



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 02:33 am BST

PDB ID : 1ZMD
Title : Crystal Structure of Human dihydrolipoamide dehydrogenase complexed to NADH
Authors : Brautigam, C.A.; Chuang, J.L.; Tomchick, D.R.; Machius, M.; Chuang, D.T.
Deposited on : 2005-05-10
Resolution : 2.08 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

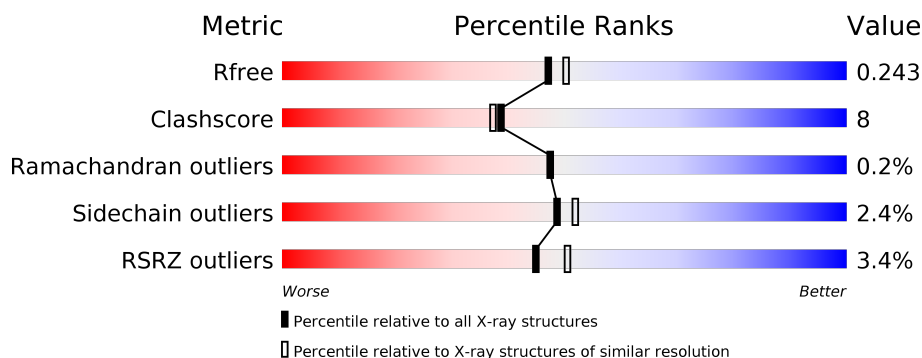
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.08 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}	130704	6189 (2.10-2.06)
Clashscore	141614	6738 (2.10-2.06)
Ramachandran outliers	138981	6663 (2.10-2.06)
Sidechain outliers	138945	6664 (2.10-2.06)
RSRZ outliers	127900	6057 (2.10-2.06)

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 on 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	474	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>17%</div> <div>.</div> </div> </div>
1	B	474	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>18%</div> <div>..</div> </div> </div>
1	C	474	<div> <div>5%</div> <div> <div></div> <div>84%</div> <div>15%</div> <div>..</div> </div> </div>
1	D	474	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>15%</div> <div>.</div> </div> </div>
1	E	474	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>14%</div> <div>.</div> </div> </div>
1	F	474	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>..</div> </div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	G	474	<div><div><div>%</div><div><div></div><div>85%</div><div>14%</div></div><div>.</div></div></div>
1	H	474	<div><div><div>2%</div><div><div></div><div>84%</div><div>14%</div></div><div>..</div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 30958 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 Dihydrolipoyl dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	3	0
			3527	2220	611	675	21			
1	B	471	Total	C	N	O	S	0	2	0
			3508	2209	606	672	21			
1	C	471	Total	C	N	O	S	0	2	0
			3508	2209	606	672	21			
1	D	472	Total	C	N	O	S	0	2	0
			3517	2214	608	674	21			
1	E	473	Total	C	N	O	S	0	2	0
			3525	2218	609	677	21			
1	F	471	Total	C	N	O	S	0	2	0
			3508	2209	606	672	21			
1	G	472	Total	C	N	O	S	0	3	0
			3527	2220	611	675	21			
1	H	471	Total	C	N	O	S	0	2	0
			3508	2209	606	672	21			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



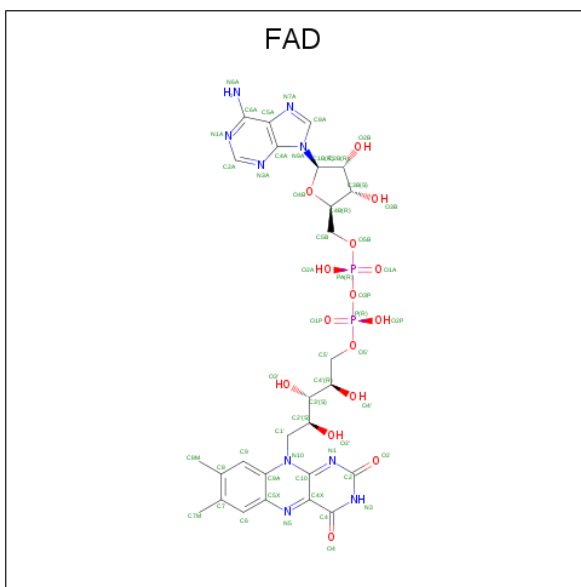
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

Continued on next page...

Continued from previous page...

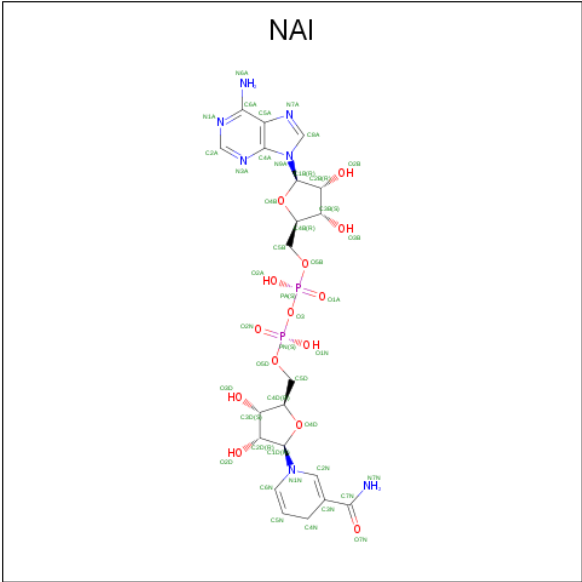
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	B	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	C	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	D	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	E	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	F	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	G	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	H	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 4 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	194	Total	O	0	0
			194	194		
5	B	183	Total	O	0	0
			183	183		
5	C	197	Total	O	0	0
			197	197		
5	D	193	Total	O	0	0
			193	193		

Continued on next page...

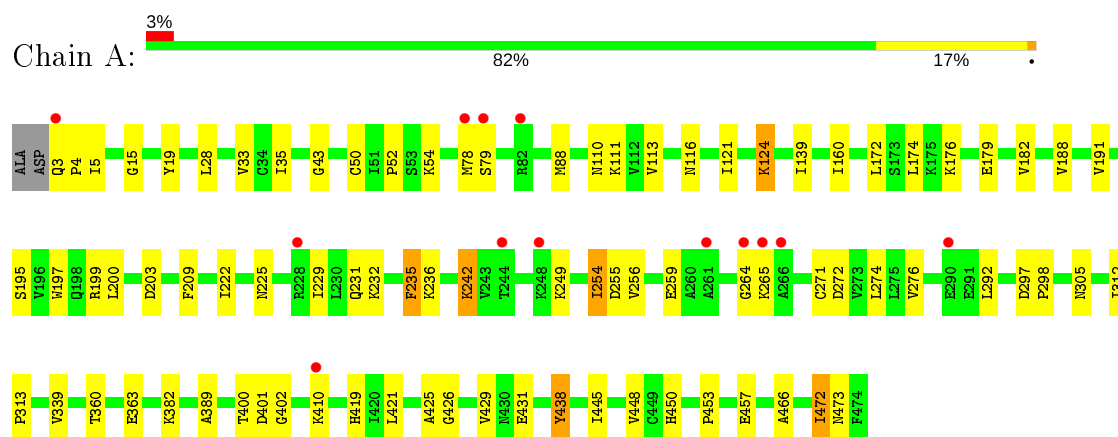
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	233	Total 233	O 233	0	0
5	F	311	Total 311	O 311	0	0
5	G	331	Total 331	O 331	0	0
5	H	257	Total 257	O 257	0	0

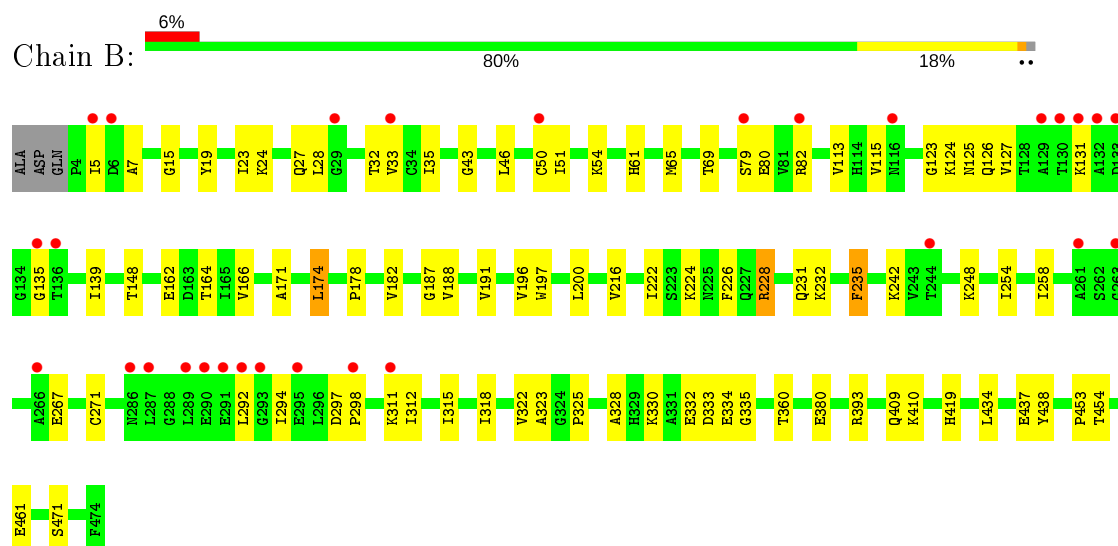
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

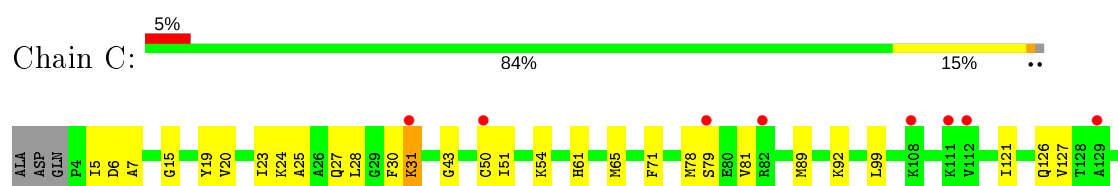
• Molecule 1: Dihydrolipoyl dehydrogenase

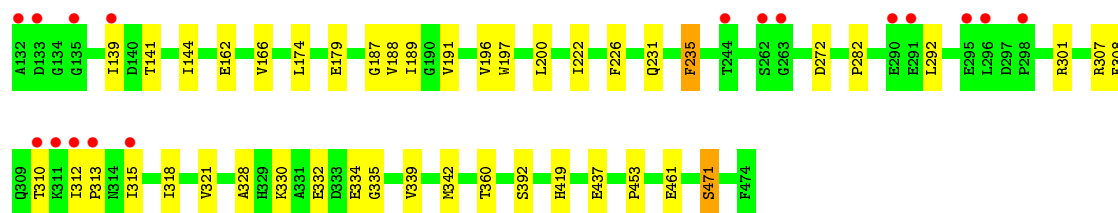


• Molecule 1: Dihydrolipoyl dehydrogenase

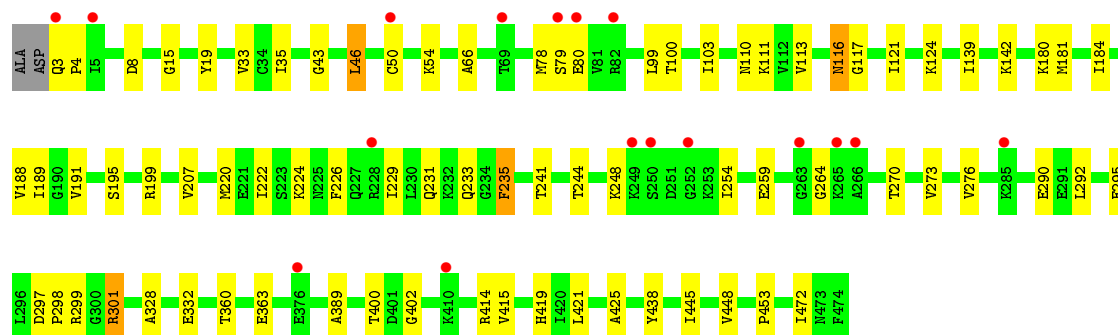
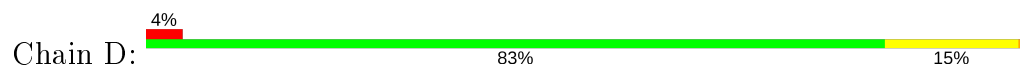


• Molecule 1: Dihydrolipoyl dehydrogenase

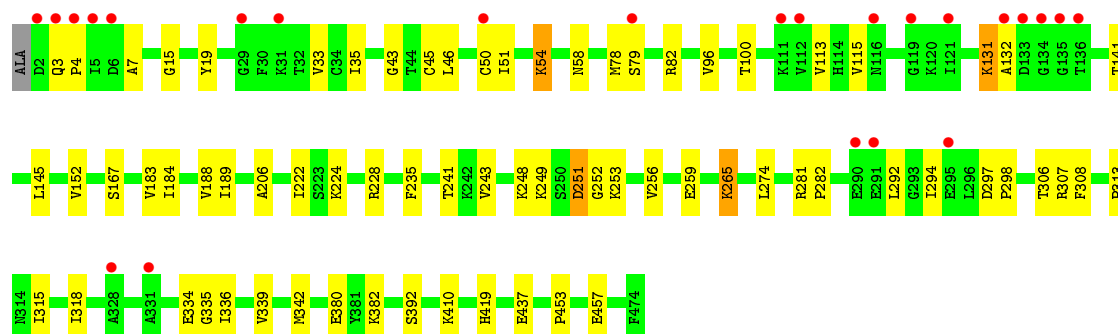
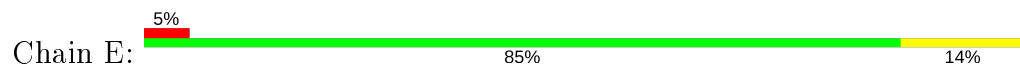




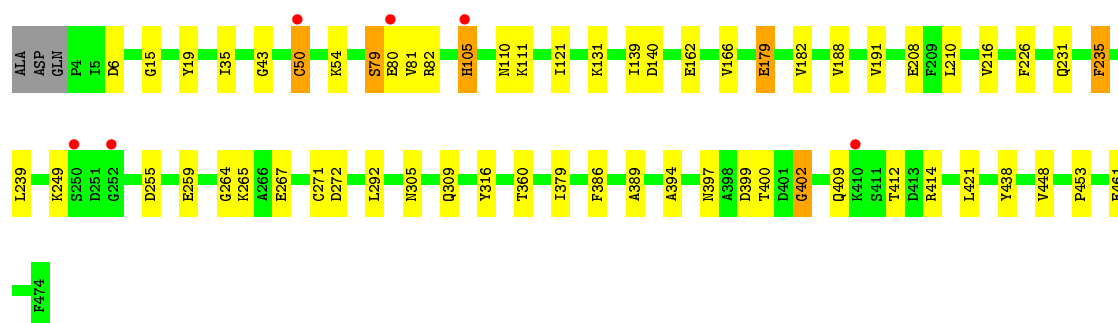
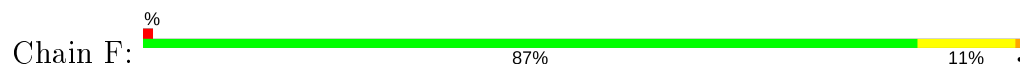
• Molecule 1: Dihydrolipoyl dehydrogenase



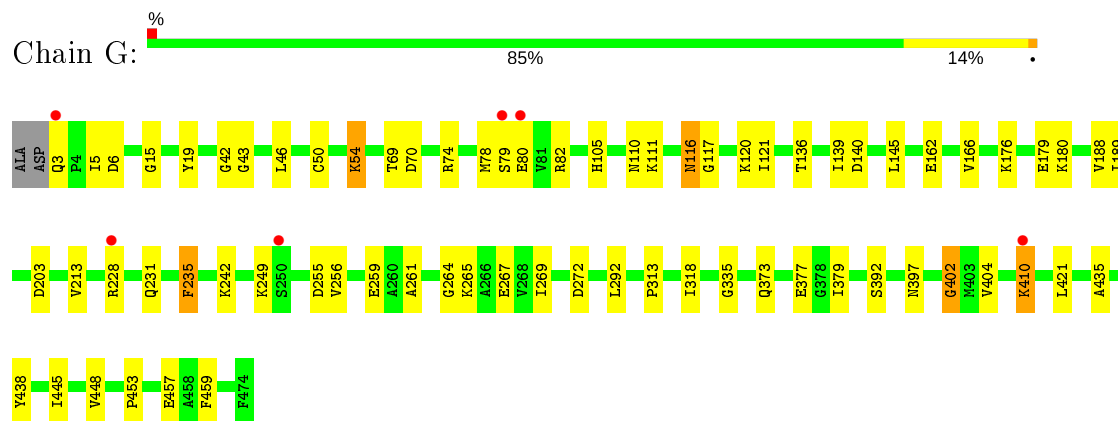
• Molecule 1: Dihydrolipoyl dehydrogenase



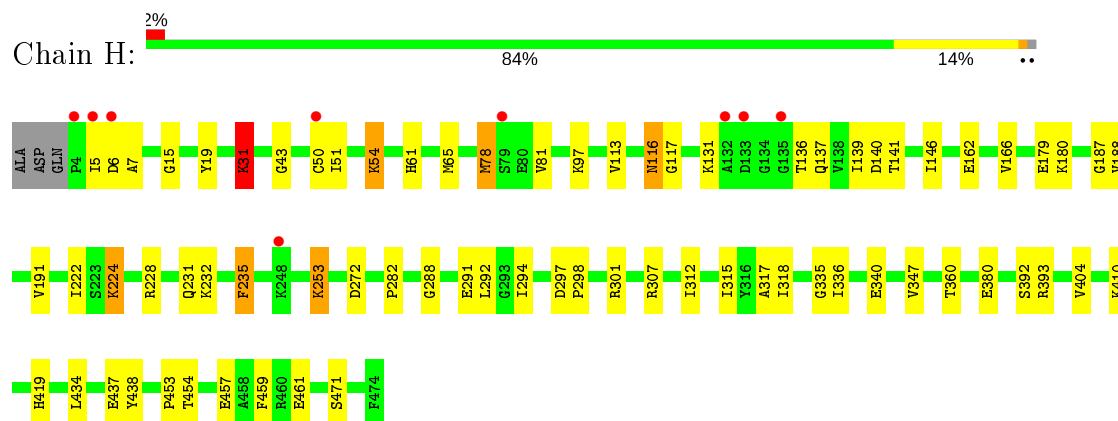
• Molecule 1: Dihydrolipoyl dehydrogenase



- Molecule 1: Dihydrolipoyl dehydrogenase



- Molecule 1: Dihydrolipoyl dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	175.87Å 210.91Å 127.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.06 – 2.08 40.77 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.7 (33.06-2.08) 99.8 (40.77-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.57 (at 2.08Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.249 0.212 , 0.243	Depositor DCC
R_{free} test set	1748 reflections (0.62%)	wwPDB-VP
Wilson B-factor (Å ²)	26.8	Xtriage
Anisotropy	0.542	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	30958	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 54.00 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.8873e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAI, SO4, FAD

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	0.55	0/3583	0.74	1/4840 (0.0%)
1	B	0.54	0/3563	0.74	0/4812
1	C	0.53	0/3563	0.74	0/4812
1	D	0.55	0/3572	0.73	0/4825
1	E	0.60	0/3580	0.78	0/4836
1	F	0.65	0/3563	0.79	3/4812 (0.1%)
1	G	0.65	0/3583	0.79	2/4840 (0.0%)
1	H	0.63	0/3563	0.78	1/4812 (0.0%)
All	All	0.59	0/28570	0.76	7/38589 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	402	GLY	N-CA-C	6.08	128.29	113.10
1	H	31	LYS	N-CA-C	-5.52	96.09	111.00
1	A	313	PRO	N-CA-C	5.30	125.88	112.10
1	F	50[A]	CYS	CA-CB-SG	5.13	123.24	114.00
1	F	50[B]	CYS	CA-CB-SG	5.13	123.24	114.00
1	G	313	PRO	N-CA-C	5.07	125.28	112.10
1	F	402	GLY	N-CA-C	5.00	125.61	113.10

There are no chirality outliers.

There are no planarity outliers.

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	3527	0	3570	63	0
1	B	3508	0	3557	82	0
1	C	3508	0	3557	53	0
1	D	3517	0	3564	53	0
1	E	3525	0	3568	59	0
1	F	3508	0	3557	42	0
1	G	3527	0	3570	55	0
1	H	3508	0	3557	58	0
2	A	15	0	0	0	0
2	B	20	0	0	1	0
2	C	20	0	0	1	0
2	D	20	0	0	0	0
2	E	20	0	0	0	0
2	F	20	0	0	1	0
2	G	20	0	0	1	0
2	H	20	0	0	0	0
3	A	53	0	31	4	0
3	B	53	0	31	5	0
3	C	53	0	31	3	0
3	D	53	0	31	4	0
3	E	53	0	31	3	0
3	F	53	0	31	4	0
3	G	53	0	31	5	0
3	H	53	0	31	2	0
4	A	44	0	27	1	0
4	B	44	0	27	4	0
4	C	44	0	27	5	0
4	D	44	0	27	2	0
4	E	44	0	27	4	0
4	F	44	0	27	1	0
4	G	44	0	27	3	0
4	H	44	0	27	3	0
5	A	194	0	0	2	0
5	B	183	0	0	2	0
5	C	197	0	0	2	0
5	D	193	0	0	0	0
5	E	233	0	0	3	0
5	F	311	0	0	1	0
5	G	331	0	0	8	0
5	H	257	0	0	4	0
All	All	30958	0	28964	437	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.

All (437) 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:C:437:GLU:HG3	1:D:448:VAL:HG22	1.39	1.03
1:G:50[B]:CYS:SG	1:H:453:PRO:HD3	2.04	0.98
1:E:453:PRO:HD3	1:F:50[B]:CYS:SG	2.03	0.97
1:C:121:ILE:HG21	1:C:292:LEU:HD11	1.48	0.94
1:E:335:GLY:O	1:E:339:VAL:HG22	1.71	0.91
1:A:50[B]:CYS:SG	1:B:453:PRO:HD3	2.12	0.90
1:G:397:ASN:HB3	5:H:2107:HOH:O	1.71	0.89
1:E:339:VAL:HG12	1:E:342:MET:HE2	1.56	0.88
1:B:5:ILE:HD11	1:B:139:ILE:HG12	1.59	0.85
5:E:2078:HOH:O	1:F:397:ASN:HB3	1.77	0.84
1:E:339:VAL:HG12	1:E:342:MET:CE	2.10	0.82
1:A:448:VAL:HG12	1:B:437:GLU:HG2	1.63	0.80
1:A:33:VAL:HG11	1:A:139:ILE:HD13	1.64	0.80
1:C:20:VAL:HG13	1:D:472:ILE:HD13	1.65	0.78
1:D:46:LEU:HD22	1:D:100:THR:HG23	1.65	0.78
1:A:179:GLU:HB3	1:A:272:ASP:OD2	1.84	0.77
1:A:259:GLU:HB2	1:A:264:GLY:O	1.85	0.76
1:B:80:GLU:HG2	1:B:82:ARG:CZ	2.16	0.74
1:C:453:PRO:HD3	1:D:50[B]:CYS:SG	2.28	0.73
1:F:179:GLU:HG2	1:F:272:ASP:OD2	1.90	0.72
1:G:176:LYS:HD3	5:G:2135:HOH:O	1.90	0.71
1:D:46:LEU:HD21	1:D:99:LEU:HB2	1.72	0.71
1:A:28:LEU:HD12	1:A:339:VAL:HG12	1.72	0.70
1:D:121:ILE:HG21	1:D:292:LEU:HD11	1.74	0.70
1:A:182:VAL:HG12	1:A:271:CYS:SG	2.31	0.70
1:B:5:ILE:CD1	1:B:139:ILE:HG12	2.21	0.69
1:B:191:VAL:HG12	1:B:360:THR:HG21	1.74	0.69
1:E:184:ILE:CD1	1:E:243:VAL:HG21	2.22	0.69
1:E:437:GLU:HG3	1:F:448:VAL:HG22	1.75	0.68
1:E:46:LEU:HA	1:E:51:ILE:HD13	1.75	0.68
1:B:33:VAL:HG12	1:B:113:VAL:CG1	2.24	0.68
1:A:160:ILE:HD11	1:A:276:VAL:HB	1.77	0.67
1:C:191:VAL:HG12	1:C:360:THR:HG21	1.77	0.67
1:E:339:VAL:HA	1:E:342:MET:HE2	1.77	0.67
1:B:115:VAL:HG21	1:B:139:ILE:HD11	1.76	0.67
1:F:216:VAL:HG22	5:F:2095:HOH:O	1.93	0.67
1:F:259:GLU:HB2	1:F:264:GLY:O	1.95	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:322:VAL:HG12	1:B:323:ALA:N	2.10	0.66
1:H:392:SER:HB2	5:H:2165:HOH:O	1.94	0.65
1:B:294:ILE:HD11	1:B:312:ILE:HD13	1.78	0.65
1:B:162:GLU:HA	1:B:166:VAL:HG12	1.79	0.64
1:B:228:ARG:CG	1:B:228:ARG:HH21	2.09	0.64
1:H:228:ARG:HG2	1:H:228:ARG:HH11	1.61	0.64
1:G:15:GLY:HA3	3:G:480:FAD:H52A	1.78	0.64
1:A:124:LYS:HD2	1:A:312:ILE:HD13	1.79	0.64
1:A:3:GLN:N	1:A:4:PRO:HD3	2.12	0.64
1:C:78:MET:HG2	1:D:78:MET:HG2	1.78	0.64
1:H:162:GLU:HA	1:H:166:VAL:HG12	1.79	0.64
1:G:50[B]:CYS:SG	1:H:453:PRO:CD	2.83	0.64
1:B:171:ALA:HA	1:B:174:LEU:HD13	1.78	0.64
1:G:242:LYS:HE3	1:G:261:ALA:HA	1.80	0.64
1:A:121:ILE:HG21	1:A:292:LEU:HD11	1.80	0.63
1:D:259:GLU:HB2	1:D:264:GLY:O	1.97	0.63
1:C:28:LEU:HD12	1:C:339:VAL:HG12	1.80	0.63
1:A:265:LYS:NZ	1:A:265:LYS:HB3	2.12	0.63
1:B:15:GLY:HA3	3:B:480:FAD:H52A	1.80	0.63
1:F:249:LYS:HD2	1:F:255:ASP:CG	2.19	0.63
1:C:78:MET:HG2	1:D:78:MET:HE2	1.81	0.62
1:B:228:ARG:HH21	1:B:228:ARG:CB	2.12	0.62
1:F:15:GLY:HA3	3:F:480:FAD:H52A	1.79	0.62
1:B:297:ASP:HB2	1:B:298:PRO:HD2	1.81	0.62
1:A:174:LEU:HD12	1:A:197:TRP:CE2	2.33	0.62
1:H:191:VAL:HG12	1:H:360:THR:HG21	1.81	0.62
1:B:228:ARG:HB2	1:B:228:ARG:HH21	1.65	0.61
1:C:79:SER:OG	1:D:79:SER:HB2	2.00	0.61
1:E:50[B]:CYS:SG	1:E:51:ILE:HD12	2.40	0.61
1:B:33:VAL:HG12	1:B:113:VAL:HG11	1.82	0.60
1:A:15:GLY:HA3	3:A:480:FAD:H52A	1.82	0.60
1:B:23:ILE:O	1:B:27:GLN:HG3	2.02	0.60
1:E:188:VAL:HG12	4:E:481:NAI:H42N	1.84	0.59
1:G:188:VAL:HG12	4:G:481:NAI:H42N	1.84	0.59
1:G:78:MET:HG2	1:H:81:VAL:HG22	1.84	0.59
1:H:188:VAL:HG12	4:H:481:NAI:H42N	1.84	0.59
1:B:164:THR:OG1	1:B:254:ILE:HD11	2.03	0.59
1:E:224:LYS:O	1:E:228:ARG:HG3	2.02	0.59
1:H:297:ASP:HB2	1:H:298:PRO:HD2	1.84	0.59
1:E:184:ILE:HD13	1:E:243:VAL:HG21	1.84	0.59
1:G:179:GLU:HB3	1:G:272:ASP:OD2	2.01	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:180:LYS:HE3	5:H:2119:HOH:O	2.02	0.59
1:F:379:ILE:HA	2:F:2021:SO4:O3	2.01	0.59
1:E:15:GLY:HA3	3:E:480:FAD:H52A	1.85	0.59
1:G:5:ILE:CG1	1:G:139:ILE:HG13	2.33	0.59
1:A:79:SER:OG	1:B:79:SER:HB2	2.02	0.58
1:A:305:ASN:HB2	1:F:438:TYR:O	2.03	0.58
1:D:363:GLU:HB2	1:D:425:ALA:HB3	1.85	0.58
1:B:228:ARG:HG2	1:B:228:ARG:NH2	2.17	0.58
1:B:46:LEU:HD22	1:B:51:ILE:HG13	1.84	0.58
1:A:203:ASP:OD1	1:A:236:LYS:NZ	2.32	0.58
1:A:35:ILE:HD11	1:A:139:ILE:HD12	1.85	0.58
1:E:251:ASP:HB3	1:E:253:LYS:H	1.68	0.57
1:G:145:LEU:HD11	1:G:318:ILE:HG12	1.86	0.57
1:A:182:VAL:HG13	1:A:274:LEU:CD1	2.34	0.57
1:B:322:VAL:HG12	1:B:323:ALA:H	1.69	0.57
1:E:453:PRO:CD	1:F:50[B]:CYS:SG	2.86	0.57
1:G:249:LYS:HD2	1:G:255:ASP:CG	2.25	0.57
1:A:28:LEU:HD12	1:A:339:VAL:CG1	2.35	0.57
1:H:179:GLU:HB3	1:H:272:ASP:OD2	2.05	0.57
1:H:253:LYS:HE2	1:H:272:ASP:OD1	2.05	0.57
1:G:15:GLY:CA	3:G:480:FAD:H52A	2.34	0.56
1:D:46:LEU:HD21	1:D:99:LEU:CB	2.35	0.56
1:C:282:PRO:HG2	1:C:301:ARG:HG2	1.86	0.56
1:G:116:ASN:HD22	1:G:117:GLY:N	2.04	0.56
1:G:70:ASP:O	1:G:74:ARG:HG3	2.05	0.56
1:A:401:ASP:HB2	5:A:2105:HOH:O	2.05	0.56
1:D:207:VAL:HG13	1:D:241:THR:HB	1.87	0.56
1:E:339:VAL:CG1	1:E:342:MET:HE2	2.34	0.56
1:D:116:ASN:HD22	1:D:117:GLY:N	2.03	0.56
1:H:139:ILE:HD12	1:H:139:ILE:N	2.21	0.56
1:C:78:MET:CG	1:D:78:MET:HE2	2.37	0.55
1:E:78:MET:HG3	1:F:81:VAL:HG22	1.87	0.55
1:A:50[B]:CYS:SG	1:B:453:PRO:CD	2.92	0.55
1:B:35:ILE:HD11	1:B:139:ILE:HD12	1.88	0.55
1:D:15:GLY:HA3	3:D:480:FAD:H52A	1.88	0.55
1:H:380:GLU:HG3	1:H:410:LYS:HD2	1.89	0.55
1:B:325:PRO:HG2	1:B:330:LYS:HG3	1.88	0.55
1:C:28:LEU:HD12	1:C:339:VAL:CG1	2.36	0.55
1:G:318:ILE:HD13	1:G:335:GLY:HA2	1.89	0.55
1:B:174:LEU:HD23	1:B:178:PRO:HD3	1.87	0.55
1:A:363:GLU:HB2	1:A:425:ALA:HB3	1.89	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:282:PRO:CG	1:C:301:ARG:HG2	2.38	0.54
1:E:292:LEU:HD21	1:E:315:ILE:HD13	1.89	0.54
1:F:210:LEU:HD11	5:G:2107:HOH:O	2.07	0.54
1:D:43:GLY:HA2	3:D:480:FAD:O3B	2.07	0.54
1:E:392:SER:HB2	5:E:2090:HOH:O	2.08	0.54
1:G:379:ILE:HA	2:G:2025:SO4:O4	2.08	0.54
1:A:191:VAL:HG12	1:A:360:THR:HG21	1.91	0.54
1:B:33:VAL:HG21	1:B:139:ILE:HG21	1.90	0.54
1:C:31:LYS:NZ	1:C:31:LYS:HB2	2.22	0.54
1:E:318:ILE:HD13	1:E:335:GLY:HA2	1.90	0.54
1:D:191:VAL:HG12	1:D:360:THR:HG21	1.89	0.53
1:C:81:VAL:HG22	1:D:78:MET:HG2	1.90	0.53
1:F:231:GLN:HA	1:F:235:PHE:O	2.07	0.53
1:H:6:ASP:HA	1:H:140:ASP:O	2.09	0.53
1:B:43:GLY:HA2	3:B:480:FAD:O3B	2.07	0.53
1:D:8:ASP:OD1	1:D:142:LYS:HD2	2.08	0.53
1:E:33:VAL:HG22	1:E:113:VAL:HB	1.91	0.53
1:E:292:LEU:HD23	1:E:294:ILE:HD12	1.91	0.53
1:C:162:GLU:HA	1:C:166:VAL:HG12	1.91	0.53
1:F:80:GLU:HG2	1:F:82:ARG:HG3	1.89	0.53
1:A:445:ILE:O	1:A:448:VAL:HG22	2.08	0.53
1:C:15:GLY:HA3	3:C:480:FAD:H52A	1.91	0.53
1:A:182:VAL:HG13	1:A:274:LEU:HD12	1.91	0.53
3:G:480:FAD:H51A	5:G:2027:HOH:O	2.08	0.53
1:C:282:PRO:HG2	1:C:301:ARG:CG	2.38	0.52
1:F:15:GLY:CA	3:F:480:FAD:H52A	2.38	0.52
1:C:43:GLY:HA2	3:C:480:FAD:O3B	2.08	0.52
1:H:50[B]:CYS:SG	1:H:51:ILE:HD13	2.49	0.52
1:B:318:ILE:HD13	1:B:335:GLY:HA2	1.92	0.52
1:D:184:ILE:HD12	1:D:276:VAL:HG22	1.91	0.52
1:E:50[B]:CYS:O	1:E:54:LYS:HD2	2.10	0.52
1:F:35:ILE:HD11	1:F:139:ILE:HD13	1.91	0.52
1:F:162:GLU:HA	1:F:166:VAL:HG12	1.90	0.52
1:A:453:PRO:HG3	1:B:50[B]:CYS:SG	2.50	0.52
1:B:328:ALA:O	1:B:332:GLU:HG3	2.10	0.52
1:E:7:ALA:O	1:E:141:THR:HA	2.09	0.52
1:B:292:LEU:HD21	1:B:315:ILE:CD1	2.39	0.52
1:H:282:PRO:HG2	1:H:301:ARG:CG	2.41	0.51
1:A:110:ASN:O	1:A:111:LYS:HB2	2.11	0.51
1:G:256:VAL:HB	1:G:269:ILE:HG13	1.93	0.51
1:E:189:ILE:HG13	4:E:481:NAI:H6N	1.92	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:311:LYS:HB3	1:B:311:LYS:NZ	2.25	0.51
1:H:318:ILE:HD13	1:H:335:GLY:HA2	1.92	0.51
1:B:294:ILE:CD1	1:B:312:ILE:HD13	2.39	0.51
1:D:195:SER:O	1:D:199:ARG:HG3	2.10	0.51
1:F:105:HIS:C	1:F:105:HIS:CD2	2.83	0.51
1:G:80:GLU:OE1	1:G:82:ARG:HD3	2.11	0.51
1:A:448:VAL:HG23	1:A:450:HIS:CE1	2.45	0.51
1:C:339:VAL:HA	1:C:342:MET:HE2	1.91	0.51
1:C:392:SER:HB2	5:C:2114:HOH:O	2.08	0.51
1:D:389:ALA:HA	1:D:400:THR:HB	1.93	0.51
1:E:292:LEU:HD21	1:E:315:ILE:CD1	2.40	0.51
1:H:453:PRO:HA	1:H:457:GLU:OE1	2.11	0.51
1:B:224:LYS:HA	1:B:224:LYS:HE2	1.92	0.51
1:C:188:VAL:HG12	4:C:481:NAI:H42N	1.92	0.51
1:G:189:ILE:HG13	4:G:481:NAI:H6N	1.92	0.51
1:A:389:ALA:HA	1:A:400:THR:HB	1.93	0.50
1:G:5:ILE:HG13	1:G:139:ILE:HG13	1.92	0.50
1:G:228:ARG:HD3	5:G:2303:HOH:O	2.11	0.50
1:A:15:GLY:CA	3:A:480:FAD:H52A	2.41	0.50
1:A:182:VAL:HG11	1:A:256:VAL:HG21	1.92	0.50
1:A:402:GLY:HA3	1:A:421:LEU:O	2.11	0.50
1:H:292:LEU:HD23	1:H:294:ILE:HD12	1.94	0.50
1:D:46:LEU:HD22	1:D:100:THR:CG2	2.39	0.50
1:H:282:PRO:CG	1:H:301:ARG:HG2	2.42	0.50
1:H:15:GLY:HA3	3:H:480:FAD:H52A	1.94	0.50
1:A:254:ILE:HD11	1:A:256:VAL:HG22	1.93	0.50
1:G:249:LYS:HD2	1:G:255:ASP:OD1	2.11	0.50
1:B:188:VAL:HG12	4:B:481:NAI:H42N	1.93	0.50
1:F:188:VAL:HG12	4:F:481:NAI:H42N	1.94	0.50
1:A:209:PHE:HD1	1:A:242:LYS:HB2	1.76	0.49
1:D:66:ALA:HB1	1:D:78:MET:HE1	1.93	0.49
1:E:43:GLY:HA2	3:E:480:FAD:O3B	2.11	0.49
1:E:51:ILE:HD12	1:E:51:ILE:N	2.26	0.49
1:G:69:THR:HG23	5:G:2314:HOH:O	2.10	0.49
1:H:188:VAL:CG1	4:H:481:NAI:H42N	2.42	0.49
1:H:307:ARG:HA	1:H:347:VAL:HG11	1.94	0.49
1:H:5:ILE:CG1	1:H:139:ILE:HG13	2.42	0.49
1:A:453:PRO:HA	1:A:457:GLU:OE2	2.12	0.49
1:B:164:THR:HG21	1:B:248:LYS:NZ	2.28	0.49
1:E:281:ARG:HB2	1:E:282:PRO:HD2	1.94	0.49
1:H:461:GLU:OE2	1:H:471:SER:HB2	2.13	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:LEU:CD1	1:A:339:VAL:HG12	2.43	0.49
1:B:61:HIS:O	1:B:65:MET:HG2	2.12	0.49
1:C:461:GLU:HB3	1:C:471:SER:OG	2.12	0.49
1:G:15:GLY:N	3:G:480:FAD:H52A	2.27	0.49
1:C:471:SER:HB2	5:C:2123:HOH:O	2.12	0.49
1:E:297:ASP:HB2	1:E:298:PRO:HD2	1.95	0.49
1:G:50[B]:CYS:SG	1:H:453:PRO:HG3	2.53	0.49
1:H:292:LEU:HG	1:H:312:ILE:CD1	2.42	0.49
1:B:131:LYS:HE3	1:B:135:GLY:O	2.13	0.49
1:C:191:VAL:HG12	1:C:360:THR:CG2	2.42	0.49
1:C:174:LEU:HB2	1:C:197:TRP:CZ2	2.48	0.49
1:G:402:GLY:HA3	1:G:421:LEU:O	2.12	0.49
1:E:50[A]:CYS:O	1:E:54:LYS:HD2	2.12	0.49
1:E:188:VAL:CG1	4:E:481:NAI:H42N	2.43	0.49
1:H:146:ILE:O	1:H:317:ALA:HA	2.13	0.49
1:B:322:VAL:CG1	1:B:323:ALA:N	2.76	0.49
1:C:51:ILE:HG21	1:C:99:LEU:HD12	1.95	0.48
1:H:380:GLU:CG	1:H:410:LYS:HD2	2.43	0.48
1:E:82:ARG:HH11	1:E:82:ARG:HG3	1.78	0.48
1:G:242:LYS:NZ	1:G:261:ALA:O	2.45	0.48
1:E:336:ILE:O	1:E:339:VAL:HG23	2.13	0.48
1:A:265:LYS:HB3	1:A:265:LYS:HZ2	1.76	0.48
1:A:43:GLY:HA2	3:A:480:FAD:O3B	2.14	0.48
1:B:322:VAL:HG13	5:B:2139:HOH:O	2.14	0.48
1:B:32:THR:O	1:B:113:VAL:HG12	2.14	0.48
1:F:182:VAL:HG23	1:F:271:CYS:HB3	1.94	0.48
1:H:116:ASN:HD22	1:H:117:GLY:N	2.10	0.48
1:C:318:ILE:HD13	1:C:335:GLY:HA2	1.94	0.48
1:E:145:LEU:HD11	1:E:318:ILE:HG23	1.96	0.48
1:D:188:VAL:HG12	4:D:481:NAI:H42N	1.96	0.48
1:H:292:LEU:HD21	1:H:315:ILE:HD13	1.96	0.48
1:F:191:VAL:HG12	1:F:360:THR:HG21	1.96	0.48
1:G:105[A]:HIS:CD2	5:G:2332:HOH:O	2.65	0.48
1:B:191:VAL:HG11	1:B:226:PHE:CZ	2.49	0.48
1:B:69:THR:O	1:B:69:THR:HG22	2.13	0.48
1:F:121:ILE:HG21	1:F:292:LEU:HD11	1.96	0.48
1:F:43:GLY:HA2	3:F:480:FAD:O3B	2.14	0.48
1:G:448:VAL:HG22	1:H:437:GLU:HG2	1.96	0.47
1:G:50[B]:CYS:O	1:G:54:LYS:HD2	2.14	0.47
1:A:297:ASP:HB2	1:A:298:PRO:CD	2.44	0.47
1:H:282:PRO:HG2	1:H:301:ARG:HG2	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:297:ASP:HB2	1:D:298:PRO:CD	2.45	0.47
1:G:110:ASN:O	1:G:111:LYS:HB2	2.15	0.47
1:H:222:ILE:HG13	1:H:419:HIS:HB3	1.95	0.47
1:C:328:ALA:O	1:C:332:GLU:HG3	2.15	0.47
1:F:265:LYS:O	1:F:267:GLU:HG3	2.14	0.47
1:D:248:LYS:HD2	1:D:254:ILE:HG13	1.97	0.47
1:E:307:ARG:HB3	1:E:307:ARG:HH11	1.79	0.47
1:F:191:VAL:HG11	1:F:226:PHE:CZ	2.50	0.47
1:G:259:GLU:HB2	1:G:264:GLY:O	2.15	0.47
1:G:6:ASP:HA	1:G:140:ASP:O	2.15	0.47
1:B:228:ARG:O	1:B:232:LYS:HG3	2.14	0.47
1:B:231:GLN:HA	1:B:235:PHE:O	2.15	0.47
1:C:196:VAL:O	1:C:200:LEU:HG	2.15	0.47
1:G:410:LYS:HD2	1:G:410:LYS:HA	1.56	0.47
1:H:50[A]:CYS:O	1:H:54:LYS:HD2	2.15	0.47
1:B:292:LEU:HD21	1:B:315:ILE:HD12	1.97	0.46
1:D:231:GLN:HA	1:D:235:PHE:O	2.16	0.46
1:F:389:ALA:HA	1:F:400:THR:HB	1.96	0.46
1:E:50[B]:CYS:SG	1:F:453:PRO:HG3	2.55	0.46
1:B:24:LYS:O	1:B:28:LEU:HG	2.15	0.46
1:C:191:VAL:HG11	1:C:226:PHE:CZ	2.51	0.46
1:C:301:ARG:NH1	2:C:2010:SO4:O4	2.49	0.46
1:C:187:GLY:HA3	4:C:481:NAI:O5B	2.16	0.46
1:E:15:GLY:CA	3:E:480:FAD:H52A	2.44	0.46
1:B:174:LEU:HB2	1:B:197:TRP:CZ2	2.51	0.46
1:D:33:VAL:HG22	1:D:113:VAL:HB	1.98	0.46
1:E:35:ILE:CD1	1:E:115:VAL:HB	2.45	0.46
1:B:409:GLN:NE2	2:B:2005:SO4:O1	2.49	0.46
1:B:434:LEU:HD11	1:B:438:TYR:HE2	1.81	0.46
1:G:43:GLY:HA2	3:G:480:FAD:O3B	2.15	0.46
1:C:189:ILE:HG13	4:C:481:NAI:H6N	1.97	0.46
1:E:3:GLN:N	1:E:4:PRO:HD3	2.30	0.46
1:D:15:GLY:CA	3:D:480:FAD:H52A	2.45	0.45
1:C:7:ALA:O	1:C:141:THR:HA	2.16	0.45
1:D:46:LEU:HD12	1:D:103:ILE:HD11	1.97	0.45
1:D:189:ILE:HG13	4:D:481:NAI:H6N	1.98	0.45
1:D:66:ALA:HB1	1:D:78:MET:CE	2.47	0.45
1:G:404:VAL:HG11	1:G:459:PHE:HA	1.98	0.45
1:A:438:TYR:O	1:F:305:ASN:HB2	2.16	0.45
1:G:318:ILE:HD13	1:G:335:GLY:CA	2.46	0.45
1:F:309:GLN:HG2	1:F:316:TYR:CE2	2.52	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:472:ILE:HG22	1:A:473:ASN:ND2	2.31	0.45
1:B:322:VAL:CG1	1:B:323:ALA:H	2.28	0.45
1:C:61:HIS:O	1:C:65:MET:HG2	2.17	0.45
1:D:121:ILE:CG2	1:D:292:LEU:HD11	2.45	0.45
1:B:33:VAL:HA	1:B:113:VAL:CG1	2.46	0.45
1:D:254:ILE:HD12	1:D:254:ILE:N	2.32	0.45
1:D:299:ARG:HB2	1:D:301:ARG:HD3	1.99	0.45
1:E:183:VAL:O	1:E:206:ALA:HA	2.17	0.45
1:E:453:PRO:HA	1:E:457:GLU:OE1	2.17	0.45
1:A:195:SER:O	1:A:199:ARG:HG3	2.16	0.45
1:G:162:GLU:HA	1:G:166:VAL:HG12	1.98	0.45
1:E:152:VAL:HG21	1:E:167:SER:HB2	1.98	0.44
1:A:174:LEU:HB2	1:A:197:TRP:CZ2	2.52	0.44
1:H:288:GLY:O	1:H:292:LEU:HB2	2.16	0.44
1:E:82:ARG:NH1	1:E:82:ARG:HG3	2.33	0.44
1:G:435:ALA:HB1	1:G:445:ILE:HD11	2.00	0.44
1:A:174:LEU:HB2	1:A:197:TRP:CH2	2.53	0.44
1:C:24:LYS:O	1:C:28:LEU:HG	2.16	0.44
1:F:402:GLY:HA3	1:F:421:LEU:O	2.17	0.44
1:G:120:LYS:NZ	5:G:2329:HOH:O	2.33	0.44
1:H:31:LYS:HE3	1:H:113:VAL:HG23	1.98	0.44
1:B:228:ARG:HG2	1:B:228:ARG:HH21	1.78	0.44
1:F:208:GLU:HB3	1:F:239:LEU:HD23	2.00	0.44
1:F:80:GLU:CD	1:F:82:ARG:HD3	2.37	0.44
1:A:52:PRO:HG2	1:A:172:LEU:HD13	2.00	0.44
1:B:123:GLY:O	1:B:125:ASN:N	2.51	0.44
1:C:5:ILE:HB	1:C:139:ILE:HD13	1.99	0.44
1:G:265:LYS:O	1:G:267:GLU:HG3	2.17	0.44
1:G:392:SER:HB2	5:G:2284:HOH:O	2.18	0.44
1:G:50[B]:CYS:SG	1:H:453:PRO:CG	3.06	0.44
1:A:225:ASN:O	1:A:229:ILE:HG13	2.18	0.44
1:B:242:LYS:HE2	1:B:242:LYS:HB3	1.72	0.44
1:D:180:LYS:HD3	1:D:270:THR:O	2.17	0.44
1:C:71:PHE:HE2	1:D:66:ALA:HB2	1.83	0.44
1:A:254:ILE:HG13	1:A:254:ILE:O	2.18	0.44
1:D:220:MET:O	1:D:224:LYS:HG3	2.17	0.44
1:F:110:ASN:O	1:F:111:LYS:HB2	2.17	0.44
1:C:179:GLU:HB3	1:C:272:ASP:OD2	2.18	0.43
1:G:242:LYS:HE3	1:G:261:ALA:O	2.18	0.43
1:A:88:MET:HE3	1:A:200:LEU:HD11	2.00	0.43
1:A:426:GLY:O	1:A:429:VAL:HG12	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:ILE:HG13	1:B:419:HIS:HB3	1.99	0.43
1:B:33:VAL:HA	1:B:113:VAL:HG13	2.00	0.43
1:C:144:ILE:HD12	1:C:315:ILE:CD1	2.48	0.43
1:E:308:PHE:CZ	1:E:334:GLU:HG2	2.53	0.43
1:H:393:ARG:HD2	1:H:454:THR:HA	2.00	0.43
1:B:131:LYS:HB2	1:B:135:GLY:O	2.18	0.43
1:C:126:GLN:HG2	1:C:127:VAL:N	2.33	0.43
1:D:191:VAL:HG11	1:D:226:PHE:CZ	2.54	0.43
1:H:5:ILE:HG13	1:H:139:ILE:HG13	2.00	0.43
1:H:43:GLY:HA2	3:H:480:FAD:O3B	2.18	0.43
1:G:231:GLN:HA	1:G:235:PHE:O	2.18	0.43
1:F:249:LYS:HD2	1:F:255:ASP:OD1	2.17	0.43
1:H:224:LYS:O	1:H:228:ARG:HG3	2.18	0.43
1:H:50[B]:CYS:O	1:H:54:LYS:HD2	2.18	0.43
1:H:61:HIS:O	1:H:65:MET:HG2	2.19	0.43
1:A:231:GLN:HA	1:A:235:PHE:O	2.18	0.43
1:A:472:ILE:HD11	1:B:24:LYS:HD3	2.00	0.43
1:H:7:ALA:O	1:H:141:THR:HA	2.19	0.43
1:B:164:THR:HG21	1:B:248:LYS:CE	2.49	0.43
1:B:330:LYS:O	1:B:334:GLU:HG3	2.19	0.43
1:D:181:MET:HG3	1:D:273:VAL:HG23	2.01	0.43
1:A:15:GLY:N	3:A:480:FAD:H52A	2.33	0.43
1:B:188:VAL:CG1	4:B:481:NAI:H42N	2.49	0.43
1:B:191:VAL:HG12	1:B:360:THR:CG2	2.45	0.42
1:D:110:ASN:O	1:D:111:LYS:HB2	2.19	0.42
1:A:176:LYS:HD3	1:A:176:LYS:HA	1.75	0.42
1:C:222:ILE:HG13	1:C:419:HIS:HB3	2.01	0.42
1:D:15:GLY:N	3:D:480:FAD:H52A	2.34	0.42
1:D:3:GLN:HA	1:D:4:PRO:HD3	1.91	0.42
1:E:249:LYS:C	1:E:251:ASP:N	2.72	0.42
1:G:453:PRO:HA	1:G:457:GLU:OE2	2.20	0.42
1:A:5:ILE:HG23	5:A:2150:HOH:O	2.19	0.42
1:B:410:LYS:HE2	1:B:410:LYS:HB2	1.85	0.42
1:E:131:LYS:O	1:E:132:ALA:HB3	2.19	0.42
1:E:382:LYS:HD3	5:E:2123:HOH:O	2.18	0.42
1:F:191:VAL:HG11	1:F:226:PHE:CE2	2.54	0.42
1:G:242:LYS:CE	1:G:261:ALA:O	2.67	0.42
1:G:242:LYS:HE3	1:G:261:ALA:CA	2.49	0.42
1:G:121:ILE:HG21	1:G:292:LEU:HD11	2.00	0.42
1:G:80:GLU:O	1:H:78:MET:HA	2.19	0.42
1:H:5:ILE:O	1:H:139:ILE:HA	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:393:ARG:HD2	1:B:454:THR:HA	2.02	0.42
1:A:33:VAL:HG22	1:A:113:VAL:HB	2.02	0.42
1:B:162:GLU:CA	1:B:166:VAL:HG12	2.48	0.42
1:B:196:VAL:O	1:B:200:LEU:HG	2.20	0.42
1:B:80:GLU:HG2	1:B:82:ARG:NH2	2.35	0.42
1:C:307:ARG:O	1:C:308:PHE:HB2	2.20	0.42
1:E:248:LYS:NZ	1:E:252:GLY:HA2	2.35	0.42
1:E:241:THR:HG23	1:E:259:GLU:O	2.20	0.42
1:F:409:GLN:HG2	1:F:412:THR:OG1	2.19	0.42
1:A:121:ILE:CG2	1:A:292:LEU:HD11	2.47	0.42
1:C:330:LYS:O	1:C:334:GLU:HG3	2.20	0.42
1:E:307:ARG:O	1:E:308:PHE:HB2	2.20	0.42
1:G:373:GLN:O	1:G:377:GLU:HG3	2.19	0.42
1:H:282:PRO:HG2	1:H:301:ARG:HG3	2.02	0.42
1:H:336:ILE:O	1:H:340:GLU:HG3	2.20	0.42
1:H:50[B]:CYS:SG	1:H:51:ILE:N	2.91	0.42
1:A:382:LYS:HD2	1:A:466:ALA:O	2.20	0.42
1:B:258:ILE:HG13	1:B:267:GLU:HB2	2.02	0.42
1:E:96:VAL:O	1:E:100:THR:HG23	2.20	0.42
1:A:188:VAL:HG12	4:A:481:NAI:H42N	2.02	0.42
1:B:312:ILE:HD12	1:B:312:ILE:N	2.34	0.42
1:B:15:GLY:CA	3:B:480:FAD:H52A	2.47	0.42
1:B:187:GLY:HA3	4:B:481:NAI:O5B	2.20	0.42
1:E:222:ILE:HG13	1:E:419:HIS:HB3	2.01	0.42
1:E:79:SER:HB2	1:F:79:SER:OG	2.20	0.42
1:G:180:LYS:HG2	1:G:203:ASP:HB3	2.02	0.42
1:H:231:GLN:HA	1:H:235:PHE:O	2.20	0.42
1:B:216:VAL:HG22	5:B:2188:HOH:O	2.19	0.41
1:D:222:ILE:HG13	1:D:419:HIS:HB3	2.01	0.41
1:D:415:VAL:HG21	1:D:445:ILE:HD12	2.02	0.41
1:E:189:ILE:HG13	4:E:481:NAI:C6N	2.49	0.41
1:B:126:GLN:HG2	1:B:127:VAL:N	2.34	0.41
1:C:292:LEU:HB3	1:C:312:ILE:HD11	2.02	0.41
1:C:25:ALA:O	1:C:30:PHE:HB2	2.20	0.41
1:D:35:ILE:HD11	1:D:139:ILE:HD13	2.01	0.41
1:H:97:LYS:HE3	5:H:2203:HOH:O	2.20	0.41
1:A:222:ILE:HG13	1:A:419:HIS:HB3	2.01	0.41
1:B:461:GLU:OE1	1:B:471:SER:HB2	2.19	0.41
1:D:402:GLY:HA3	1:D:421:LEU:O	2.20	0.41
1:H:434:LEU:HD11	1:H:438:TYR:HE2	1.84	0.41
1:E:256:VAL:HG21	1:E:274:LEU:HD13	2.02	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:15:GLY:N	3:F:480:FAD:H52A	2.36	0.41
1:D:229:ILE:O	1:D:233:GLN:HG3	2.21	0.41
1:A:182:VAL:HG13	1:A:274:LEU:HD13	2.02	0.41
1:C:188:VAL:CG1	4:C:481:NAI:H42N	2.50	0.41
1:H:228:ARG:HH11	1:H:228:ARG:CG	2.26	0.41
1:B:148:THR:O	3:B:480:FAD:H8A	2.21	0.41
1:C:310:THR:HG23	1:C:315:ILE:O	2.21	0.41
3:B:480:FAD:N5	4:B:481:NAI:H4N	2.36	0.41
1:D:328:ALA:O	1:D:332:GLU:HG3	2.20	0.41
1:E:265:LYS:O	1:E:265:LYS:HG3	2.21	0.41
1:H:404:VAL:HG11	1:H:459:PHE:HA	2.03	0.41
1:H:136:THR:CG2	1:H:137:GLN:N	2.84	0.41
1:H:187:GLY:HA3	4:H:481:NAI:O5B	2.21	0.41
1:B:182:VAL:HG23	1:B:271:CYS:HB3	2.02	0.41
1:E:45[B]:CYS:SG	1:E:46:LEU:HD12	2.61	0.41
1:C:23:ILE:O	1:C:27:GLN:HG3	2.22	0.40
1:C:89:MET:O	1:C:92:LYS:HB3	2.21	0.40
1:D:244:THR:HG21	1:D:259:GLU:OE1	2.21	0.40
1:F:386:PHE:CD2	1:F:461:GLU:HB3	2.56	0.40
1:F:394:ALA:HB1	1:F:399:ASP:O	2.20	0.40
1:B:7:ALA:HB2	1:B:33:VAL:HG13	2.02	0.40
1:E:410:LYS:HE2	1:E:410:LYS:HB2	1.81	0.40
1:G:189:ILE:HG13	4:G:481:NAI:C6N	2.51	0.40
1:C:231:GLN:HA	1:C:235:PHE:O	2.21	0.40
1:G:42:GLY:O	1:G:46:LEU:HD23	2.21	0.40
1:A:182:VAL:CG1	1:A:274:LEU:HD13	2.51	0.40
1:F:6:ASP:HA	1:F:140:ASP:O	2.20	0.40
1:A:249:LYS:HD3	1:A:255:ASP:OD2	2.22	0.40
1:B:311:LYS:HZ3	1:B:311:LYS:HB3	1.87	0.40
3:C:480:FAD:N5	4:C:481:NAI:H4N	2.36	0.40
1:C:50[B]:CYS:SG	1:D:453:PRO:HD3	2.62	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	473/474 (100%)	455 (96%)	18 (4%)	0	100	100
1	B	471/474 (99%)	451 (96%)	19 (4%)	1 (0%)	47	47
1	C	471/474 (99%)	452 (96%)	18 (4%)	1 (0%)	47	47
1	D	472/474 (100%)	455 (96%)	16 (3%)	1 (0%)	47	47
1	E	473/474 (100%)	455 (96%)	15 (3%)	3 (1%)	25	20
1	F	471/474 (99%)	459 (98%)	11 (2%)	1 (0%)	47	47
1	G	473/474 (100%)	456 (96%)	16 (3%)	1 (0%)	47	47
1	H	471/474 (99%)	457 (97%)	14 (3%)	0	100	100
All	All	3775/3792 (100%)	3640 (96%)	127 (3%)	8 (0%)	47	47

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	79	SER
1	B	124	LYS
1	E	131	LYS
1	F	79	SER
1	D	80	GLU
1	E	265	LYS
1	C	313	PRO
1	E	313	PRO

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	375/373 (100%)	362 (96%)	13 (4%)	36	36
1	B	373/373 (100%)	366 (98%)	7 (2%)	57	61
1	C	373/373 (100%)	366 (98%)	7 (2%)	57	61
1	D	374/373 (100%)	363 (97%)	11 (3%)	42	44

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	375/373 (100%)	368 (98%)	7 (2%)	57	61
1	F	373/373 (100%)	366 (98%)	7 (2%)	57	61
1	G	375/373 (100%)	366 (98%)	9 (2%)	49	52
1	H	373/373 (100%)	362 (97%)	11 (3%)	42	44
All	All	2991/2984 (100%)	2919 (98%)	72 (2%)	49	52

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	TYR
1	A	54	LYS
1	A	78	MET
1	A	116	ASN
1	A	124	LYS
1	A	232	LYS
1	A	235	PHE
1	A	242	LYS
1	A	254	ILE
1	A	410	LYS
1	A	431	GLU
1	A	438	TYR
1	A	472	ILE
1	B	19	TYR
1	B	54	LYS
1	B	174	LEU
1	B	228	ARG
1	B	235	PHE
1	B	333	ASP
1	B	380	GLU
1	C	6	ASP
1	C	19	TYR
1	C	31	LYS
1	C	54	LYS
1	C	235	PHE
1	C	321	VAL
1	C	471	SER
1	D	19	TYR
1	D	46	LEU
1	D	54	LYS
1	D	116	ASN
1	D	124	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	235	PHE
1	D	290	GLU
1	D	295	GLU
1	D	301	ARG
1	D	414	ARG
1	D	438	TYR
1	E	19	TYR
1	E	54	LYS
1	E	58	ASN
1	E	235	PHE
1	E	251	ASP
1	E	306	THR
1	E	380	GLU
1	F	19	TYR
1	F	54	LYS
1	F	105	HIS
1	F	131	LYS
1	F	179	GLU
1	F	235	PHE
1	F	414	ARG
1	G	3	GLN
1	G	19	TYR
1	G	54	LYS
1	G	116	ASN
1	G	136	THR
1	G	213	VAL
1	G	235	PHE
1	G	410	LYS
1	G	438	TYR
1	H	19	TYR
1	H	31	LYS
1	H	54	LYS
1	H	78	MET
1	H	116	ASN
1	H	131	LYS
1	H	224	LYS
1	H	232	LYS
1	H	235	PHE
1	H	253	LYS
1	H	291	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	473	ASN
1	B	67	HIS
1	B	109	GLN
1	C	58	ASN
1	C	137	GLN
1	D	116	ASN
1	E	225	ASN
1	E	231	GLN
1	G	116	ASN
1	G	137	GLN
1	H	67	HIS
1	H	114	HIS
1	H	116	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

47 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	$\# Z > 2$
2	SO4	E	2017	-	4,4,4	0.62	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	FAD	B	480	-	51,58,58	1.78	9 (17%)	60,89,89	2.08	10 (16%)
4	NAI	F	481	-	42,48,48	1.41	5 (11%)	47,73,73	1.35	5 (10%)
2	SO4	B	2006	-	4,4,4	0.63	0	6,6,6	0.05	0
2	SO4	A	2002	-	4,4,4	0.62	0	6,6,6	0.06	0
3	FAD	F	480	-	51,58,58	1.84	11 (21%)	60,89,89	2.12	13 (21%)
3	FAD	G	480	-	51,58,58	1.72	10 (19%)	60,89,89	1.93	11 (18%)
2	SO4	D	2014	-	4,4,4	0.61	0	6,6,6	0.06	0
4	NAI	A	481	-	42,48,48	1.41	5 (11%)	47,73,73	1.37	5 (10%)
2	SO4	E	2016	-	4,4,4	0.62	0	6,6,6	0.04	0
4	NAI	E	481	-	42,48,48	1.40	5 (11%)	47,73,73	1.30	5 (10%)
2	SO4	B	2004	-	4,4,4	0.61	0	6,6,6	0.05	0
2	SO4	D	2011	-	4,4,4	0.62	0	6,6,6	0.07	0
2	SO4	E	2015	-	4,4,4	0.60	0	6,6,6	0.07	0
2	SO4	A	2000	-	4,4,4	0.62	0	6,6,6	0.04	0
2	SO4	G	2026	-	4,4,4	0.61	0	6,6,6	0.05	0
2	SO4	F	2020	-	4,4,4	0.63	0	6,6,6	0.06	0
3	FAD	C	480	-	51,58,58	2.02	9 (17%)	60,89,89	2.07	9 (15%)
4	NAI	G	481	-	42,48,48	1.41	5 (11%)	47,73,73	1.32	5 (10%)
4	NAI	C	481	-	42,48,48	1.40	5 (11%)	47,73,73	1.30	5 (10%)
2	SO4	A	2001	-	4,4,4	0.62	0	6,6,6	0.07	0
2	SO4	F	2021	-	4,4,4	0.62	0	6,6,6	0.05	0
2	SO4	H	2030	-	4,4,4	0.61	0	6,6,6	0.07	0
2	SO4	H	2028	-	4,4,4	0.62	0	6,6,6	0.05	0
2	SO4	F	2022	-	4,4,4	0.63	0	6,6,6	0.05	0
2	SO4	D	2012	-	4,4,4	0.62	0	6,6,6	0.06	0
2	SO4	G	2025	-	4,4,4	0.63	0	6,6,6	0.04	0
2	SO4	C	2008	-	4,4,4	0.62	0	6,6,6	0.05	0
4	NAI	H	481	-	42,48,48	1.39	5 (11%)	47,73,73	1.32	5 (10%)
2	SO4	C	2010	-	4,4,4	0.63	0	6,6,6	0.06	0
2	SO4	G	2024	-	4,4,4	0.64	0	6,6,6	0.09	0
2	SO4	B	2005	-	4,4,4	0.63	0	6,6,6	0.10	0
2	SO4	H	2027	-	4,4,4	0.62	0	6,6,6	0.05	0
3	FAD	H	480	-	51,58,58	1.71	10 (19%)	60,89,89	2.20	10 (16%)
4	NAI	D	481	-	42,48,48	1.42	5 (11%)	47,73,73	1.36	5 (10%)
2	SO4	D	2013	-	4,4,4	0.62	0	6,6,6	0.05	0
2	SO4	G	2023	-	4,4,4	0.61	0	6,6,6	0.07	0
2	SO4	B	2003	-	4,4,4	0.61	0	6,6,6	0.07	0
2	SO4	C	2009	-	4,4,4	0.64	0	6,6,6	0.12	0
4	NAI	B	481	-	42,48,48	1.40	5 (11%)	47,73,73	1.32	5 (10%)
2	SO4	C	2007	-	4,4,4	0.62	0	6,6,6	0.05	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	F	2019	-	4,4,4	0.62	0	6,6,6	0.07	0
2	SO4	H	2029	-	4,4,4	0.63	0	6,6,6	0.08	0
3	FAD	E	480	-	51,58,58	1.99	12 (23%)	60,89,89	2.10	9 (15%)
3	FAD	A	480	-	51,58,58	1.74	10 (19%)	60,89,89	1.97	9 (15%)
3	FAD	D	480	-	51,58,58	1.78	11 (21%)	60,89,89	2.03	8 (13%)
2	SO4	E	2018	-	4,4,4	0.62	0	6,6,6	0.06	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
3	FAD	B	480	-	-	3/30/50/50	0/6/6/6
3	FAD	C	480	-	-	4/30/50/50	0/6/6/6
3	FAD	F	480	-	-	2/30/50/50	0/6/6/6
3	FAD	A	480	-	-	3/30/50/50	0/6/6/6
3	FAD	D	480	-	-	4/30/50/50	0/6/6/6
3	FAD	G	480	-	-	1/30/50/50	0/6/6/6
3	FAD	E	480	-	-	2/30/50/50	0/6/6/6
3	FAD	H	480	-	-	1/30/50/50	0/6/6/6
4	NAI	C	481	-	-	3/25/72/72	0/5/5/5
4	NAI	F	481	-	-	2/25/72/72	0/5/5/5
4	NAI	E	481	-	-	3/25/72/72	0/5/5/5
4	NAI	D	481	-	-	2/25/72/72	0/5/5/5
4	NAI	H	481	-	-	4/25/72/72	0/5/5/5
4	NAI	B	481	-	-	4/25/72/72	0/5/5/5
4	NAI	A	481	-	-	2/25/72/72	0/5/5/5
4	NAI	G	481	-	-	2/25/72/72	0/5/5/5

All (122) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	480	FAD	C4X-N5	6.21	1.42	1.33
3	C	480	FAD	C4X-C10	5.79	1.44	1.38
3	E	480	FAD	C4X-C10	5.79	1.44	1.38
3	E	480	FAD	C4X-N5	5.55	1.41	1.33
3	E	480	FAD	C10-N1	5.46	1.40	1.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	481	NAI	C4N-C3N	-5.31	1.39	1.49
4	A	481	NAI	C4N-C3N	-5.30	1.39	1.49
4	G	481	NAI	C4N-C3N	-5.28	1.39	1.49
4	B	481	NAI	C4N-C3N	-5.27	1.39	1.49
4	F	481	NAI	C4N-C3N	-5.27	1.39	1.49
4	E	481	NAI	C4N-C3N	-5.21	1.39	1.49
4	C	481	NAI	C4N-C3N	-5.16	1.39	1.49
3	H	480	FAD	C10-N1	5.12	1.39	1.33
4	H	481	NAI	C4N-C3N	-5.12	1.39	1.49
3	F	480	FAD	C4X-C10	5.08	1.43	1.38
3	C	480	FAD	C10-N1	5.04	1.39	1.33
3	A	480	FAD	C4X-C10	5.01	1.43	1.38
3	A	480	FAD	C4X-N5	4.90	1.40	1.33
3	B	480	FAD	C4X-N5	4.78	1.40	1.33
3	D	480	FAD	C4X-N5	4.78	1.40	1.33
3	D	480	FAD	C4X-C10	4.77	1.43	1.38
3	F	480	FAD	C4X-N5	4.73	1.40	1.33
3	G	480	FAD	C4-N3	4.61	1.41	1.33
3	B	480	FAD	C10-N1	4.60	1.39	1.33
3	G	480	FAD	C4X-N5	4.33	1.39	1.33
3	C	480	FAD	C4-N3	4.30	1.40	1.33
3	C	480	FAD	C9A-N10	4.30	1.44	1.38
3	F	480	FAD	O4B-C1B	4.28	1.47	1.41
3	H	480	FAD	C4X-N5	4.19	1.39	1.33
3	G	480	FAD	C10-N1	4.17	1.38	1.33
3	A	480	FAD	C4-N3	4.03	1.40	1.33
3	E	480	FAD	C4-N3	3.89	1.39	1.33
3	B	480	FAD	C4X-C10	3.87	1.42	1.38
3	G	480	FAD	C2A-N3A	3.81	1.38	1.32
3	F	480	FAD	C2A-N3A	3.80	1.38	1.32
3	D	480	FAD	C4-N3	3.79	1.39	1.33
3	H	480	FAD	C4X-C10	3.76	1.42	1.38
3	F	480	FAD	C10-N1	3.73	1.38	1.33
3	C	480	FAD	C5X-N5	3.72	1.41	1.35
3	B	480	FAD	C4-N3	3.72	1.39	1.33
4	D	481	NAI	C6N-C5N	3.67	1.39	1.33
3	B	480	FAD	C5X-N5	3.63	1.41	1.35
4	E	481	NAI	C6N-C5N	3.62	1.39	1.33
4	F	481	NAI	C6N-C5N	3.62	1.39	1.33
4	B	481	NAI	C6N-C5N	3.59	1.39	1.33
4	C	481	NAI	C4N-C5N	-3.58	1.39	1.48
4	G	481	NAI	C6N-C5N	3.57	1.39	1.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	480	FAD	C5B-C4B	3.56	1.62	1.51
3	C	480	FAD	C2A-N3A	3.53	1.37	1.32
4	H	481	NAI	C6N-C5N	3.53	1.39	1.33
4	F	481	NAI	C4N-C5N	-3.52	1.39	1.48
4	A	481	NAI	C4N-C5N	-3.52	1.39	1.48
4	C	481	NAI	C6N-C5N	3.51	1.39	1.33
4	D	481	NAI	C4N-C5N	-3.50	1.39	1.48
4	B	481	NAI	C4N-C5N	-3.50	1.39	1.48
4	A	481	NAI	C6N-C5N	3.49	1.39	1.33
4	G	481	NAI	C4N-C5N	-3.48	1.39	1.48
3	F	480	FAD	C4-N3	3.46	1.39	1.33
4	H	481	NAI	C4N-C5N	-3.46	1.39	1.48
3	A	480	FAD	C10-N1	3.46	1.37	1.33
3	E	480	FAD	C9A-N10	3.46	1.43	1.38
3	G	480	FAD	O4B-C1B	3.46	1.45	1.41
4	E	481	NAI	C4N-C5N	-3.42	1.40	1.48
3	H	480	FAD	C9A-N10	3.34	1.43	1.38
3	D	480	FAD	C9A-N10	3.34	1.43	1.38
3	D	480	FAD	C5X-N5	3.34	1.40	1.35
3	B	480	FAD	C2A-N3A	3.32	1.37	1.32
3	G	480	FAD	C4X-C10	3.32	1.42	1.38
3	D	480	FAD	C10-N1	3.32	1.37	1.33
3	A	480	FAD	C9A-N10	3.30	1.43	1.38
3	D	480	FAD	O4B-C1B	3.23	1.45	1.41
3	E	480	FAD	C2A-N3A	3.18	1.37	1.32
3	F	480	FAD	C5B-C4B	3.12	1.61	1.51
3	E	480	FAD	C5X-N5	3.04	1.40	1.35
3	H	480	FAD	C4-N3	2.99	1.38	1.33
3	F	480	FAD	C9A-N10	2.85	1.42	1.38
3	B	480	FAD	C9A-N10	2.81	1.42	1.38
3	E	480	FAD	C1'-N10	2.81	1.51	1.48
3	H	480	FAD	C5B-C4B	2.78	1.60	1.51
3	A	480	FAD	C2A-N1A	2.69	1.38	1.33
3	H	480	FAD	C5X-N5	2.68	1.39	1.35
3	B	480	FAD	O4B-C1B	2.67	1.44	1.41
3	B	480	FAD	C2A-N1A	2.66	1.38	1.33
3	D	480	FAD	C5B-C4B	2.66	1.59	1.51
3	H	480	FAD	O4B-C1B	2.65	1.44	1.41
3	A	480	FAD	C2A-N3A	2.65	1.36	1.32
3	A	480	FAD	C5X-N5	2.64	1.39	1.35
3	E	480	FAD	C5B-C4B	2.61	1.59	1.51
3	H	480	FAD	C2A-N3A	2.58	1.36	1.32

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	480	FAD	C2A-N1A	2.57	1.38	1.33
3	H	480	FAD	C2A-N1A	2.56	1.38	1.33
4	G	481	NAI	O4B-C1B	2.55	1.44	1.41
3	A	480	FAD	O4B-C1B	2.51	1.44	1.41
4	F	481	NAI	O4B-C1B	2.51	1.44	1.41
4	H	481	NAI	O4B-C1B	2.49	1.44	1.41
3	C	480	FAD	C2A-N1A	2.45	1.38	1.33
3	F	480	FAD	C5X-N5	2.45	1.39	1.35
4	A	481	NAI	O4B-C1B	2.43	1.44	1.41
3	G	480	FAD	C2A-N1A	2.41	1.38	1.33
3	E	480	FAD	O4B-C1B	2.39	1.44	1.41
3	F	480	FAD	C1'-N10	2.37	1.50	1.48
3	D	480	FAD	C9-C8	2.37	1.43	1.37
3	F	480	FAD	P-O2P	-2.35	1.44	1.55
4	C	481	NAI	O4B-C1B	2.32	1.44	1.41
4	D	481	NAI	O4B-C1B	2.29	1.44	1.41
4	B	481	NAI	O4B-C1B	2.27	1.44	1.41
4	C	481	NAI	C7N-C3N	2.26	1.53	1.48
4	E	481	NAI	O4B-C1B	2.25	1.44	1.41
3	C	480	FAD	O4B-C1B	2.24	1.44	1.41
3	D	480	FAD	C2A-N3A	2.23	1.35	1.32
4	E	481	NAI	C7N-C3N	2.22	1.53	1.48
3	A	480	FAD	C5B-C4B	2.19	1.58	1.51
4	B	481	NAI	C7N-C3N	2.17	1.53	1.48
4	A	481	NAI	C7N-C3N	2.17	1.53	1.48
4	G	481	NAI	C7N-C3N	2.15	1.53	1.48
3	G	480	FAD	C9A-N10	2.14	1.41	1.38
4	D	481	NAI	C7N-C3N	2.14	1.53	1.48
4	F	481	NAI	C7N-C3N	2.10	1.53	1.48
3	D	480	FAD	C2A-N1A	2.08	1.37	1.33
3	G	480	FAD	C5X-N5	2.05	1.38	1.35
4	H	481	NAI	C7N-C3N	2.04	1.53	1.48
3	E	480	FAD	C9A-C5X	2.02	1.46	1.42

All (119) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	480	FAD	C4-N3-C2	11.91	125.20	115.14
3	E	480	FAD	C4-N3-C2	11.57	124.91	115.14
3	C	480	FAD	C4-N3-C2	11.17	124.58	115.14
3	B	480	FAD	C4-N3-C2	11.17	124.57	115.14
3	F	480	FAD	C4-N3-C2	10.90	124.35	115.14

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	480	FAD	C4-N3-C2	10.29	123.83	115.14
3	A	480	FAD	C4-N3-C2	10.18	123.74	115.14
3	G	480	FAD	C4-N3-C2	9.68	123.31	115.14
3	D	480	FAD	O4B-C1B-C2B	-5.13	99.42	106.93
3	A	480	FAD	O4B-C1B-C2B	-5.09	99.49	106.93
4	B	481	NAI	N3A-C2A-N1A	-5.00	120.86	128.68
4	D	481	NAI	N3A-C2A-N1A	-5.00	120.87	128.68
4	H	481	NAI	N3A-C2A-N1A	-4.99	120.88	128.68
4	A	481	NAI	N3A-C2A-N1A	-4.97	120.91	128.68
3	F	480	FAD	C4X-C4-N3	-4.95	116.66	123.43
4	F	481	NAI	N3A-C2A-N1A	-4.88	121.05	128.68
3	C	480	FAD	C4X-C4-N3	-4.87	116.77	123.43
4	C	481	NAI	N3A-C2A-N1A	-4.82	121.14	128.68
4	G	481	NAI	N3A-C2A-N1A	-4.81	121.16	128.68
4	E	481	NAI	N3A-C2A-N1A	-4.72	121.29	128.68
3	B	480	FAD	C4X-C4-N3	-4.67	117.05	123.43
3	D	480	FAD	C4X-C4-N3	-4.63	117.10	123.43
4	A	481	NAI	O4B-C1B-C2B	-4.43	100.44	106.93
3	H	480	FAD	C4X-C4-N3	-4.38	117.44	123.43
3	G	480	FAD	C4X-N5-C5X	4.32	121.09	116.77
3	H	480	FAD	C4X-N5-C5X	4.31	121.08	116.77
3	B	480	FAD	O4B-C1B-C2B	-4.21	100.78	106.93
3	H	480	FAD	C1'-N10-C9A	4.18	121.59	118.29
3	A	480	FAD	C4X-N5-C5X	4.18	120.95	116.77
3	A	480	FAD	C4X-C4-N3	-4.18	117.72	123.43
3	G	480	FAD	C4X-C4-N3	-4.14	117.77	123.43
3	E	480	FAD	C4X-C4-N3	-4.09	117.83	123.43
3	H	480	FAD	O4B-C1B-C2B	-4.01	101.06	106.93
3	D	480	FAD	C1'-N10-C9A	3.97	121.42	118.29
3	F	480	FAD	C4X-N5-C5X	3.96	120.73	116.77
3	E	480	FAD	C4-C4X-C10	-3.91	117.36	119.95
3	E	480	FAD	C1'-N10-C9A	3.83	121.31	118.29
4	G	481	NAI	O4B-C1B-C2B	-3.82	101.35	106.93
4	D	481	NAI	O4B-C1B-C2B	-3.81	101.36	106.93
3	C	480	FAD	C1'-N10-C9A	3.79	121.28	118.29
3	G	480	FAD	O4B-C1B-C2B	-3.76	101.43	106.93
3	B	480	FAD	C4X-N5-C5X	3.75	120.52	116.77
3	D	480	FAD	C4X-N5-C5X	3.69	120.45	116.77
3	F	480	FAD	O4B-C1B-C2B	-3.68	101.55	106.93
3	H	480	FAD	C4-C4X-C10	-3.65	117.54	119.95
3	E	480	FAD	O4B-C1B-C2B	-3.63	101.62	106.93
3	C	480	FAD	C4X-N5-C5X	3.61	120.37	116.77

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	480	FAD	C4X-N5-C5X	3.50	120.27	116.77
3	C	480	FAD	O4B-C1B-C2B	-3.43	101.92	106.93
3	B	480	FAD	C1'-N10-C9A	3.34	120.92	118.29
3	C	480	FAD	C4-C4X-C10	-3.23	117.81	119.95
3	B	480	FAD	C4-C4X-C10	-3.18	117.84	119.95
4	E	481	NAI	O4B-C1B-C2B	-3.05	102.46	106.93
4	F	481	NAI	O4B-C1B-C2B	-2.97	102.59	106.93
4	D	481	NAI	PN-O3-PA	-2.93	122.78	132.83
3	F	480	FAD	C4-C4X-C10	-2.92	118.02	119.95
3	F	480	FAD	C1'-N10-C9A	2.91	120.58	118.29
4	C	481	NAI	O4B-C1B-C2B	-2.90	102.69	106.93
4	F	481	NAI	O4D-C1D-N1N	2.87	113.67	108.06
3	D	480	FAD	C4-C4X-C10	-2.84	118.07	119.95
4	F	481	NAI	PN-O3-PA	-2.84	123.09	132.83
4	H	481	NAI	O4D-C1D-N1N	2.80	113.52	108.06
4	B	481	NAI	O4B-C1B-C2B	-2.78	102.86	106.93
3	F	480	FAD	N3A-C2A-N1A	-2.77	124.35	128.68
4	B	481	NAI	O4D-C1D-N1N	2.72	113.38	108.06
3	H	480	FAD	C4-C4X-N5	2.70	121.69	118.60
4	F	481	NAI	C3N-C2N-N1N	-2.69	119.25	123.10
3	E	480	FAD	C5A-C6A-N6A	2.68	124.43	120.35
3	G	480	FAD	PA-O5B-C5B	2.67	137.32	121.68
3	E	480	FAD	C4-C4X-N5	2.66	121.64	118.60
4	B	481	NAI	PN-O3-PA	-2.66	123.70	132.83
4	G	481	NAI	C3N-C2N-N1N	-2.65	119.31	123.10
4	E	481	NAI	PN-O3-PA	-2.64	123.78	132.83
4	G	481	NAI	PN-O3-PA	-2.59	123.92	132.83
4	C	481	NAI	O4D-C1D-N1N	2.58	113.10	108.06
3	B	480	FAD	C4-C4X-N5	2.57	121.53	118.60
3	F	480	FAD	C4-C4X-N5	2.55	121.52	118.60
4	C	481	NAI	PN-O3-PA	-2.55	124.07	132.83
4	D	481	NAI	O4D-C1D-N1N	2.55	113.04	108.06
3	H	480	FAD	N3A-C2A-N1A	-2.53	124.72	128.68
4	A	481	NAI	PN-O3-PA	-2.52	124.19	132.83
4	A	481	NAI	C3N-C2N-N1N	-2.51	119.52	123.10
4	H	481	NAI	C3N-C2N-N1N	-2.50	119.53	123.10
4	D	481	NAI	C3N-C2N-N1N	-2.50	119.53	123.10
3	D	480	FAD	N3A-C2A-N1A	-2.50	124.78	128.68
3	C	480	FAD	C4-C4X-N5	2.48	121.43	118.60
3	F	480	FAD	C5X-C9A-N10	2.47	119.51	117.72
4	C	481	NAI	C3N-C2N-N1N	-2.47	119.57	123.10
3	A	480	FAD	C1'-N10-C9A	2.47	120.23	118.29

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	481	NAI	C3N-C2N-N1N	-2.45	119.60	123.10
3	C	480	FAD	C5A-C6A-N6A	2.45	124.07	120.35
4	H	481	NAI	PN-O3-PA	-2.43	124.48	132.83
3	F	480	FAD	PA-O5B-C5B	2.42	135.90	121.68
4	H	481	NAI	O4B-C1B-C2B	-2.42	103.38	106.93
3	B	480	FAD	N3A-C2A-N1A	-2.42	124.89	128.68
4	E	481	NAI	C3N-C2N-N1N	-2.41	119.66	123.10
4	G	481	NAI	O4D-C1D-N1N	2.40	112.76	108.06
4	E	481	NAI	O4D-C1D-N1N	2.39	112.73	108.06
3	G	480	FAD	C4-C4X-C10	-2.35	118.40	119.95
3	F	480	FAD	C5A-C6A-N6A	2.33	123.89	120.35
3	G	480	FAD	C1'-N10-C9A	2.32	120.12	118.29
3	G	480	FAD	N3A-C2A-N1A	-2.32	125.06	128.68
4	A	481	NAI	O4D-C1D-N1N	2.31	112.57	108.06
3	D	480	FAD	C4-C4X-N5	2.31	121.23	118.60
3	B	480	FAD	P-O3P-PA	2.24	140.50	132.83
3	F	480	FAD	P-O3P-PA	2.21	140.40	132.83
3	A	480	FAD	C4-C4X-C10	-2.20	118.49	119.95
3	C	480	FAD	N3A-C2A-N1A	-2.15	125.31	128.68
3	G	480	FAD	O4B-C4B-C5B	2.13	116.39	109.37
3	H	480	FAD	C5A-C6A-N6A	2.13	123.58	120.35
3	A	480	FAD	N3A-C2A-N1A	-2.12	125.36	128.68
3	H	480	FAD	PA-O5B-C5B	2.11	134.08	121.68
3	A	480	FAD	P-O3P-PA	2.10	140.03	132.83
3	A	480	FAD	C4-C4X-N5	2.07	120.97	118.60
3	F	480	FAD	O4B-C4B-C5B	2.07	116.19	109.37
3	G	480	FAD	C4-C4X-N5	2.06	120.95	118.60
3	B	480	FAD	C5A-C6A-N6A	2.05	123.47	120.35
3	G	480	FAD	C5A-C6A-N6A	2.05	123.47	120.35
3	E	480	FAD	N3A-C2A-N1A	-2.04	125.49	128.68

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	481	NAI	O4D-C1D-N1N-C2N
4	A	481	NAI	O4D-C1D-N1N-C2N
4	E	481	NAI	O4D-C1D-N1N-C2N
4	C	481	NAI	O4D-C1D-N1N-C2N
4	H	481	NAI	O4D-C1D-N1N-C2N
4	D	481	NAI	O4D-C1D-N1N-C2N
4	B	481	NAI	O4D-C1D-N1N-C2N

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
4	G	481	NAI	O4D-C1D-N1N-C2N
4	H	481	NAI	PA-O3-PN-O2N
4	B	481	NAI	PA-O3-PN-O2N
3	F	480	FAD	O4B-C4B-C5B-O5B
3	F	480	FAD	PA-O3P-P-O5'
3	G	480	FAD	PA-O3P-P-O5'
3	B	480	FAD	PA-O3P-P-O5'
3	C	480	FAD	PA-O3P-P-O5'
3	H	480	FAD	PA-O3P-P-O5'
3	A	480	FAD	PA-O3P-P-O5'
3	D	480	FAD	PA-O3P-P-O5'
3	D	480	FAD	O4B-C4B-C5B-O5B
4	H	481	NAI	PA-O3-PN-O1N
4	B	481	NAI	PA-O3-PN-O1N
4	E	481	NAI	O4B-C4B-C5B-O5B
3	C	480	FAD	O4B-C4B-C5B-O5B
4	C	481	NAI	O4B-C4B-C5B-O5B
4	E	481	NAI	PA-O3-PN-O1N
3	B	480	FAD	P-O3P-PA-O1A
3	C	480	FAD	P-O3P-PA-O1A
4	C	481	NAI	PA-O3-PN-O1N
3	A	480	FAD	P-O3P-PA-O1A
3	D	480	FAD	P-O3P-PA-O1A
3	E	480	FAD	PA-O3P-P-O5'
3	E	480	FAD	O4B-C4B-C5B-O5B
4	F	481	NAI	O4B-C4B-C5B-O5B
4	A	481	NAI	O4B-C4B-C5B-O5B
4	G	481	NAI	O4B-C4B-C5B-O5B
4	D	481	NAI	O4B-C4B-C5B-O5B
4	B	481	NAI	O4B-C4B-C5B-O5B
3	B	480	FAD	P-O3P-PA-O2A
3	C	480	FAD	P-O3P-PA-O2A
3	A	480	FAD	P-O3P-PA-O2A
3	D	480	FAD	P-O3P-PA-O2A
4	H	481	NAI	O4B-C4B-C5B-O5B

There are no ring outliers.

20 monomers are involved in 55 short contacts:

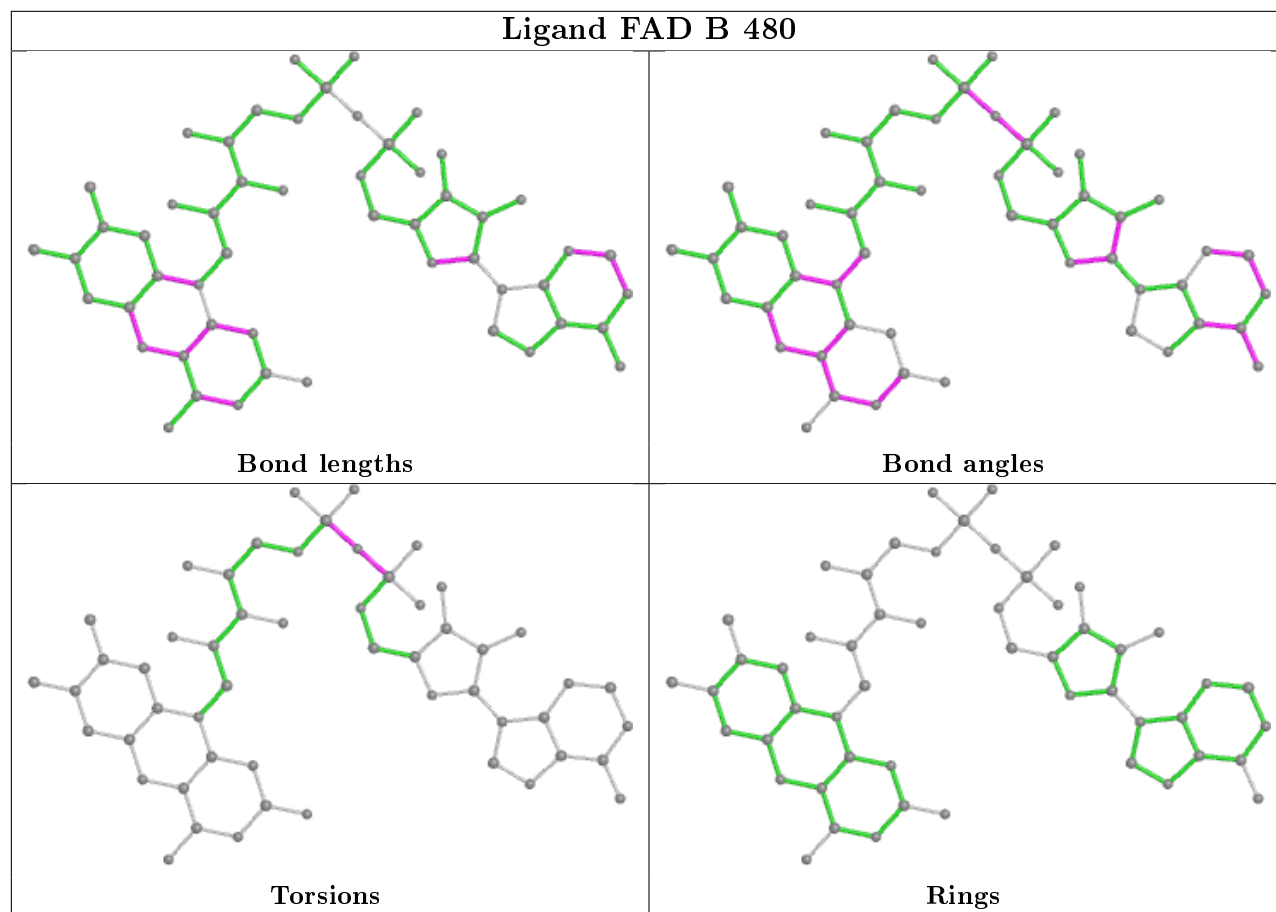
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	480	FAD	5	0
4	F	481	NAI	1	0

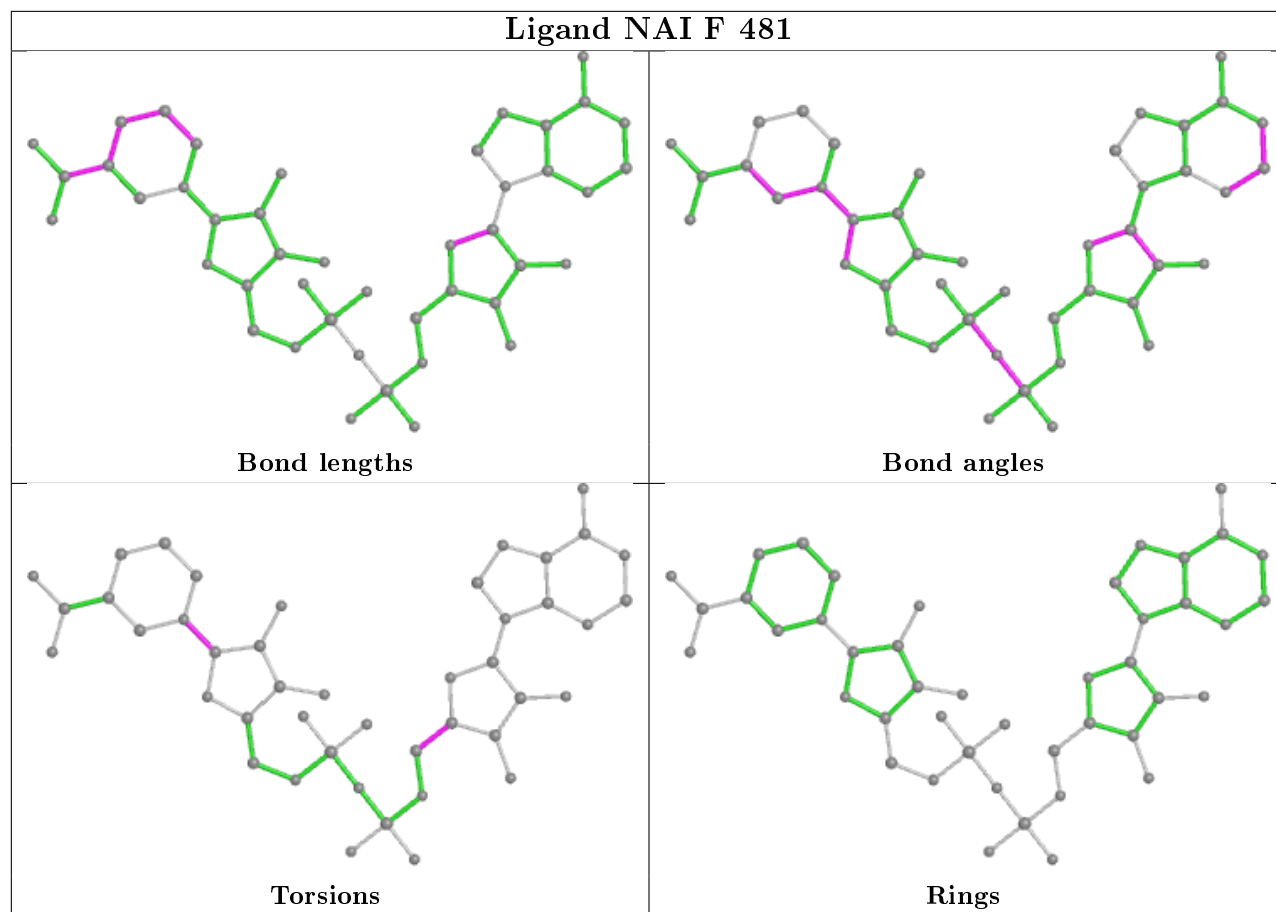
Continued on next page...

Continued from previous page...

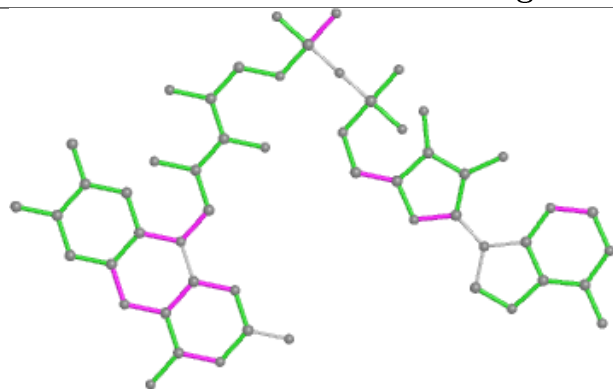
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	480	FAD	4	0
3	G	480	FAD	5	0
4	A	481	NAI	1	0
4	E	481	NAI	4	0
3	C	480	FAD	3	0
4	G	481	NAI	3	0
4	C	481	NAI	5	0
2	F	2021	SO4	1	0
2	G	2025	SO4	1	0
4	H	481	NAI	3	0
2	C	2010	SO4	1	0
2	B	2005	SO4	1	0
3	H	480	FAD	2	0
4	D	481	NAI	2	0
4	B	481	NAI	4	0
3	E	480	FAD	3	0
3	A	480	FAD	4	0
3	D	480	FAD	4	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.

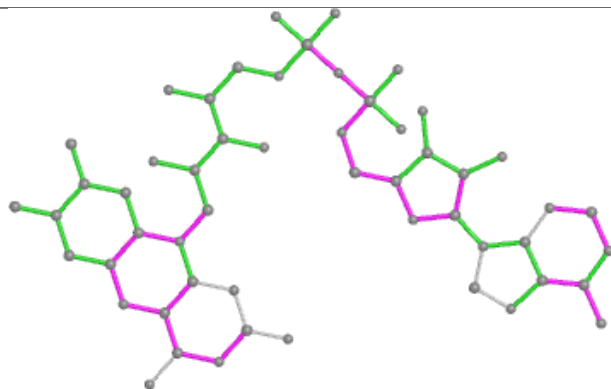




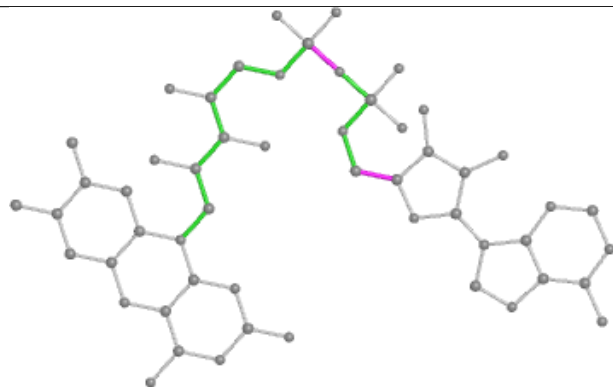
Ligand FAD F 480



Bond lengths



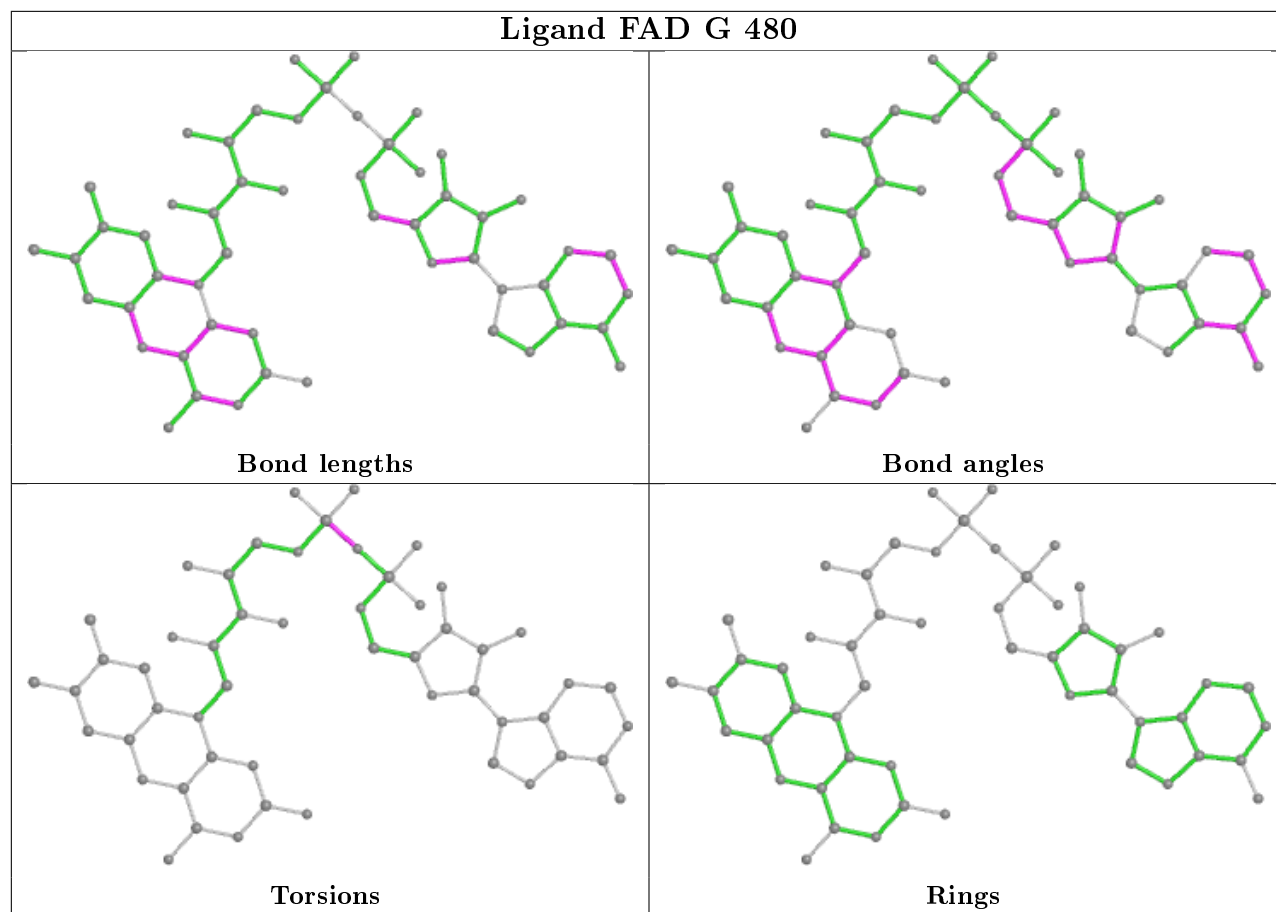
Bond angles

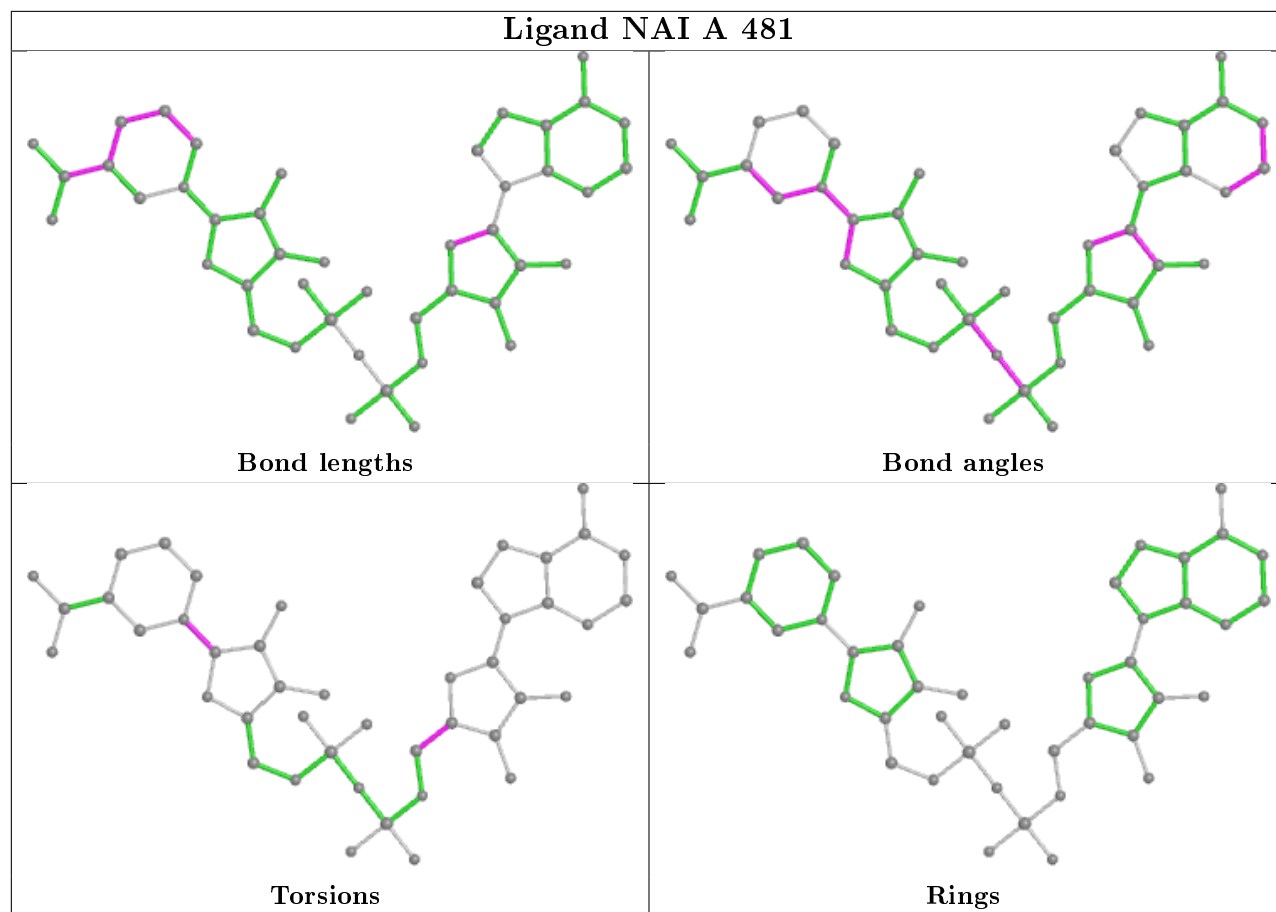


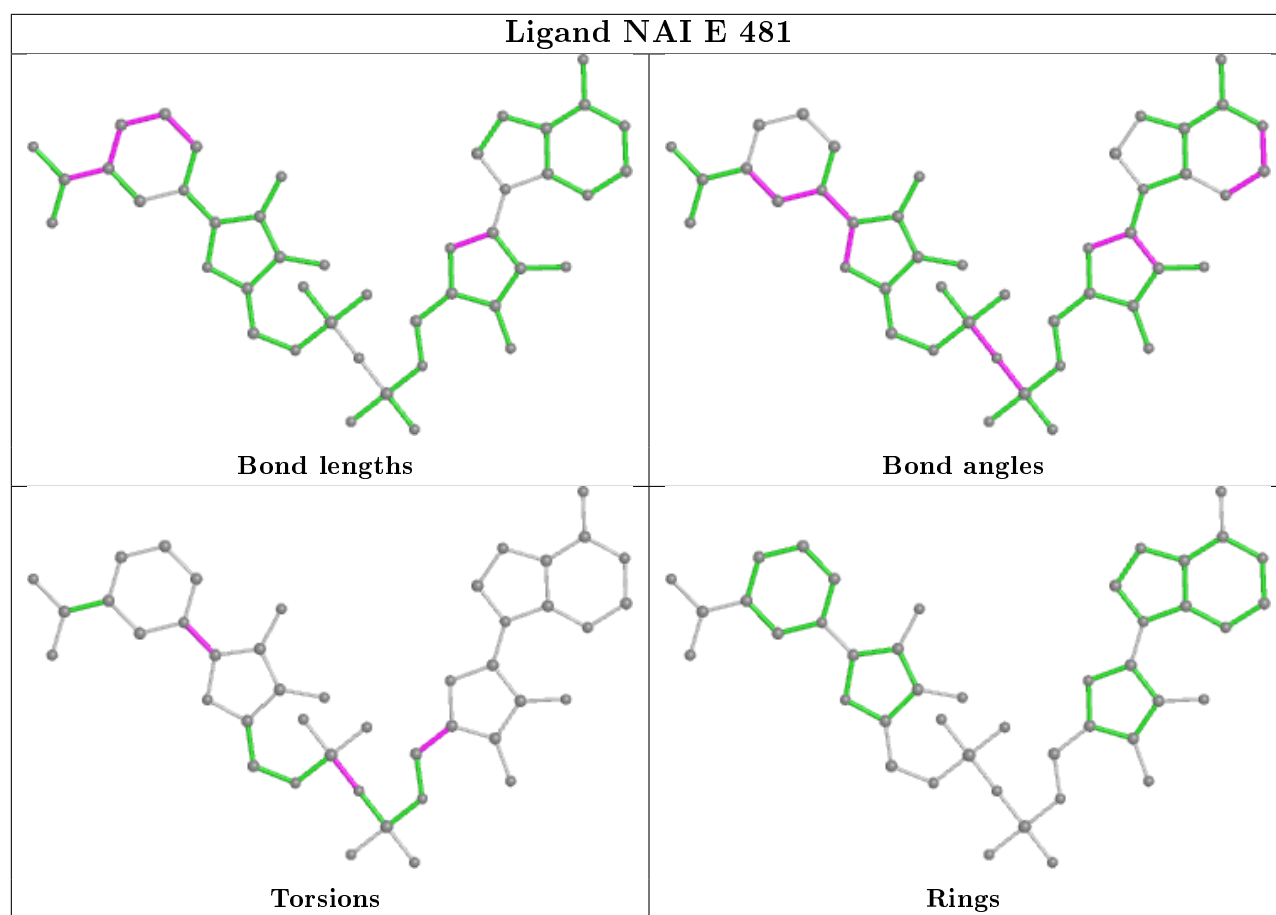
Torsions

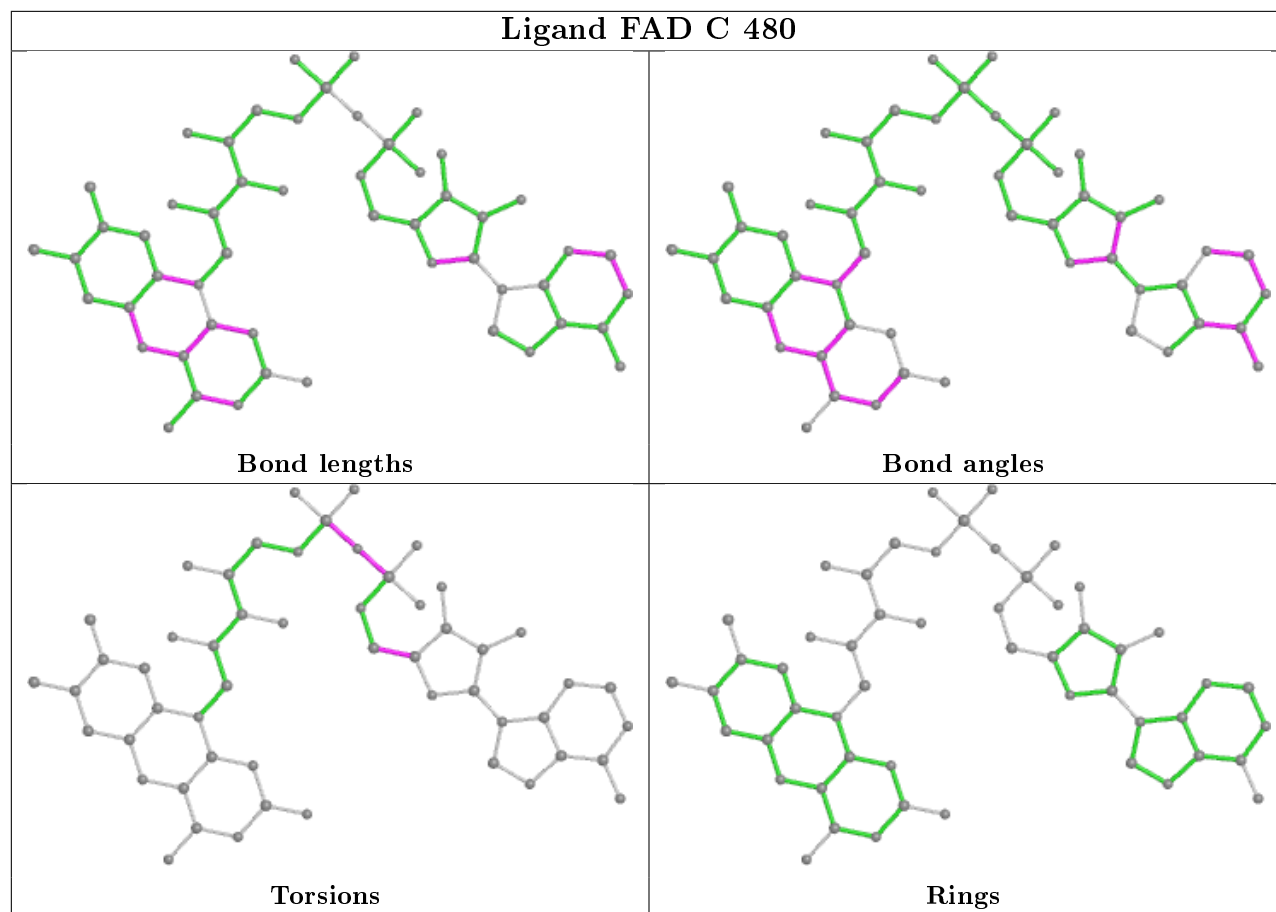


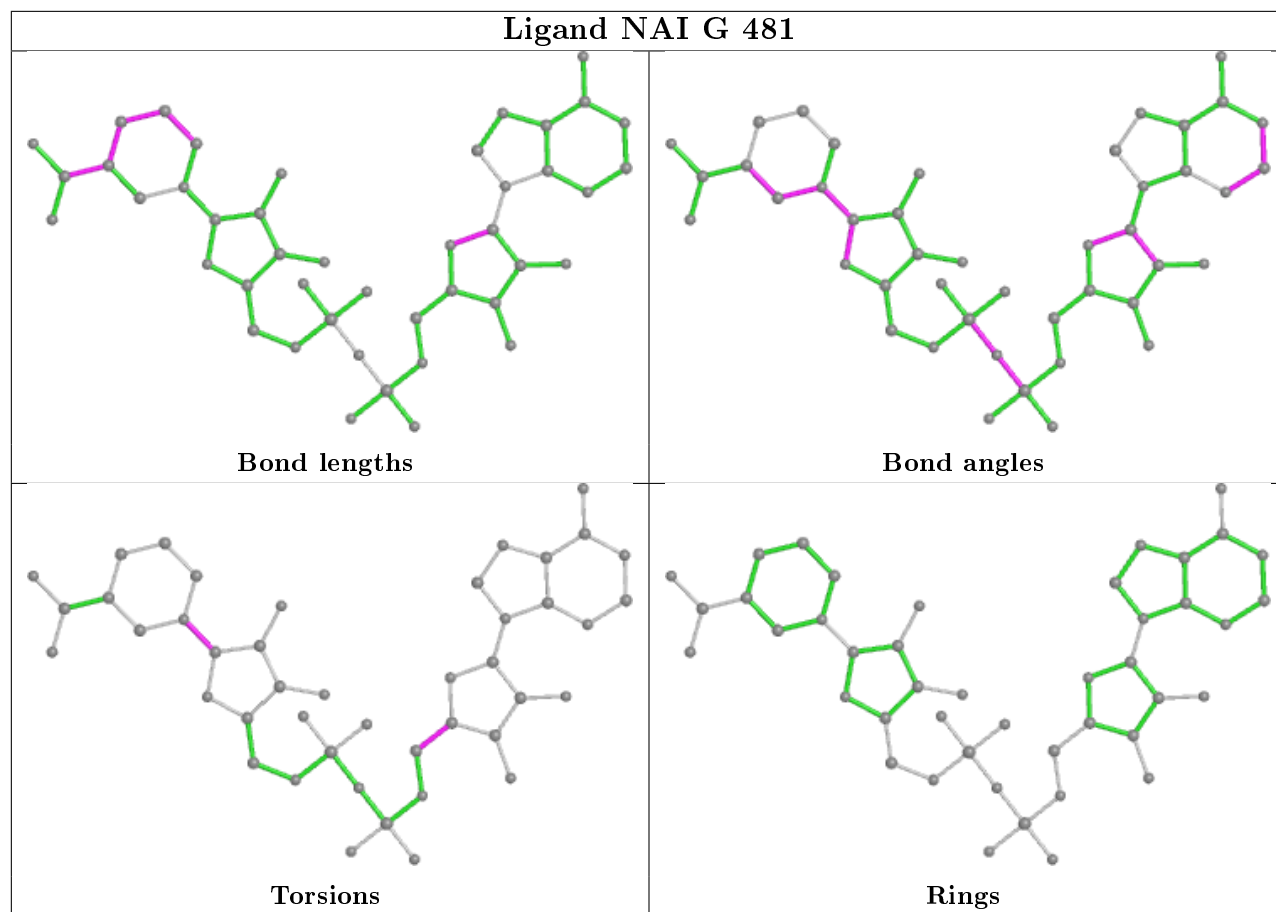
Rings

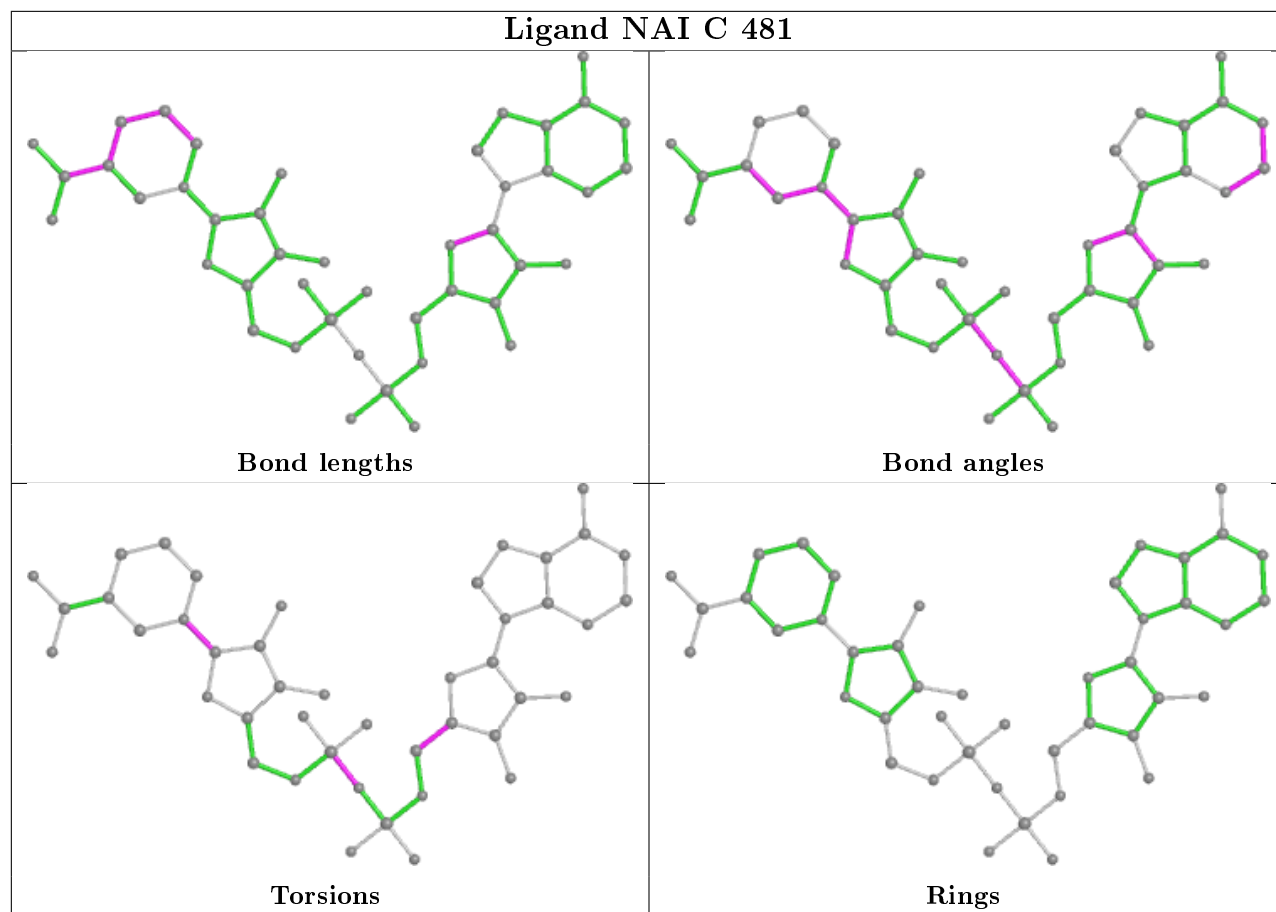


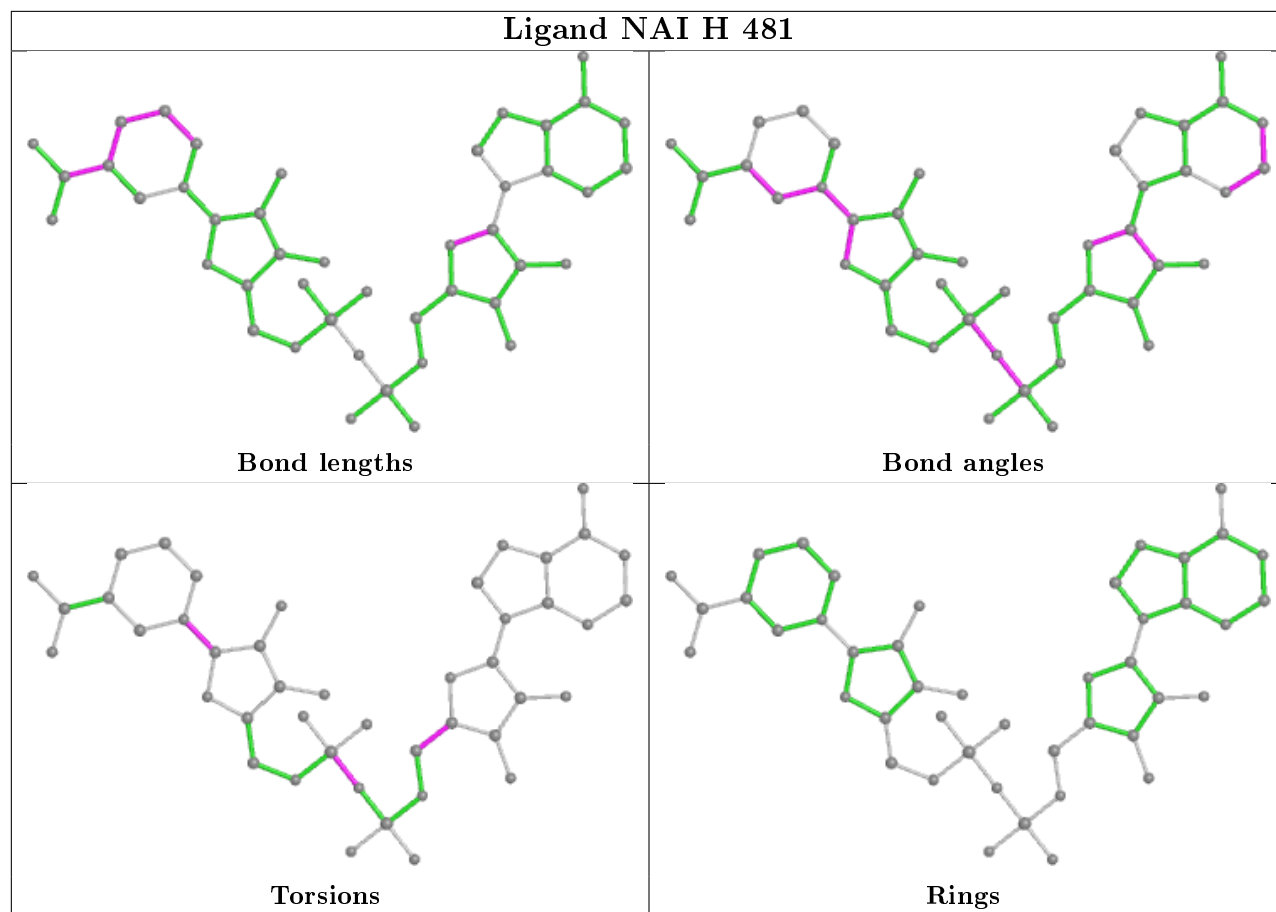


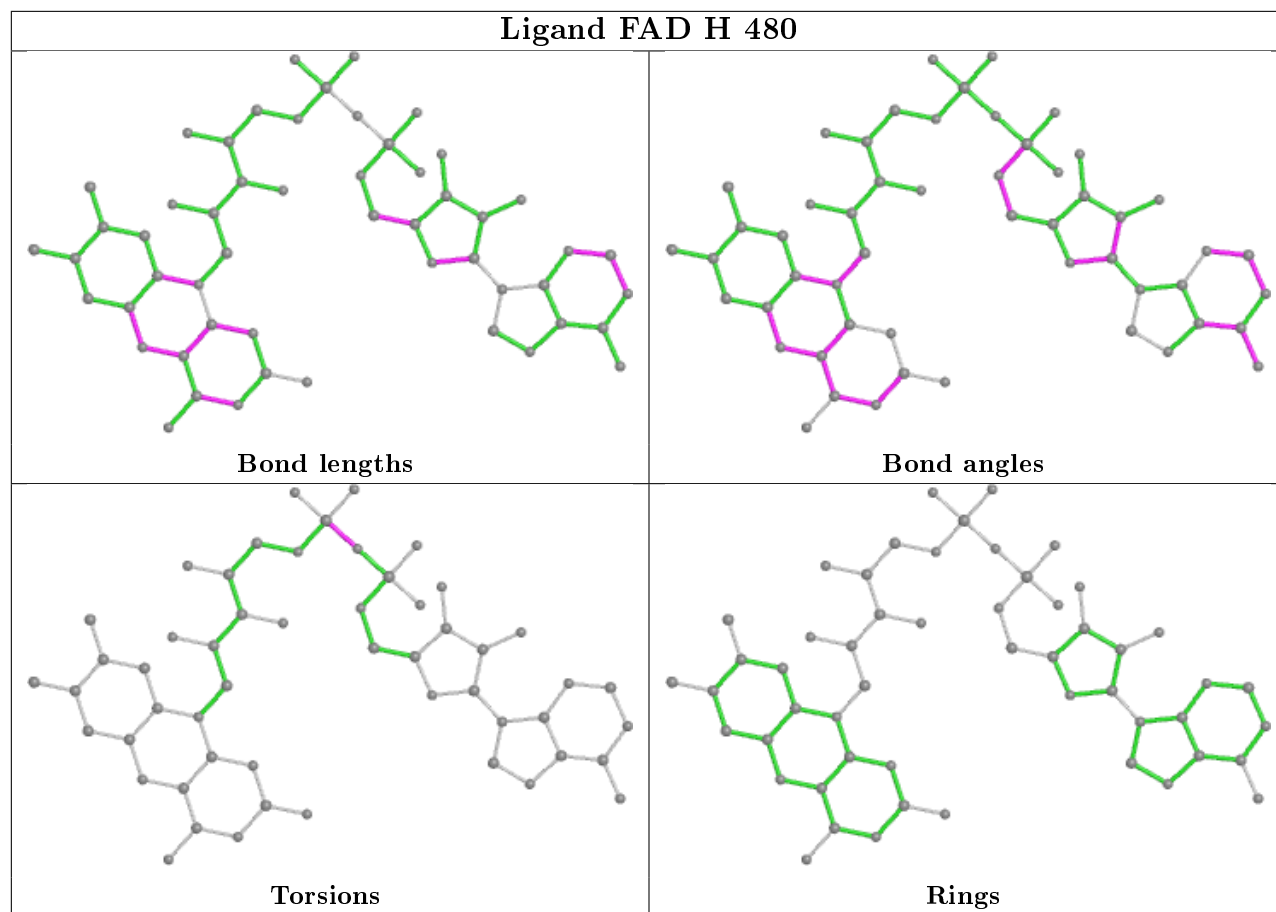


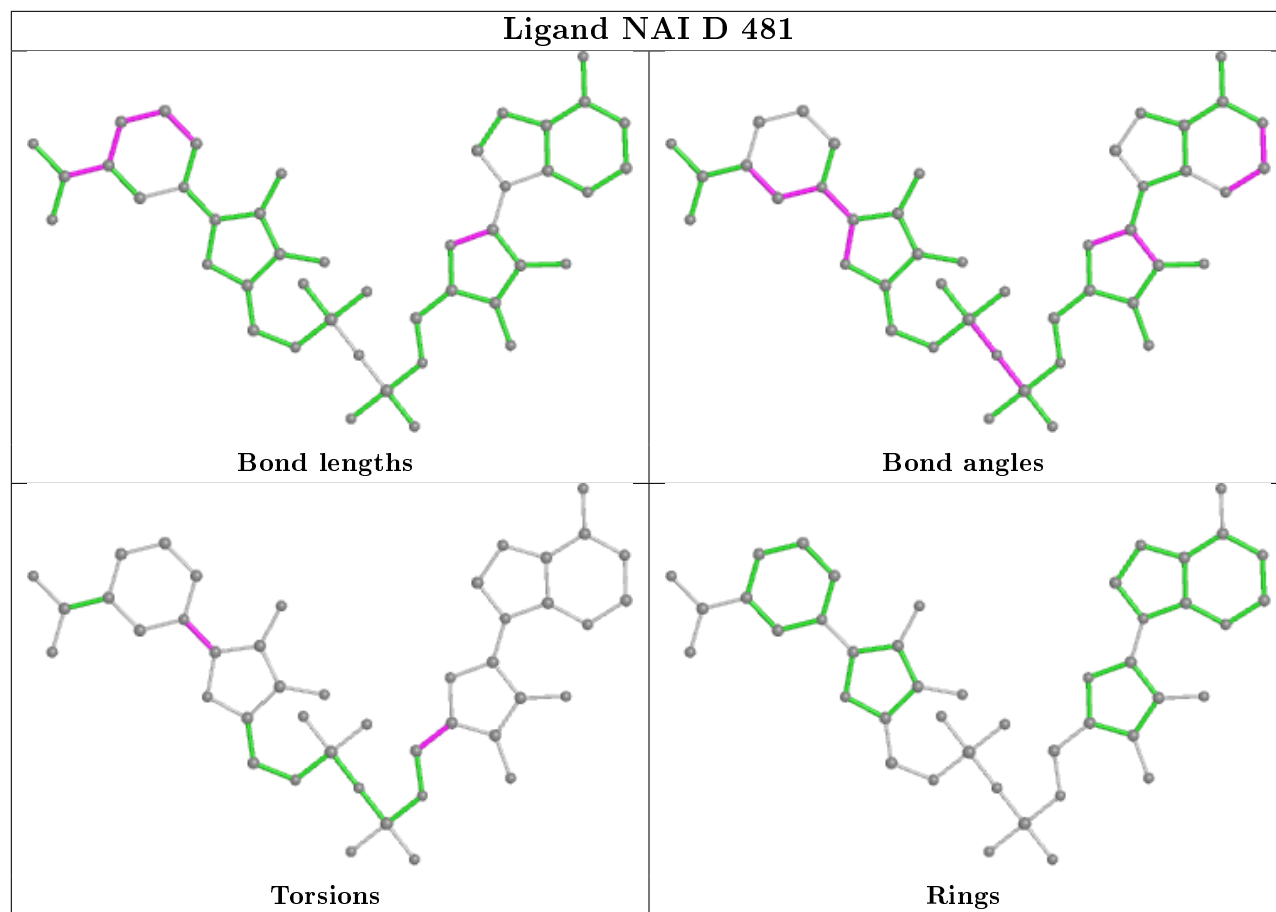


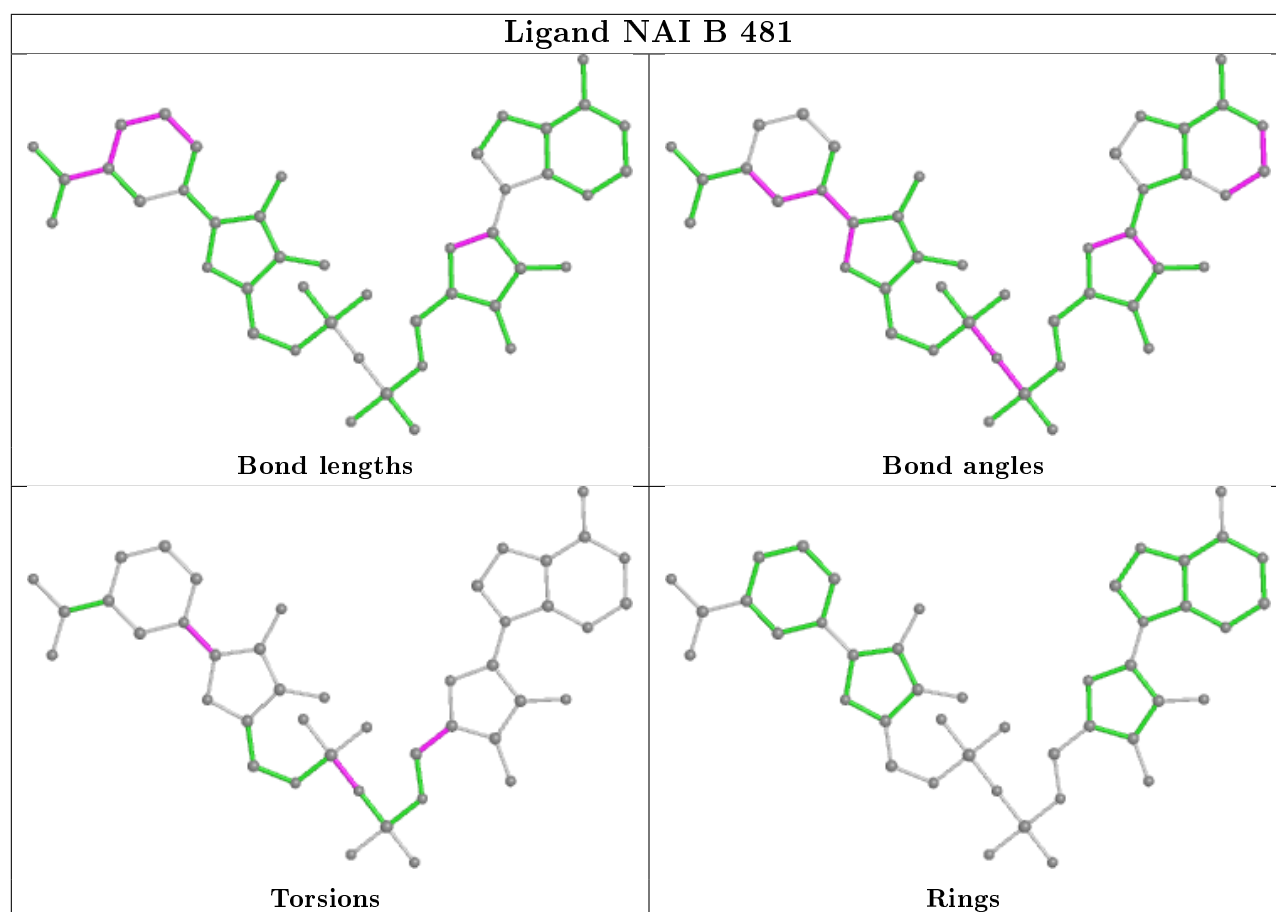


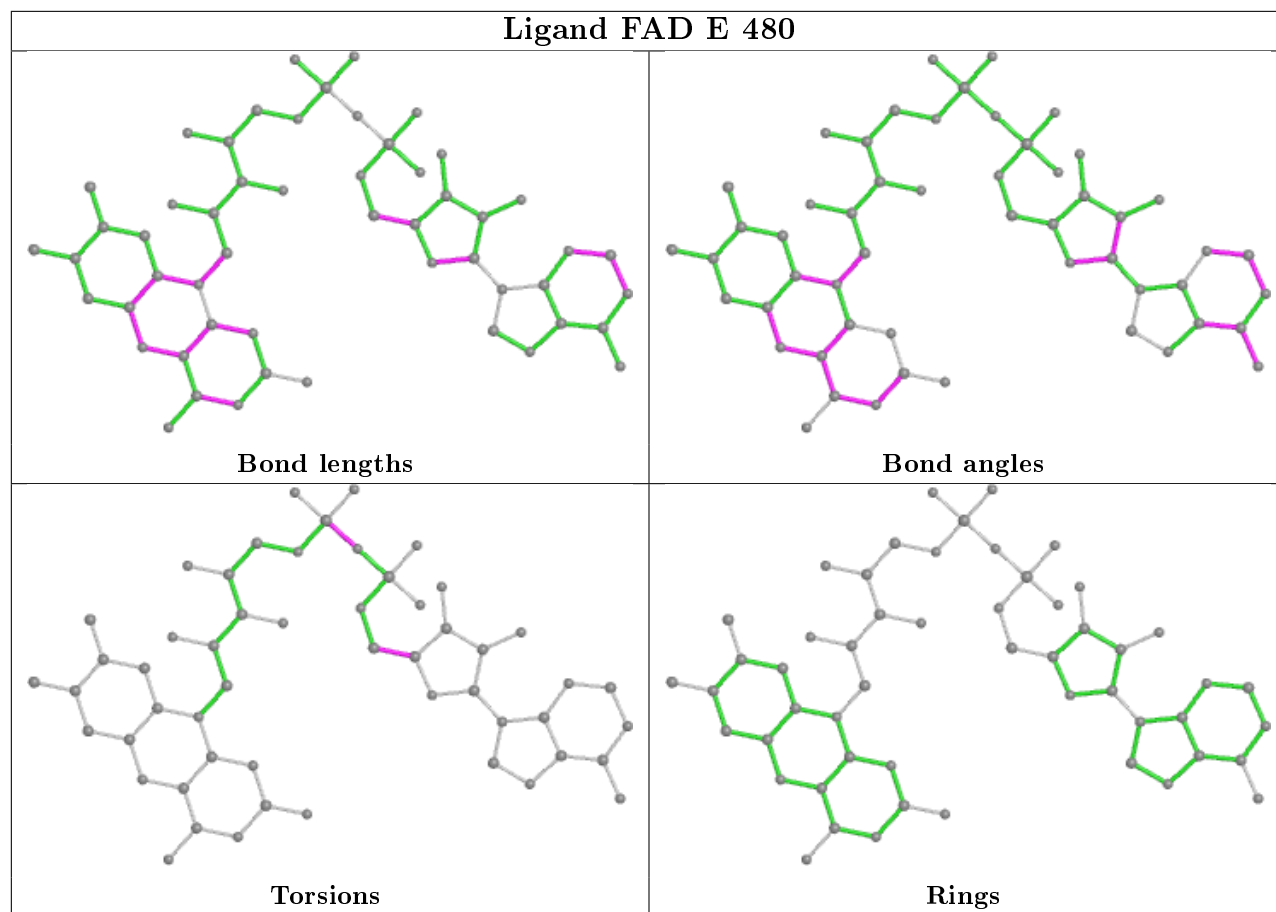


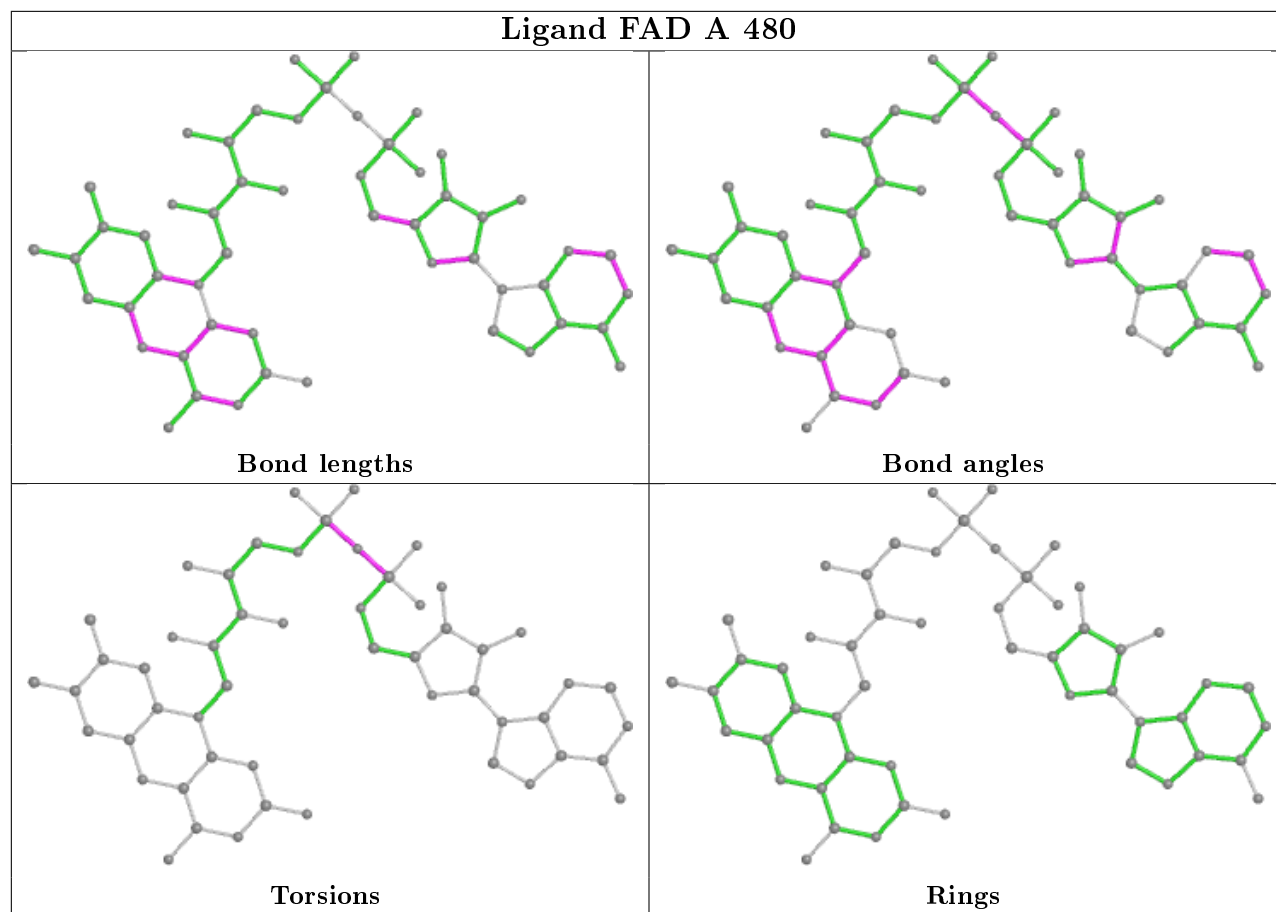


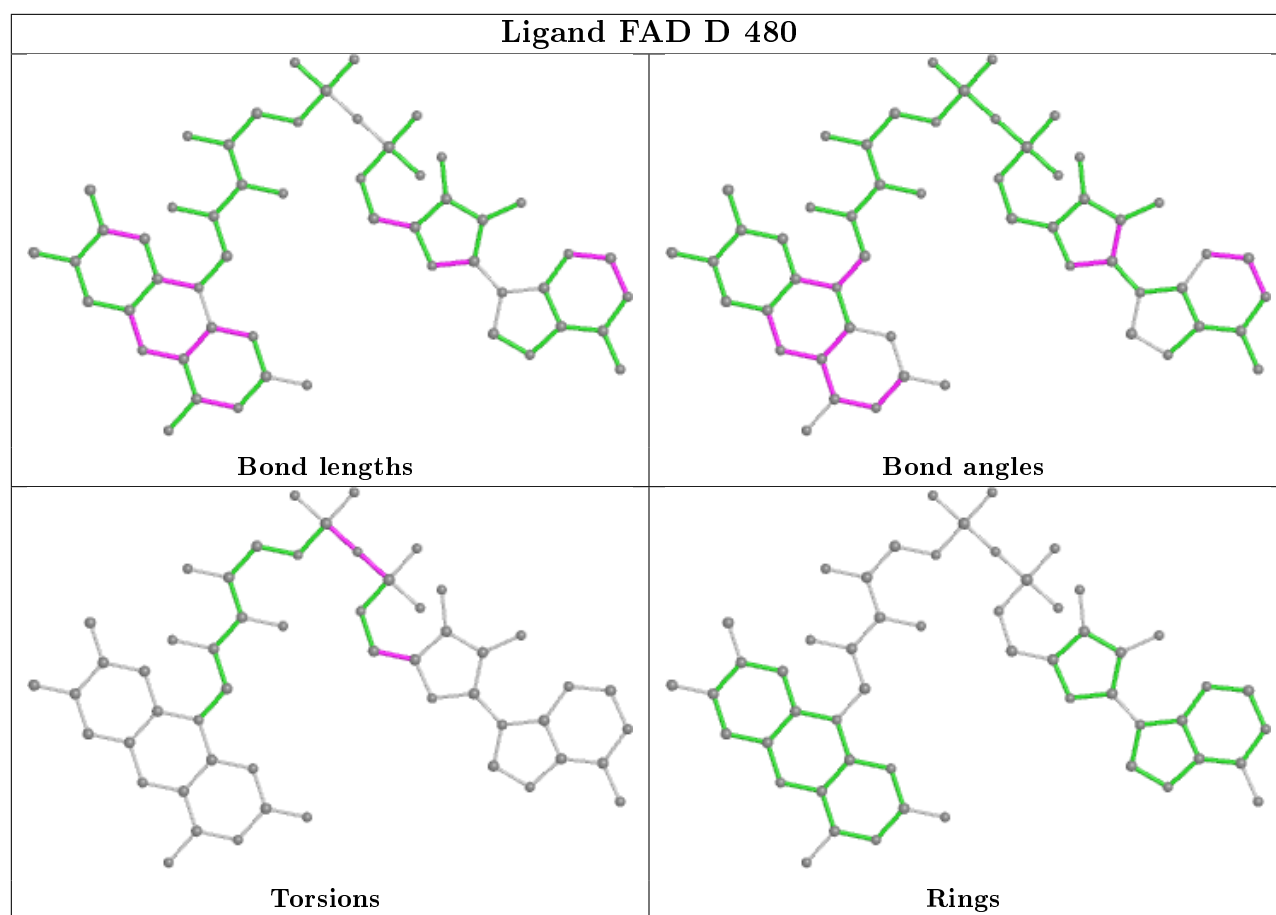












5.7 Other polymers [i](#)

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	472/474 (99%)	0.04	13 (2%)	53	58	18, 33, 53, 69	0
1	B	471/474 (99%)	0.27	29 (6%)	20	24	17, 36, 64, 77	0
1	C	471/474 (99%)	0.23	25 (5%)	26	31	15, 35, 63, 76	0
1	D	472/474 (99%)	-0.00	17 (3%)	42	48	17, 31, 51, 62	0
1	E	473/474 (99%)	0.16	24 (5%)	28	32	15, 31, 56, 85	0
1	F	471/474 (99%)	-0.14	6 (1%)	77	79	14, 25, 43, 62	0
1	G	472/474 (99%)	-0.14	6 (1%)	77	79	15, 25, 41, 59	0
1	H	471/474 (99%)	-0.08	9 (1%)	66	70	13, 29, 47, 67	0
All	All	3773/3792 (99%)	0.04	129 (3%)	45	50	13, 30, 55, 85	0

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	2	ASP	6.6
1	C	132	ALA	5.8
1	E	79	SER	5.1
1	H	79	SER	4.8
1	B	5	ILE	4.6
1	F	250	SER	4.5
1	G	250	SER	4.3
1	C	135	GLY	4.3
1	B	135	GLY	4.3
1	B	133	ASP	4.2
1	B	129	ALA	4.1
1	H	133	ASP	4.0
1	B	244	THR	4.0
1	B	50[A]	CYS	3.9
1	C	79	SER	3.9
1	D	79	SER	3.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	5	ILE	3.7
1	C	31	LYS	3.6
1	C	295	GLU	3.6
1	D	252	GLY	3.6
1	H	5	ILE	3.5
1	H	50[A]	CYS	3.4
1	E	132	ALA	3.4
1	C	133	ASP	3.3
1	B	29	GLY	3.3
1	A	79	SER	3.3
1	B	311	LYS	3.3
1	B	298	PRO	3.3
1	B	132	ALA	3.3
1	E	3	GLN	3.2
1	B	291	GLU	3.2
1	C	312	ILE	3.2
1	C	310	THR	3.2
1	C	263	GLY	3.1
1	H	132	ALA	3.1
1	D	228	ARG	3.1
1	D	265	LYS	3.1
1	B	263	GLY	3.1
1	D	3	GLN	3.1
1	C	111	LYS	3.1
1	E	133	ASP	3.1
1	D	50[A]	CYS	3.0
1	C	291	GLU	3.0
1	F	80	GLU	3.0
1	F	252	GLY	2.9
1	C	50[A]	CYS	2.9
1	F	105	HIS	2.9
1	E	135	GLY	2.9
1	B	292	LEU	2.9
1	A	261	ALA	2.8
1	E	291	GLU	2.8
1	E	134	GLY	2.7
1	D	82	ARG	2.7
1	C	315	ILE	2.7
1	B	136	THR	2.7
1	B	266	ALA	2.7
1	B	33	VAL	2.7
1	F	50[A]	CYS	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	G	80	GLU	2.6
1	B	261	ALA	2.6
1	B	130	THR	2.6
1	B	287	LEU	2.6
1	E	4	PRO	2.6
1	D	250	SER	2.6
1	B	290	GLU	2.6
1	G	3	GLN	2.6
1	C	82	ARG	2.6
1	C	129	ALA	2.6
1	E	119	GLY	2.6
1	A	248	LYS	2.5
1	B	131	LYS	2.5
1	C	298	PRO	2.5
1	A	244	THR	2.5
1	A	410	LYS	2.5
1	D	266	ALA	2.5
1	E	290	GLU	2.4
1	D	5	ILE	2.4
1	E	121	ILE	2.4
1	B	286	ASN	2.4
1	D	69	THR	2.4
1	A	264	GLY	2.4
1	C	244	THR	2.4
1	E	6	ASP	2.4
1	D	263	GLY	2.3
1	B	289	LEU	2.3
1	B	295	GLU	2.3
1	B	6	ASP	2.3
1	A	290	GLU	2.3
1	B	79	SER	2.3
1	B	293	GLY	2.3
1	C	296	LEU	2.3
1	A	228	ARG	2.3
1	B	82	ARG	2.3
1	D	80	GLU	2.3
1	E	328	ALA	2.3
1	A	265	LYS	2.3
1	A	266	ALA	2.2
1	A	82	ARG	2.2
1	H	248	LYS	2.2
1	B	116	ASN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	78	MET	2.2
1	E	295	GLU	2.2
1	C	108	LYS	2.2
1	E	31	LYS	2.2
1	A	3	GLN	2.2
1	G	79	SER	2.2
1	D	249	LYS	2.2
1	E	111	LYS	2.2
1	C	290	GLU	2.2
1	G	228	ARG	2.2
1	C	311	LYS	2.1
1	D	410	LYS	2.1
1	E	50[A]	CYS	2.1
1	F	410	LYS	2.1
1	H	6	ASP	2.1
1	D	376	GLU	2.1
1	E	29	GLY	2.1
1	C	313	PRO	2.1
1	D	285	LYS	2.1
1	C	112	VAL	2.1
1	E	331	ALA	2.1
1	G	410	LYS	2.1
1	H	135	GLY	2.1
1	C	262	SER	2.1
1	E	136	THR	2.1
1	C	139	ILE	2.0
1	E	112	VAL	2.0
1	H	4	PRO	2.0
1	E	116	ASN	2.0

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 carbohydrates 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
2	SO4	C	2008	5/5	0.75	0.22	105,105,105,106	0
2	SO4	G	2025	5/5	0.83	0.35	104,104,104,105	0
2	SO4	E	2016	5/5	0.85	0.28	106,106,106,106	0
2	SO4	F	2020	5/5	0.85	0.15	56,58,59,60	0
2	SO4	G	2024	5/5	0.85	0.14	55,56,56,57	0
2	SO4	D	2012	5/5	0.86	0.15	69,69,70,70	0
2	SO4	H	2028	5/5	0.86	0.21	95,95,96,96	0
2	SO4	B	2004	5/5	0.87	0.15	103,103,103,103	0
2	SO4	B	2005	5/5	0.87	0.25	53,55,56,56	0
2	SO4	A	2002	5/5	0.88	0.15	81,82,82,83	0
2	SO4	D	2013	5/5	0.88	0.31	102,102,103,103	0
2	SO4	D	2014	5/5	0.89	0.18	69,70,71,71	0
2	SO4	A	2001	5/5	0.89	0.15	68,68,69,69	0
4	NAI	A	481	44/44	0.89	0.14	32,46,55,59	0
2	SO4	G	2026	5/5	0.90	0.10	79,79,80,80	0
4	NAI	D	481	44/44	0.91	0.15	39,46,54,57	0
2	SO4	H	2030	5/5	0.91	0.13	67,68,68,69	0
2	SO4	C	2009	5/5	0.91	0.16	49,50,53,53	0
4	NAI	B	481	44/44	0.91	0.12	31,37,42,43	0
3	FAD	C	480	53/53	0.92	0.14	27,32,44,45	0
4	NAI	G	481	44/44	0.92	0.14	25,36,48,50	0
4	NAI	C	481	44/44	0.92	0.13	31,37,42,45	0
4	NAI	E	481	44/44	0.92	0.14	26,37,42,43	0
3	FAD	B	480	53/53	0.92	0.14	27,32,43,44	0
2	SO4	E	2017	5/5	0.93	0.14	60,60,62,63	0
4	NAI	F	481	44/44	0.93	0.13	27,35,38,40	0
3	FAD	E	480	53/53	0.93	0.14	24,29,37,38	0
2	SO4	F	2022	5/5	0.94	0.13	79,79,80,80	0
4	NAI	H	481	44/44	0.94	0.11	26,35,38,38	0
2	SO4	C	2010	5/5	0.94	0.13	70,71,71,71	0
2	SO4	E	2018	5/5	0.94	0.14	69,69,70,71	0
2	SO4	H	2029	5/5	0.94	0.16	53,54,56,56	0
2	SO4	B	2006	5/5	0.94	0.15	77,77,77,78	0
3	FAD	A	480	53/53	0.94	0.13	18,25,28,31	0
2	SO4	F	2021	5/5	0.95	0.25	84,84,85,85	0
3	FAD	H	480	53/53	0.95	0.14	17,23,32,33	0
3	FAD	G	480	53/53	0.95	0.14	14,19,22,25	0

Continued on next page...

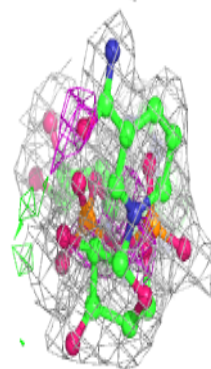
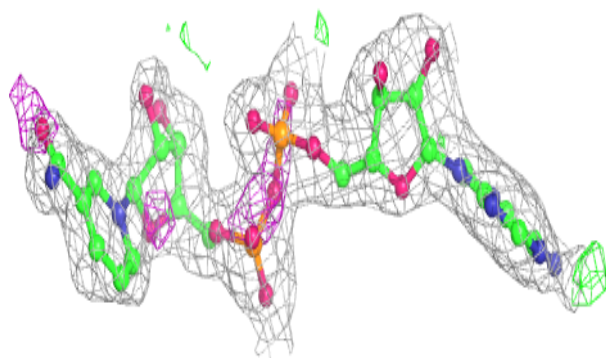
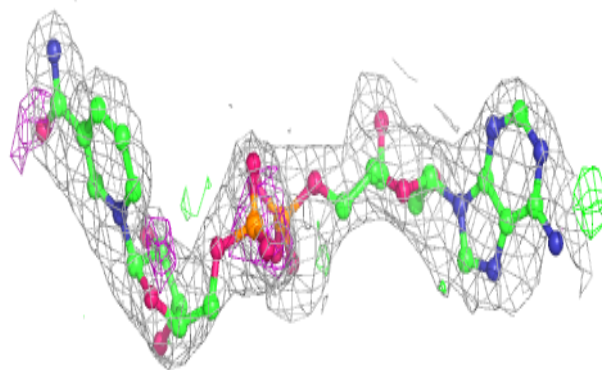
Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	H	2027	5/5	0.96	0.19	56,57,59,59	0
3	FAD	F	480	53/53	0.96	0.14	13,17,22,26	0
2	SO4	E	2015	5/5	0.96	0.11	54,54,55,57	0
3	FAD	D	480	53/53	0.96	0.12	18,22,26,27	0
2	SO4	C	2007	5/5	0.97	0.17	69,70,71,71	0
2	SO4	F	2019	5/5	0.97	0.16	61,63,63,63	0
2	SO4	G	2023	5/5	0.97	0.12	69,69,70,70	0
2	SO4	B	2003	5/5	0.98	0.16	65,66,66,67	0
2	SO4	A	2000	5/5	0.98	0.14	57,58,60,60	0
2	SO4	D	2011	5/5	0.98	0.17	57,59,60,61	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.

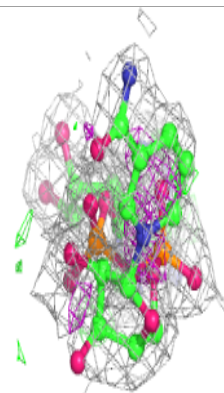
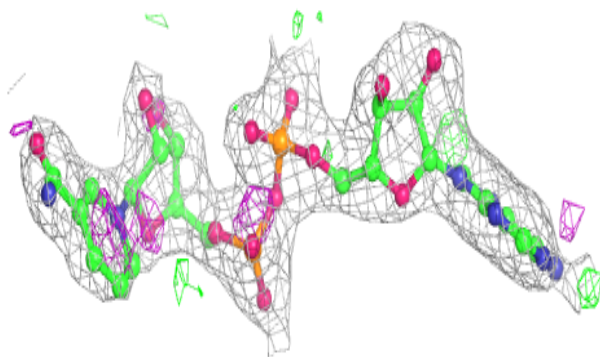
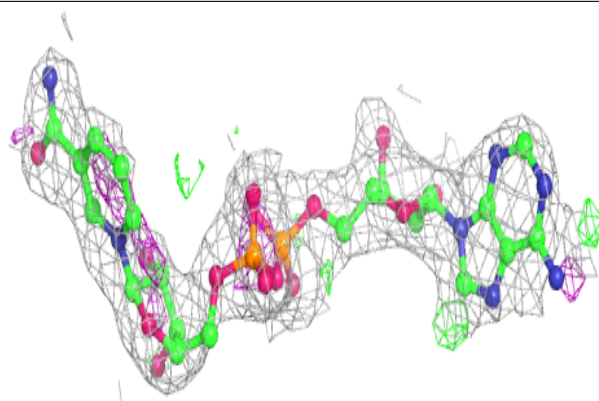
Electron density around NAI A 481:

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

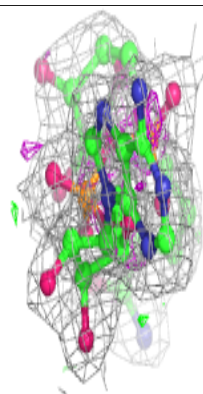
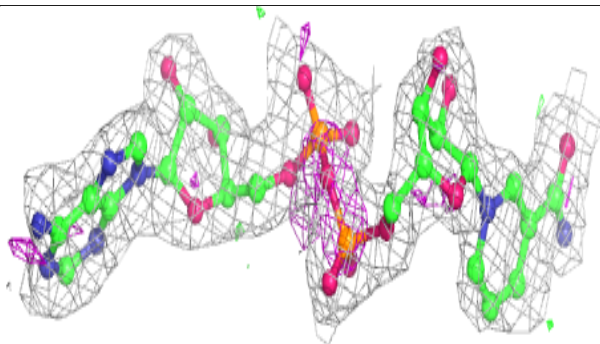
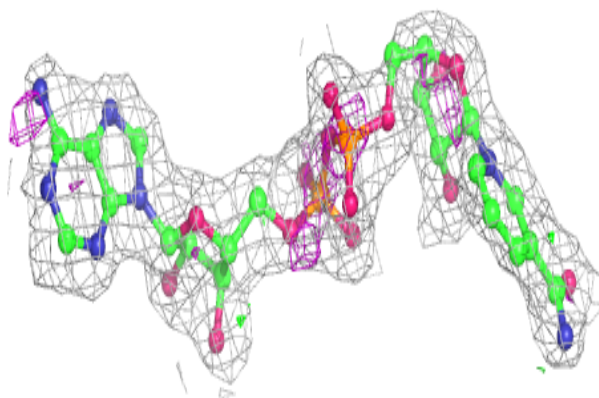


Electron density around NAI D 481:

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

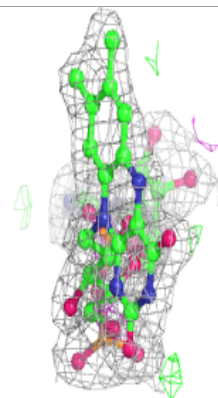
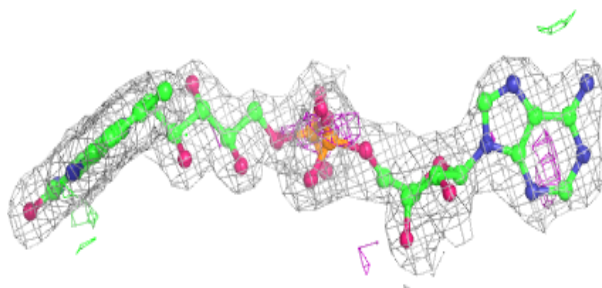
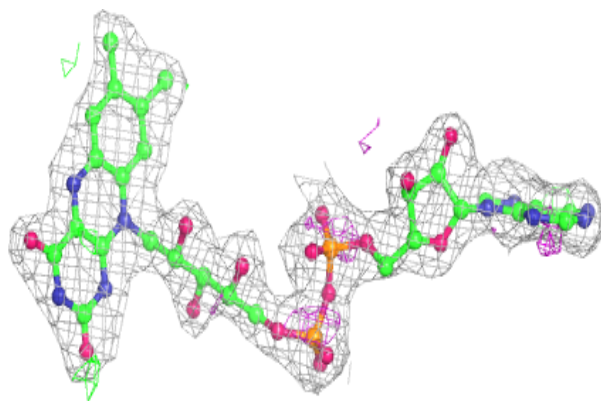
**Electron density around NAI B 481:**

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

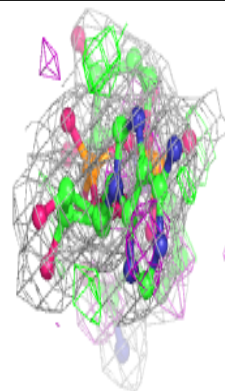
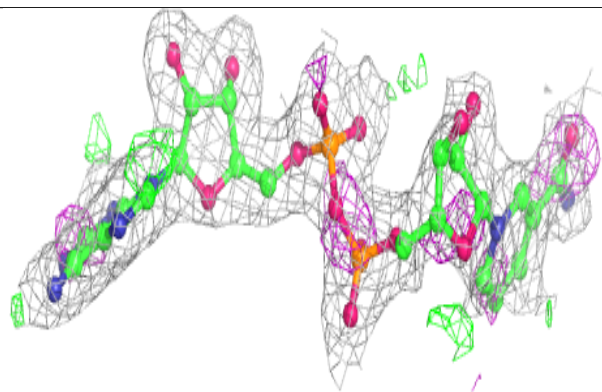
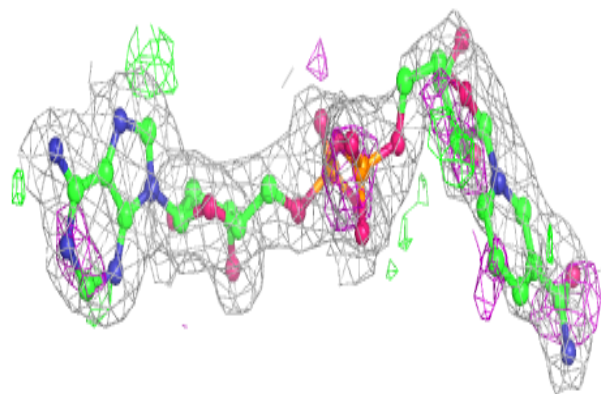


Electron density around FAD C 480:

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

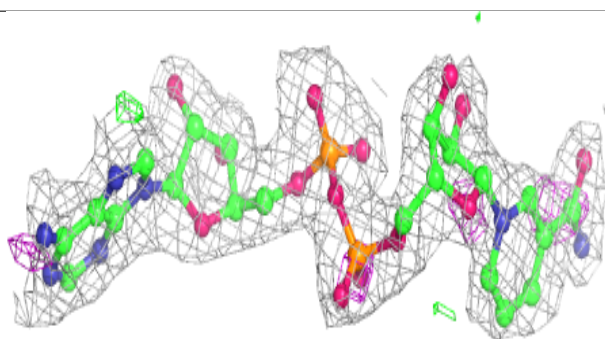
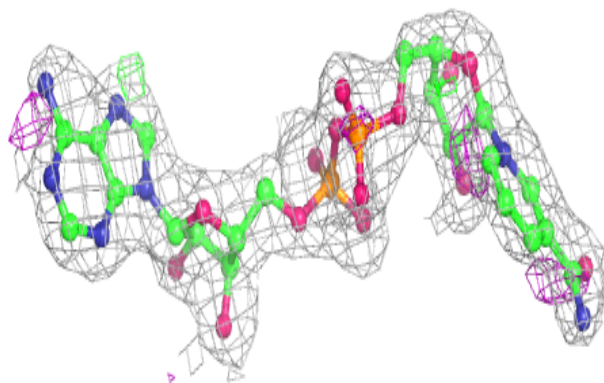
**Electron density around NAI G 481:**

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

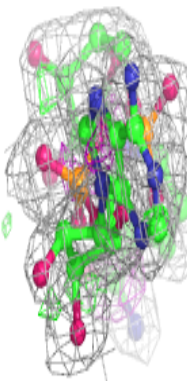
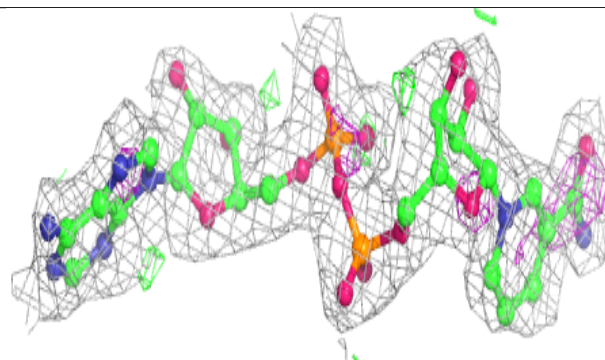
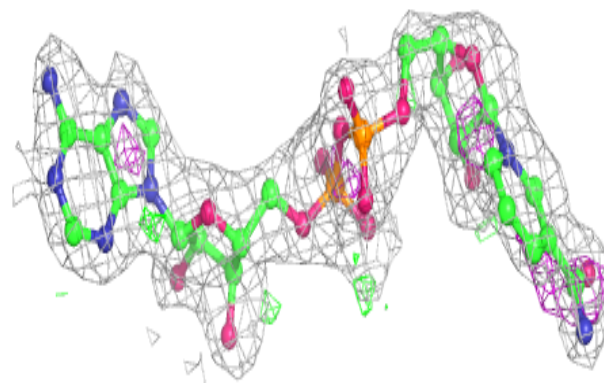


Electron density around NAI C 481:

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

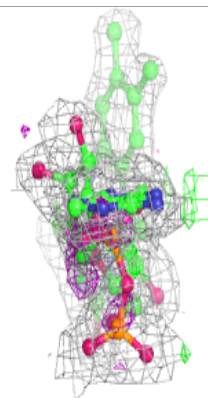
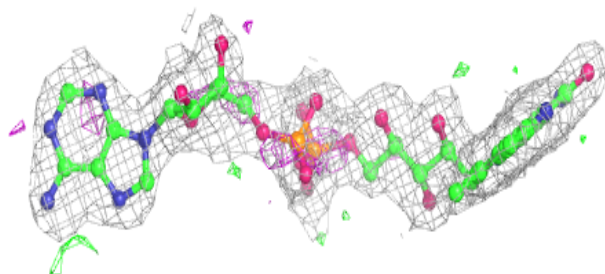
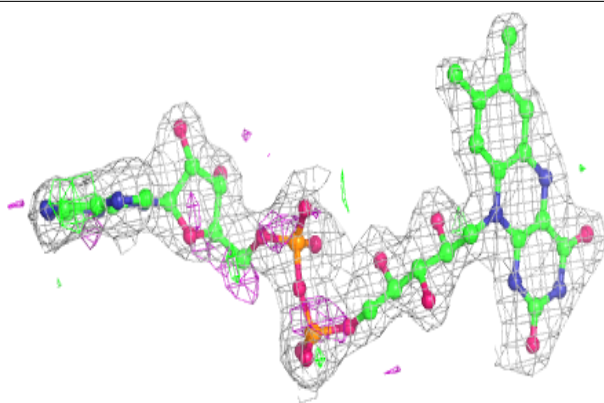
**Electron density around NAI E 481:**

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

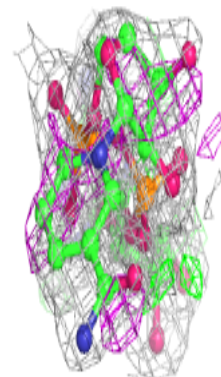
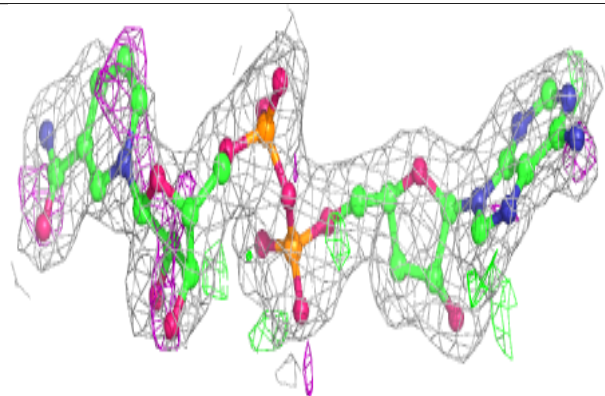
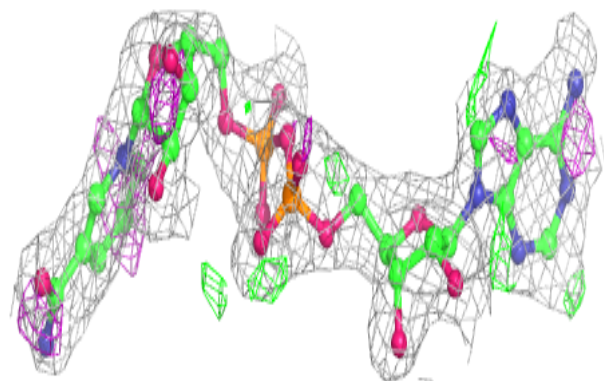


Electron density around FAD B 480:

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

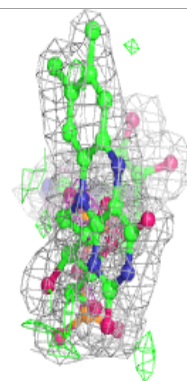
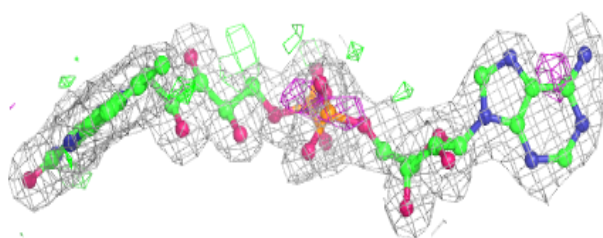
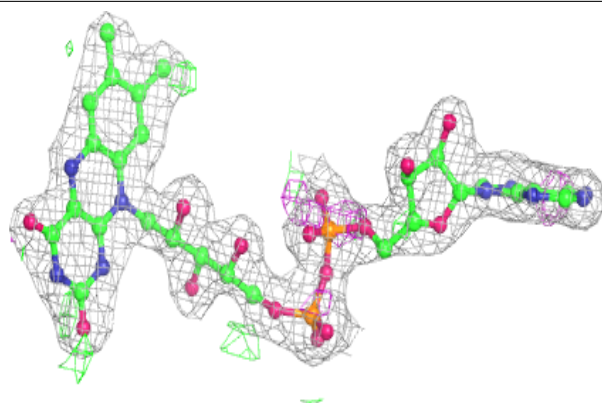
**Electron density around NAI F 481:**

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

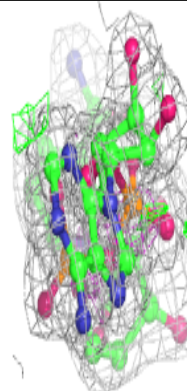
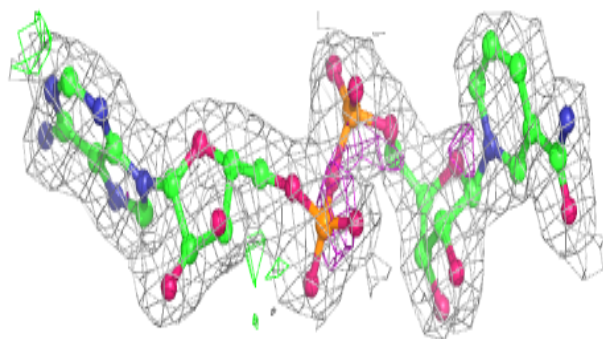
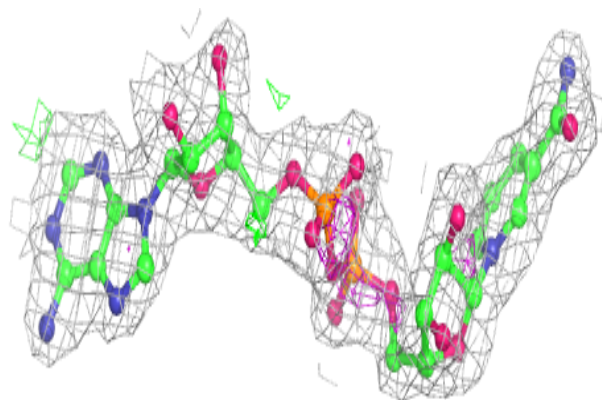


Electron density around FAD E 480:

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

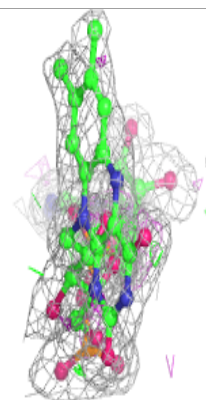
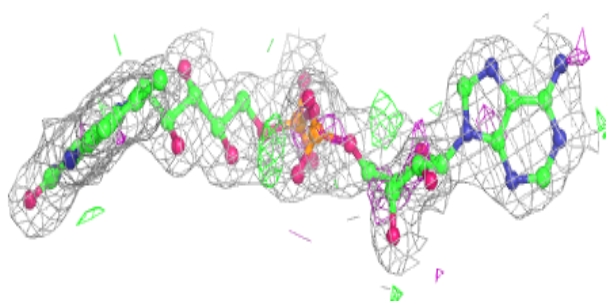
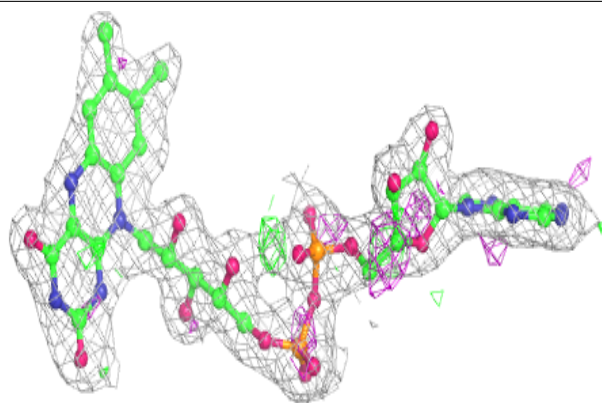
**Electron density around NAI H 481:**

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

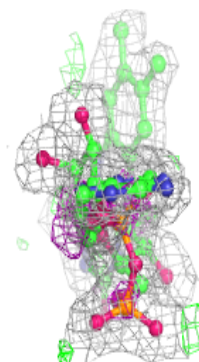
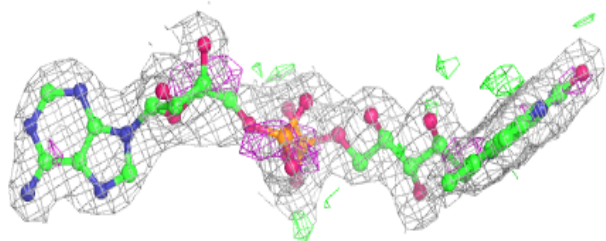
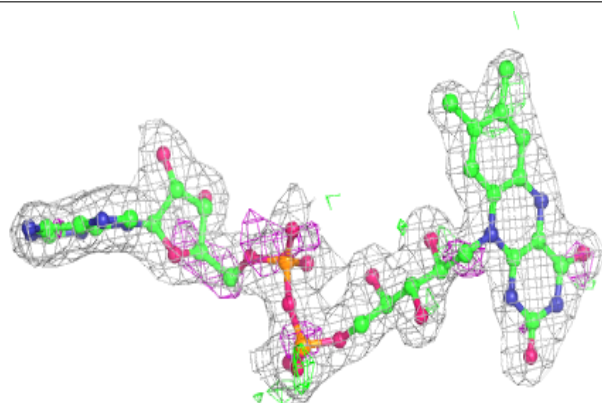


Electron density around FAD A 480:

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

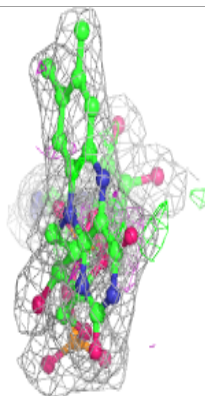
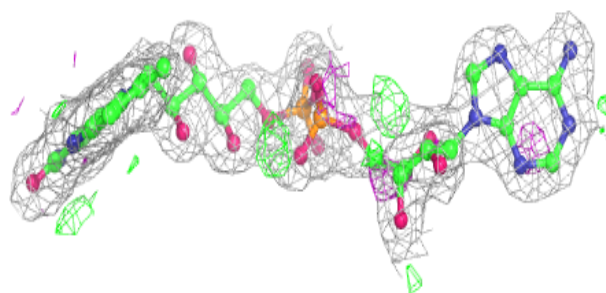
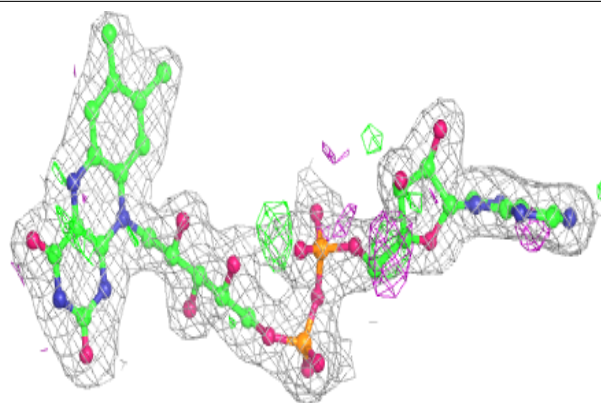
**Electron density around FAD H 480:**

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

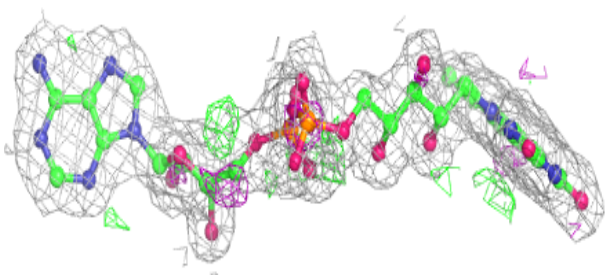
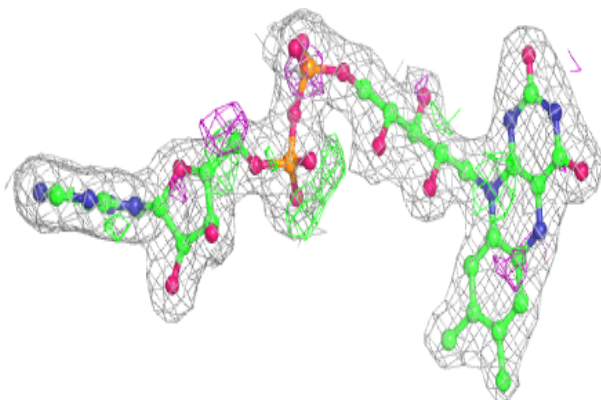


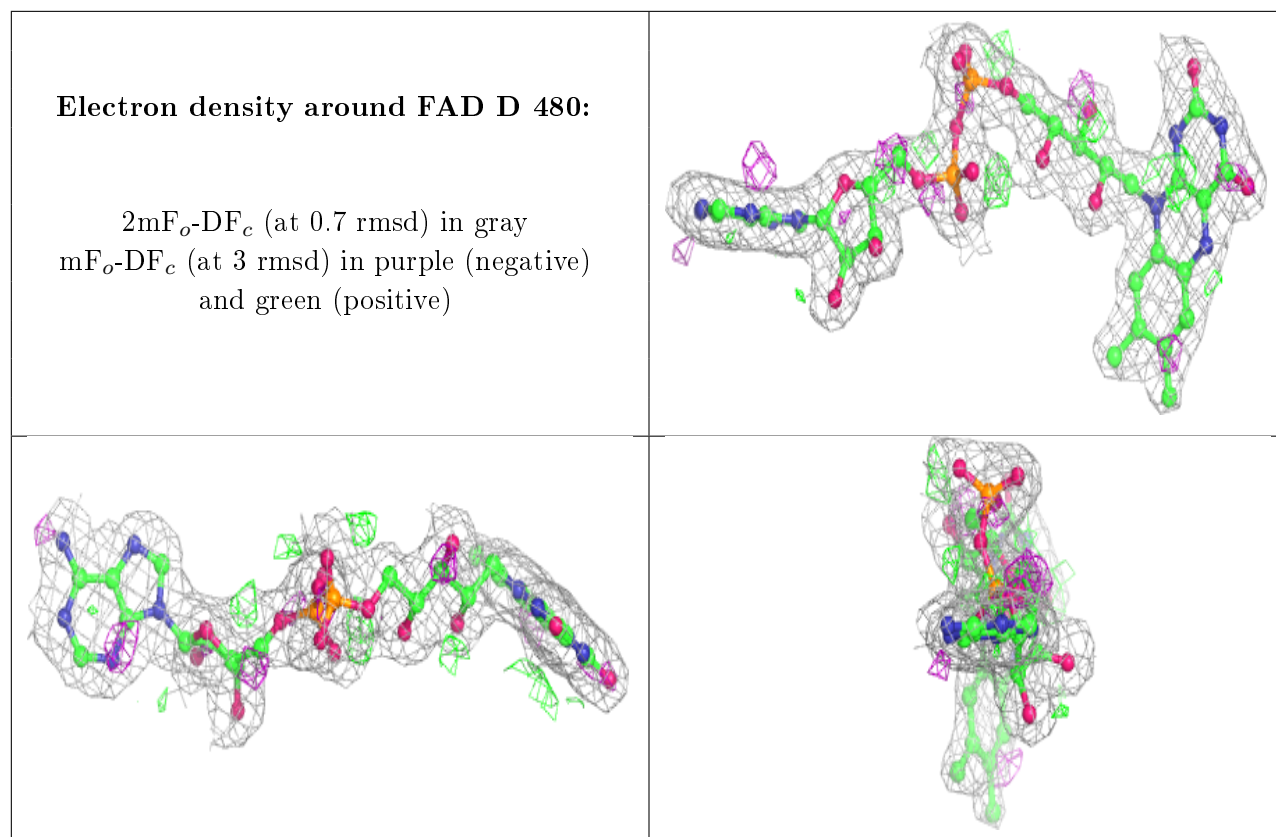
Electron density around FAD G 480:

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

**Electron density around FAD F 480:**

$2mF_o-DF_c$ (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.