Q7A6D8
A	-15	SER	-	expression tag	UNP Q7A6D8
A	-14	ASP	-	expression tag	UNP Q7A6D8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	TYR	-	expression tag	UNP Q7A6D8
A	-12	ASP	-	expression tag	UNP Q7A6D8
A	-11	ILE	-	expression tag	UNP Q7A6D8
A	-10	PRO	-	expression tag	UNP Q7A6D8
A	-9	THR	-	expression tag	UNP Q7A6D8
A	-8	THR	-	expression tag	UNP Q7A6D8
A	-7	GLU	-	expression tag	UNP Q7A6D8
A	-6	ASN	-	expression tag	UNP Q7A6D8
A	-5	LEU	-	expression tag	UNP Q7A6D8
A	-4	TYR	-	expression tag	UNP Q7A6D8
A	-3	PHE	-	expression tag	UNP Q7A6D8
A	-2	GLN	-	expression tag	UNP Q7A6D8
A	-1	GLY	-	expression tag	UNP Q7A6D8
A	0	ALA	-	expression tag	UNP Q7A6D8
A	2	VAL	LEU	engineered mutation	UNP Q7A6D8
B	-25	MET	-	expression tag	UNP Q7A6D8
B	-24	LYS	-	expression tag	UNP Q7A6D8
B	-23	HIS	-	expression tag	UNP Q7A6D8
B	-22	HIS	-	expression tag	UNP Q7A6D8
B	-21	HIS	-	expression tag	UNP Q7A6D8
B	-20	HIS	-	expression tag	UNP Q7A6D8
B	-19	HIS	-	expression tag	UNP Q7A6D8
B	-18	HIS	-	expression tag	UNP Q7A6D8
B	-17	PRO	-	expression tag	UNP Q7A6D8
B	-16	MET	-	expression tag	UNP Q7A6D8
B	-15	SER	-	expression tag	UNP Q7A6D8
B	-14	ASP	-	expression tag	UNP Q7A6D8
B	-13	TYR	-	expression tag	UNP Q7A6D8
B	-12	ASP	-	expression tag	UNP Q7A6D8
B	-11	ILE	-	expression tag	UNP Q7A6D8
B	-10	PRO	-	expression tag	UNP Q7A6D8
B	-9	THR	-	expression tag	UNP Q7A6D8
B	-8	THR	-	expression tag	UNP Q7A6D8
B	-7	GLU	-	expression tag	UNP Q7A6D8
B	-6	ASN	-	expression tag	UNP Q7A6D8
B	-5	LEU	-	expression tag	UNP Q7A6D8
B	-4	TYR	-	expression tag	UNP Q7A6D8
B	-3	PHE	-	expression tag	UNP Q7A6D8
B	-2	GLN	-	expression tag	UNP Q7A6D8
B	-1	GLY	-	expression tag	UNP Q7A6D8
B	0	ALA	-	expression tag	UNP Q7A6D8
B	2	VAL	LEU	engineered mutation	UNP Q7A6D8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-25	MET	-	expression tag	UNP Q7A6D8
C	-24	LYS	-	expression tag	UNP Q7A6D8
C	-23	HIS	-	expression tag	UNP Q7A6D8
C	-22	HIS	-	expression tag	UNP Q7A6D8
C	-21	HIS	-	expression tag	UNP Q7A6D8
C	-20	HIS	-	expression tag	UNP Q7A6D8
C	-19	HIS	-	expression tag	UNP Q7A6D8
C	-18	HIS	-	expression tag	UNP Q7A6D8
C	-17	PRO	-	expression tag	UNP Q7A6D8
C	-16	MET	-	expression tag	UNP Q7A6D8
C	-15	SER	-	expression tag	UNP Q7A6D8
C	-14	ASP	-	expression tag	UNP Q7A6D8
C	-13	TYR	-	expression tag	UNP Q7A6D8
C	-12	ASP	-	expression tag	UNP Q7A6D8
C	-11	ILE	-	expression tag	UNP Q7A6D8
C	-10	PRO	-	expression tag	UNP Q7A6D8
C	-9	THR	-	expression tag	UNP Q7A6D8
C	-8	THR	-	expression tag	UNP Q7A6D8
C	-7	GLU	-	expression tag	UNP Q7A6D8
C	-6	ASN	-	expression tag	UNP Q7A6D8
C	-5	LEU	-	expression tag	UNP Q7A6D8
C	-4	TYR	-	expression tag	UNP Q7A6D8
C	-3	PHE	-	expression tag	UNP Q7A6D8
C	-2	GLN	-	expression tag	UNP Q7A6D8
C	-1	GLY	-	expression tag	UNP Q7A6D8
C	0	ALA	-	expression tag	UNP Q7A6D8
C	2	VAL	LEU	engineered mutation	UNP Q7A6D8
D	-25	MET	-	expression tag	UNP Q7A6D8
D	-24	LYS	-	expression tag	UNP Q7A6D8
D	-23	HIS	-	expression tag	UNP Q7A6D8
D	-22	HIS	-	expression tag	UNP Q7A6D8
D	-21	HIS	-	expression tag	UNP Q7A6D8
D	-20	HIS	-	expression tag	UNP Q7A6D8
D	-19	HIS	-	expression tag	UNP Q7A6D8
D	-18	HIS	-	expression tag	UNP Q7A6D8
D	-17	PRO	-	expression tag	UNP Q7A6D8
D	-16	MET	-	expression tag	UNP Q7A6D8
D	-15	SER	-	expression tag	UNP Q7A6D8
D	-14	ASP	-	expression tag	UNP Q7A6D8
D	-13	TYR	-	expression tag	UNP Q7A6D8
D	-12	ASP	-	expression tag	UNP Q7A6D8
D	-11	ILE	-	expression tag	UNP Q7A6D8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-10	PRO	-	expression tag	UNP Q7A6D8
D	-9	THR	-	expression tag	UNP Q7A6D8
D	-8	THR	-	expression tag	UNP Q7A6D8
D	-7	GLU	-	expression tag	UNP Q7A6D8
D	-6	ASN	-	expression tag	UNP Q7A6D8
D	-5	LEU	-	expression tag	UNP Q7A6D8
D	-4	TYR	-	expression tag	UNP Q7A6D8
D	-3	PHE	-	expression tag	UNP Q7A6D8
D	-2	GLN	-	expression tag	UNP Q7A6D8
D	-1	GLY	-	expression tag	UNP Q7A6D8
D	0	ALA	-	expression tag	UNP Q7A6D8
D	2	VAL	LEU	engineered mutation	UNP Q7A6D8
E	-25	MET	-	expression tag	UNP Q7A6D8
E	-24	LYS	-	expression tag	UNP Q7A6D8
E	-23	HIS	-	expression tag	UNP Q7A6D8
E	-22	HIS	-	expression tag	UNP Q7A6D8
E	-21	HIS	-	expression tag	UNP Q7A6D8
E	-20	HIS	-	expression tag	UNP Q7A6D8
E	-19	HIS	-	expression tag	UNP Q7A6D8
E	-18	HIS	-	expression tag	UNP Q7A6D8
E	-17	PRO	-	expression tag	UNP Q7A6D8
E	-16	MET	-	expression tag	UNP Q7A6D8
E	-15	SER	-	expression tag	UNP Q7A6D8
E	-14	ASP	-	expression tag	UNP Q7A6D8
E	-13	TYR	-	expression tag	UNP Q7A6D8
E	-12	ASP	-	expression tag	UNP Q7A6D8
E	-11	ILE	-	expression tag	UNP Q7A6D8
E	-10	PRO	-	expression tag	UNP Q7A6D8
E	-9	THR	-	expression tag	UNP Q7A6D8
E	-8	THR	-	expression tag	UNP Q7A6D8
E	-7	GLU	-	expression tag	UNP Q7A6D8
E	-6	ASN	-	expression tag	UNP Q7A6D8
E	-5	LEU	-	expression tag	UNP Q7A6D8
E	-4	TYR	-	expression tag	UNP Q7A6D8
E	-3	PHE	-	expression tag	UNP Q7A6D8
E	-2	GLN	-	expression tag	UNP Q7A6D8
E	-1	GLY	-	expression tag	UNP Q7A6D8
E	0	ALA	-	expression tag	UNP Q7A6D8
E	2	VAL	LEU	engineered mutation	UNP Q7A6D8
F	-25	MET	-	expression tag	UNP Q7A6D8
F	-24	LYS	-	expression tag	UNP Q7A6D8
F	-23	HIS	-	expression tag	UNP Q7A6D8

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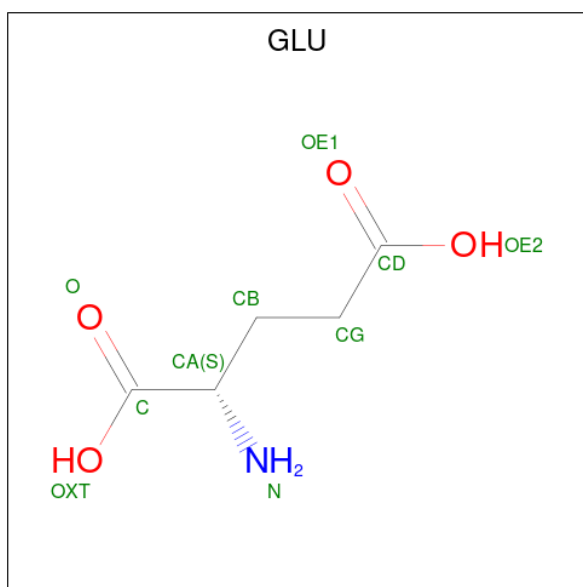
Chain	Residue	Modelled	Actual	Comment	Reference
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F	-21	HIS	-	expression tag	UNP Q7A6D8
F	-20	HIS	-	expression tag	UNP Q7A6D8
F	-19	HIS	-	expression tag	UNP Q7A6D8
F	-18	HIS	-	expression tag	UNP Q7A6D8
F	-17	PRO	-	expression tag	UNP Q7A6D8
F	-16	MET	-	expression tag	UNP Q7A6D8
F	-15	SER	-	expression tag	UNP Q7A6D8
F	-14	ASP	-	expression tag	UNP Q7A6D8
F	-13	TYR	-	expression tag	UNP Q7A6D8
F	-12	ASP	-	expression tag	UNP Q7A6D8
F	-11	ILE	-	expression tag	UNP Q7A6D8
F	-10	PRO	-	expression tag	UNP Q7A6D8
F	-9	THR	-	expression tag	UNP Q7A6D8
F	-8	THR	-	expression tag	UNP Q7A6D8
F	-7	GLU	-	expression tag	UNP Q7A6D8
F	-6	ASN	-	expression tag	UNP Q7A6D8
F	-5	LEU	-	expression tag	UNP Q7A6D8
F	-4	TYR	-	expression tag	UNP Q7A6D8
F	-3	PHE	-	expression tag	UNP Q7A6D8
F	-2	GLN	-	expression tag	UNP Q7A6D8
F	-1	GLY	-	expression tag	UNP Q7A6D8
F	0	ALA	-	expression tag	UNP Q7A6D8
F	2	VAL	LEU	engineered mutation	UNP Q7A6D8
G	-25	MET	-	expression tag	UNP Q7A6D8
G	-24	LYS	-	expression tag	UNP Q7A6D8
G	-23	HIS	-	expression tag	UNP Q7A6D8
G	-22	HIS	-	expression tag	UNP Q7A6D8
G	-21	HIS	-	expression tag	UNP Q7A6D8
G	-20	HIS	-	expression tag	UNP Q7A6D8
G	-19	HIS	-	expression tag	UNP Q7A6D8
G	-18	HIS	-	expression tag	UNP Q7A6D8
G	-17	PRO	-	expression tag	UNP Q7A6D8
G	-16	MET	-	expression tag	UNP Q7A6D8
G	-15	SER	-	expression tag	UNP Q7A6D8
G	-14	ASP	-	expression tag	UNP Q7A6D8
G	-13	TYR	-	expression tag	UNP Q7A6D8
G	-12	ASP	-	expression tag	UNP Q7A6D8
G	-11	ILE	-	expression tag	UNP Q7A6D8
G	-10	PRO	-	expression tag	UNP Q7A6D8
G	-9	THR	-	expression tag	UNP Q7A6D8
G	-8	THR	-	expression tag	UNP Q7A6D8

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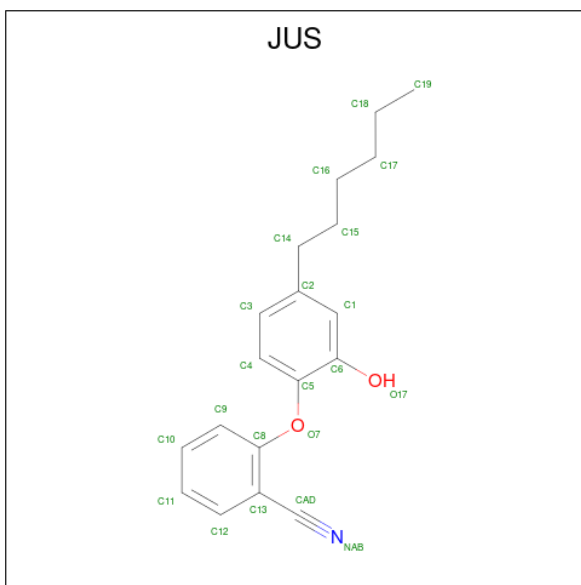
Chain	Residue	Modelled	Actual	Comment	Reference
G	-7	GLU	-	expression tag	UNP Q7A6D8
G	-6	ASN	-	expression tag	UNP Q7A6D8
G	-5	LEU	-	expression tag	UNP Q7A6D8
G	-4	TYR	-	expression tag	UNP Q7A6D8
G	-3	PHE	-	expression tag	UNP Q7A6D8
G	-2	GLN	-	expression tag	UNP Q7A6D8
G	-1	GLY	-	expression tag	UNP Q7A6D8
G	0	ALA	-	expression tag	UNP Q7A6D8
G	2	VAL	LEU	engineered mutation	UNP Q7A6D8
H	-25	MET	-	expression tag	UNP Q7A6D8
H	-24	LYS	-	expression tag	UNP Q7A6D8
H	-23	HIS	-	expression tag	UNP Q7A6D8
H	-22	HIS	-	expression tag	UNP Q7A6D8
H	-21	HIS	-	expression tag	UNP Q7A6D8
H	-20	HIS	-	expression tag	UNP Q7A6D8
H	-19	HIS	-	expression tag	UNP Q7A6D8
H	-18	HIS	-	expression tag	UNP Q7A6D8
H	-17	PRO	-	expression tag	UNP Q7A6D8
H	-16	MET	-	expression tag	UNP Q7A6D8
H	-15	SER	-	expression tag	UNP Q7A6D8
H	-14	ASP	-	expression tag	UNP Q7A6D8
H	-13	TYR	-	expression tag	UNP Q7A6D8
H	-12	ASP	-	expression tag	UNP Q7A6D8
H	-11	ILE	-	expression tag	UNP Q7A6D8
H	-10	PRO	-	expression tag	UNP Q7A6D8
H	-9	THR	-	expression tag	UNP Q7A6D8
H	-8	THR	-	expression tag	UNP Q7A6D8
H	-7	GLU	-	expression tag	UNP Q7A6D8
H	-6	ASN	-	expression tag	UNP Q7A6D8
H	-5	LEU	-	expression tag	UNP Q7A6D8
H	-4	TYR	-	expression tag	UNP Q7A6D8
H	-3	PHE	-	expression tag	UNP Q7A6D8
H	-2	GLN	-	expression tag	UNP Q7A6D8
H	-1	GLY	-	expression tag	UNP Q7A6D8
H	0	ALA	-	expression tag	UNP Q7A6D8
H	2	VAL	LEU	engineered mutation	UNP Q7A6D8

- Molecule 2 is GLUTAMIC ACID (CCD ID: GLU) (formula: $C_5H_9NO_4$).



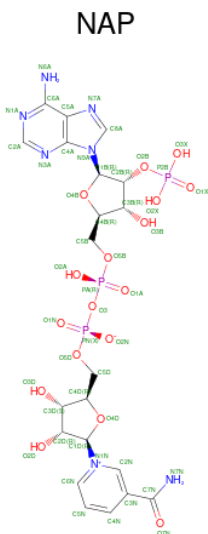
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	5	1	4		
2	B	1	Total	C	N	O	0	0
			10	5	1	4		
2	C	1	Total	C	N	O	0	0
			10	5	1	4		
2	E	1	Total	C	N	O	0	0
			10	5	1	4		
2	F	1	Total	C	N	O	0	0
			10	5	1	4		
2	G	1	Total	C	N	O	0	0
			10	5	1	4		

- Molecule 3 is 2-(2-CYANOPHENOXY)-5-HEXYLPHENOL (CCD ID: JUS) (formula: C₁₉H₂₁NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			22	19	1	2		
3	B	1	Total	C	N	O	0	0
			22	19	1	2		
3	C	1	Total	C	N	O	0	0
			22	19	1	2		
3	D	1	Total	C	N	O	0	0
			22	19	1	2		
3	E	1	Total	C	N	O	0	0
			22	19	1	2		
3	F	1	Total	C	N	O	0	0
			22	19	1	2		
3	G	1	Total	C	N	O	0	0
			22	19	1	2		
3	H	1	Total	C	N	O	0	0
			22	19	1	2		

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	C	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	D	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	E	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	F	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	G	1	Total 48	C 21	N 7	O 17	P 3	0	0
4	H	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	170	Total O 170 170	0	0
5	B	135	Total O 135 135	0	0
5	C	130	Total O 130 130	0	0
5	D	99	Total O 99 99	0	0

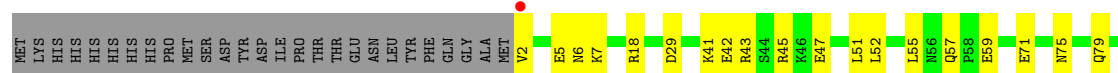
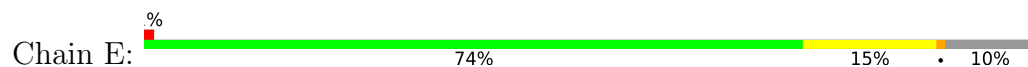
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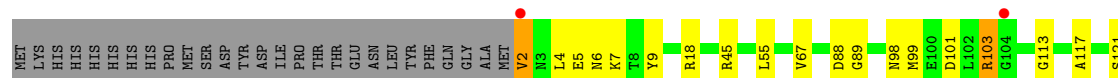
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	144	Total 144	O 144	0	0
5	F	130	Total 130	O 130	0	0
5	G	125	Total 125	O 125	0	0
5	H	105	Total 105	O 105	0	0



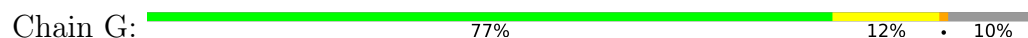
• Molecule 1: ENOYL-[ACYL-CARRIER-PROTEIN] REDUCTASE [NADPH]



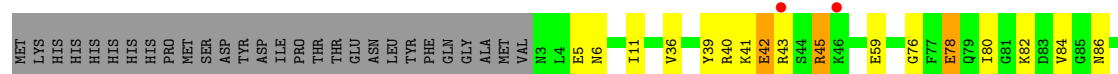
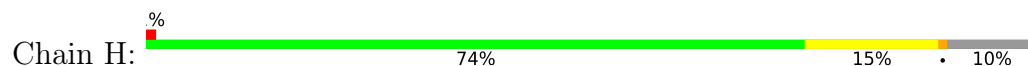
• Molecule 1: ENOYL-[ACYL-CARRIER-PROTEIN] REDUCTASE [NADPH]



• Molecule 1: ENOYL-[ACYL-CARRIER-PROTEIN] REDUCTASE [NADPH]



• Molecule 1: ENOYL-[ACYL-CARRIER-PROTEIN] REDUCTASE [NADPH]



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	89.92Å 94.60Å 94.27Å 97.92° 111.96° 97.31°	Depositor
Resolution (Å)	40.90 – 2.25 40.90 – 2.25	Depositor EDS
% Data completeness (in resolution range)	97.5 (40.90-2.25) 97.5 (40.90-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.151 , 0.206 0.174 , 0.226	Depositor DCC
R_{free} test set	6528 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.574	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17807	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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: JUS, NAP

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	0.68	0/2091	1.09	0/2813
1	B	0.66	0/2087	1.02	0/2809
1	C	0.64	0/2039	1.07	1/2746 (0.0%)
1	D	0.69	0/2001	1.09	2/2696 (0.1%)
1	E	0.66	0/2098	1.09	1/2823 (0.0%)
1	F	0.67	0/2078	1.04	0/2797
1	G	0.67	1/2028 (0.0%)	1.07	3/2732 (0.1%)
1	H	0.63	0/2012	1.11	6/2710 (0.2%)
All	All	0.66	1/16434 (0.0%)	1.07	13/22126 (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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	153	ALA	CA-CB	-7.49	1.43	1.53

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	99	MET	N-CA-C	6.55	119.26	111.33
1	H	165	ALA	N-CA-C	-5.81	104.95	111.28
1	H	179	GLY	O-C-N	5.74	123.14	121.07
1	H	94	ILE	N-CA-C	5.73	116.54	108.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	H	191	GLY	CA-C-N	-5.68	113.85	119.99

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	2	VAL	Peptide

5.2 Too-close contacts [i](#)

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	2052	0	2071	64	0
1	B	2051	0	2067	31	0
1	C	2000	0	2022	33	0
1	D	1971	0	1981	26	0
1	E	2059	0	2080	47	0
1	F	2045	0	2061	50	0
1	G	1992	0	2009	28	0
1	H	1979	0	1994	26	0
2	A	10	0	5	15	0
2	B	10	0	5	7	0
2	C	10	0	5	2	0
2	E	10	0	5	7	0
2	F	10	0	5	12	0
2	G	10	0	5	2	0
3	A	22	0	20	3	0
3	B	22	0	21	2	0
3	C	22	0	20	4	0
3	D	22	0	21	0	0
3	E	22	0	20	1	0
3	F	22	0	20	1	0
3	G	22	0	20	2	0
3	H	22	0	20	2	0
4	A	48	0	25	1	0
4	B	48	0	25	0	0
4	C	48	0	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	48	0	25	0	0
4	E	48	0	25	1	0
4	F	48	0	25	0	0
4	G	48	0	25	0	0
4	H	48	0	25	3	0
5	A	170	0	0	15	0
5	B	135	0	0	3	0
5	C	130	0	0	6	0
5	D	99	0	0	2	0
5	E	144	0	0	10	0
5	F	130	0	0	6	0
5	G	125	0	0	2	0
5	H	105	0	0	6	0
All	All	17807	0	16677	285	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 9.

The worst 5 of 285 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:E:103[A]:ARG:NH1	2:E:1001:GLU:OE2	1.77	1.16
1:F:103[B]:ARG:HH12	2:F:1001:GLU:HG2	1.07	1.16
1:F:103[B]:ARG:NH1	2:F:1001:GLU:HG2	1.62	1.15
1:F:103[B]:ARG:HH12	2:F:1001:GLU:CG	1.62	1.11
2:F:1001:GLU:O	2:F:1001:GLU:HG3	1.51	1.06

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	265/282 (94%)	250 (94%)	12 (4%)	3 (1%)	11	8
1	B	265/282 (94%)	254 (96%)	11 (4%)	0	100	100
1	C	259/282 (92%)	246 (95%)	13 (5%)	0	100	100
1	D	255/282 (90%)	237 (93%)	17 (7%)	1 (0%)	30	31
1	E	266/282 (94%)	255 (96%)	10 (4%)	1 (0%)	30	31
1	F	264/282 (94%)	249 (94%)	15 (6%)	0	100	100
1	G	258/282 (92%)	245 (95%)	13 (5%)	0	100	100
1	H	256/282 (91%)	238 (93%)	18 (7%)	0	100	100
All	All	2088/2256 (93%)	1974 (94%)	109 (5%)	5 (0%)	43	50

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	148	LEU
1	A	99[A]	MET
1	A	99[B]	MET
1	A	148	LEU
1	D	156	ASN

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	219/234 (94%)	213 (97%)	6 (3%)	39	49
1	B	219/234 (94%)	208 (95%)	11 (5%)	22	24
1	C	214/234 (92%)	202 (94%)	12 (6%)	19	20
1	D	210/234 (90%)	200 (95%)	10 (5%)	23	25
1	E	220/234 (94%)	208 (94%)	12 (6%)	19	20
1	F	218/234 (93%)	207 (95%)	11 (5%)	22	24
1	G	213/234 (91%)	205 (96%)	8 (4%)	29	36
1	H	211/234 (90%)	199 (94%)	12 (6%)	18	19

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1724/1872 (92%)	1642 (95%)	82 (5%)	23	25

5 of 82 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	206	THR
1	H	11	ILE
1	F	209	LYS
1	G	112	GLU
1	H	59	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	170	ASN
1	G	6	ASN
1	H	253	HIS
1	G	3	ASN
1	G	156	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](#)

22 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	JUS	H	1257	-	23,23,23	0.58	0	29,29,29	0.90	2 (6%)
4	NAP	F	1258	-	50,52,52	1.57	4 (8%)	71,80,80	1.88	15 (21%)
2	GLU	G	1001	-	8,9,9	1.28	1 (12%)	8,11,11	1.06	0
3	JUS	E	1257	-	23,23,23	0.64	0	29,29,29	0.97	3 (10%)
2	GLU	A	1001	-	8,9,9	1.27	1 (12%)	8,11,11	1.05	0
3	JUS	B	1257	-	23,23,23	0.71	0	29,29,29	0.91	2 (6%)
3	JUS	A	1257	-	23,23,23	0.70	0	29,29,29	1.05	2 (6%)
2	GLU	E	1001	-	8,9,9	1.06	0	8,11,11	1.32	2 (25%)
3	JUS	D	1257	-	23,23,23	0.64	0	29,29,29	1.01	2 (6%)
3	JUS	G	1257	-	23,23,23	0.51	0	29,29,29	1.08	2 (6%)
2	GLU	F	1001	-	8,9,9	1.33	1 (12%)	8,11,11	0.91	0
4	NAP	E	1258	-	50,52,52	1.51	4 (8%)	71,80,80	1.93	15 (21%)
2	GLU	C	1001	-	8,9,9	1.03	0	8,11,11	1.54	2 (25%)
4	NAP	B	1258	-	50,52,52	1.46	4 (8%)	71,80,80	1.98	18 (25%)
4	NAP	A	1258	-	50,52,52	1.58	6 (12%)	71,80,80	1.98	18 (25%)
4	NAP	H	1258	-	50,52,52	1.70	6 (12%)	71,80,80	1.92	17 (23%)
4	NAP	G	1258	-	50,52,52	1.56	7 (14%)	71,80,80	1.78	17 (23%)
4	NAP	C	1258	-	50,52,52	1.60	5 (10%)	71,80,80	1.98	16 (22%)
3	JUS	F	1257	-	23,23,23	0.58	0	29,29,29	1.14	1 (3%)
4	NAP	D	1258	-	50,52,52	1.72	5 (10%)	71,80,80	1.80	13 (18%)
3	JUS	C	1257	-	23,23,23	0.61	0	29,29,29	0.83	1 (3%)
2	GLU	B	1001	-	8,9,9	1.19	0	8,11,11	1.41	1 (12%)

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	JUS	H	1257	-	-	4/12/12/12	0/2/2/2
4	NAP	F	1258	-	-	6/35/67/67	0/5/5/5
2	GLU	G	1001	-	-	7/9/9/9	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	JUS	E	1257	-	-	3/12/12/12	0/2/2/2
2	GLU	A	1001	-	-	6/9/9/9	-
3	JUS	B	1257	-	-	2/12/12/12	0/2/2/2
3	JUS	A	1257	-	-	2/12/12/12	0/2/2/2
2	GLU	E	1001	-	-	8/9/9/9	-
3	JUS	D	1257	-	-	1/12/12/12	0/2/2/2
3	JUS	G	1257	-	-	2/12/12/12	0/2/2/2
2	GLU	F	1001	-	-	4/9/9/9	-
4	NAP	E	1258	-	-	6/35/67/67	0/5/5/5
2	GLU	C	1001	-	-	8/9/9/9	-
4	NAP	B	1258	-	-	7/35/67/67	0/5/5/5
4	NAP	A	1258	-	-	6/35/67/67	0/5/5/5
4	NAP	H	1258	-	-	5/35/67/67	0/5/5/5
4	NAP	G	1258	-	-	8/35/67/67	0/5/5/5
4	NAP	C	1258	-	-	5/35/67/67	0/5/5/5
3	JUS	F	1257	-	-	3/12/12/12	0/2/2/2
4	NAP	D	1258	-	-	7/35/67/67	0/5/5/5
3	JUS	C	1257	-	-	2/12/12/12	0/2/2/2
2	GLU	B	1001	-	-	4/9/9/9	-

The worst 5 of 44 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1258	NAP	O7N-C7N	9.20	1.41	1.24
4	D	1258	NAP	O7N-C7N	9.06	1.41	1.24
4	F	1258	NAP	O7N-C7N	8.60	1.40	1.24
4	C	1258	NAP	O7N-C7N	8.56	1.40	1.24
4	G	1258	NAP	O7N-C7N	8.17	1.39	1.24

The worst 5 of 149 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1258	NAP	N3A-C2A-N1A	-6.58	118.61	128.58
4	D	1258	NAP	N3A-C2A-N1A	-6.48	118.78	128.58
4	F	1258	NAP	N3A-C2A-N1A	-6.33	119.00	128.58
4	C	1258	NAP	N3A-C2A-N1A	-6.20	119.19	128.58
4	A	1258	NAP	N3A-C2A-N1A	-6.18	119.23	128.58

There are no chirality outliers.

5 of 106 torsion outliers are listed below:

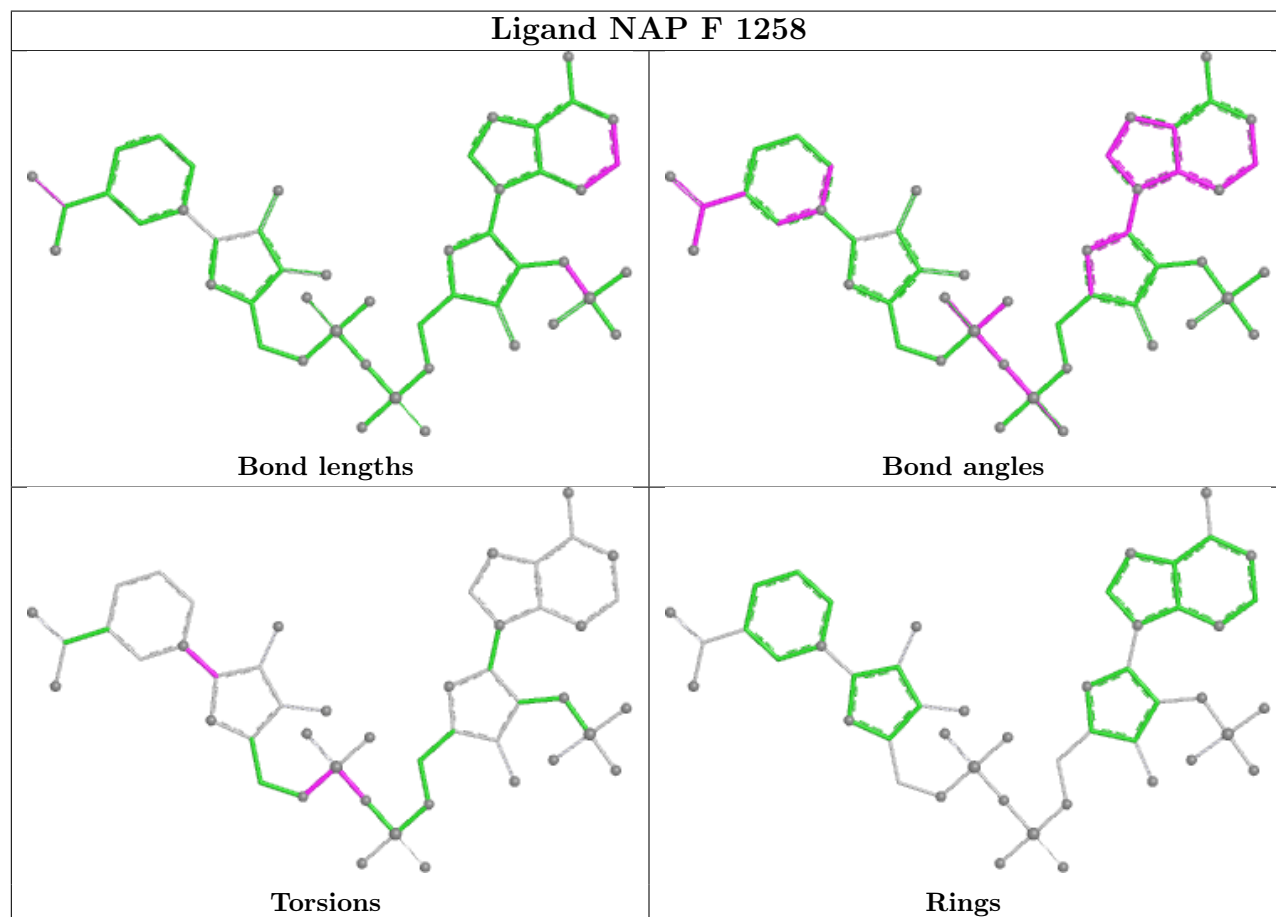
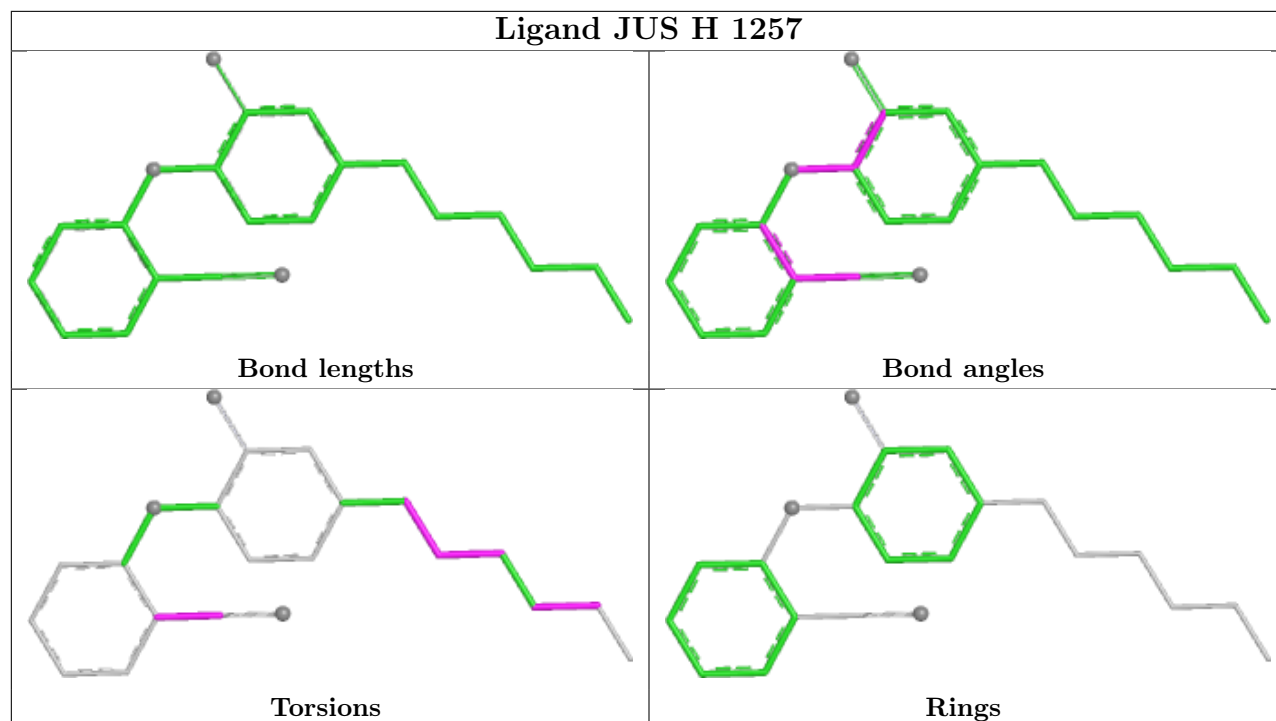
Mol	Chain	Res	Type	Atoms
2	A	1001	GLU	O-C-CA-N
2	B	1001	GLU	N-CA-CB-CG
2	B	1001	GLU	C-CA-CB-CG
2	C	1001	GLU	N-CA-CB-CG
2	E	1001	GLU	O-C-CA-N

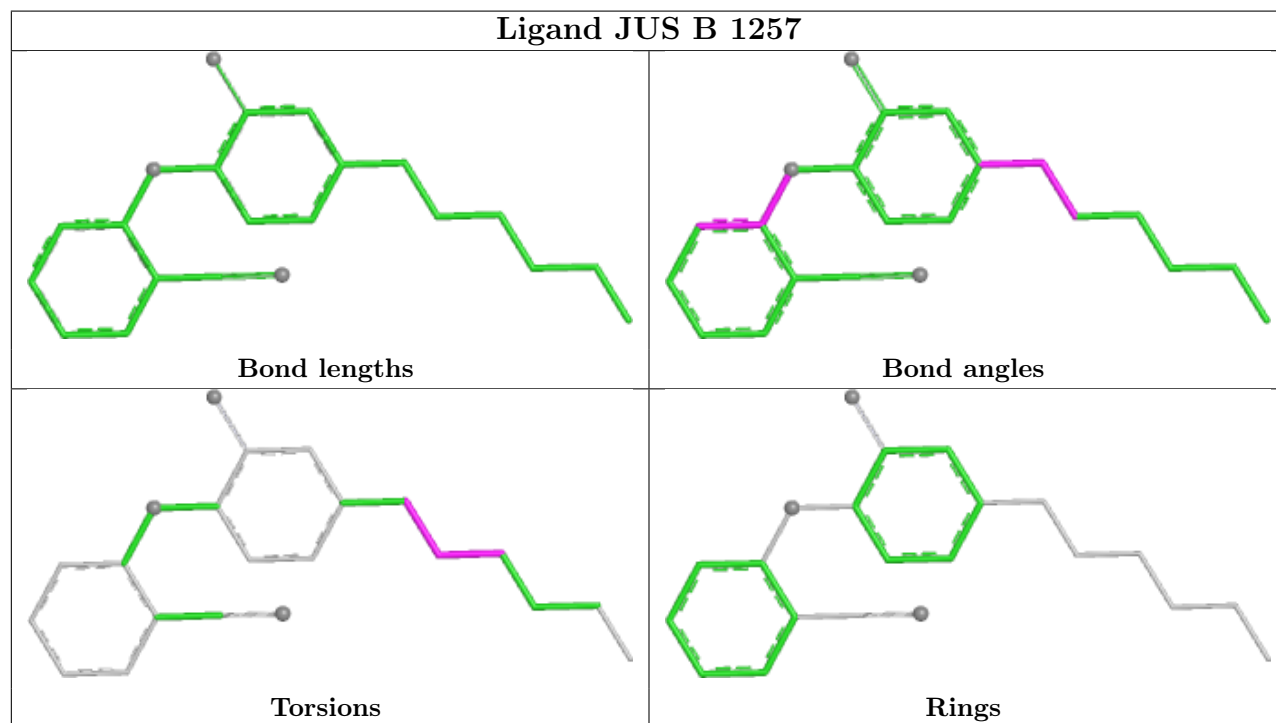
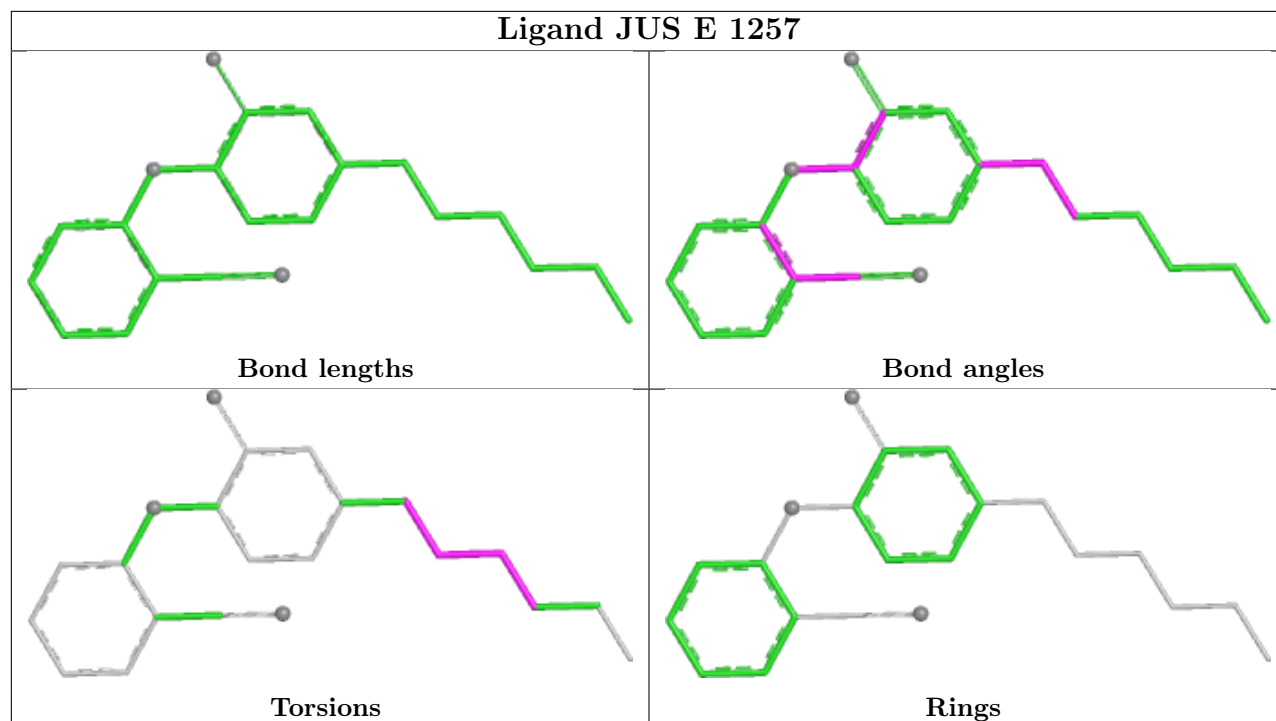
There are no ring outliers.

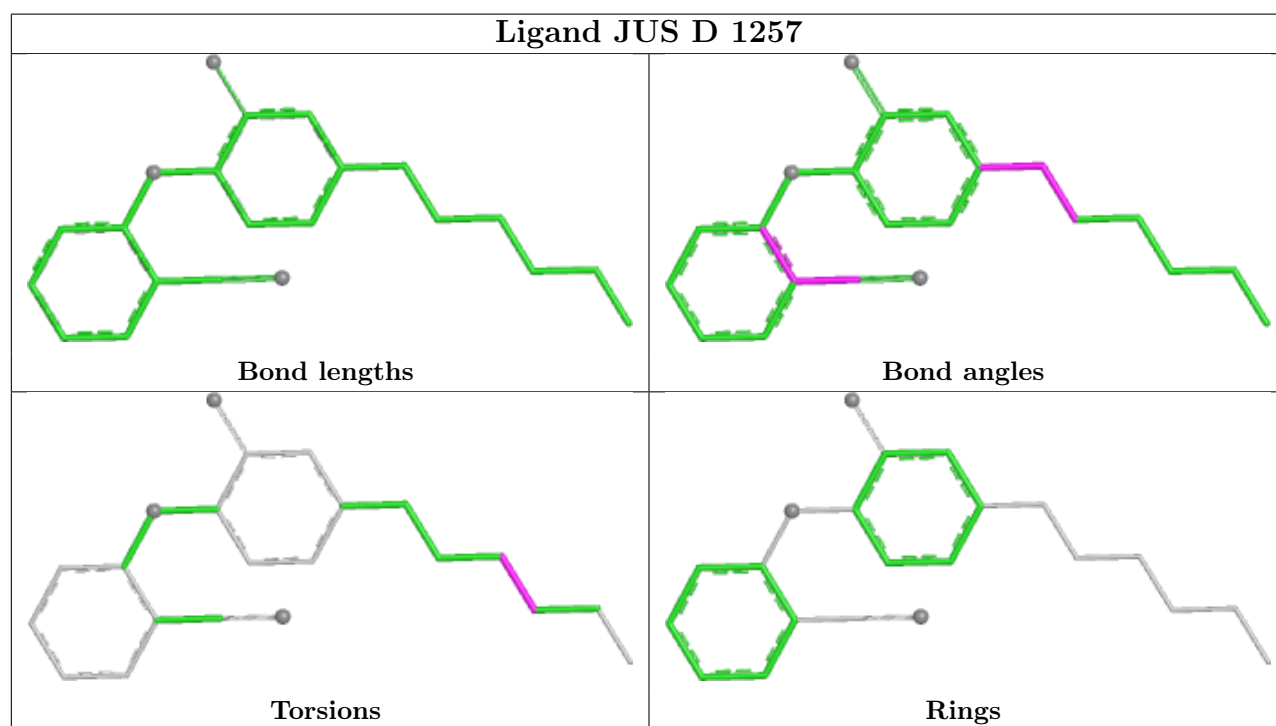
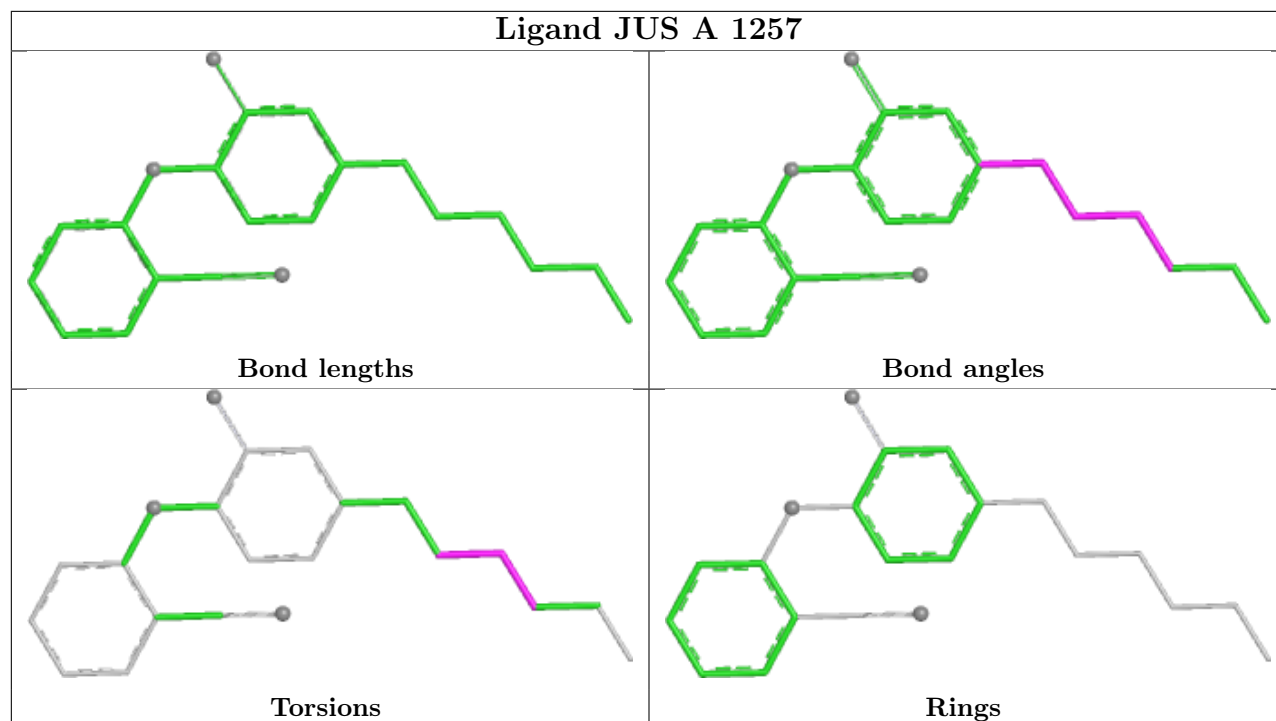
17 monomers are involved in 65 short contacts:

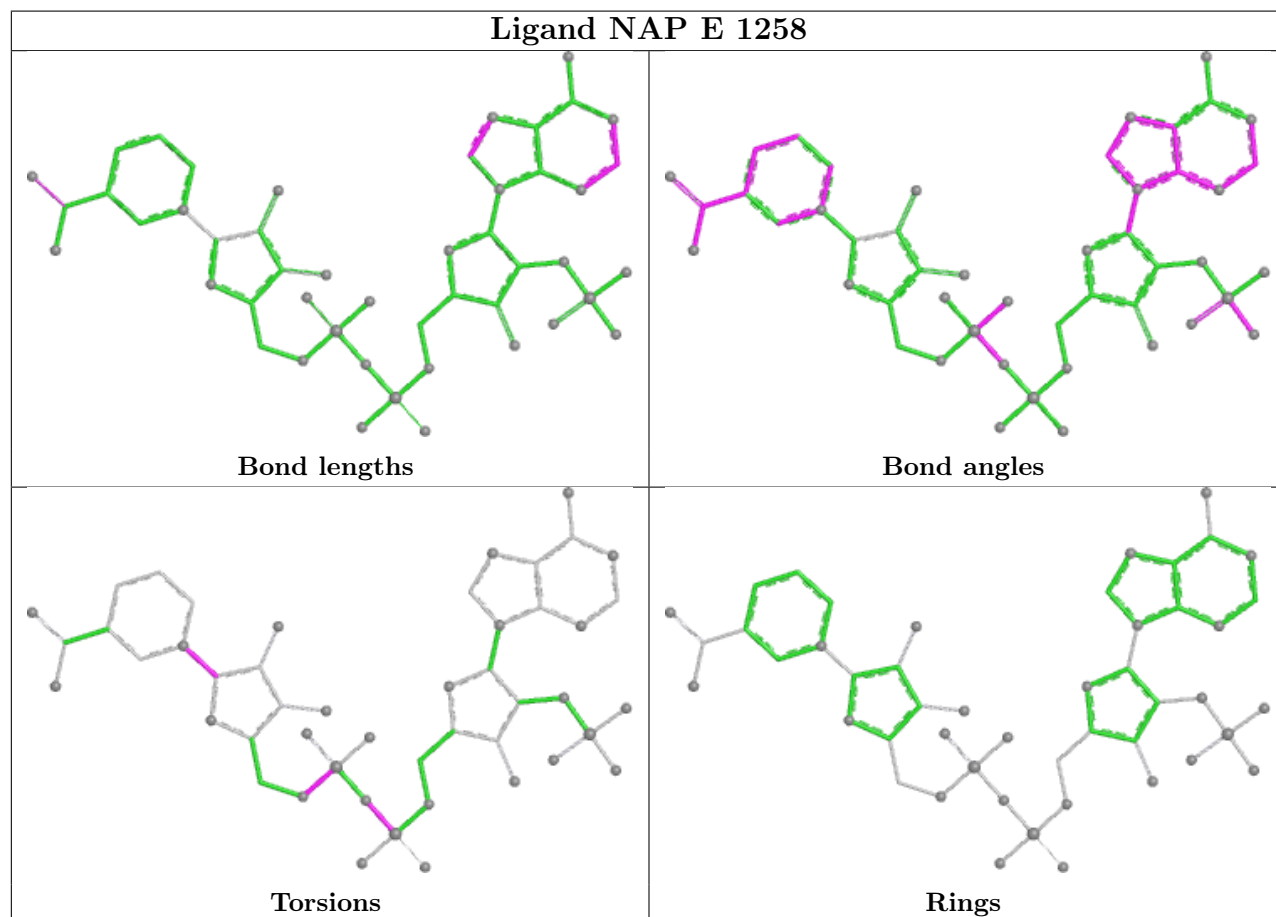
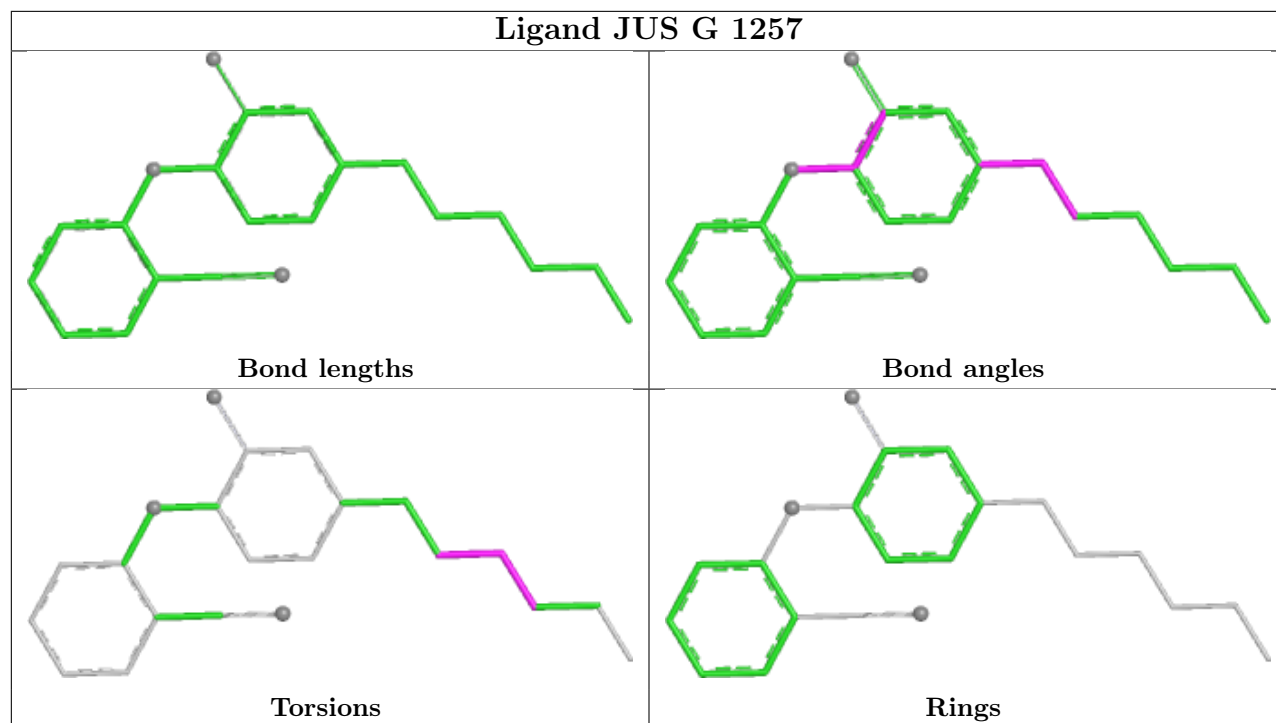
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1257	JUS	2	0
2	G	1001	GLU	2	0
3	E	1257	JUS	1	0
2	A	1001	GLU	15	0
3	B	1257	JUS	2	0
3	A	1257	JUS	3	0
2	E	1001	GLU	7	0
3	G	1257	JUS	2	0
2	F	1001	GLU	12	0
4	E	1258	NAP	1	0
2	C	1001	GLU	2	0
4	A	1258	NAP	1	0
4	H	1258	NAP	3	0
4	C	1258	NAP	1	0
3	F	1257	JUS	1	0
3	C	1257	JUS	4	0
2	B	1001	GLU	7	0

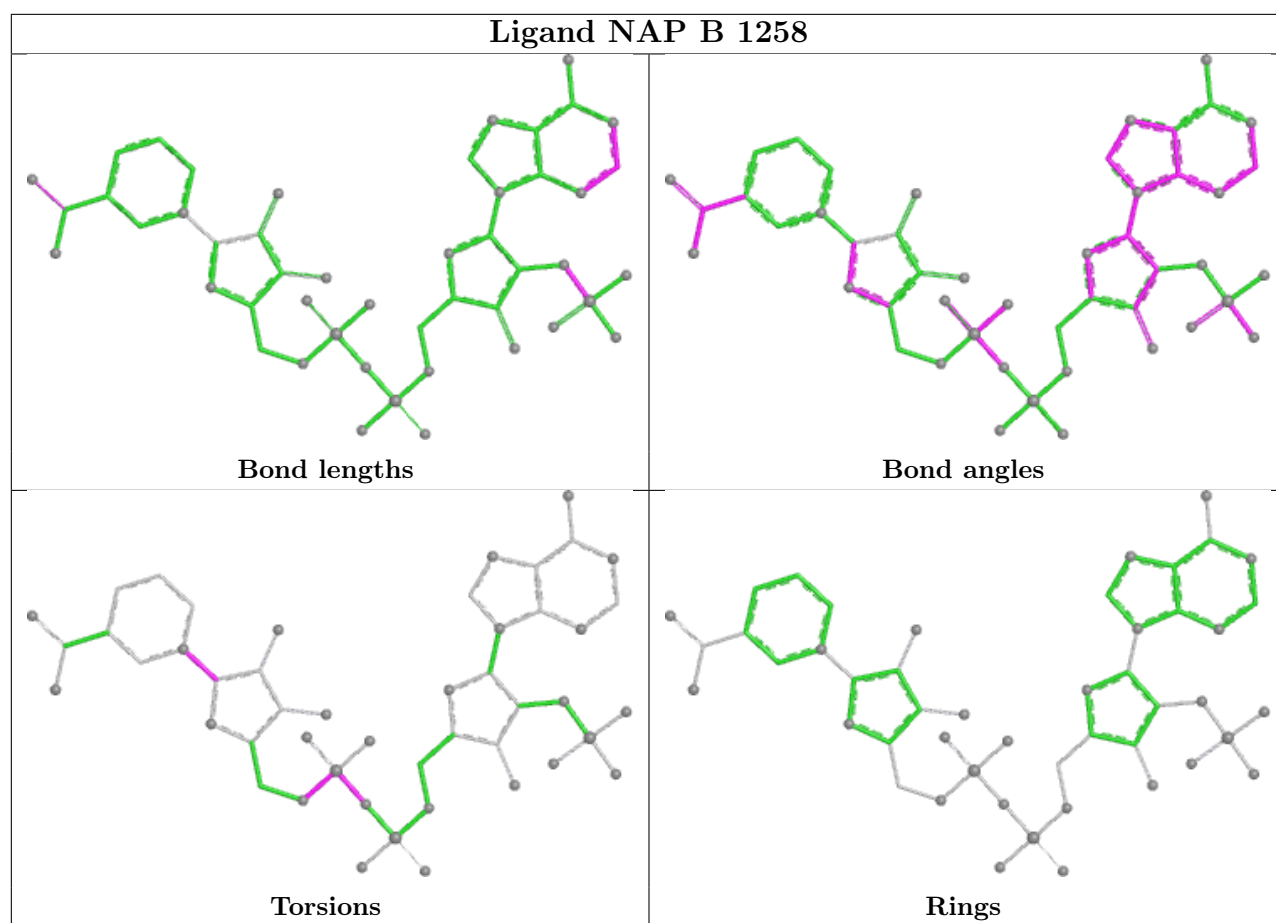
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

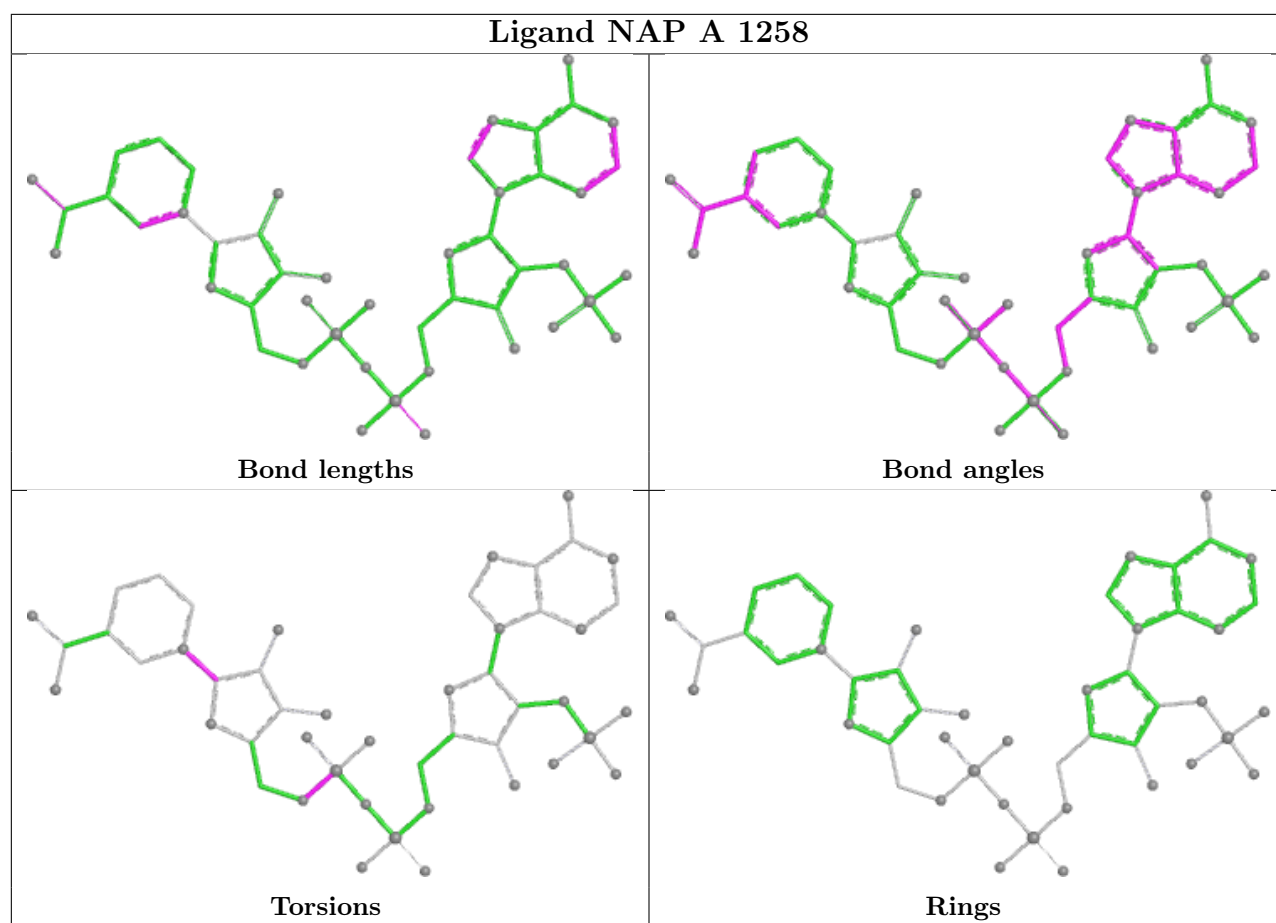


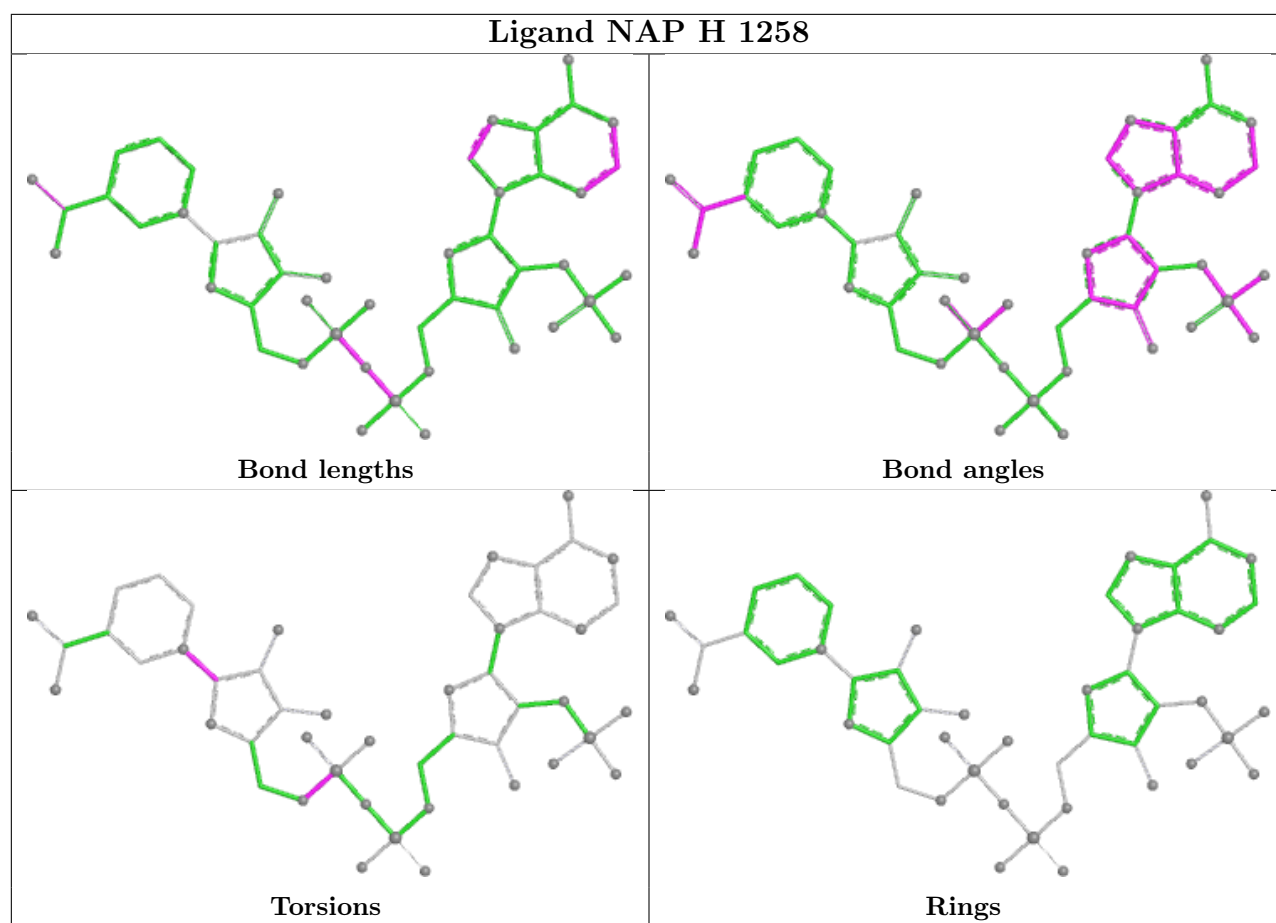


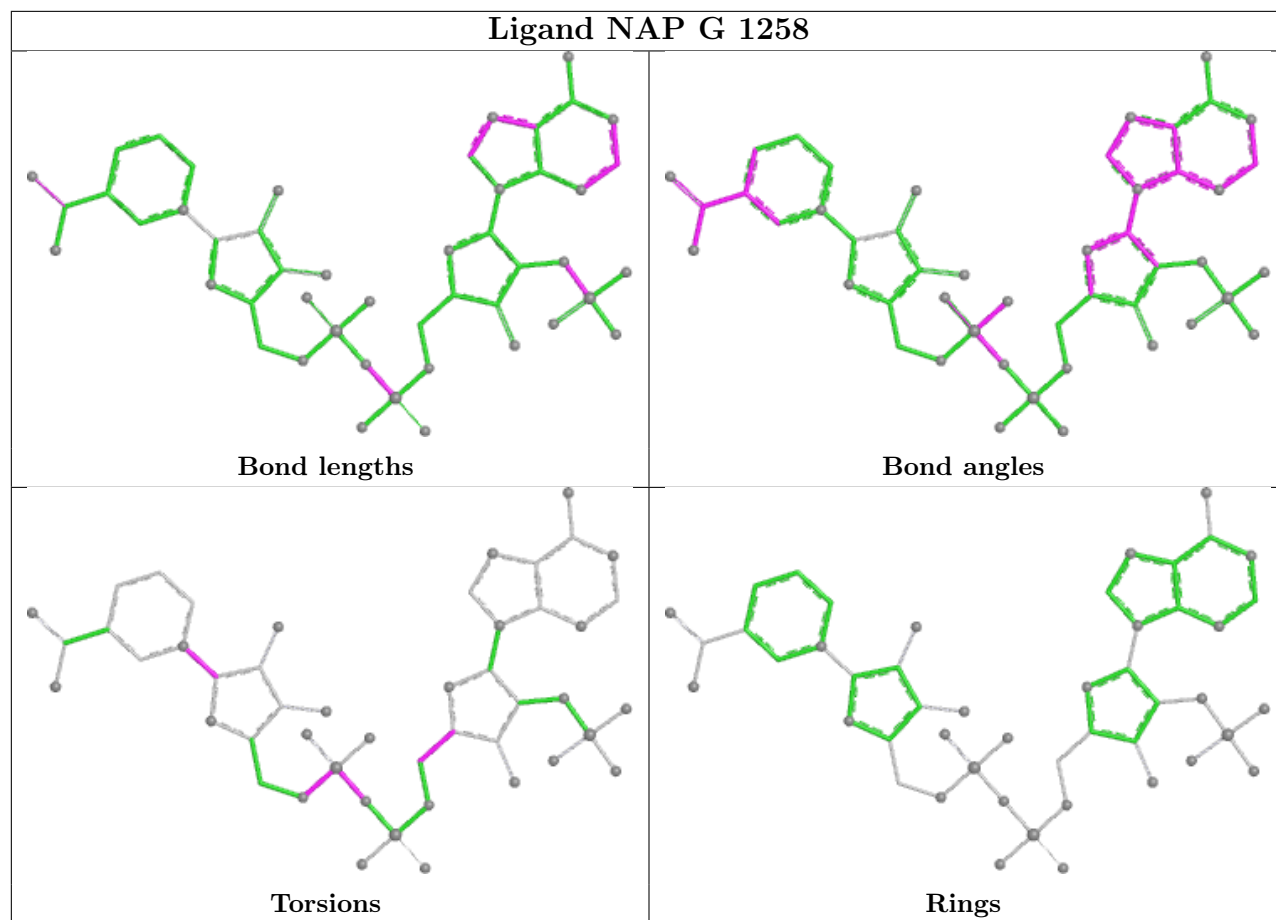


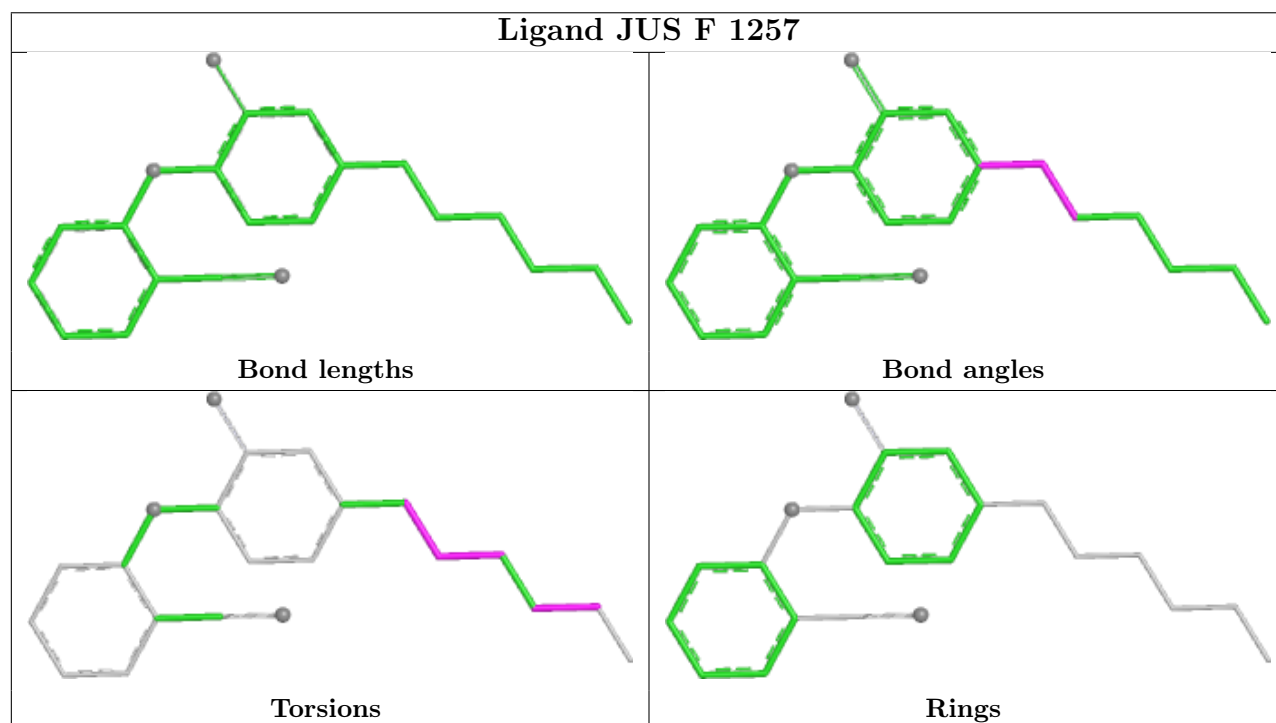
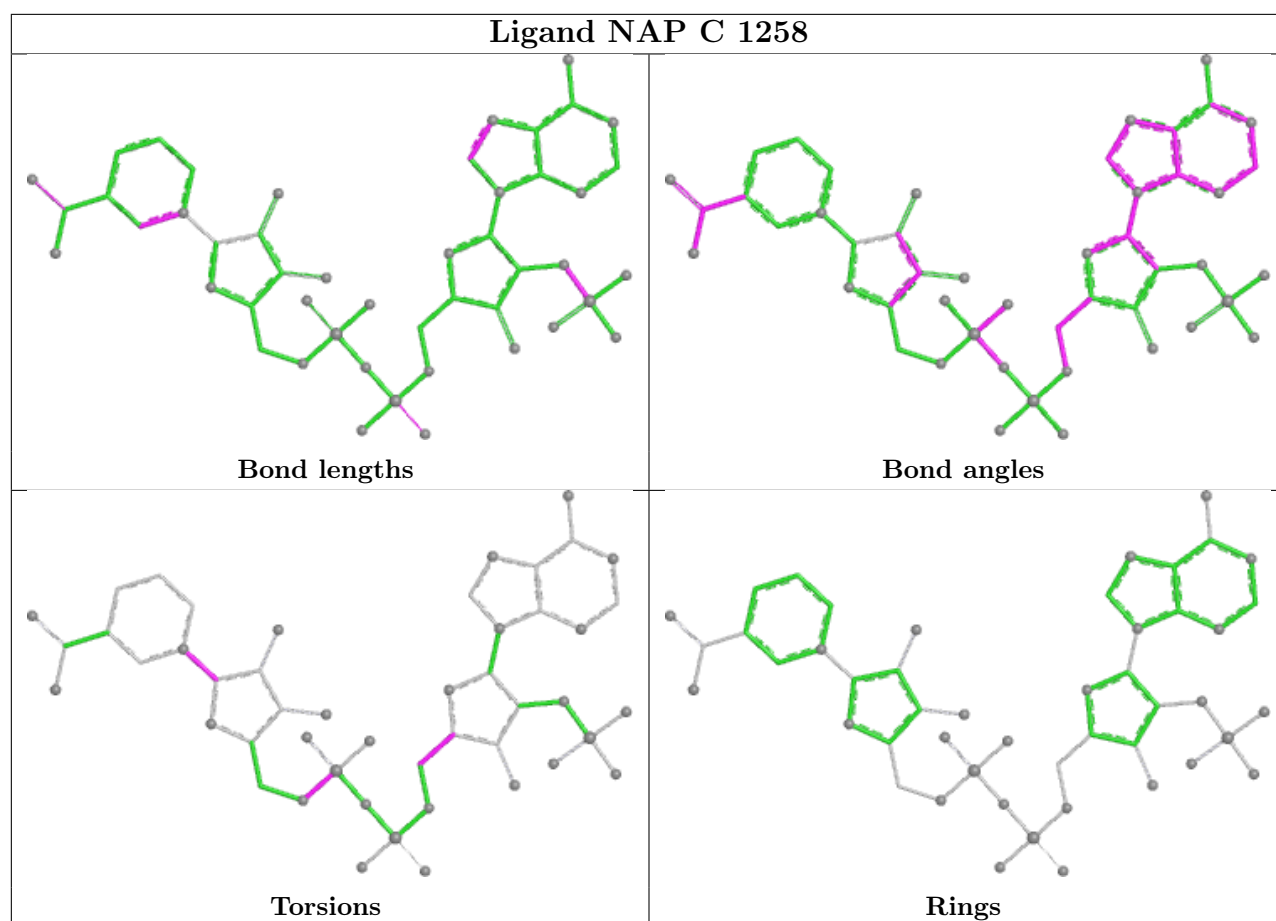




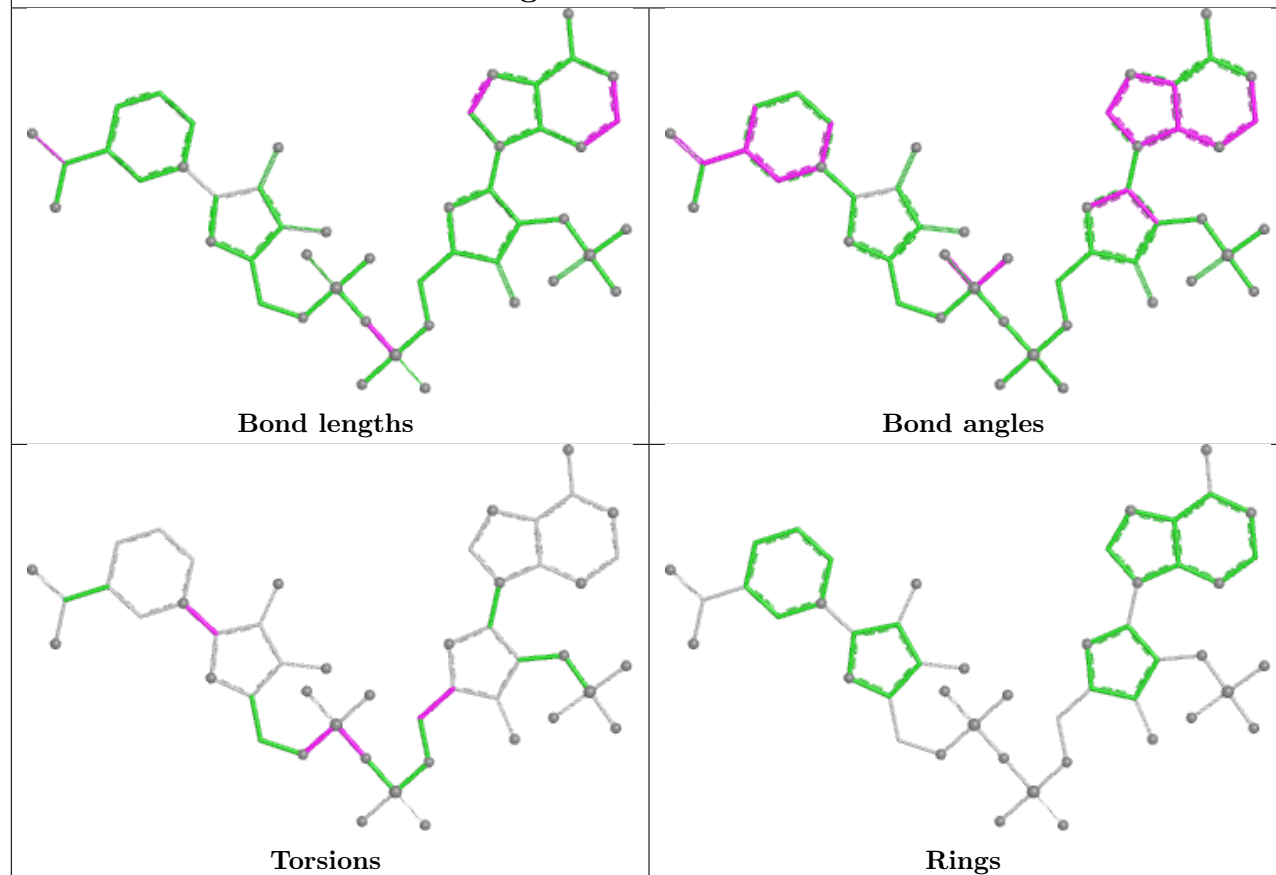




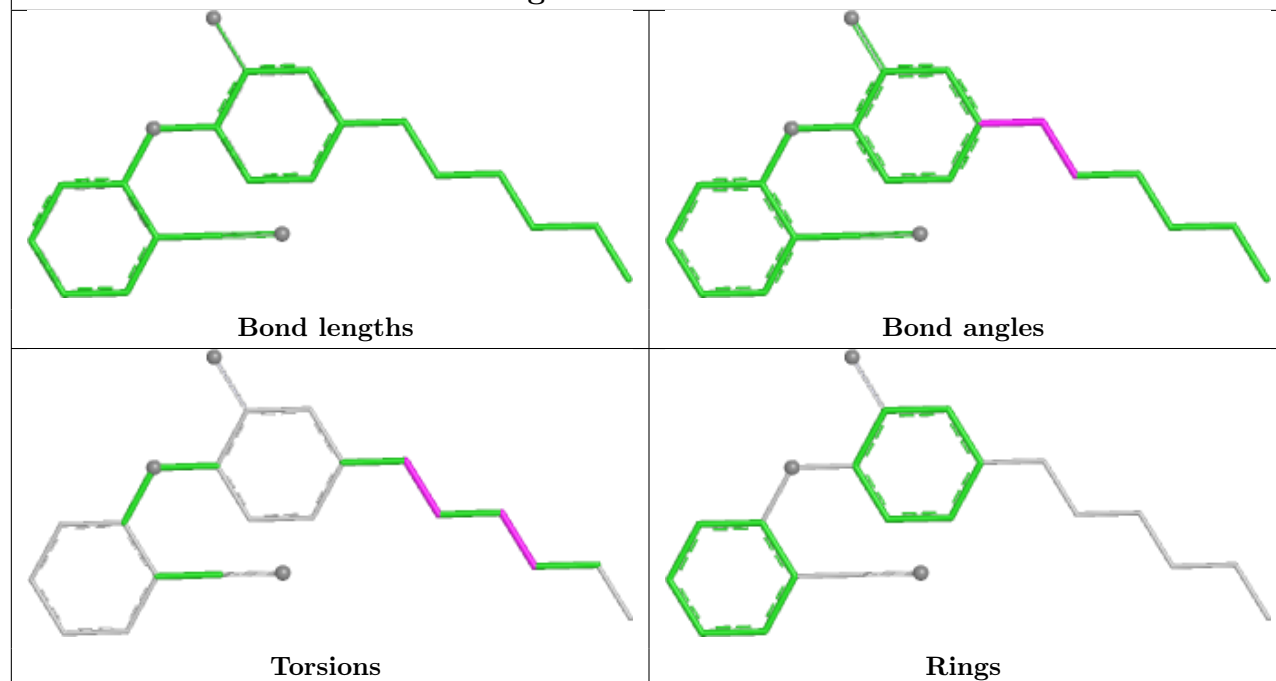




Ligand NAP D 1258



Ligand JUS C 1257



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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	254/282 (90%)	-0.63	1 (0%) 88 89	13, 28, 51, 64	13 (5%)
1	B	255/282 (90%)	-0.58	3 (1%) 76 77	12, 31, 56, 74	12 (4%)
1	C	255/282 (90%)	-0.51	1 (0%) 88 89	14, 33, 56, 74	6 (2%)
1	D	254/282 (90%)	-0.47	0 100 100	13, 33, 63, 84	3 (1%)
1	E	255/282 (90%)	-0.62	2 (0%) 82 84	12, 27, 53, 75	13 (5%)
1	F	255/282 (90%)	-0.57	2 (0%) 82 84	12, 30, 58, 71	11 (4%)
1	G	255/282 (90%)	-0.50	1 (0%) 88 89	14, 33, 58, 70	5 (1%)
1	H	254/282 (90%)	-0.39	4 (1%) 70 71	12, 35, 67, 95	4 (1%)
All	All	2037/2256 (90%)	-0.53	14 (0%) 84 85	12, 32, 59, 95	67 (3%)

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	104[A]	GLY	4.1
1	F	2	VAL	4.0
1	G	2	VAL	3.2
1	C	2	VAL	3.2
1	E	2	VAL	3.2

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

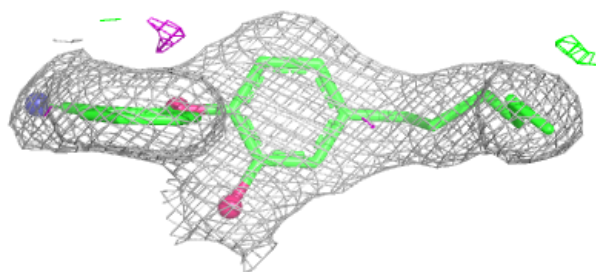
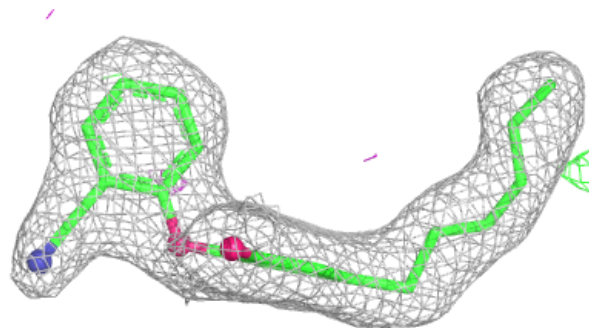
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	GLU	B	1001	10/10	0.63	0.19	70,72,78,78	0
2	GLU	E	1001	10/10	0.73	0.17	81,83,85,85	0
2	GLU	A	1001	10/10	0.74	0.15	72,74,77,78	0
2	GLU	G	1001	10/10	0.76	0.17	78,80,82,85	0
2	GLU	C	1001	10/10	0.77	0.14	73,74,79,80	0
2	GLU	F	1001	10/10	0.83	0.14	71,77,79,80	0
3	JUS	C	1257	22/22	0.96	0.07	26,30,33,36	0
3	JUS	H	1257	22/22	0.96	0.07	31,37,40,42	0
4	NAP	H	1258	48/48	0.96	0.07	26,37,55,58	0
3	JUS	D	1257	22/22	0.97	0.07	30,33,38,41	0
3	JUS	E	1257	22/22	0.97	0.07	22,28,33,34	0
3	JUS	F	1257	22/22	0.97	0.06	20,28,33,37	0
3	JUS	G	1257	22/22	0.97	0.07	23,31,37,40	0
3	JUS	B	1257	22/22	0.97	0.06	21,25,37,41	0
4	NAP	D	1258	48/48	0.97	0.06	26,35,52,55	0
3	JUS	A	1257	22/22	0.97	0.07	19,28,34,41	0
4	NAP	C	1258	48/48	0.98	0.06	23,30,35,42	0
4	NAP	A	1258	48/48	0.98	0.05	21,28,32,33	0
4	NAP	E	1258	48/48	0.98	0.05	18,27,33,40	0
4	NAP	G	1258	48/48	0.98	0.06	25,31,36,41	0
4	NAP	B	1258	48/48	0.98	0.05	25,31,37,40	0
4	NAP	F	1258	48/48	0.99	0.05	23,31,36,39	0

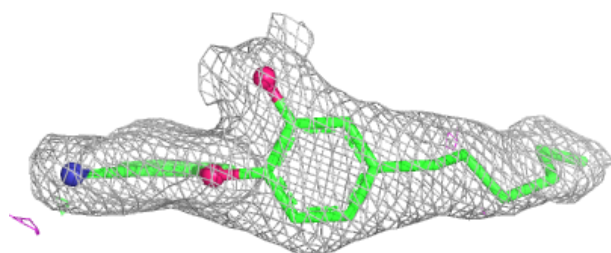
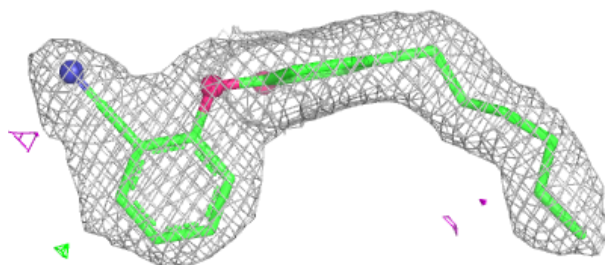
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around JUS C 1257:

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

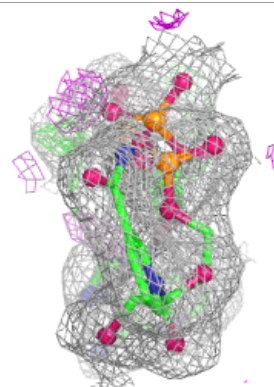
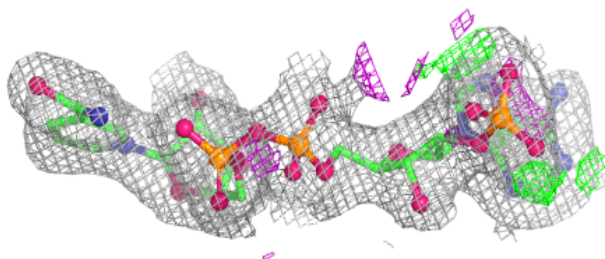
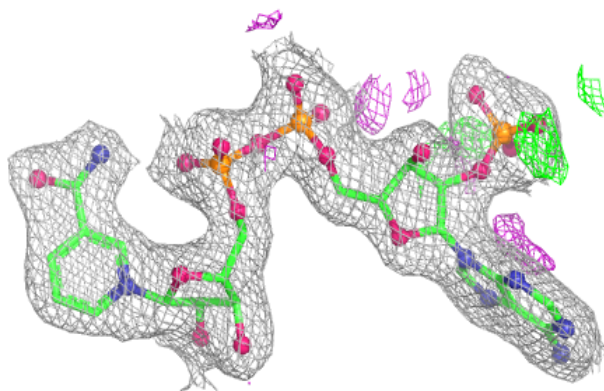
**Electron density around JUS H 1257:**

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

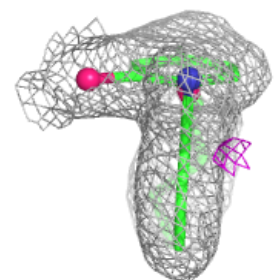
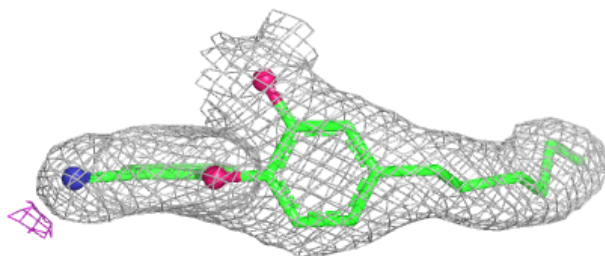
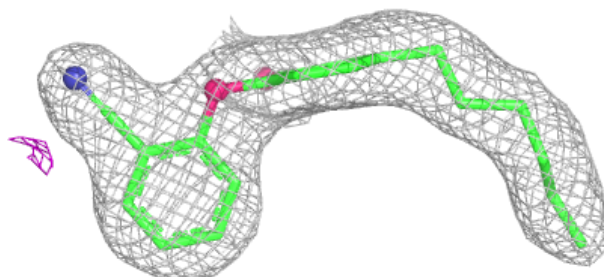


Electron density around NAP H 1258:

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

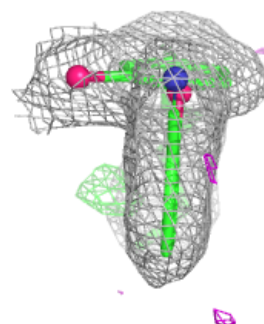
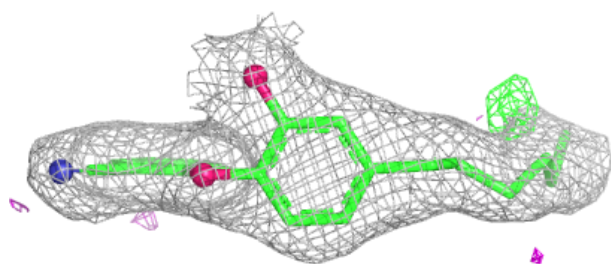
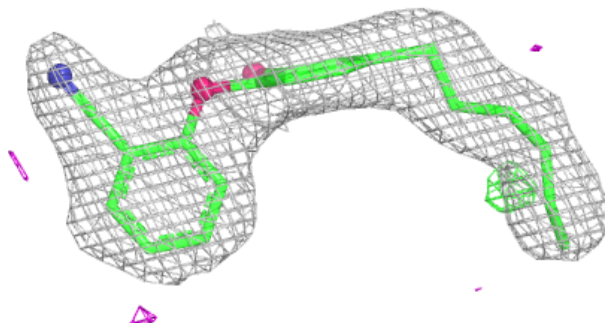
**Electron density around JUS D 1257:**

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

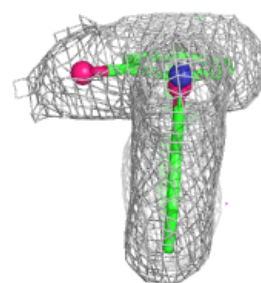
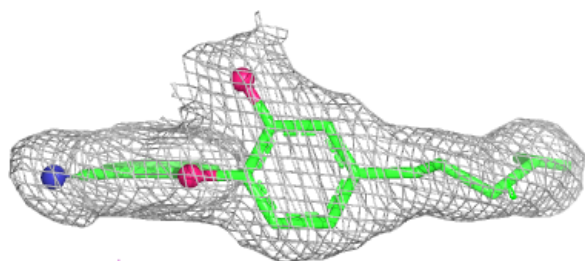
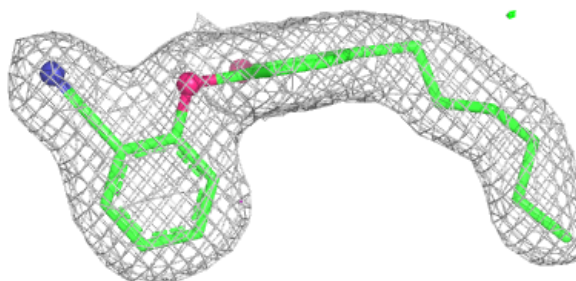


Electron density around JUS E 1257:

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

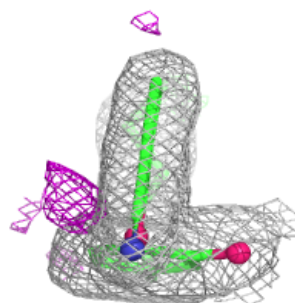
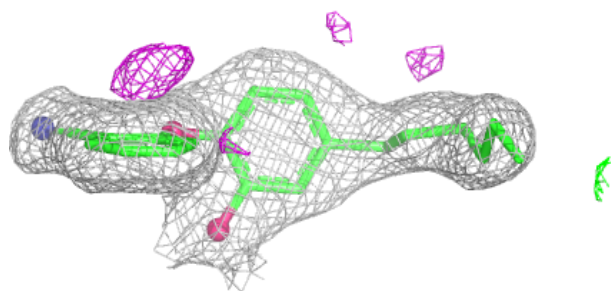
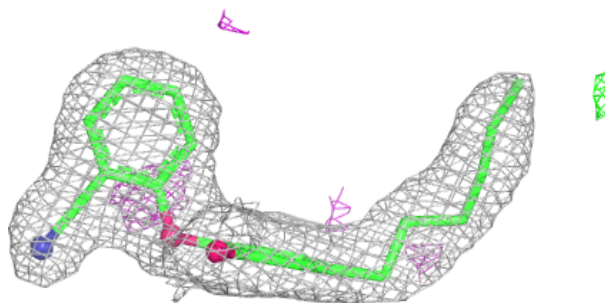
**Electron density around JUS F 1257:**

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and green (positive)

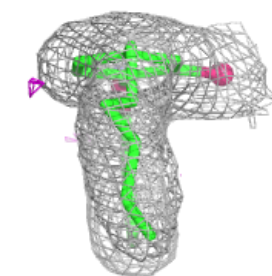
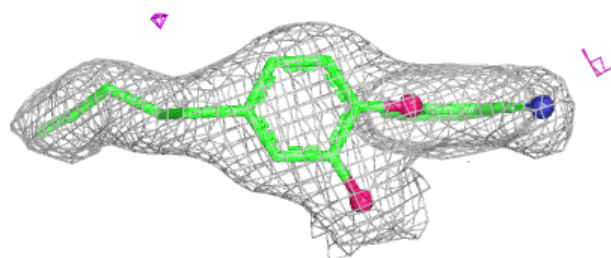
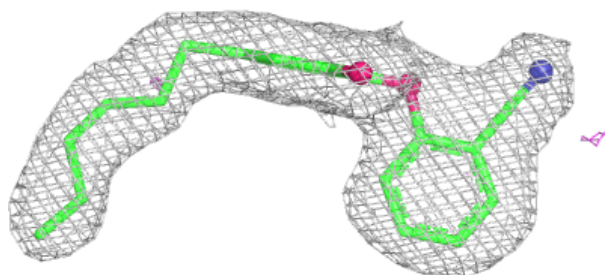


Electron density around JUS G 1257:

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

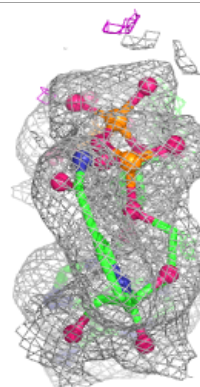
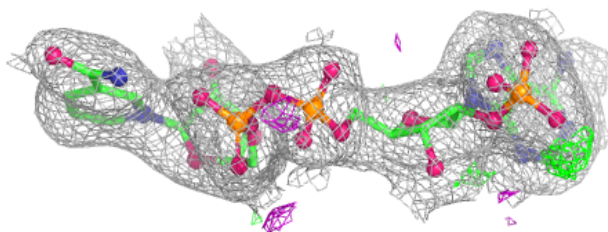
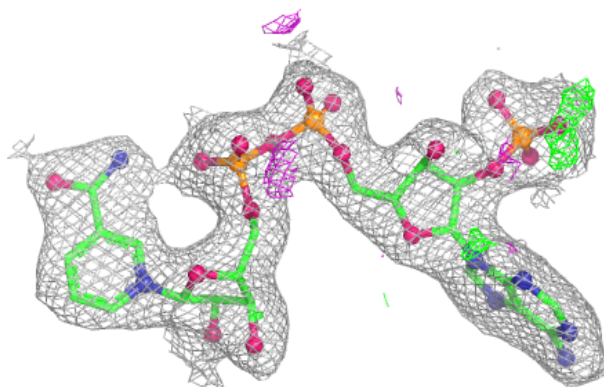
**Electron density around JUS B 1257:**

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

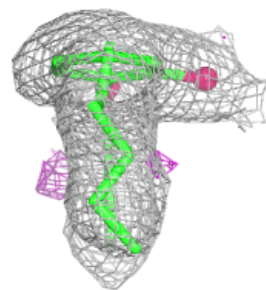
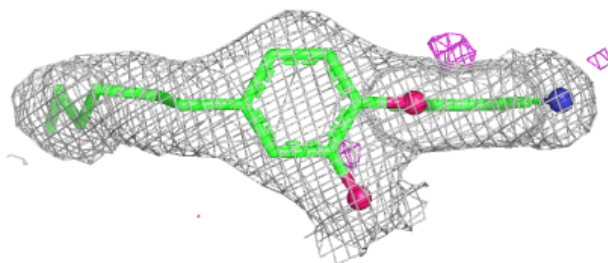
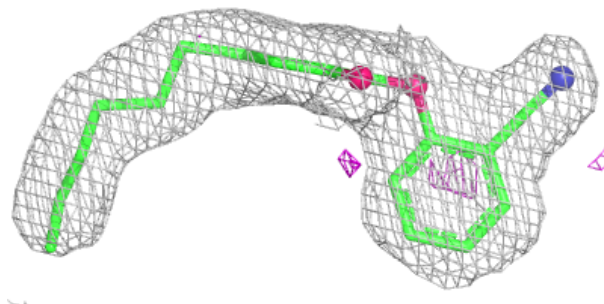


Electron density around NAP D 1258:

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

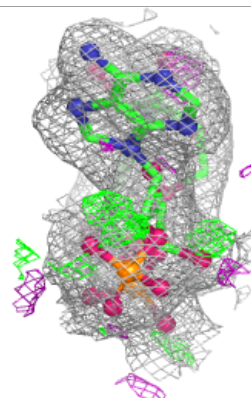
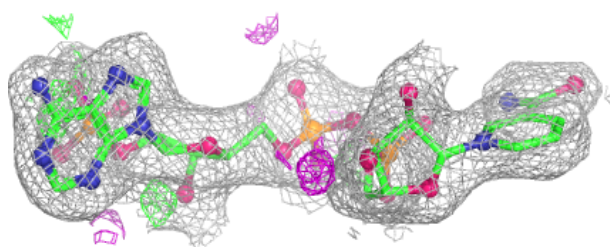
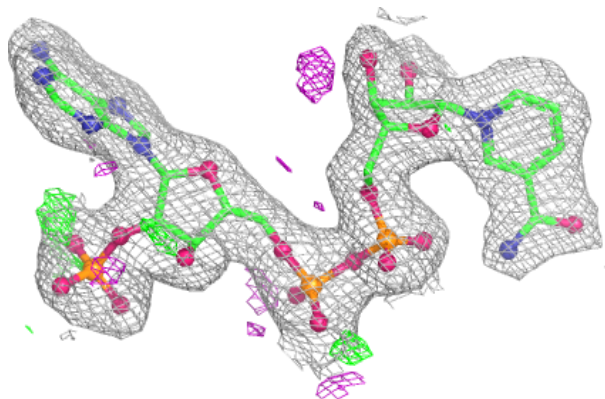
**Electron density around JUS A 1257:**

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

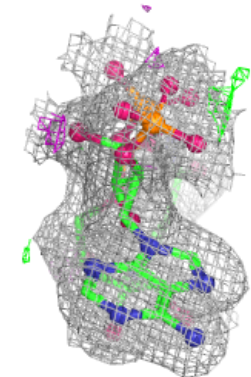
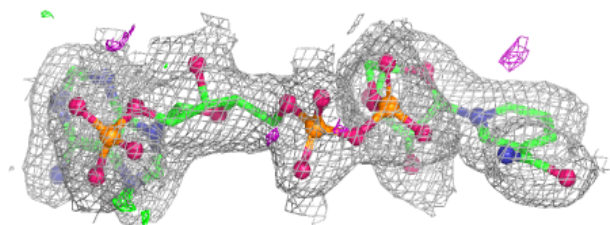
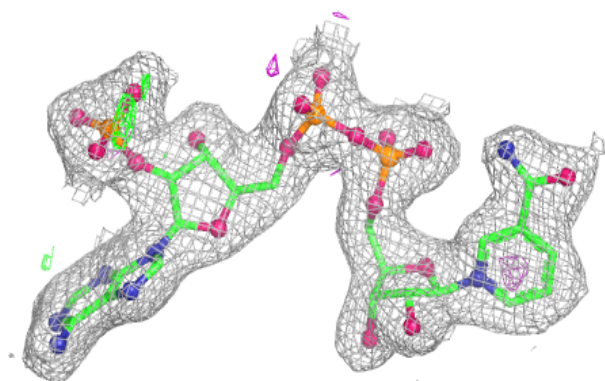


Electron density around NAP C 1258:

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

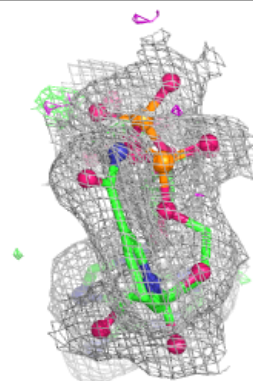
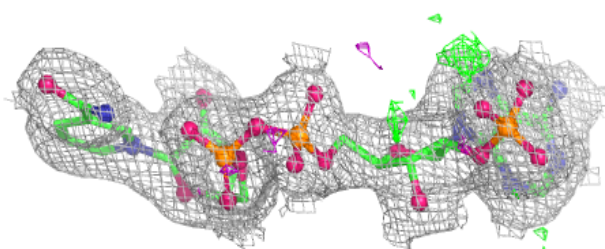
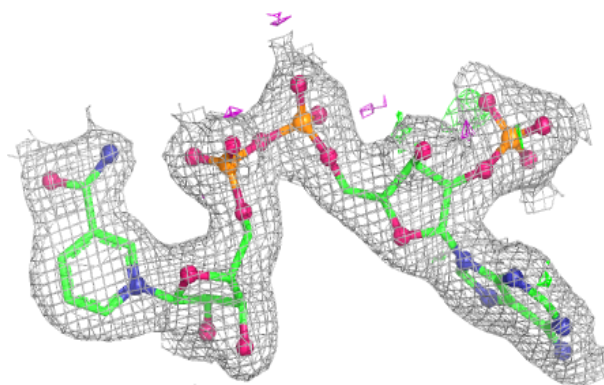
**Electron density around NAP A 1258:**

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

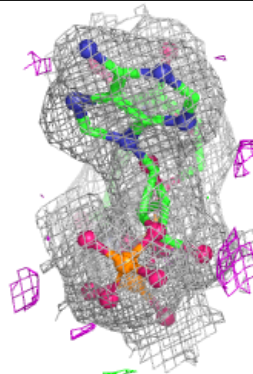
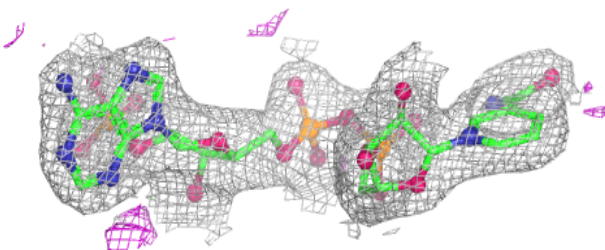
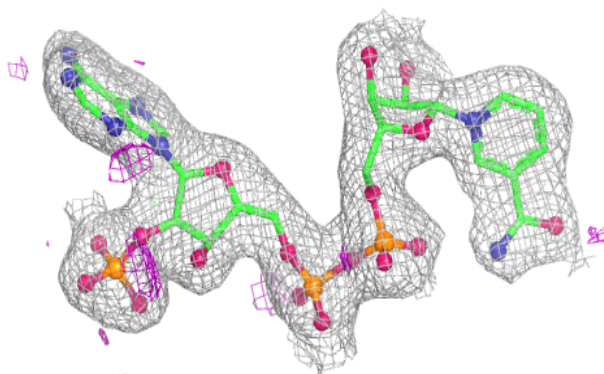


Electron density around NAP E 1258:

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

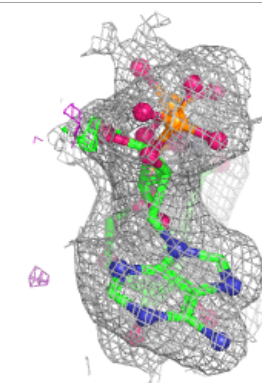
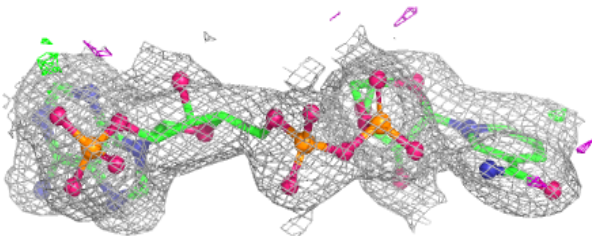
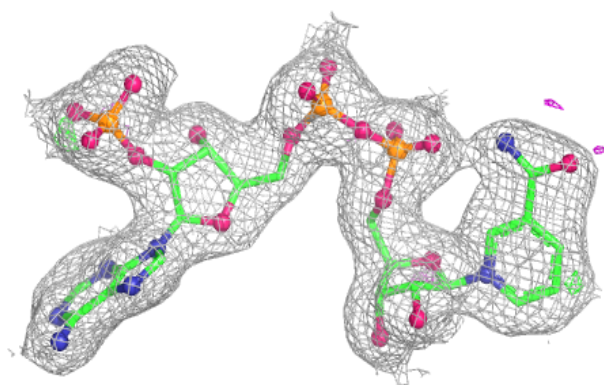
**Electron density around NAP G 1258:**

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

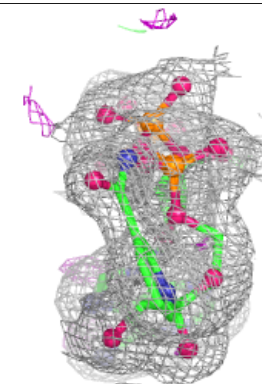
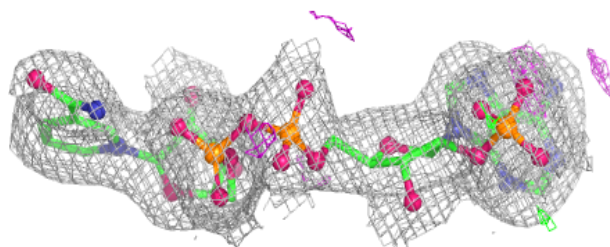
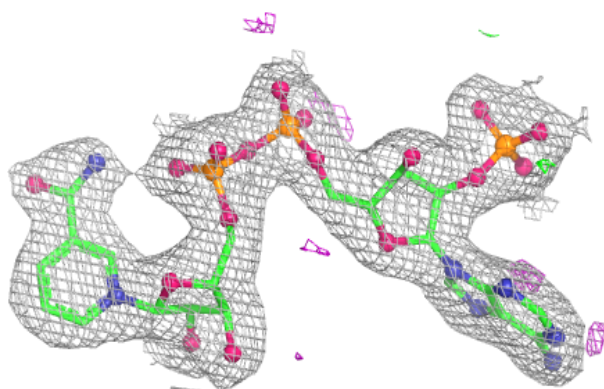


Electron density around NAP B 1258:

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

**Electron density around NAP F 1258:**

$2mF_o-DF_c$ (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.