



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:11 PM UTC

PDB ID : 3UKL / pdb_00003ukl
Title : Crystal structure of UDP-galactopyranose mutase from *Aspergillus fumigatus* in complex with UDP
Authors : Van Straaten, K.E.; Sanders, D.A.R.
Deposited on : 2011-11-09
Resolution : 2.63 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

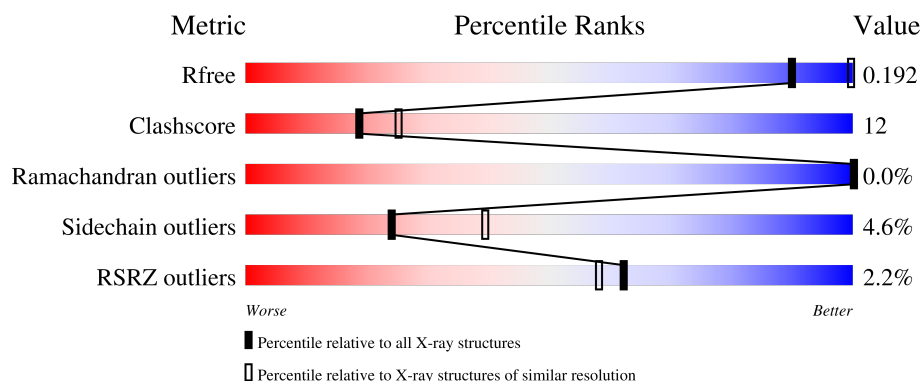
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.63 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.



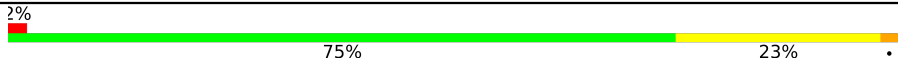
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	510	3% 76% 23% .
1	B	510	2% 78% 20% .
1	C	510	2% 78% 21% .
1	D	510	3% 75% 23% .
1	E	510	2% 78% 20% .

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Mol	Chain	Length	Quality of chain
1	F	510	
1	G	510	
1	H	510	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 33747 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.

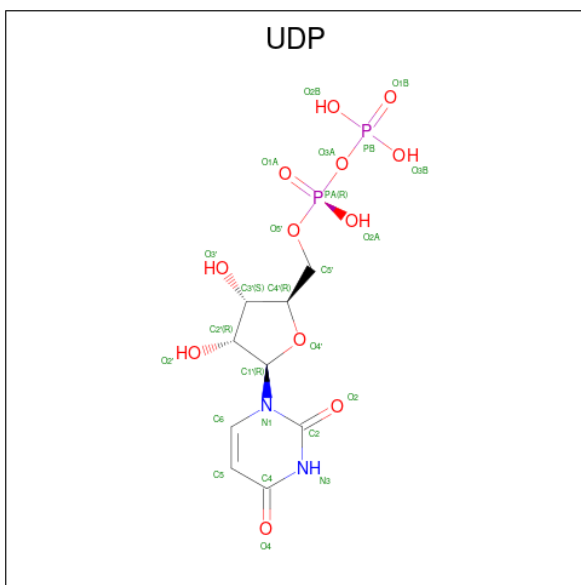
- Molecule 1 is a protein called UDP-galactopyranose mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	510	Total	C	N	O	S	0	0	0
			4004	2540	687	756	21			
1	B	510	Total	C	N	O	S	0	1	0
			4015	2549	688	757	21			
1	C	510	Total	C	N	O	S	0	1	0
			4015	2546	691	757	21			
1	D	510	Total	C	N	O	S	0	0	0
			4004	2540	687	756	21			
1	E	510	Total	C	N	O	S	0	1	0
			4015	2549	688	757	21			
1	F	510	Total	C	N	O	S	0	0	0
			4004	2540	687	756	21			
1	G	510	Total	C	N	O	S	0	0	0
			4004	2540	687	756	21			
1	H	510	Total	C	N	O	S	0	1	0
			4015	2546	691	757	21			

There are 8 discrepancies between the modelled and reference sequences:

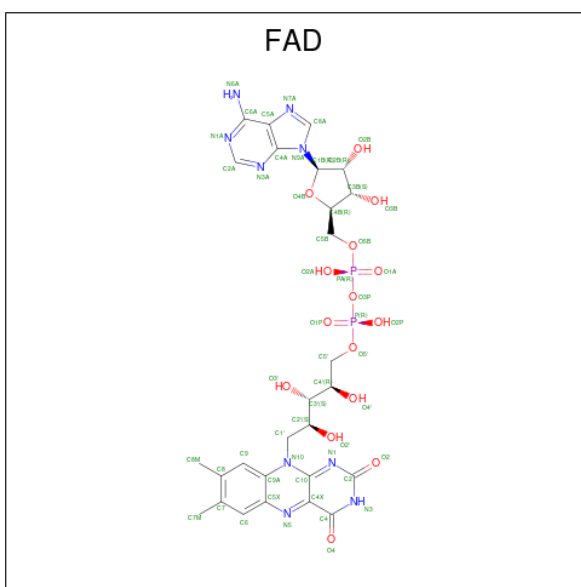
Chain	Residue	Modelled	Actual	Comment	Reference
A	511	LEU	-	expression tag	UNP Q4W1X2
B	511	LEU	-	expression tag	UNP Q4W1X2
C	511	LEU	-	expression tag	UNP Q4W1X2
D	511	LEU	-	expression tag	UNP Q4W1X2
E	511	LEU	-	expression tag	UNP Q4W1X2
F	511	LEU	-	expression tag	UNP Q4W1X2
G	511	LEU	-	expression tag	UNP Q4W1X2
H	511	LEU	-	expression tag	UNP Q4W1X2

- Molecule 2 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	B	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	C	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	D	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	E	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	F	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	G	1	Total 25	C 9	N 2	O 12	P 2	0	0
2	H	1	Total 25	C 9	N 2	O 12	P 2	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: 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 CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Cl 2 2	0	0
4	B	1	Total Cl 1 1	0	0
4	C	1	Total Cl 1 1	0	0
4	E	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	2	Total	Cl	0	0
			2	2		
4	H	1	Total	Cl	0	0
			1	1		

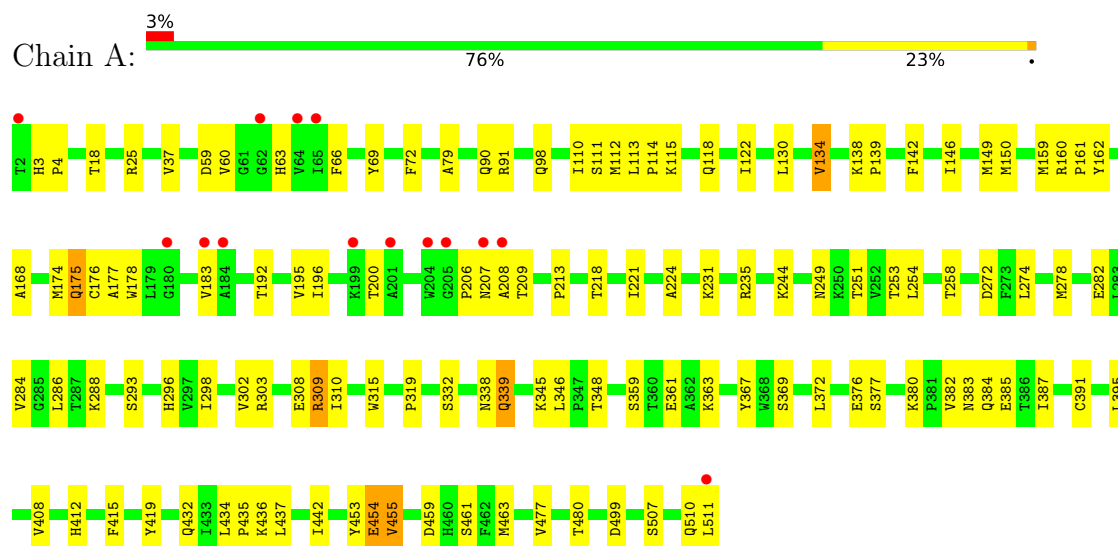
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	148	Total	O	0	0
			148	148		
5	B	151	Total	O	0	0
			151	151		
5	C	155	Total	O	0	0
			155	155		
5	D	140	Total	O	0	0
			140	140		
5	E	125	Total	O	0	0
			125	125		
5	F	122	Total	O	0	0
			122	122		
5	G	100	Total	O	0	0
			100	100		
5	H	98	Total	O	0	0
			98	98		

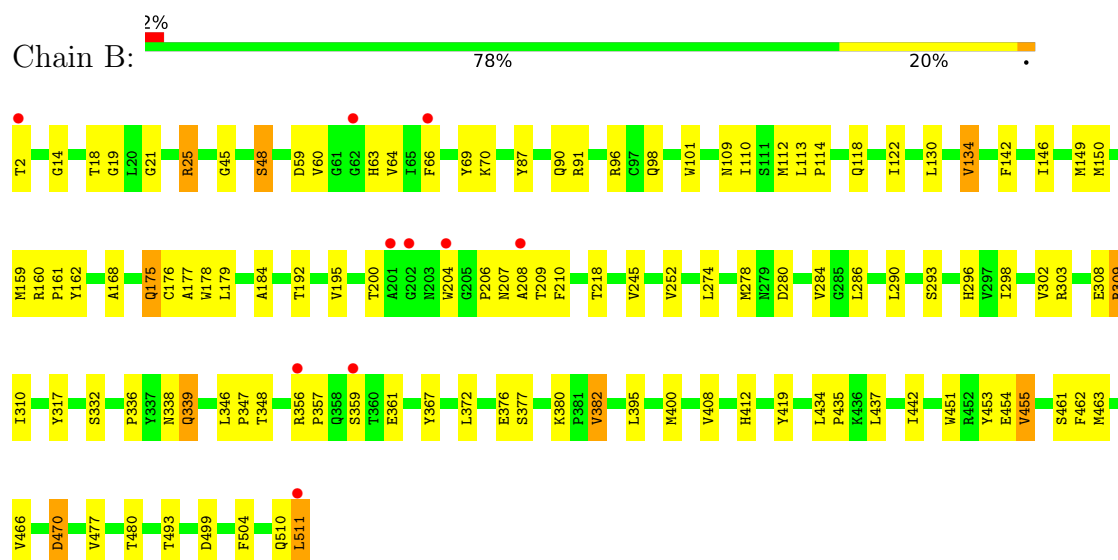
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

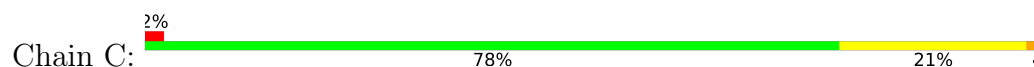
• Molecule 1: UDP-galactopyranose mutase

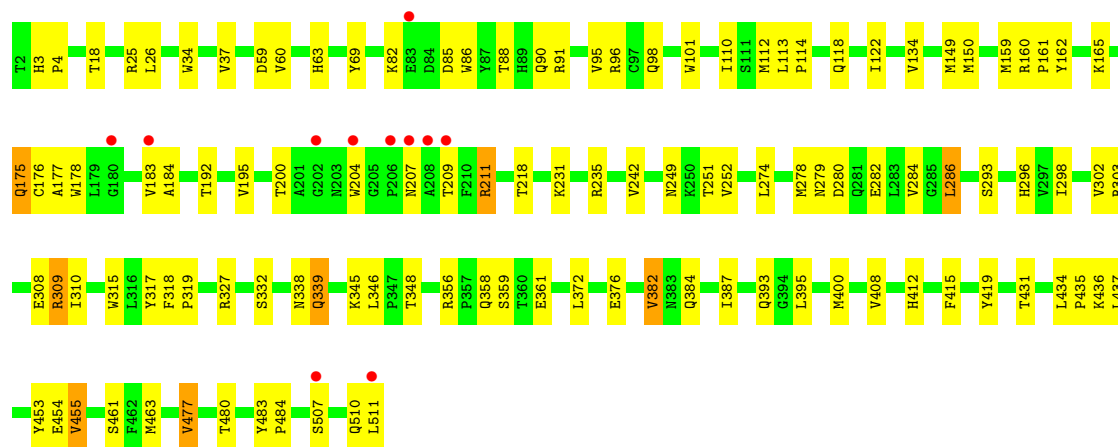


• Molecule 1: UDP-galactopyranose mutase

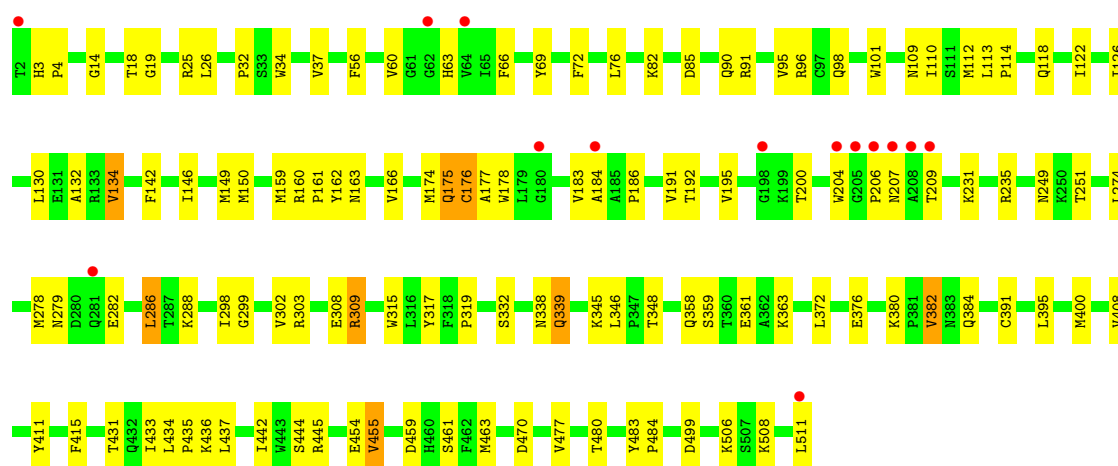
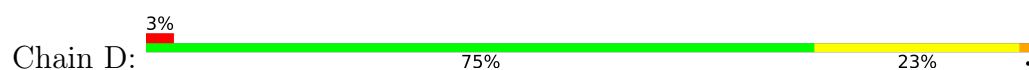


• Molecule 1: UDP-galactopyranose mutase

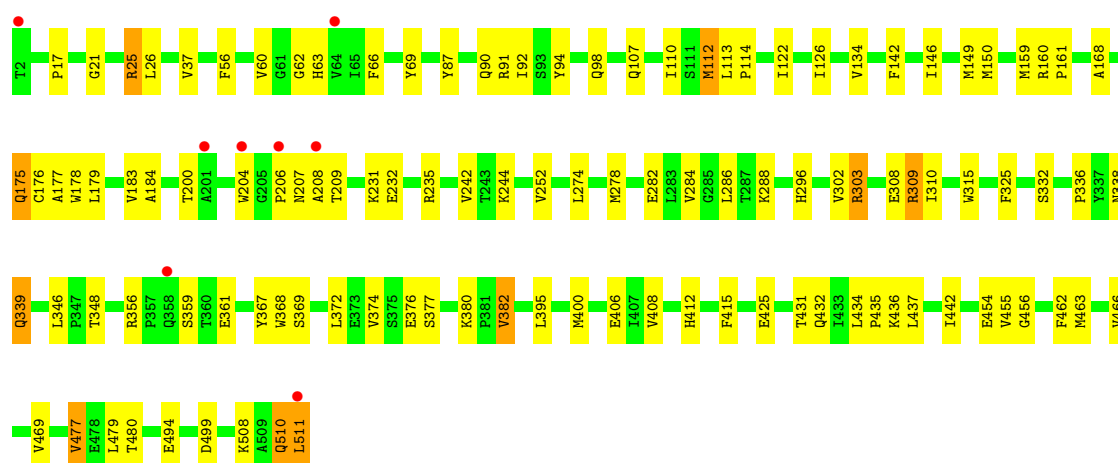
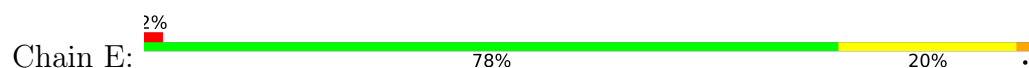




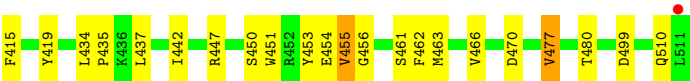
• Molecule 1: UDP-galactopyranose mutase



• Molecule 1: UDP-galactopyranose mutase



• Molecule 1: UDP-galactopyranose mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	71.44Å 129.20Å 172.82Å 90.07° 84.22° 82.24°	Depositor
Resolution (Å)	48.00 – 2.63 48.00 – 2.63	Depositor EDS
% Data completeness (in resolution range)	91.3 (48.00-2.63) 91.3 (48.00-2.63)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, R_{free}	0.185 , 0.231 (Not available) , 0.192	Depositor DCC
R_{free} test set	8265 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	33747	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: FAD, CL, UDP

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	0.32	0/4103	0.71	0/5576
1	B	0.32	0/4115	0.72	0/5592
1	C	0.32	0/4114	0.73	0/5590
1	D	0.32	0/4103	0.73	0/5576
1	E	0.33	0/4115	0.72	0/5592
1	F	0.32	0/4103	0.72	0/5576
1	G	0.30	0/4103	0.72	0/5576
1	H	0.30	0/4114	0.72	2/5590 (0.0%)
All	All	0.32	0/32870	0.72	2/44668 (0.0%)

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	E	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	80	LEU	CA-C-N	5.25	124.75	119.19
1	H	80	LEU	C-N-CA	5.25	124.75	119.19

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	303	ARG	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4004	0	3925	95	0
1	B	4015	0	3933	91	0
1	C	4015	0	3937	92	0
1	D	4004	0	3925	96	0
1	E	4015	0	3933	104	0
1	F	4004	0	3925	102	0
1	G	4004	0	3925	103	0
1	H	4015	0	3937	90	0
2	A	25	0	11	3	0
2	B	25	0	11	2	0
2	C	25	0	11	4	0
2	D	25	0	11	3	0
2	E	25	0	11	2	0
2	F	25	0	11	2	0
2	G	25	0	11	3	0
2	H	25	0	11	2	0
3	A	53	0	30	1	0
3	B	53	0	30	1	0
3	C	53	0	30	1	0
3	D	53	0	30	1	0
3	E	53	0	30	1	0
3	F	53	0	30	1	0
3	G	53	0	30	2	0
3	H	53	0	30	3	0
4	A	2	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
4	F	2	0	0	0	0
4	H	1	0	0	0	0
5	A	148	0	0	3	0
5	B	151	0	0	5	0
5	C	155	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	140	0	0	6	0
5	E	125	0	0	5	0
5	F	122	0	0	3	0
5	G	100	0	0	7	0
5	H	98	0	0	7	0
All	All	33747	0	31768	764	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 12.

All (764) 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:86:TRP:CG	1:C:211[A]:ARG:HH22	1.71	1.07
1:F:296:HIS:HD2	1:F:412:HIS:HE1	1.15	0.91
1:C:150:MET:HE1	1:C:159:MET:HE3	1.56	0.87
1:D:150:MET:HE1	1:D:159:MET:HE3	1.56	0.86
1:C:332:SER:HA	1:C:339:GLN:NE2	1.91	0.85
1:E:296:HIS:HD2	1:E:412:HIS:HE1	1.23	0.85
1:H:174:MET:SD	5:H:577:HOH:O	2.35	0.84
1:H:150:MET:HE1	1:H:159:MET:HE3	1.59	0.83
1:C:86:TRP:CG	1:C:211[A]:ARG:NH2	2.47	0.82
1:C:296:HIS:HD2	1:C:412:HIS:HE1	1.25	0.82
1:B:150:MET:HE1	1:B:159:MET:HE3	1.60	0.81
1:A:296:HIS:HD2	1:A:412:HIS:HE1	1.27	0.81
1:B:296:HIS:HD2	1:B:412:HIS:HE1	1.30	0.80
1:C:274:LEU:HG	1:C:278:MET:HE3	1.63	0.79
1:F:98:GLN:NE2	1:F:114:PRO:HG2	1.96	0.79
1:E:150:MET:HE1	1:E:159:MET:HE3	1.63	0.79
1:A:150:MET:HE1	1:A:159:MET:HE3	1.64	0.78
1:D:178:TRP:HB2	1:D:454:GLU:HG3	1.64	0.78
1:F:150:MET:HE1	1:F:159:MET:HE3	1.65	0.77
1:B:511:LEU:HD12	1:B:511:LEU:H	1.48	0.77
1:D:98:GLN:NE2	1:D:114:PRO:HG2	2.00	0.75
1:H:332:SER:HA	1:H:339:GLN:NE2	2.01	0.75
1:F:296:HIS:HD2	1:F:412:HIS:CE1	2.03	0.75
1:A:159:MET:HE1	2:A:800:UDP:O2	1.87	0.75
1:A:380:LYS:NZ	5:A:522:HOH:O	2.21	0.73
1:F:274:LEU:HG	1:F:278:MET:HE3	1.71	0.73
1:G:150:MET:HE1	1:G:159:MET:HE3	1.69	0.73
1:G:274:LEU:HG	1:G:278:MET:HE3	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:332:SER:HA	1:B:339:GLN:NE2	2.04	0.72
1:D:332:SER:HA	1:D:339:GLN:NE2	2.05	0.72
1:G:178:TRP:HB2	1:G:454:GLU:HG3	1.71	0.72
1:G:98:GLN:NE2	1:G:114:PRO:HG2	2.05	0.72
1:G:175:GLN:HG2	1:G:176:CYS:N	2.05	0.72
1:A:178:TRP:HB2	1:A:454:GLU:HG3	1.69	0.71
1:F:296:HIS:HE1	1:F:376:GLU:OE2	1.73	0.71
1:F:178:TRP:HB2	1:F:454:GLU:HG3	1.72	0.71
1:C:356:ARG:NH1	5:C:610:HOH:O	2.20	0.71
1:E:98:GLN:NE2	1:E:114:PRO:HG2	2.06	0.71
1:E:296:HIS:CD2	1:E:412:HIS:HE1	2.08	0.71
1:G:346:LEU:HD13	1:G:408:VAL:HG11	1.72	0.71
1:E:296:HIS:HD2	1:E:412:HIS:CE1	2.08	0.71
1:A:282:GLU:OE2	1:A:436:LYS:NZ	2.23	0.70
1:C:159:MET:HE1	2:C:800:UDP:O2	1.90	0.70
1:C:278:MET:HE1	1:C:437:LEU:HD22	1.72	0.70
1:F:455:VAL:HG13	1:F:455:VAL:O	1.91	0.70
1:G:332:SER:HA	1:G:339:GLN:NE2	2.06	0.70
1:E:380:LYS:NZ	5:E:885:HOH:O	2.23	0.70
1:A:303:ARG:NH2	1:A:346:LEU:O	2.23	0.70
1:B:178:TRP:HB2	1:B:454:GLU:HG3	1.73	0.70
1:F:296:HIS:CD2	1:F:412:HIS:HE1	2.04	0.70
1:B:98:GLN:NE2	1:B:114:PRO:HG2	2.06	0.69
1:D:175:GLN:HG2	1:D:176:CYS:N	2.07	0.69
1:E:178:TRP:HB2	1:E:454:GLU:HG3	1.73	0.69
1:E:346:LEU:HD13	1:E:408:VAL:HG11	1.75	0.69
1:A:332:SER:HA	1:A:339:GLN:NE2	2.07	0.69
1:C:178:TRP:HB2	1:C:454:GLU:HG3	1.74	0.69
1:E:274:LEU:HG	1:E:278:MET:HE3	1.75	0.69
1:B:175:GLN:HG2	1:B:176:CYS:N	2.07	0.69
1:B:455:VAL:HG13	1:B:455:VAL:O	1.93	0.69
1:E:175:GLN:HE21	1:E:178:TRP:HD1	1.39	0.69
1:H:159:MET:HE1	2:H:800:UDP:O2	1.93	0.69
1:F:278:MET:HE1	1:F:437:LEU:HD22	1.73	0.69
1:A:338:ASN:HB2	1:A:339:GLN:NE2	2.08	0.68
1:F:175:GLN:HG2	1:F:176:CYS:N	2.09	0.68
1:F:332:SER:HA	1:F:339:GLN:NE2	2.07	0.68
1:C:86:TRP:CD2	1:C:211[A]:ARG:NH2	2.62	0.68
1:H:178:TRP:HB2	1:H:454:GLU:HG3	1.76	0.68
1:B:380:LYS:NZ	5:B:883:HOH:O	2.27	0.67
1:C:296:HIS:HD2	1:C:412:HIS:CE1	2.11	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:25:ARG:NH2	1:H:499:ASP:OD2	2.26	0.67
1:C:455:VAL:O	1:C:455:VAL:HG13	1.94	0.67
1:C:296:HIS:HE1	1:C:376:GLU:OE2	1.78	0.67
1:H:332:SER:HA	1:H:339:GLN:HE21	1.59	0.67
1:B:296:HIS:CD2	1:B:412:HIS:HE1	2.13	0.67
1:D:178:TRP:HA	1:D:454:GLU:HG2	1.77	0.67
1:H:455:VAL:O	1:H:455:VAL:HG13	1.95	0.67
1:E:282:GLU:OE2	1:E:436:LYS:NZ	2.29	0.66
1:F:332:SER:HB3	1:F:369:SER:H	1.60	0.66
1:G:455:VAL:HG13	1:G:455:VAL:O	1.94	0.66
1:A:63:HIS:N	5:A:967:HOH:O	2.20	0.66
1:H:91:ARG:HD3	1:H:207:ASN:HB2	1.77	0.66
1:E:303:ARG:NH1	1:E:406:GLU:OE2	2.27	0.66
1:G:69:TYR:CG	1:G:463:MET:HG3	2.30	0.66
1:A:296:HIS:HD2	1:A:412:HIS:CE1	2.11	0.66
1:A:98:GLN:NE2	1:A:114:PRO:HG2	2.11	0.66
1:C:86:TRP:CB	1:C:211[A]:ARG:HH22	2.08	0.66
1:E:332:SER:HA	1:E:339:GLN:NE2	2.10	0.66
1:G:380:LYS:NZ	5:G:887:HOH:O	2.28	0.66
1:H:98:GLN:NE2	1:H:114:PRO:HG2	2.11	0.65
1:B:25:ARG:NH2	1:B:499:ASP:OD2	2.27	0.65
1:E:455:VAL:HG13	1:E:455:VAL:O	1.96	0.65
1:D:274:LEU:HG	1:D:278:MET:HE3	1.78	0.65
1:A:296:HIS:CD2	1:A:412:HIS:HE1	2.12	0.65
1:E:338:ASN:HB2	1:E:339:GLN:NE2	2.12	0.65
1:G:303:ARG:NH2	1:G:346:LEU:O	2.30	0.65
1:H:372:LEU:HD21	1:H:395:LEU:HD21	1.79	0.65
1:C:98:GLN:NE2	1:C:114:PRO:HG2	2.12	0.65
1:A:178:TRP:HA	1:A:454:GLU:HG2	1.78	0.65
1:E:175:GLN:HG2	1:E:176:CYS:N	2.12	0.65
1:B:356:ARG:HG2	1:B:357:PRO:HD2	1.79	0.64
1:A:134:VAL:HG22	1:D:134:VAL:HG22	1.77	0.64
1:B:296:HIS:HD2	1:B:412:HIS:CE1	2.13	0.64
1:G:296:HIS:HD2	1:G:412:HIS:HE1	1.45	0.64
1:A:274:LEU:HG	1:A:278:MET:HE3	1.79	0.64
1:G:338:ASN:HB2	1:G:339:GLN:HE21	1.62	0.64
1:C:175:GLN:HG2	1:C:176:CYS:N	2.12	0.64
1:H:303:ARG:NH2	1:H:346:LEU:O	2.28	0.64
1:C:274:LEU:CG	1:C:278:MET:HE3	2.27	0.64
1:B:178:TRP:HA	1:B:454:GLU:HG2	1.80	0.64
1:F:338:ASN:HB2	1:F:339:GLN:NE2	2.13	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:380:LYS:NZ	5:D:884:HOH:O	2.30	0.63
1:G:309:ARG:NH1	1:G:400:MET:O	2.31	0.63
1:C:178:TRP:HA	1:C:454:GLU:HG2	1.80	0.63
1:E:303:ARG:NH2	1:E:346:LEU:O	2.30	0.63
1:H:178:TRP:HA	1:H:454:GLU:HG2	1.80	0.63
1:A:118:GLN:NE2	1:A:195:VAL:HG13	2.14	0.63
1:B:274:LEU:HD22	5:B:520:HOH:O	1.99	0.63
1:D:455:VAL:HG21	1:D:480:THR:HG23	1.81	0.63
1:B:309:ARG:NH1	1:B:400:MET:O	2.32	0.63
1:G:178:TRP:HA	1:G:454:GLU:HG2	1.80	0.63
1:C:282:GLU:OE2	1:C:436:LYS:NZ	2.32	0.63
1:C:303:ARG:NH2	1:C:346:LEU:O	2.29	0.63
1:C:296:HIS:CD2	1:C:412:HIS:HE1	2.12	0.62
1:D:372:LEU:HD21	1:D:395:LEU:HD21	1.81	0.62
1:E:309:ARG:HG2	1:E:310:ILE:HD12	1.81	0.62
1:A:175:GLN:HG2	1:A:176:CYS:N	2.13	0.62
1:H:309:ARG:HG2	1:H:310:ILE:HD12	1.81	0.62
1:B:122:ILE:HD12	1:B:192:THR:HG22	1.81	0.62
1:D:455:VAL:CG2	1:D:480:THR:HG23	2.29	0.62
1:G:296:HIS:CD2	1:G:412:HIS:HE1	2.18	0.62
1:H:32:PRO:HD2	5:H:516:HOH:O	2.00	0.62
1:H:60:VAL:HG13	1:H:415:PHE:CE2	2.34	0.62
1:A:338:ASN:HB2	1:A:339:GLN:HE21	1.64	0.62
1:A:372:LEU:HD21	1:A:395:LEU:HD21	1.82	0.62
1:D:384:GLN:HG3	5:D:880:HOH:O	1.99	0.62
1:A:296:HIS:HE1	1:A:376:GLU:OE2	1.81	0.62
1:B:66[B]:PHE:CD2	1:B:66[B]:PHE:N	2.67	0.62
1:G:159:MET:HE1	2:G:800:UDP:O2	1.99	0.62
1:C:63:HIS:N	5:C:964:HOH:O	2.32	0.62
1:A:359:SER:OG	1:A:361:GLU:HG2	2.00	0.61
1:F:178:TRP:HA	1:F:454:GLU:HG2	1.81	0.61
1:G:346:LEU:HD13	1:G:408:VAL:CG1	2.29	0.61
1:D:338:ASN:HB2	1:D:339:GLN:NE2	2.15	0.61
1:C:69:TYR:CG	1:C:463:MET:HG3	2.36	0.61
1:D:91:ARG:HG3	1:D:315:TRP:CH2	2.36	0.61
1:F:159:MET:HE1	2:F:800:UDP:O2	2.00	0.61
1:A:69:TYR:CG	1:A:463:MET:HG3	2.35	0.61
1:D:122:ILE:HD12	1:D:192:THR:HG22	1.82	0.61
1:H:122:ILE:HD12	1:H:192:THR:HG22	1.83	0.61
1:B:274:LEU:HG	1:B:278:MET:HE3	1.81	0.61
1:C:122:ILE:HD12	1:C:192:THR:HG22	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:60:VAL:HG13	1:D:415:PHE:CE2	2.36	0.61
1:B:296:HIS:HE1	1:B:376:GLU:OE2	1.84	0.60
1:B:338:ASN:HB2	1:B:339:GLN:NE2	2.16	0.60
1:F:309:ARG:NH1	1:F:400:MET:O	2.34	0.60
1:D:338:ASN:HB2	1:D:339:GLN:HE21	1.65	0.60
1:E:232:GLU:HG2	1:G:232:GLU:HG3	1.83	0.60
1:E:510:GLN:HE22	1:G:506:LYS:HG2	1.67	0.60
1:C:332:SER:HA	1:C:339:GLN:HE21	1.64	0.60
1:D:91:ARG:HD3	1:D:207:ASN:CB	2.32	0.60
1:A:455:VAL:O	1:A:455:VAL:HG13	2.01	0.60
1:E:63:HIS:N	5:E:966:HOH:O	2.33	0.60
1:D:359:SER:OG	1:D:361:GLU:HG2	2.02	0.60
1:D:37:VAL:HG12	1:D:235:ARG:HB3	1.83	0.59
1:D:455:VAL:HG13	1:D:455:VAL:O	2.01	0.59
1:G:242:VAL:HG13	1:G:252:VAL:HG13	1.84	0.59
1:F:284:VAL:HG12	1:F:288:LYS:HE2	1.84	0.59
1:G:274:LEU:CG	1:G:278:MET:HE3	2.32	0.59
1:C:91:ARG:HG3	1:C:315:TRP:CH2	2.37	0.59
1:D:159:MET:HE1	2:D:800:UDP:O2	2.03	0.59
1:E:25:ARG:NH2	1:E:499:ASP:OD2	2.31	0.59
1:E:296:HIS:HE1	1:E:376:GLU:OE2	1.86	0.59
1:E:278:MET:HE1	1:E:437:LEU:HD22	1.84	0.59
1:H:175:GLN:HG2	1:H:176:CYS:N	2.18	0.59
1:A:274:LEU:CG	1:A:278:MET:HE3	2.33	0.59
1:F:346:LEU:HD13	1:F:408:VAL:HG11	1.85	0.59
1:B:348:THR:N	5:B:545:HOH:O	2.36	0.58
1:F:37:VAL:HG12	1:F:235:ARG:HB3	1.83	0.58
1:A:122:ILE:HD12	1:A:192:THR:HG22	1.85	0.58
1:D:69:TYR:CG	1:D:463:MET:HG3	2.37	0.58
1:H:118:GLN:NE2	1:H:195:VAL:HG13	2.18	0.58
1:D:274:LEU:CG	1:D:278:MET:HE3	2.33	0.58
1:D:118:GLN:NE2	1:D:195:VAL:HG13	2.18	0.58
1:E:178:TRP:HA	1:E:454:GLU:HG2	1.85	0.58
1:G:175:GLN:HG2	1:G:177:ALA:H	1.68	0.58
1:H:91:ARG:HD3	1:H:207:ASN:CB	2.33	0.58
1:D:303:ARG:NH2	1:D:346:LEU:O	2.32	0.58
1:D:431:THR:O	1:D:435:PRO:HG2	2.03	0.58
1:E:183:VAL:HG22	2:E:800:UDP:H1'	1.86	0.58
1:F:338:ASN:HB2	1:F:339:GLN:HE21	1.67	0.58
1:G:309:ARG:HG2	1:G:310:ILE:HD12	1.86	0.58
1:H:69:TYR:CG	1:H:463:MET:HG3	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:91:ARG:HG3	1:H:315:TRP:CH2	2.39	0.58
1:C:372:LEU:HD21	1:C:395:LEU:HD21	1.85	0.58
1:G:338:ASN:HB2	1:G:339:GLN:NE2	2.17	0.58
1:B:142:PHE:O	1:B:146:ILE:HG13	2.03	0.57
1:F:367:TYR:HE1	1:F:408:VAL:HG21	1.67	0.57
1:F:274:LEU:CG	1:F:278:MET:HE3	2.34	0.57
1:H:359:SER:OG	1:H:361:GLU:HG2	2.05	0.57
1:B:112:MET:CE	1:B:200:THR:HG22	2.34	0.57
1:H:112:MET:CE	1:H:200:THR:HG22	2.34	0.57
1:F:69:TYR:CG	1:F:463:MET:HG3	2.39	0.57
1:B:159:MET:HE1	2:B:800:UDP:O2	2.05	0.57
1:D:130:LEU:HD23	1:F:130:LEU:HD23	1.87	0.57
1:E:332:SER:HB3	1:E:369:SER:H	1.70	0.57
1:C:150:MET:CE	1:C:159:MET:HE3	2.33	0.56
1:C:384:GLN:HG3	5:C:879:HOH:O	2.05	0.56
1:D:183:VAL:HG22	2:D:800:UDP:H1'	1.87	0.56
1:D:109:ASN:HA	5:D:1045:HOH:O	2.03	0.56
1:A:91:ARG:HH11	1:A:207:ASN:HB2	1.69	0.56
1:C:88:THR:HG23	1:C:211[A]:ARG:HG2	1.85	0.56
1:C:507:SER:O	1:C:511:LEU:HD13	2.05	0.56
1:F:376:GLU:HB2	1:F:382:VAL:CG1	2.35	0.56
1:H:175:GLN:HG2	1:H:177:ALA:H	1.71	0.56
1:A:60:VAL:HG12	1:A:60:VAL:O	2.05	0.56
1:B:142:PHE:CE1	1:B:179:LEU:HD23	2.40	0.56
1:C:175:GLN:HG2	1:C:177:ALA:H	1.70	0.56
1:A:18:THR:OG1	1:A:461:SER:HB3	2.06	0.56
1:B:175:GLN:HG2	1:B:177:ALA:H	1.70	0.56
1:E:455:VAL:CG2	1:E:480:THR:HG23	2.36	0.56
1:A:244:LYS:HB2	1:A:253:THR:HB	1.88	0.56
1:C:18:THR:OG1	1:C:461:SER:HB3	2.05	0.56
1:D:508:LYS:HA	1:D:511:LEU:HD12	1.87	0.56
1:G:25:ARG:NH2	1:G:499:ASP:OD2	2.34	0.56
1:F:359:SER:OG	1:F:361:GLU:HG2	2.05	0.56
1:A:346:LEU:HD13	1:A:408:VAL:HG11	1.88	0.55
1:G:122:ILE:HD12	1:G:192:THR:HG22	1.88	0.55
1:H:274:LEU:HG	1:H:278:MET:HE3	1.88	0.55
1:A:455:VAL:CG2	1:A:480:THR:HG23	2.37	0.55
1:B:109:ASN:HA	5:B:602:HOH:O	2.06	0.55
1:E:175:GLN:HE21	1:E:178:TRP:CD1	2.23	0.55
1:E:359:SER:OG	1:E:361:GLU:HG2	2.05	0.55
1:C:86:TRP:HB2	1:C:211[A]:ARG:HH12	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:367:TYR:HE1	1:E:408:VAL:HG21	1.72	0.55
1:F:112:MET:CE	1:F:200:THR:HG22	2.36	0.55
1:A:254:LEU:HD12	1:A:258:THR:HB	1.88	0.55
1:D:91:ARG:HD3	1:D:207:ASN:HB2	1.87	0.55
1:A:507:SER:O	1:A:511:LEU:HD13	2.05	0.55
1:E:338:ASN:HB2	1:E:339:GLN:HE21	1.71	0.55
1:F:183:VAL:HG22	2:F:800:UDP:H1'	1.88	0.55
1:F:142:PHE:HD2	1:F:174:MET:HE3	1.70	0.55
1:H:112:MET:HE3	1:H:200:THR:HG22	1.89	0.55
1:A:112:MET:CE	1:A:200:THR:HG22	2.36	0.55
1:F:303:ARG:NH2	1:F:346:LEU:O	2.35	0.55
1:B:274:LEU:CG	1:B:278:MET:HE3	2.36	0.54
1:B:278:MET:HE1	1:B:437:LEU:HD22	1.88	0.54
1:E:274:LEU:CG	1:E:278:MET:HE3	2.37	0.54
1:F:309:ARG:HG2	1:F:310:ILE:HD12	1.89	0.54
1:H:142:PHE:HD2	1:H:174:MET:HE3	1.72	0.54
1:A:278:MET:HE1	1:A:437:LEU:HD22	1.88	0.54
1:C:37:VAL:HG12	1:C:235:ARG:HB3	1.89	0.54
1:D:69:TYR:CD1	1:D:463:MET:HG3	2.42	0.54
1:E:159:MET:HE1	2:E:800:UDP:O2	2.06	0.54
1:G:372:LEU:HD21	1:G:395:LEU:HD21	1.89	0.54
1:A:91:ARG:NH1	1:A:207:ASN:HB2	2.21	0.54
1:F:348:THR:N	5:F:520:HOH:O	2.39	0.54
1:F:110:ILE:HG23	1:F:113:LEU:HD12	1.90	0.54
1:F:434:LEU:HB2	1:F:435:PRO:HD3	1.89	0.54
1:C:91:ARG:HD3	1:C:207:ASN:HB2	1.90	0.54
1:D:112:MET:HE1	1:D:200:THR:HG22	1.90	0.54
1:F:310:ILE:HG12	1:F:368:TRP:CZ3	2.43	0.54
1:B:332:SER:HA	1:B:339:GLN:HE21	1.70	0.54
1:F:332:SER:HB3	1:F:369:SER:N	2.21	0.54
1:H:434:LEU:HB2	1:H:435:PRO:HD3	1.88	0.54
1:A:175:GLN:HG2	1:A:177:ALA:H	1.72	0.53
1:D:434:LEU:HB2	1:D:435:PRO:HD3	1.90	0.53
1:F:91:ARG:HD3	1:F:207:ASN:HB2	1.90	0.53
1:H:278:MET:HE1	1:H:437:LEU:HD22	1.90	0.53
1:C:110:ILE:HG23	1:C:113:LEU:HD12	1.89	0.53
1:D:112:MET:CE	1:D:200:THR:HG22	2.37	0.53
1:F:249:ASN:O	1:F:251:THR:HG23	2.09	0.53
1:G:26:LEU:HG	1:G:469:VAL:HG21	1.90	0.53
1:C:309:ARG:HG2	1:C:310:ILE:HD12	1.89	0.53
1:D:278:MET:HE1	1:D:437:LEU:HD22	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:69:TYR:CD1	1:G:463:MET:HG3	2.43	0.53
1:A:91:ARG:HG3	1:A:315:TRP:CH2	2.44	0.53
1:G:146:ILE:HD13	1:G:159:MET:HB3	1.91	0.53
1:A:91:ARG:HD3	1:A:207:ASN:HB2	1.91	0.53
1:A:162:TYR:HB2	1:A:319:PRO:HB3	1.90	0.53
1:A:278:MET:HE2	1:A:442:ILE:HD13	1.91	0.53
1:E:456:GLY:HA2	5:E:908:HOH:O	2.09	0.53
1:E:508:LYS:O	1:E:511:LEU:HD12	2.09	0.53
1:C:112:MET:CE	1:C:200:THR:HG22	2.39	0.53
1:C:86:TRP:CB	1:C:211[A]:ARG:HH12	2.22	0.53
1:E:112:MET:CE	1:E:200:THR:HG22	2.39	0.53
1:F:91:ARG:HG3	1:F:315:TRP:CH2	2.44	0.53
1:G:274:LEU:O	1:G:278:MET:HG3	2.08	0.53
1:B:69:TYR:CG	1:B:463:MET:HG3	2.44	0.52
1:C:69:TYR:CD1	1:C:463:MET:HG3	2.44	0.52
1:C:118:GLN:NE2	1:C:195:VAL:HG13	2.24	0.52
1:C:455:VAL:O	1:C:455:VAL:CG1	2.57	0.52
1:E:376:GLU:HB2	1:E:382:VAL:CG1	2.39	0.52
1:F:98:GLN:HE22	1:F:114:PRO:HG2	1.72	0.52
1:B:118:GLN:NE2	1:B:195:VAL:HG13	2.24	0.52
1:B:376:GLU:HB2	1:B:382:VAL:CG1	2.40	0.52
1:E:110:ILE:HG23	1:E:113:LEU:HD12	1.92	0.52
1:G:494:GLU:HG2	1:H:477:VAL:HA	1.91	0.52
1:D:142:PHE:HD2	1:D:174:MET:HE3	1.75	0.52
1:B:455:VAL:O	1:B:455:VAL:CG1	2.58	0.52
1:E:510:GLN:OE1	1:G:510:GLN:NE2	2.43	0.52
1:F:332:SER:HA	1:F:339:GLN:HE21	1.73	0.52
1:G:91:ARG:HD3	1:G:207:ASN:HB2	1.92	0.52
1:B:112:MET:HE3	1:B:200:THR:HG22	1.91	0.52
1:C:60:VAL:HG13	1:C:415:PHE:CE2	2.45	0.52
1:E:309:ARG:NH1	1:E:400:MET:O	2.43	0.52
1:B:18:THR:OG1	1:B:461:SER:HB3	2.10	0.52
1:E:346:LEU:HD13	1:E:408:VAL:CG1	2.40	0.52
1:G:296:HIS:HE1	1:G:376:GLU:OE2	1.93	0.52
1:D:175:GLN:HG2	1:D:177:ALA:H	1.75	0.51
1:B:434:LEU:HB2	1:B:435:PRO:HD3	1.93	0.51
1:C:26:LEU:HD13	1:C:34:TRP:HB2	1.93	0.51
1:C:162:TYR:HB2	1:C:319:PRO:HB3	1.92	0.51
1:E:66[A]:PHE:CE2	1:E:206:PRO:HB3	2.44	0.51
1:E:175:GLN:HG2	1:E:177:ALA:H	1.74	0.51
1:F:118:GLN:NE2	1:F:195:VAL:HG13	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:56:PHE:CE1	1:H:339:GLN:HB3	2.45	0.51
1:C:359:SER:OG	1:C:361:GLU:HG2	2.10	0.51
1:E:69:TYR:CG	1:E:463:MET:HG3	2.45	0.51
1:G:453:TYR:C	1:G:455:VAL:H	2.19	0.51
1:A:112:MET:HE1	1:A:200:THR:HG22	1.91	0.51
1:H:25:ARG:HH11	1:H:470:ASP:CG	2.18	0.51
1:A:159:MET:HE1	2:A:800:UDP:C2	2.45	0.51
1:A:434:LEU:HB2	1:A:435:PRO:HD3	1.93	0.51
1:C:434:LEU:HB2	1:C:435:PRO:HD3	1.92	0.51
1:B:359:SER:OG	1:B:361:GLU:HG2	2.11	0.51
1:C:165:LYS:NZ	1:C:318:PHE:O	2.38	0.51
1:E:332:SER:HB3	1:E:369:SER:N	2.26	0.51
1:H:338:ASN:HB2	1:H:339:GLN:NE2	2.25	0.51
1:E:175:GLN:NE2	1:E:178:TRP:HD1	2.06	0.51
1:E:372:LEU:HD21	1:E:395:LEU:HD21	1.91	0.51
1:F:372:LEU:HD21	1:F:395:LEU:HD21	1.91	0.51
1:C:393:GLN:HG2	5:C:551:HOH:O	2.10	0.51
1:D:278:MET:HE2	1:D:442:ILE:HD13	1.92	0.51
1:E:168:ALA:HB1	1:E:377:SER:OG	2.11	0.51
1:F:278:MET:HE1	1:F:437:LEU:CD2	2.41	0.51
1:C:278:MET:HE1	1:C:437:LEU:CD2	2.39	0.51
1:B:25:ARG:HH11	1:B:470:ASP:CG	2.18	0.51
1:C:91:ARG:HD3	1:C:207:ASN:CB	2.40	0.51
1:C:338:ASN:HB2	1:C:339:GLN:NE2	2.25	0.51
1:G:310:ILE:HG12	1:G:368:TRP:CZ3	2.46	0.51
1:H:332:SER:HB3	1:H:369:SER:H	1.75	0.51
1:E:175:GLN:HE22	1:E:178:TRP:HB3	1.76	0.50
1:H:332:SER:HB3	1:H:369:SER:N	2.25	0.50
1:G:504:PHE:CD2	1:H:32:PRO:HG3	2.45	0.50
1:G:107:GLN:HB2	1:G:186:PRO:HG3	1.92	0.50
1:G:180:GLY:N	5:G:526:HOH:O	2.25	0.50
1:D:162:TYR:HB2	1:D:319:PRO:HB3	1.93	0.50
1:E:60:VAL:HG12	1:E:60:VAL:O	2.11	0.50
1:F:455:VAL:CG2	1:F:480:THR:HG23	2.42	0.50
1:G:359:SER:OG	1:G:361:GLU:HG2	2.11	0.50
1:H:447:ARG:NH1	5:H:874:HOH:O	2.33	0.50
1:D:110:ILE:HG23	1:D:113:LEU:HD12	1.92	0.50
1:F:455:VAL:O	1:F:455:VAL:CG1	2.58	0.50
1:H:146:ILE:HD13	1:H:159:MET:HB3	1.92	0.50
1:G:112:MET:CE	1:G:200:THR:HG22	2.42	0.50
1:D:63:HIS:N	5:D:965:HOH:O	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:VAL:HG22	1:F:134:VAL:HG22	1.94	0.50
1:B:338:ASN:HB2	1:B:339:GLN:HE21	1.76	0.50
1:B:372:LEU:HD21	1:B:395:LEU:HD21	1.92	0.49
1:D:146:ILE:HD13	1:D:159:MET:HB3	1.93	0.49
1:E:107:GLN:HE21	1:E:184:ALA:HB3	1.76	0.49
1:A:196:ILE:HD11	1:B:122:ILE:HG13	1.94	0.49
1:E:91:ARG:HD3	1:E:207:ASN:HB2	1.93	0.49
1:F:142:PHE:O	1:F:146:ILE:HG13	2.12	0.49
1:F:162:TYR:HH	1:F:317:TYR:HD1	1.59	0.49
1:B:168:ALA:HB1	1:B:377:SER:OG	2.12	0.49
1:D:91:ARG:HH11	1:D:207:ASN:HB2	1.78	0.49
1:G:69:TYR:CD2	1:G:463:MET:HG3	2.47	0.49
1:G:118:GLN:NE2	1:G:195:VAL:HG13	2.27	0.49
1:A:455:VAL:HG22	1:A:480:THR:HG23	1.93	0.49
1:E:455:VAL:HG21	1:E:480:THR:HG23	1.95	0.49
1:A:60:VAL:HG13	1:A:415:PHE:CE2	2.47	0.49
1:A:91:ARG:HD3	1:A:207:ASN:CB	2.43	0.49
1:C:82:LYS:HB2	1:C:85:ASP:OD1	2.13	0.49
1:H:455:VAL:O	1:H:455:VAL:CG1	2.61	0.49
1:B:146:ILE:HD13	1:B:159:MET:HB3	1.94	0.49
1:B:278:MET:HE2	1:B:442:ILE:HD13	1.94	0.49
1:H:91:ARG:HH11	1:H:207:ASN:HB2	1.78	0.49
1:E:160:ARG:N	1:E:161:PRO:HD2	2.28	0.49
1:E:278:MET:HE2	1:E:442:ILE:HD13	1.94	0.49
1:F:146:ILE:HD13	1:F:159:MET:HB3	1.94	0.49
1:F:282:GLU:OE2	1:F:436:LYS:NZ	2.39	0.49
1:H:160:ARG:N	1:H:161:PRO:HD2	2.28	0.49
1:D:346:LEU:HD13	1:D:408:VAL:HG11	1.94	0.49
1:F:112:MET:HE1	1:F:200:THR:HG22	1.94	0.49
1:B:87:TYR:CE2	1:B:336:PRO:HD2	2.48	0.48
1:D:18:THR:OG1	1:D:461:SER:HB3	2.13	0.48
1:D:66:PHE:CE2	1:D:206:PRO:HB3	2.49	0.48
1:D:282:GLU:OE2	1:D:436:LYS:NZ	2.42	0.48
1:F:14:GLY:O	1:F:19:GLY:HA3	2.13	0.48
1:F:313:LYS:HB3	1:F:316:LEU:HD21	1.95	0.48
1:B:25:ARG:HD2	1:B:466:VAL:HG13	1.95	0.48
1:C:455:VAL:CG2	1:C:480:THR:HG23	2.43	0.48
1:D:60:VAL:HG12	1:D:60:VAL:O	2.13	0.48
1:H:66:PHE:CE2	1:H:206:PRO:HB3	2.47	0.48
1:A:249:ASN:O	1:A:251:THR:HG23	2.12	0.48
1:B:336:PRO:HD3	1:C:356:ARG:NH1	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:91:ARG:HH11	1:C:207:ASN:HB2	1.77	0.48
1:E:142:PHE:O	1:E:146:ILE:HG13	2.13	0.48
1:G:455:VAL:O	1:G:455:VAL:CG1	2.59	0.48
1:D:25:ARG:HH11	1:D:470:ASP:CG	2.22	0.48
1:D:91:ARG:HD3	1:D:207:ASN:HB3	1.94	0.48
1:E:91:ARG:HD3	1:E:207:ASN:CB	2.43	0.48
1:G:447:ARG:NH1	5:G:611:HOH:O	2.31	0.48
1:G:122:ILE:O	1:G:126:ILE:HG13	2.14	0.48
1:C:453:TYR:C	1:C:455:VAL:H	2.21	0.48
1:F:56:PHE:CE1	1:F:339:GLN:HB3	2.47	0.48
1:G:453:TYR:C	1:G:455:VAL:N	2.71	0.48
1:A:309:ARG:HG2	1:A:310:ILE:HD12	1.95	0.48
1:D:66:PHE:CD2	1:D:206:PRO:HB3	2.49	0.48
1:D:303:ARG:NH1	1:D:348:THR:OG1	2.46	0.48
1:G:296:HIS:HD2	1:G:412:HIS:CE1	2.28	0.48
1:H:274:LEU:CG	1:H:278:MET:HE3	2.43	0.48
1:A:183:VAL:HG22	2:A:800:UDP:H1'	1.94	0.48
1:F:346:LEU:HD12	1:F:367:TYR:CZ	2.49	0.48
1:H:66:PHE:CD2	1:H:206:PRO:HB3	2.48	0.48
1:A:69:TYR:CD1	1:A:463:MET:HG3	2.49	0.48
1:E:91:ARG:HG3	1:E:315:TRP:CH2	2.48	0.48
1:F:25:ARG:NH2	1:F:499:ASP:OD2	2.36	0.48
1:B:178:TRP:HB2	1:B:454:GLU:CG	2.43	0.47
1:E:284:VAL:O	1:E:288:LYS:HG3	2.14	0.47
1:B:504:PHE:CD2	1:D:32:PRO:HB3	2.49	0.47
1:E:477:VAL:HG22	1:E:479:LEU:HG	1.95	0.47
1:F:114:PRO:O	1:F:118:GLN:HG3	2.14	0.47
1:G:149:MET:O	1:G:186:PRO:HD2	2.15	0.47
1:D:98:GLN:HE21	1:D:114:PRO:HG2	1.79	0.47
1:E:175:GLN:NE2	1:E:178:TRP:CD1	2.82	0.47
1:E:244:LYS:HG2	5:E:590:HOH:O	2.14	0.47
1:E:510:GLN:O	1:E:511:LEU:C	2.57	0.47
1:G:159:MET:HE1	2:G:800:UDP:C2	2.49	0.47
1:H:278:MET:HE2	1:H:442:ILE:HD13	1.95	0.47
1:F:160:ARG:N	1:F:161:PRO:HD2	2.30	0.47
1:G:332:SER:HB3	1:G:369:SER:H	1.80	0.47
1:C:26:LEU:HD13	1:C:34:TRP:CB	2.45	0.47
1:H:60:VAL:HG12	1:H:60:VAL:O	2.14	0.47
1:A:59:ASP:OD2	1:A:218:THR:HG23	2.15	0.47
1:C:59:ASP:OD2	1:C:218:THR:HG23	2.14	0.47
1:H:150:MET:CE	1:H:159:MET:HE3	2.37	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:LEU:HD23	1:B:130:LEU:HD23	1.97	0.47
1:A:142:PHE:HD2	1:A:174:MET:HE3	1.80	0.47
1:B:278:MET:HE1	1:B:437:LEU:CD2	2.44	0.47
1:B:303:ARG:NH2	1:B:346:LEU:O	2.44	0.47
1:D:63:HIS:CE1	3:D:601:FAD:C9A	2.98	0.47
1:D:160:ARG:N	1:D:161:PRO:HD2	2.30	0.47
1:F:347:PRO:C	5:F:520:HOH:O	2.58	0.47
1:F:384:GLN:O	1:F:387:ILE:HG22	2.13	0.47
1:G:60:VAL:HG13	1:G:415:PHE:CE2	2.49	0.47
1:E:310:ILE:HG12	1:E:368:TRP:CZ3	2.50	0.47
1:F:163:ASN:HA	1:F:166:VAL:HG12	1.97	0.47
1:G:447:ARG:NH2	5:G:537:HOH:O	2.47	0.47
1:A:150:MET:CE	1:A:159:MET:HE3	2.38	0.47
1:A:332:SER:HB3	1:A:369:SER:N	2.30	0.47
1:B:455:VAL:CG2	1:B:480:THR:HG23	2.45	0.47
1:C:183:VAL:HG22	2:C:800:UDP:H1'	1.97	0.47
1:E:87:TYR:CE2	1:E:336:PRO:HD2	2.50	0.47
1:A:110:ILE:HG23	1:A:113:LEU:HD12	1.97	0.47
1:B:347:PRO:C	5:B:545:HOH:O	2.58	0.47
1:F:60:VAL:HG13	1:F:415:PHE:CE2	2.50	0.47
1:F:280:ASP:O	1:F:284:VAL:HG23	2.15	0.47
1:G:46:LEU:HB2	3:G:607:FAD:O4'	2.15	0.47
1:D:309:ARG:NH1	1:D:400:MET:O	2.48	0.46
1:E:325:PHE:HA	1:E:374:VAL:HG22	1.97	0.46
1:B:96:ARG:HB2	1:B:101:TRP:CZ3	2.50	0.46
1:C:160:ARG:N	1:C:161:PRO:HD2	2.30	0.46
1:E:26:LEU:HG	1:E:469:VAL:HG21	1.97	0.46
1:F:212:PHE:CG	1:F:213:PRO:HD2	2.50	0.46
1:A:332:SER:HA	1:A:339:GLN:HE21	1.79	0.46
1:E:66[B]:PHE:CD2	1:E:66[B]:PHE:N	2.82	0.46
1:H:18:THR:OG1	1:H:461:SER:HB3	2.16	0.46
1:B:293:SER:OG	1:B:419:TYR:HB2	2.15	0.46
1:E:112:MET:HE3	1:E:200:THR:HG22	1.98	0.46
1:F:60:VAL:HG12	1:F:60:VAL:O	2.14	0.46
1:F:69:TYR:CD1	1:F:463:MET:HG3	2.51	0.46
1:G:110:ILE:HG23	1:G:113:LEU:HD12	1.97	0.46
1:A:367:TYR:HE1	1:A:408:VAL:HG21	1.80	0.46
1:E:91:ARG:HH11	1:E:207:ASN:HB2	1.79	0.46
1:E:184:ALA:HB2	1:E:204:TRP:CD2	2.50	0.46
1:F:367:TYR:CE1	1:F:408:VAL:HG21	2.49	0.46
1:G:282:GLU:OE2	1:G:436:LYS:NZ	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:37:VAL:HG12	1:H:235:ARG:HB3	1.97	0.46
1:H:367:TYR:HE1	1:H:408:VAL:HG21	1.79	0.46
1:H:182:ARG:HG2	5:H:864:HOH:O	2.16	0.46
1:B:290:LEU:HD13	1:B:451:TRP:CD2	2.51	0.46
1:E:91:ARG:NH1	1:E:207:ASN:HB2	2.30	0.46
1:G:263:LYS:NZ	5:G:945:HOH:O	2.47	0.46
1:H:92:ILE:HG22	1:H:94:TYR:HE2	1.81	0.46
1:A:346:LEU:HD13	1:A:408:VAL:CG1	2.46	0.46
1:B:25:ARG:NH1	1:B:470:ASP:OD1	2.47	0.46
1:B:66[A]:PHE:CE2	1:B:206:PRO:HB3	2.51	0.46
1:C:60:VAL:HG12	1:C:60:VAL:O	2.16	0.46
1:C:242:VAL:HG13	1:C:252:VAL:HG13	1.97	0.46
1:D:98:GLN:HE22	1:D:114:PRO:HG2	1.78	0.46
1:F:91:ARG:HH11	1:F:207:ASN:HB2	1.81	0.46
1:G:431:THR:O	1:G:435:PRO:HG2	2.16	0.46
1:D:184:ALA:HB2	1:D:204:TRP:CD2	2.51	0.45
1:E:56:PHE:CE1	1:E:339:GLN:HB3	2.51	0.45
1:F:184:ALA:HB2	1:F:204:TRP:CD2	2.51	0.45
1:G:456:GLY:HA2	5:G:516:HOH:O	2.16	0.45
1:F:175:GLN:HG2	1:F:177:ALA:H	1.81	0.45
1:G:278:MET:HE1	1:G:437:LEU:HD22	1.97	0.45
1:G:495:ARG:O	1:G:496:ARG:HD2	2.16	0.45
1:A:298:ILE:HB	1:A:372:LEU:HB2	1.97	0.45
1:B:453:TYR:C	1:B:455:VAL:H	2.23	0.45
3:B:604:FAD:H9	3:B:604:FAD:H1'1	1.44	0.45
1:G:37:VAL:HG12	1:G:235:ARG:HB3	1.97	0.45
1:G:168:ALA:HB1	1:G:377:SER:OG	2.16	0.45
1:A:284:VAL:HG12	1:A:288:LYS:HE3	1.99	0.45
1:E:431:THR:O	1:E:435:PRO:HG2	2.17	0.45
1:F:278:MET:HE2	1:F:442:ILE:HD13	1.99	0.45
1:G:432:GLN:C	1:G:435:PRO:HD2	2.41	0.45
1:E:17:PRO:O	1:E:462:PHE:HA	2.16	0.45
1:F:87:TYR:CE2	1:F:336:PRO:HD2	2.51	0.45
1:G:280:ASP:O	1:G:284:VAL:HG23	2.16	0.45
1:H:63:HIS:CE1	3:H:605:FAD:C9A	2.99	0.45
1:B:69:TYR:CD1	1:B:463:MET:HG3	2.51	0.45
1:B:298:ILE:HB	1:B:372:LEU:HB2	1.97	0.45
1:C:298:ILE:HB	1:C:372:LEU:HB2	1.98	0.45
1:E:142:PHE:CE1	1:E:179:LEU:HD23	2.52	0.45
1:E:150:MET:CE	1:E:159:MET:HE3	2.40	0.45
1:G:160:ARG:N	1:G:161:PRO:HD2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:280:ASP:O	1:C:284:VAL:HG23	2.17	0.45
1:H:79:ALA:HB1	1:H:224:ALA:CB	2.47	0.45
1:B:63:HIS:HB2	1:B:218:THR:HG21	1.99	0.45
1:B:110:ILE:HG23	1:B:113:LEU:HD12	1.98	0.45
1:B:511:LEU:HD12	1:B:511:LEU:N	2.24	0.45
1:E:37:VAL:HG12	1:E:235:ARG:HB3	1.99	0.45
1:E:278:MET:HE1	1:E:437:LEU:CD2	2.47	0.45
1:A:213:PRO:HG2	1:A:218:THR:HA	1.99	0.45
1:G:101:TRP:CH2	1:G:316:LEU:HD13	2.52	0.45
1:H:453:TYR:C	1:H:455:VAL:N	2.75	0.45
1:H:456:GLY:O	3:H:605:FAD:O3'	2.35	0.45
1:A:146:ILE:HD13	1:A:159:MET:HB3	1.99	0.45
1:B:59:ASP:OD2	1:B:218:THR:HG23	2.16	0.45
1:C:86:TRP:CB	1:C:211[A]:ARG:NH2	2.76	0.45
1:D:25:ARG:O	1:D:25:ARG:HG3	2.16	0.45
1:D:91:ARG:NH1	1:D:207:ASN:HB2	2.32	0.45
1:D:95:VAL:HG22	1:D:317:TYR:HB2	1.98	0.45
1:D:150:MET:CE	1:D:159:MET:HE3	2.39	0.45
1:E:207:ASN:O	1:E:208:ALA:C	2.60	0.45
1:F:17:PRO:O	1:F:462:PHE:HA	2.17	0.45
1:A:332:SER:HB3	1:A:369:SER:H	1.81	0.44
1:C:303:ARG:NH1	1:C:348:THR:OG1	2.50	0.44
1:D:159:MET:HE1	2:D:800:UDP:C2	2.51	0.44
1:D:298:ILE:HB	1:D:372:LEU:HB2	1.99	0.44
1:E:178:TRP:HB2	1:E:454:GLU:CG	2.45	0.44
1:F:456:GLY:HA2	5:F:518:HOH:O	2.17	0.44
1:H:183:VAL:HG22	2:H:800:UDP:H1'	1.99	0.44
1:H:274:LEU:O	1:H:278:MET:HG3	2.17	0.44
1:A:432:GLN:HG2	5:A:571:HOH:O	2.16	0.44
1:B:339:GLN:N	1:B:339:GLN:CD	2.75	0.44
1:D:483:TYR:N	1:D:484:PRO:HD3	2.33	0.44
1:G:66:PHE:CE2	1:G:206:PRO:HB3	2.52	0.44
1:G:142:PHE:O	1:G:146:ILE:HG13	2.16	0.44
1:A:66:PHE:CD2	1:A:206:PRO:HB3	2.52	0.44
1:B:45:GLY:O	1:B:48:SER:OG	2.35	0.44
1:D:444:SER:O	1:D:445:ARG:HD3	2.18	0.44
1:A:25:ARG:NH2	1:A:499:ASP:OD2	2.42	0.44
1:C:211[A]:ARG:HB3	1:C:211[A]:ARG:CZ	2.47	0.44
1:D:345:LYS:HA	1:D:363:LYS:O	2.17	0.44
1:E:92:ILE:HG22	1:E:94:TYR:HE2	1.83	0.44
1:F:150:MET:CE	1:F:159:MET:HE3	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:69:TYR:CD2	1:H:463:MET:HG3	2.52	0.44
1:H:184:ALA:HB2	1:H:204:TRP:CD2	2.53	0.44
1:A:160:ARG:N	1:A:161:PRO:HD2	2.32	0.44
1:D:163:ASN:HA	1:D:166:VAL:HG12	2.00	0.44
1:F:18:THR:OG1	1:F:461:SER:HB3	2.17	0.44
1:F:168:ALA:HB1	1:F:377:SER:OG	2.18	0.44
1:G:66:PHE:CD2	1:G:206:PRO:HB3	2.53	0.44
1:C:376:GLU:HB2	1:C:382:VAL:CG1	2.48	0.44
1:E:60:VAL:HG13	1:E:415:PHE:CE2	2.53	0.44
1:F:207:ASN:O	1:F:208:ALA:C	2.60	0.44
1:G:91:ARG:HG3	1:G:315:TRP:CH2	2.52	0.44
1:H:278:MET:HE1	1:H:437:LEU:CD2	2.48	0.44
1:A:384:GLN:O	1:A:387:ILE:HG22	2.18	0.44
1:D:82:LYS:HB2	1:D:85:ASP:OD1	2.18	0.44
1:F:299:GLY:HA3	1:F:411:TYR:HB3	2.00	0.44
1:G:311:GLY:HA3	5:G:984:HOH:O	2.17	0.44
1:B:159:MET:HE1	2:B:800:UDP:C2	2.53	0.44
1:C:453:TYR:C	1:C:455:VAL:N	2.73	0.44
1:D:122:ILE:HD11	1:D:195:VAL:HG21	1.99	0.44
1:H:109:ASN:HA	5:H:1046:HOH:O	2.17	0.44
1:H:294:SER:HB2	1:H:376:GLU:HB3	2.00	0.44
1:H:338:ASN:HB2	1:H:339:GLN:HE21	1.82	0.44
1:H:453:TYR:C	1:H:455:VAL:H	2.24	0.44
1:A:213:PRO:HG3	1:A:221:ILE:HD11	1.99	0.43
1:A:303:ARG:NH1	1:A:348:THR:OG1	2.50	0.43
1:D:299:GLY:HA3	1:D:411:TYR:HB3	1.99	0.43
1:E:332:SER:CB	1:E:369:SER:H	2.30	0.43
1:H:290:LEU:HD13	1:H:451:TRP:CD2	2.53	0.43
1:A:72:PHE:CD1	1:A:459:ASP:HA	2.53	0.43
1:A:207:ASN:O	1:A:208:ALA:C	2.61	0.43
1:C:346:LEU:HD13	1:C:408:VAL:HG11	2.01	0.43
1:E:25:ARG:HD2	1:E:466:VAL:HG13	2.01	0.43
1:E:122:ILE:O	1:E:126:ILE:HG13	2.17	0.43
1:G:184:ALA:HB2	1:G:204:TRP:CD2	2.54	0.43
1:A:37:VAL:HG12	1:A:235:ARG:HB3	2.01	0.43
1:A:453:TYR:C	1:A:455:VAL:H	2.25	0.43
1:B:367:TYR:HE1	1:B:408:VAL:HG21	1.82	0.43
1:H:31:GLY:N	5:H:516:HOH:O	2.34	0.43
1:H:207:ASN:O	1:H:208:ALA:C	2.61	0.43
1:H:303:ARG:NH1	1:H:348:THR:OG1	2.51	0.43
1:A:3:HIS:HA	1:A:4:PRO:HD3	1.88	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:602:FAD:H1'1	3:A:602:FAD:H9	1.41	0.43
1:D:346:LEU:HD13	1:D:408:VAL:CG1	2.48	0.43
1:A:453:TYR:C	1:A:455:VAL:N	2.75	0.43
1:B:14:GLY:O	1:B:19:GLY:HA3	2.17	0.43
1:E:339:GLN:N	1:E:339:GLN:CD	2.77	0.43
1:F:388:LEU:O	1:F:392:ILE:HG13	2.19	0.43
1:G:332:SER:HA	1:G:339:GLN:HE21	1.81	0.43
1:B:346:LEU:HD13	1:B:408:VAL:HG11	2.00	0.43
1:E:303:ARG:NH1	1:E:348:THR:OG1	2.51	0.43
1:H:91:ARG:NH1	1:H:207:ASN:HB2	2.33	0.43
1:A:178:TRP:CA	1:A:454:GLU:HG2	2.47	0.43
1:F:339:GLN:N	1:F:339:GLN:CD	2.77	0.43
1:H:162:TYR:HB2	1:H:319:PRO:HB3	1.99	0.43
1:B:453:TYR:C	1:B:455:VAL:N	2.75	0.43
1:C:3:HIS:HA	1:C:4:PRO:HD3	1.85	0.43
1:H:150:MET:HG2	1:H:186:PRO:HG3	2.00	0.43
1:B:184:ALA:HB2	1:B:204:TRP:CD2	2.54	0.43
1:D:249:ASN:O	1:D:251:THR:HG23	2.19	0.43
1:F:91:ARG:HD3	1:F:207:ASN:CB	2.48	0.43
1:G:142:PHE:HB3	1:G:171:THR:O	2.19	0.43
1:H:455:VAL:CG2	1:H:480:THR:HG23	2.48	0.43
1:F:69:TYR:CD2	1:F:463:MET:HG3	2.54	0.43
1:F:274:LEU:O	1:F:278:MET:HG3	2.18	0.43
1:H:293:SER:OG	1:H:419:TYR:HB2	2.19	0.43
1:A:69:TYR:CD2	1:A:463:MET:HG3	2.53	0.42
1:A:345:LYS:HA	1:A:363:LYS:O	2.19	0.42
1:C:159:MET:HE1	2:C:800:UDP:C2	2.53	0.42
1:G:98:GLN:HE22	1:G:114:PRO:HG2	1.83	0.42
1:G:213:PRO:HG3	1:G:218:THR:HA	2.00	0.42
1:G:432:GLN:O	1:G:435:PRO:HD2	2.18	0.42
1:H:104:TYR:O	1:H:202:GLY:HA2	2.18	0.42
1:B:98:GLN:HE21	1:B:114:PRO:HG2	1.83	0.42
1:B:160:ARG:N	1:B:161:PRO:HD2	2.34	0.42
1:B:207:ASN:O	1:B:208:ALA:C	2.62	0.42
1:D:56:PHE:CE1	1:D:339:GLN:HB3	2.54	0.42
1:D:96:ARG:HB2	1:D:101:TRP:CZ3	2.55	0.42
1:D:274:LEU:HD22	5:D:1014:HOH:O	2.18	0.42
1:G:165:LYS:NZ	1:G:318:PHE:O	2.50	0.42
1:C:309:ARG:NH1	1:C:400:MET:O	2.52	0.42
1:E:62:GLY:HA2	5:E:966:HOH:O	2.19	0.42
1:E:232:GLU:HG3	1:G:232:GLU:HG2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:432:GLN:O	1:E:435:PRO:HD2	2.19	0.42
1:A:178:TRP:HB2	1:A:454:GLU:CG	2.43	0.42
1:A:293:SER:OG	1:A:419:TYR:HB2	2.18	0.42
1:C:455:VAL:HG22	1:C:480:THR:HG23	2.01	0.42
1:D:132:ALA:C	1:D:134:VAL:H	2.27	0.42
1:E:175:GLN:O	1:E:178:TRP:CD1	2.73	0.42
1:F:82:LYS:HB2	1:F:85:ASP:OD1	2.19	0.42
1:H:462:PHE:O	1:H:466:VAL:HG23	2.19	0.42
1:B:91:ARG:HD3	1:B:207:ASN:HB2	2.01	0.42
1:G:384:GLN:O	1:G:387:ILE:HG22	2.20	0.42
1:B:178:TRP:CA	1:B:454:GLU:HG2	2.47	0.42
1:C:184:ALA:HB2	1:C:204:TRP:CD2	2.55	0.42
1:F:286:LEU:HB3	1:F:433:ILE:HG12	2.02	0.42
1:G:150:MET:CE	1:G:159:MET:HE3	2.44	0.42
1:H:175:GLN:O	1:H:178:TRP:CD1	2.72	0.42
1:D:26:LEU:HD13	1:D:34:TRP:HB2	2.02	0.42
1:D:72:PHE:CD1	1:D:459:ASP:HA	2.54	0.42
1:D:150:MET:HG2	1:D:186:PRO:HG3	2.00	0.42
1:E:69:TYR:CD1	1:E:463:MET:HG3	2.54	0.42
3:E:606:FAD:H1'1	3:E:606:FAD:H9	1.65	0.42
1:H:256:ASP:OD2	1:H:256:ASP:C	2.62	0.42
1:C:431:THR:O	1:C:435:PRO:HG2	2.20	0.42
1:E:66[A]:PHE:CD2	1:E:206:PRO:HB3	2.55	0.42
1:D:286:LEU:HD12	1:D:286:LEU:HA	1.88	0.42
1:D:508:LYS:HD2	5:D:541:HOH:O	2.19	0.42
1:E:232:GLU:CG	1:G:232:GLU:HG3	2.49	0.42
1:F:91:ARG:NH1	1:F:207:ASN:HB2	2.35	0.42
1:F:376:GLU:HB2	1:F:382:VAL:HG11	2.02	0.42
3:H:605:FAD:HM71	3:H:605:FAD:HM83	1.84	0.42
1:A:272:ASP:OD1	1:A:272:ASP:N	2.53	0.41
1:A:383:ASN:OD1	1:A:385:GLU:HG2	2.19	0.41
1:B:2:THR:HG23	1:B:2:THR:O	2.20	0.41
1:C:98:GLN:HE21	1:C:114:PRO:HG2	1.85	0.41
1:C:249:ASN:O	1:C:251:THR:HG23	2.20	0.41
1:E:146:ILE:HD13	1:E:159:MET:HB3	2.01	0.41
1:F:284:VAL:O	1:F:288:LYS:HG3	2.19	0.41
1:G:242:VAL:HG13	1:G:252:VAL:CG1	2.50	0.41
1:A:66:PHE:CE2	1:A:206:PRO:HB3	2.56	0.41
1:D:178:TRP:CB	1:D:454:GLU:HG3	2.43	0.41
1:D:286:LEU:HB3	1:D:433:ILE:HG12	2.01	0.41
1:F:245:VAL:O	1:F:278:MET:HA	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:183:VAL:HG22	2:G:800:UDP:H1'	2.02	0.41
1:G:212:PHE:CG	1:G:213:PRO:HD2	2.55	0.41
1:H:213:PRO:HG3	1:H:218:THR:HA	2.03	0.41
1:H:325:PHE:HA	1:H:374:VAL:HG22	2.03	0.41
1:A:115:LYS:HA	1:A:118:GLN:OE1	2.19	0.41
1:A:339:GLN:CD	1:A:339:GLN:N	2.78	0.41
1:B:245:VAL:HG22	1:B:252:VAL:HG22	2.01	0.41
1:F:244:LYS:HB2	1:F:253:THR:HB	2.02	0.41
1:F:284:VAL:CG1	1:F:288:LYS:HE2	2.50	0.41
1:G:245:VAL:O	1:G:278:MET:HA	2.21	0.41
1:G:332:SER:CB	1:G:369:SER:H	2.33	0.41
1:C:112:MET:HE1	1:C:200:THR:HG22	2.01	0.41
1:C:345:LYS:NZ	5:C:528:HOH:O	2.52	0.41
1:F:213:PRO:HG3	1:F:218:THR:HA	2.01	0.41
1:G:434:LEU:HB2	1:G:435:PRO:HD3	2.01	0.41
1:H:115:LYS:HA	1:H:118:GLN:OE1	2.20	0.41
1:D:455:VAL:HG22	1:D:480:THR:HG23	2.03	0.41
1:E:425:GLU:N	1:E:425:GLU:OE2	2.51	0.41
1:F:55:GLY:O	1:F:340:PRO:HD3	2.20	0.41
1:G:14:GLY:O	1:G:19:GLY:HA3	2.19	0.41
1:H:249:ASN:O	1:H:251:THR:HG23	2.20	0.41
1:C:286:LEU:HD12	1:C:286:LEU:HA	1.92	0.41
1:G:98:GLN:HE21	1:G:114:PRO:HG2	1.83	0.41
1:H:25:ARG:NH1	1:H:470:ASP:OD1	2.47	0.41
1:C:293:SER:OG	1:C:419:TYR:HB2	2.20	0.41
1:D:3:HIS:HA	1:D:4:PRO:HD3	1.85	0.41
1:D:122:ILE:O	1:D:126:ILE:HG13	2.20	0.41
1:E:21:GLY:HA2	1:E:462:PHE:CE1	2.56	0.41
1:E:66[B]:PHE:CE2	1:E:207:ASN:OD1	2.74	0.41
1:F:3:HIS:HA	1:F:4:PRO:HD3	1.88	0.41
3:F:603:FAD:HM71	3:F:603:FAD:HM83	1.83	0.41
1:G:112:MET:HE3	1:G:200:THR:HG22	2.03	0.41
1:G:507:SER:O	1:G:511:LEU:HD13	2.20	0.41
1:B:162:TYR:HH	1:B:317:TYR:HD1	1.65	0.41
1:D:25:ARG:NH2	1:D:499:ASP:OD2	2.42	0.41
1:D:376:GLU:HB2	1:D:382:VAL:CG1	2.51	0.41
1:F:66:PHE:CD2	1:F:206:PRO:HB3	2.56	0.41
1:G:286:LEU:HD12	1:G:286:LEU:HA	1.84	0.41
1:G:392:ILE:HD13	1:G:407:ILE:CD1	2.51	0.41
3:G:607:FAD:HM83	3:G:607:FAD:HM71	1.84	0.41
1:H:111:SER:HB3	1:H:195:VAL:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:380:LYS:NZ	5:H:888:HOH:O	2.54	0.41
1:A:138:LYS:HA	1:A:139:PRO:HD3	1.95	0.41
1:A:168:ALA:HB1	1:A:377:SER:OG	2.21	0.41
1:A:274:LEU:O	1:A:278:MET:HG3	2.21	0.41
1:B:280:ASP:O	1:B:284:VAL:HG23	2.21	0.41
1:C:96:ARG:HB2	1:C:101:TRP:CZ3	2.55	0.41
1:C:327:ARG:NH1	2:C:800:UDP:O3B	2.53	0.41
1:C:384:GLN:O	1:C:387:ILE:HG22	2.21	0.41
1:C:477:VAL:HA	1:E:494:GLU:HG2	2.03	0.41
1:G:367:TYR:HE1	1:G:408:VAL:HG21	1.85	0.41
1:H:122:ILE:O	1:H:126:ILE:HG13	2.21	0.41
1:H:286:LEU:HD12	1:H:286:LEU:HA	1.83	0.41
1:A:79:ALA:HB1	1:A:224:ALA:CB	2.51	0.41
1:A:372:LEU:CD1	1:A:391:CYS:HB3	2.51	0.41
1:B:64:VAL:HG13	1:B:210:PHE:CG	2.55	0.41
1:C:63:HIS:CE1	3:C:600:FAD:C9A	3.04	0.41
1:C:483:TYR:N	1:C:484:PRO:HD3	2.36	0.41
1:F:332:SER:CB	1:F:369:SER:H	2.30	0.41
1:B:70:LYS:HB2	1:B:493:THR:HA	2.01	0.40
1:F:76:LEU:HD22	1:F:221:ILE:CG2	2.51	0.40
1:D:191:VAL:O	1:D:195:VAL:HG23	2.21	0.40
1:D:372:LEU:CD1	1:D:391:CYS:HB3	2.50	0.40
1:G:142:PHE:HD2	1:G:174:MET:HE3	1.85	0.40
1:G:272:ASP:OD1	1:G:273:PHE:N	2.54	0.40
1:B:455:VAL:HG22	1:B:480:THR:HG23	2.02	0.40
1:E:284:VAL:HG12	1:E:288:LYS:HE3	2.01	0.40
1:G:114:PRO:O	1:G:118:GLN:HG3	2.21	0.40
1:H:298:ILE:HG13	1:H:391:CYS:SG	2.61	0.40
1:B:21:GLY:HA2	1:B:462:PHE:CE1	2.57	0.40
1:B:60:VAL:HG12	1:B:60:VAL:O	2.22	0.40
1:D:14:GLY:O	1:D:19:GLY:HA3	2.21	0.40
1:F:104:TYR:O	1:F:202:GLY:HA2	2.21	0.40
1:G:18:THR:OG1	1:G:461:SER:HB3	2.20	0.40
1:G:287:THR:HG22	1:G:433:ILE:HD13	2.02	0.40
1:H:339:GLN:CD	1:H:339:GLN:N	2.79	0.40
1:C:95:VAL:HG22	1:C:317:TYR:HB2	2.04	0.40
1:E:242:VAL:HG13	1:E:252:VAL:HG13	2.03	0.40
1:E:434:LEU:HB2	1:E:435:PRO:HD3	2.03	0.40
1:F:91:ARG:HG3	1:F:315:TRP:CZ2	2.56	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	508/510 (100%)	491 (97%)	17 (3%)	0	100	100
1	B	509/510 (100%)	491 (96%)	18 (4%)	0	100	100
1	C	509/510 (100%)	490 (96%)	19 (4%)	0	100	100
1	D	508/510 (100%)	488 (96%)	20 (4%)	0	100	100
1	E	509/510 (100%)	494 (97%)	15 (3%)	0	100	100
1	F	508/510 (100%)	490 (96%)	17 (3%)	1 (0%)	43	58
1	G	508/510 (100%)	487 (96%)	21 (4%)	0	100	100
1	H	509/510 (100%)	489 (96%)	19 (4%)	1 (0%)	43	58
All	All	4068/4080 (100%)	3920 (96%)	146 (4%)	2 (0%)	100	100

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	208	ALA
1	H	208	ALA

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	432/432 (100%)	415 (96%)	17 (4%)	28	47
1	B	433/432 (100%)	414 (96%)	19 (4%)	25	41
1	C	433/432 (100%)	413 (95%)	20 (5%)	24	39

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	432/432 (100%)	412 (95%)	20 (5%)	24	39
1	E	433/432 (100%)	415 (96%)	18 (4%)	26	44
1	F	432/432 (100%)	412 (95%)	20 (5%)	24	39
1	G	432/432 (100%)	411 (95%)	21 (5%)	22	37
1	H	433/432 (100%)	409 (94%)	24 (6%)	19	32
All	All	3460/3456 (100%)	3301 (95%)	159 (5%)	24	39

All (159) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	GLN
1	A	111	SER
1	A	134	VAL
1	A	149	MET
1	A	175	GLN
1	A	209	THR
1	A	231	LYS
1	A	286	LEU
1	A	302	VAL
1	A	308	GLU
1	A	309	ARG
1	A	339	GLN
1	A	382	VAL
1	A	454	GLU
1	A	455	VAL
1	A	477	VAL
1	A	510	GLN
1	B	25	ARG
1	B	48	SER
1	B	90	GLN
1	B	134	VAL
1	B	149	MET
1	B	175	GLN
1	B	209	THR
1	B	286	LEU
1	B	302	VAL
1	B	308	GLU
1	B	309	ARG
1	B	310	ILE
1	B	339	GLN

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Mol	Chain	Res	Type
1	B	382	VAL
1	B	455	VAL
1	B	470	ASP
1	B	477	VAL
1	B	510	GLN
1	B	511	LEU
1	C	25	ARG
1	C	90	GLN
1	C	134	VAL
1	C	149	MET
1	C	175	GLN
1	C	209	THR
1	C	211[A]	ARG
1	C	211[B]	ARG
1	C	231	LYS
1	C	279	ASN
1	C	286	LEU
1	C	302	VAL
1	C	308	GLU
1	C	309	ARG
1	C	339	GLN
1	C	358	GLN
1	C	382	VAL
1	C	455	VAL
1	C	477	VAL
1	C	510	GLN
1	D	76	LEU
1	D	90	GLN
1	D	134	VAL
1	D	149	MET
1	D	175	GLN
1	D	176	CYS
1	D	209	THR
1	D	231	LYS
1	D	279	ASN
1	D	286	LEU
1	D	288	LYS
1	D	302	VAL
1	D	308	GLU
1	D	309	ARG
1	D	339	GLN
1	D	358	GLN

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Mol	Chain	Res	Type
1	D	382	VAL
1	D	455	VAL
1	D	477	VAL
1	D	506	LYS
1	E	25	ARG
1	E	90	GLN
1	E	112	MET
1	E	134	VAL
1	E	149	MET
1	E	175	GLN
1	E	209	THR
1	E	231	LYS
1	E	286	LEU
1	E	302	VAL
1	E	308	GLU
1	E	309	ARG
1	E	339	GLN
1	E	356	ARG
1	E	382	VAL
1	E	477	VAL
1	E	510	GLN
1	E	511	LEU
1	F	25	ARG
1	F	90	GLN
1	F	134	VAL
1	F	149	MET
1	F	175	GLN
1	F	176	CYS
1	F	209	THR
1	F	231	LYS
1	F	279	ASN
1	F	286	LEU
1	F	302	VAL
1	F	309	ARG
1	F	339	GLN
1	F	356	ARG
1	F	358	GLN
1	F	382	VAL
1	F	455	VAL
1	F	477	VAL
1	F	510	GLN
1	F	511	LEU

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Mol	Chain	Res	Type
1	G	25	ARG
1	G	90	GLN
1	G	134	VAL
1	G	149	MET
1	G	175	GLN
1	G	176	CYS
1	G	179	LEU
1	G	209	THR
1	G	231	LYS
1	G	286	LEU
1	G	302	VAL
1	G	309	ARG
1	G	339	GLN
1	G	356	ARG
1	G	358	GLN
1	G	382	VAL
1	G	393	GLN
1	G	455	VAL
1	G	477	VAL
1	G	506	LYS
1	G	510	GLN
1	H	25	ARG
1	H	48	SER
1	H	76	LEU
1	H	90	GLN
1	H	134	VAL
1	H	149	MET
1	H	175	GLN
1	H	176	CYS
1	H	209	THR
1	H	231	LYS
1	H	244	LYS
1	H	252	VAL
1	H	286	LEU
1	H	302	VAL
1	H	308	GLU
1	H	309	ARG
1	H	339	GLN
1	H	358	GLN
1	H	382	VAL
1	H	393	GLN
1	H	450	SER

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Mol	Chain	Res	Type
1	H	455	VAL
1	H	477	VAL
1	H	510	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	98	GLN
1	A	296	HIS
1	A	339	GLN
1	A	393	GLN
1	A	412	HIS
1	A	458	GLN
1	B	98	GLN
1	B	296	HIS
1	B	339	GLN
1	B	393	GLN
1	B	412	HIS
1	C	98	GLN
1	C	249	ASN
1	C	296	HIS
1	C	339	GLN
1	C	412	HIS
1	C	458	GLN
1	D	27	ASN
1	D	98	GLN
1	D	246	ASN
1	D	339	GLN
1	D	393	GLN
1	D	458	GLN
1	E	98	GLN
1	E	175	GLN
1	E	296	HIS
1	E	339	GLN
1	E	393	GLN
1	E	412	HIS
1	E	458	GLN
1	E	482	ASN
1	F	98	GLN
1	F	296	HIS
1	F	339	GLN

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Mol	Chain	Res	Type
1	F	393	GLN
1	F	412	HIS
1	G	98	GLN
1	G	296	HIS
1	G	339	GLN
1	G	412	HIS
1	G	458	GLN
1	G	482	ASN
1	H	27	ASN
1	H	98	GLN
1	H	339	GLN
1	H	393	GLN
1	H	458	GLN
1	H	482	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 24 ligands modelled in this entry, 8 are monoatomic - leaving 16 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	FAD	G	607	-	58,58,58	3.30	17 (29%)	85,89,89	1.91	23 (27%)
2	UDP	G	800	-	25,26,26	1.00	0	38,40,40	1.51	5 (13%)
3	FAD	F	603	-	58,58,58	3.29	18 (31%)	85,89,89	1.93	24 (28%)
2	UDP	F	800	-	25,26,26	0.97	0	38,40,40	1.50	5 (13%)
2	UDP	H	800	-	25,26,26	0.99	1 (4%)	38,40,40	1.56	6 (15%)
3	FAD	B	604	-	58,58,58	3.25	17 (29%)	85,89,89	1.94	22 (25%)
2	UDP	C	800	-	25,26,26	0.98	0	38,40,40	1.58	7 (18%)
3	FAD	A	602	-	58,58,58	3.28	16 (27%)	85,89,89	1.98	26 (30%)
3	FAD	C	600	-	58,58,58	3.31	17 (29%)	85,89,89	1.90	24 (28%)
2	UDP	D	800	-	25,26,26	1.03	1 (4%)	38,40,40	1.55	6 (15%)
3	FAD	D	601	-	58,58,58	3.29	16 (27%)	85,89,89	1.96	25 (29%)
3	FAD	E	606	-	58,58,58	3.28	17 (29%)	85,89,89	1.95	23 (27%)
2	UDP	B	800	-	25,26,26	1.00	1 (4%)	38,40,40	1.50	5 (13%)
2	UDP	A	800	-	25,26,26	0.98	0	38,40,40	1.60	6 (15%)
2	UDP	E	800	-	25,26,26	0.99	0	38,40,40	1.48	5 (13%)
3	FAD	H	605	-	58,58,58	3.29	16 (27%)	85,89,89	2.02	24 (28%)

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	G	607	-	-	11/34/50/50	0/6/6/6
2	UDP	G	800	-	-	5/16/32/32	0/2/2/2
3	FAD	F	603	-	-	11/34/50/50	0/6/6/6
2	UDP	F	800	-	-	4/16/32/32	0/2/2/2
2	UDP	H	800	-	-	3/16/32/32	0/2/2/2
3	FAD	B	604	-	-	9/34/50/50	0/6/6/6
2	UDP	C	800	-	-	3/16/32/32	0/2/2/2
3	FAD	A	602	-	-	9/34/50/50	0/6/6/6
3	FAD	C	600	-	-	7/34/50/50	0/6/6/6
2	UDP	D	800	-	-	4/16/32/32	0/2/2/2
3	FAD	D	601	-	-	14/34/50/50	0/6/6/6
3	FAD	E	606	-	-	8/34/50/50	0/6/6/6
2	UDP	B	800	-	-	4/16/32/32	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	UDP	A	800	-	-	4/16/32/32	0/2/2/2
2	UDP	E	800	-	-	6/16/32/32	0/2/2/2
3	FAD	H	605	-	-	10/34/50/50	0/6/6/6

All (137) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	603	FAD	PA-O3P	-12.79	1.45	1.59
3	C	600	FAD	PA-O3P	-12.77	1.45	1.59
3	E	606	FAD	PA-O3P	-12.71	1.45	1.59
3	B	604	FAD	PA-O3P	-12.69	1.45	1.59
3	D	601	FAD	PA-O3P	-12.52	1.46	1.59
3	G	607	FAD	PA-O3P	-12.42	1.46	1.59
3	H	605	FAD	PA-O3P	-12.20	1.46	1.59
3	A	602	FAD	PA-O3P	-11.85	1.46	1.59
3	D	601	FAD	O4-C4	10.61	1.43	1.23
3	A	602	FAD	O4-C4	10.53	1.43	1.23
3	G	607	FAD	O4-C4	10.50	1.43	1.23
3	F	603	FAD	O4-C4	10.49	1.43	1.23
3	H	605	FAD	O4-C4	10.45	1.43	1.23
3	E	606	FAD	O4-C4	10.44	1.43	1.23
3	C	600	FAD	O4-C4	10.42	1.43	1.23
3	B	604	FAD	O4-C4	10.42	1.43	1.23
3	A	602	FAD	P-O3P	9.37	1.69	1.59
3	H	605	FAD	P-O3P	9.22	1.69	1.59
3	G	607	FAD	P-O3P	8.96	1.69	1.59
3	C	600	FAD	P-O3P	8.72	1.68	1.59
3	D	601	FAD	P-O3P	8.66	1.68	1.59
3	F	603	FAD	P-O3P	8.57	1.68	1.59
3	E	606	FAD	P-O3P	8.45	1.68	1.59
3	G	607	FAD	O2'-C2'	-8.44	1.25	1.43
3	D	601	FAD	O2'-C2'	-8.32	1.25	1.43
3	C	600	FAD	O2'-C2'	-8.30	1.26	1.43
3	E	606	FAD	O2'-C2'	-8.26	1.26	1.43
3	B	604	FAD	O2'-C2'	-8.19	1.26	1.43
3	F	603	FAD	O2'-C2'	-8.18	1.26	1.43
3	H	605	FAD	O2'-C2'	-8.16	1.26	1.43
3	A	602	FAD	O2'-C2'	-8.13	1.26	1.43
3	B	604	FAD	P-O3P	8.05	1.68	1.59
3	G	607	FAD	O2-C2	7.21	1.38	1.24
3	F	603	FAD	O2-C2	7.20	1.38	1.24
3	B	604	FAD	O2-C2	7.17	1.38	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	605	FAD	O2-C2	7.12	1.38	1.24
3	D	601	FAD	O2-C2	7.11	1.38	1.24
3	C	600	FAD	O2-C2	7.10	1.38	1.24
3	E	606	FAD	O2-C2	7.10	1.38	1.24
3	A	602	FAD	O2-C2	7.00	1.38	1.24
3	H	605	FAD	C2B-C3B	-4.85	1.40	1.53
3	B	604	FAD	C2B-C3B	-4.81	1.40	1.53
3	C	600	FAD	C1'-C2'	-4.78	1.46	1.52
3	A	602	FAD	C2B-C3B	-4.78	1.40	1.53
3	G	607	FAD	C2B-C3B	-4.70	1.40	1.53
3	E	606	FAD	C2B-C3B	-4.66	1.40	1.53
3	C	600	FAD	C2B-C3B	-4.65	1.40	1.53
3	G	607	FAD	C1'-C2'	-4.62	1.46	1.52
3	E	606	FAD	C1'-C2'	-4.61	1.46	1.52
3	F	603	FAD	C1'-C2'	-4.60	1.46	1.52
3	F	603	FAD	C2B-C3B	-4.60	1.40	1.53
3	H	605	FAD	C1'-C2'	-4.58	1.46	1.52
3	D	601	FAD	C2B-C3B	-4.55	1.41	1.53
3	D	601	FAD	C1'-C2'	-4.44	1.46	1.52
3	B	604	FAD	C1'-C2'	-4.40	1.46	1.52
3	A	602	FAD	C1'-C2'	-4.28	1.46	1.52
3	A	602	FAD	C5'-C4'	-4.09	1.46	1.51
3	H	605	FAD	C5'-C4'	-4.04	1.46	1.51
3	A	602	FAD	C9A-N10	-3.98	1.34	1.41
3	D	601	FAD	C5'-C4'	-3.85	1.46	1.51
3	C	600	FAD	C5'-C4'	-3.76	1.46	1.51
3	F	603	FAD	C5'-C4'	-3.70	1.46	1.51
3	E	606	FAD	C5'-C4'	-3.59	1.46	1.51
3	G	607	FAD	C5'-C4'	-3.58	1.46	1.51
3	E	606	FAD	C9A-N10	-3.56	1.35	1.41
3	B	604	FAD	C9A-N10	-3.52	1.35	1.41
3	F	603	FAD	C9A-N10	-3.50	1.35	1.41
3	C	600	FAD	C9A-N10	-3.49	1.35	1.41
3	H	605	FAD	C9A-N10	-3.44	1.35	1.41
3	G	607	FAD	C9A-N10	-3.37	1.35	1.41
3	B	604	FAD	C5'-C4'	-3.36	1.47	1.51
3	D	601	FAD	C9A-N10	-3.22	1.35	1.41
3	E	606	FAD	C6A-N6A	3.02	1.41	1.34
3	H	605	FAD	C6A-N6A	3.00	1.41	1.34
3	B	604	FAD	C6A-N6A	2.97	1.41	1.34
3	G	607	FAD	C6A-N6A	2.96	1.41	1.34
3	A	602	FAD	PA-O1A	-2.95	1.40	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	600	FAD	PA-O1A	-2.90	1.40	1.50
3	C	600	FAD	C6A-N6A	2.90	1.41	1.34
3	A	602	FAD	C6A-N6A	2.88	1.41	1.34
3	E	606	FAD	PA-O1A	-2.87	1.40	1.50
3	D	601	FAD	C6A-N6A	2.86	1.41	1.34
3	D	601	FAD	PA-O1A	-2.85	1.40	1.50
3	F	603	FAD	C6A-N6A	2.83	1.41	1.34
3	F	603	FAD	PA-O1A	-2.77	1.41	1.50
3	B	604	FAD	PA-O1A	-2.77	1.41	1.50
3	H	605	FAD	PA-O1A	-2.74	1.41	1.50
3	D	601	FAD	C7M-C7	2.71	1.56	1.51
3	G	607	FAD	PA-O1A	-2.68	1.41	1.50
3	C	600	FAD	C7M-C7	2.67	1.56	1.51
3	F	603	FAD	C7M-C7	2.64	1.55	1.51
3	B	604	FAD	C7M-C7	2.58	1.55	1.51
3	A	602	FAD	C7M-C7	2.57	1.55	1.51
3	H	605	FAD	C7M-C7	2.55	1.55	1.51
3	G	607	FAD	C7M-C7	2.55	1.55	1.51
3	A	602	FAD	C3B-C4B	-2.53	1.46	1.53
3	C	600	FAD	C3B-C4B	-2.47	1.46	1.53
3	D	601	FAD	C3B-C4B	-2.47	1.46	1.53
3	E	606	FAD	C7M-C7	2.47	1.55	1.51
3	B	604	FAD	C4-N3	-2.44	1.34	1.38
3	E	606	FAD	O4B-C4B	-2.37	1.39	1.45
3	E	606	FAD	C3B-C4B	-2.36	1.47	1.53
3	F	603	FAD	O4B-C4B	-2.36	1.39	1.45
3	C	600	FAD	O4B-C4B	-2.35	1.39	1.45
3	C	600	FAD	C2'-C3'	-2.34	1.49	1.53
3	D	601	FAD	O4B-C4B	-2.34	1.39	1.45
3	B	604	FAD	C3B-C4B	-2.34	1.47	1.53
3	G	607	FAD	O4B-C4B	-2.34	1.39	1.45
2	D	800	UDP	PA-O3A	2.33	1.62	1.59
3	F	603	FAD	C3B-C4B	-2.32	1.47	1.53
3	H	605	FAD	C3B-C4B	-2.30	1.47	1.53
3	A	602	FAD	O4B-C4B	-2.29	1.39	1.45
3	E	606	FAD	C2'-C3'	-2.29	1.49	1.53
3	A	602	FAD	C4-N3	-2.27	1.34	1.38
3	B	604	FAD	O4B-C4B	-2.26	1.40	1.45
3	F	603	FAD	C4-N3	-2.26	1.34	1.38
3	D	601	FAD	C4-N3	-2.25	1.34	1.38
3	H	605	FAD	C2'-C3'	-2.25	1.49	1.53
3	H	605	FAD	O4B-C4B	-2.24	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	606	FAD	C5A-N7A	-2.22	1.35	1.39
3	F	603	FAD	C5A-N7A	-2.20	1.35	1.39
3	G	607	FAD	C3B-C4B	-2.19	1.47	1.53
3	D	601	FAD	C2'-C3'	-2.15	1.49	1.53
3	G	607	FAD	C2'-C3'	-2.14	1.49	1.53
3	C	600	FAD	C4-N3	-2.13	1.34	1.38
3	A	602	FAD	C2'-C3'	-2.12	1.49	1.53
3	E	606	FAD	C4-N3	-2.12	1.34	1.38
3	G	607	FAD	C5A-N7A	-2.12	1.35	1.39
3	F	603	FAD	C2'-C3'	-2.11	1.49	1.53
3	C	600	FAD	C5A-N7A	-2.11	1.35	1.39
3	G	607	FAD	C4-N3	-2.10	1.34	1.38
3	H	605	FAD	C5A-N7A	-2.08	1.35	1.39
2	B	800	UDP	PA-O3A	2.01	1.61	1.59
3	F	603	FAD	C8A-N9A	-2.01	1.34	1.37
3	B	604	FAD	C5A-N7A	-2.01	1.35	1.39
3	B	604	FAD	C9A-C5X	-2.01	1.38	1.41
2	H	800	UDP	C6-C5	2.00	1.39	1.35

All (236) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	800	UDP	C4-N3-C2	-5.70	119.54	126.61
2	C	800	UDP	C4-N3-C2	-5.50	119.78	126.61
2	H	800	UDP	C4-N3-C2	-5.31	120.02	126.61
2	D	800	UDP	C4-N3-C2	-5.27	120.06	126.61
3	H	605	FAD	O4B-C1B-N9A	5.25	118.17	108.09
3	B	604	FAD	O4B-C1B-N9A	5.23	118.13	108.09
2	F	800	UDP	C4-N3-C2	-5.20	120.16	126.61
2	G	800	UDP	C4-N3-C2	-5.15	120.22	126.61
3	E	606	FAD	C5A-C4A-N3A	-5.13	119.66	126.72
3	D	601	FAD	C5A-C4A-N3A	-5.04	119.78	126.72
2	E	800	UDP	C4-N3-C2	-5.03	120.37	126.61
3	A	602	FAD	C5A-C4A-N3A	-5.02	119.80	126.72
2	B	800	UDP	C4-N3-C2	-4.95	120.47	126.61
3	F	603	FAD	C5A-C4A-N3A	-4.93	119.93	126.72
3	B	604	FAD	C5A-C4A-N3A	-4.91	119.95	126.72
3	A	602	FAD	O4B-C1B-N9A	4.87	117.44	108.09
3	C	600	FAD	C5A-C4A-N3A	-4.81	120.10	126.72
3	H	605	FAD	C5A-C4A-N3A	-4.74	120.19	126.72
3	G	607	FAD	C5A-C4A-N3A	-4.71	120.23	126.72
3	H	605	FAD	C4-C4X-N5	4.67	124.66	118.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	607	FAD	N9A-C8A-N7A	-4.51	107.53	113.94
3	C	600	FAD	N9A-C8A-N7A	-4.46	107.61	113.94
3	A	602	FAD	N3A-C2A-N1A	-4.41	121.91	128.58
3	F	603	FAD	N9A-C8A-N7A	-4.39	107.71	113.94
3	C	600	FAD	C4A-N9A-C8A	4.32	110.27	105.74
3	C	600	FAD	N3A-C2A-N1A	-4.30	122.08	128.58
3	H	605	FAD	N3A-C2A-N1A	-4.28	122.11	128.58
3	E	606	FAD	O4B-C1B-N9A	4.28	116.30	108.09
3	B	604	FAD	N3A-C2A-N1A	-4.27	122.11	128.58
2	A	800	UDP	N3-C2-N1	4.23	120.40	114.89
3	A	602	FAD	N9A-C8A-N7A	-4.23	107.93	113.94
3	D	601	FAD	N3A-C2A-N1A	-4.22	122.19	128.58
3	D	601	FAD	N9A-C8A-N7A	-4.22	107.95	113.94
3	E	606	FAD	N9A-C8A-N7A	-4.21	107.96	113.94
3	H	605	FAD	N9A-C8A-N7A	-4.21	107.97	113.94
3	E	606	FAD	C2'-C1'-N10	4.21	130.08	110.20
2	E	800	UDP	N3-C2-N1	4.20	120.35	114.89
3	E	606	FAD	N3A-C4A-N9A	4.19	134.30	127.17
3	C	600	FAD	C2'-C1'-N10	4.16	129.85	110.20
3	F	603	FAD	N3A-C2A-N1A	-4.15	122.30	128.58
3	E	606	FAD	N3A-C2A-N1A	-4.15	122.30	128.58
2	H	800	UDP	N3-C2-N1	4.14	120.28	114.89
3	D	601	FAD	O4B-C1B-N9A	4.09	115.95	108.09
3	G	607	FAD	O4B-C1B-N9A	4.08	115.93	108.09
2	C	800	UDP	N3-C2-N1	4.07	120.19	114.89
2	D	800	UDP	N3-C2-N1	4.06	120.18	114.89
3	G	607	FAD	C4A-N9A-C8A	4.05	109.99	105.74
3	C	600	FAD	N3A-C4A-N9A	4.03	134.01	127.17
3	G	607	FAD	N3A-C2A-N1A	-4.03	122.49	128.58
3	D	601	FAD	N3A-C4A-N9A	4.01	133.98	127.17
3	G	607	FAD	C5A-N7A-C8A	4.00	109.74	103.45
2	G	800	UDP	N3-C2-N1	4.00	120.10	114.89
3	B	604	FAD	C2'-C1'-N10	4.00	129.09	110.20
2	F	800	UDP	N3-C2-N1	3.97	120.06	114.89
3	B	604	FAD	N9A-C8A-N7A	-3.96	108.31	113.94
3	H	605	FAD	O2'-C2'-C3'	3.94	118.48	109.25
3	F	603	FAD	C4A-N9A-C8A	3.93	109.87	105.74
3	D	601	FAD	C4A-N9A-C8A	3.93	109.86	105.74
3	F	603	FAD	C5A-N7A-C8A	3.92	109.60	103.45
3	A	602	FAD	N3A-C4A-N9A	3.91	133.82	127.17
3	H	605	FAD	C4A-N9A-C8A	3.90	109.84	105.74
3	C	600	FAD	O4B-C1B-N9A	3.87	115.53	108.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	605	FAD	C2'-C1'-N10	3.87	128.49	110.20
2	B	800	UDP	N3-C2-N1	3.86	119.91	114.89
3	H	605	FAD	N3A-C4A-N9A	3.86	133.73	127.17
3	F	603	FAD	N3A-C4A-N9A	3.85	133.72	127.17
3	E	606	FAD	C4A-N9A-C8A	3.84	109.78	105.74
3	C	600	FAD	C4-C4X-N5	3.82	123.49	118.21
3	D	601	FAD	C2'-C1'-N10	3.82	128.23	110.20
3	A	602	FAD	C4A-N9A-C8A	3.81	109.74	105.74
3	A	602	FAD	C5A-N7A-C8A	3.79	109.41	103.45
3	E	606	FAD	C5A-N7A-C8A	3.77	109.37	103.45
3	C	600	FAD	C5A-N7A-C8A	3.76	109.36	103.45
3	G	607	FAD	C4-C4X-N5	3.75	123.38	118.21
3	G	607	FAD	C4A-C5A-N7A	-3.71	106.34	110.58
3	B	604	FAD	N3A-C4A-N9A	3.68	133.43	127.17
3	F	603	FAD	O2'-C2'-C3'	3.67	117.84	109.25
3	D	601	FAD	C5A-N7A-C8A	3.67	109.22	103.45
3	G	607	FAD	N3A-C4A-N9A	3.65	133.37	127.17
3	F	603	FAD	C2'-C1'-N10	3.64	127.42	110.20
3	H	605	FAD	C5A-N7A-C8A	3.62	109.14	103.45
3	F	603	FAD	C4A-C5A-N7A	-3.60	106.47	110.58
3	E	606	FAD	O5'-C5'-C4'	3.60	118.97	109.36
3	A	602	FAD	C2'-C1'-N10	3.60	127.20	110.20
3	B	604	FAD	C5A-N7A-C8A	3.58	109.08	103.45
3	F	603	FAD	C4-C4X-N5	3.57	123.14	118.21
3	A	602	FAD	C4A-C5A-N7A	-3.54	106.53	110.58
3	G	607	FAD	C2'-C1'-N10	3.53	126.88	110.20
3	D	601	FAD	C4-C4X-N5	3.50	123.05	118.21
3	B	604	FAD	C4A-C5A-N7A	-3.50	106.58	110.58
3	B	604	FAD	C4A-N9A-C8A	3.49	109.41	105.74
3	B	604	FAD	C4-N3-C2	-3.49	119.45	125.64
3	E	606	FAD	C4-N3-C2	-3.45	119.52	125.64
3	A	602	FAD	C4-N3-C2	-3.44	119.53	125.64
3	D	601	FAD	O5'-C5'-C4'	3.43	118.52	109.36
3	A	602	FAD	C2A-N3A-C4A	3.43	120.20	111.83
3	G	607	FAD	O2'-C2'-C3'	3.39	117.19	109.25
3	D	601	FAD	C4A-C5A-N7A	-3.36	106.74	110.58
3	F	603	FAD	O4B-C1B-N9A	3.35	114.52	108.09
3	H	605	FAD	C4-N3-C2	-3.34	119.72	125.64
3	F	603	FAD	C4-N3-C2	-3.33	119.72	125.64
3	D	601	FAD	C4-N3-C2	-3.33	119.74	125.64
3	H	605	FAD	C4X-C10-N10	3.29	121.19	116.48
3	E	606	FAD	C4-C4X-N5	3.28	122.73	118.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	604	FAD	O5'-C5'-C4'	3.27	118.09	109.36
3	E	606	FAD	C2A-N3A-C4A	3.27	119.81	111.83
3	C	600	FAD	C4-N3-C2	-3.27	119.84	125.64
3	A	602	FAD	C4-C4X-N5	3.27	122.72	118.21
3	E	606	FAD	C4A-C5A-N7A	-3.26	106.85	110.58
3	G	607	FAD	C4-N3-C2	-3.25	119.88	125.64
3	F	603	FAD	C2A-N3A-C4A	3.24	119.76	111.83
3	C	600	FAD	C4A-C5A-N7A	-3.24	106.87	110.58
3	D	601	FAD	C2A-N3A-C4A	3.23	119.72	111.83
3	D	601	FAD	O2'-C2'-C3'	3.22	116.78	109.25
3	E	606	FAD	O4-C4-C4X	-3.22	118.04	126.53
3	B	604	FAD	C2A-N3A-C4A	3.21	119.67	111.83
3	H	605	FAD	O5'-C5'-C4'	3.18	117.84	109.36
3	H	605	FAD	C2A-N3A-C4A	3.16	119.56	111.83
3	H	605	FAD	C4A-C5A-N7A	-3.16	106.97	110.58
3	C	600	FAD	C2A-N3A-C4A	3.16	119.54	111.83
3	H	605	FAD	O5B-C5B-C4B	3.11	119.58	108.99
3	G	607	FAD	O5'-C5'-C4'	3.11	117.66	109.36
3	G	607	FAD	C2A-N3A-C4A	3.10	119.40	111.83
2	C	800	UDP	C5-C4-N3	3.08	119.12	114.80
3	A	602	FAD	O2'-C2'-C3'	3.07	116.43	109.25
3	B	604	FAD	C4A-N9A-C1B	-3.04	119.52	126.63
3	A	602	FAD	O5B-C5B-C4B	3.04	119.34	108.99
3	H	605	FAD	C4A-N9A-C1B	-3.03	119.54	126.63
2	G	800	UDP	C5-C4-N3	3.03	119.04	114.80
3	G	607	FAD	C4A-N9A-C1B	-3.03	119.55	126.63
2	H	800	UDP	C5-C4-N3	3.02	119.03	114.80
2	B	800	UDP	C5-C4-N3	3.01	119.02	114.80
3	E	606	FAD	O5B-C5B-C4B	3.01	119.25	108.99
2	F	800	UDP	C5-C4-N3	3.01	119.01	114.80
3	D	601	FAD	C4X-C10-N10	2.99	120.76	116.48
3	H	605	FAD	C4X-C4-N3	2.98	120.84	113.25
3	B	604	FAD	O4-C4-C4X	-2.98	118.67	126.53
3	E	606	FAD	C4X-C4-N3	2.97	120.81	113.25
2	D	800	UDP	C5-C4-N3	2.97	118.95	114.80
3	A	602	FAD	C4X-C4-N3	2.95	120.76	113.25
2	A	800	UDP	C5-C4-N3	2.94	118.91	114.80
3	C	600	FAD	O2'-C2'-C3'	2.93	116.12	109.25
3	F	603	FAD	C4X-C4-N3	2.93	120.72	113.25
2	E	800	UDP	C5-C4-N3	2.93	118.91	114.80
3	F	603	FAD	O4-C4-C4X	-2.93	118.80	126.53
3	G	607	FAD	C4X-C4-N3	2.92	120.69	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	607	FAD	O4-C4-C4X	-2.91	118.85	126.53
3	B	604	FAD	C4X-C4-N3	2.91	120.66	113.25
3	F	603	FAD	C4A-N9A-C1B	-2.90	119.86	126.63
3	C	600	FAD	C4A-N9A-C1B	-2.88	119.89	126.63
3	C	600	FAD	C4X-C4-N3	2.86	120.53	113.25
3	H	605	FAD	C10-C4X-N5	-2.86	118.97	124.81
3	H	605	FAD	O4-C4-C4X	-2.86	118.99	126.53
3	F	603	FAD	O5'-C5'-C4'	2.84	116.94	109.36
3	D	601	FAD	O5B-C5B-C4B	2.84	118.67	108.99
3	D	601	FAD	O4-C4-C4X	-2.83	119.07	126.53
3	A	602	FAD	C4A-N9A-C1B	-2.82	120.03	126.63
2	B	800	UDP	O2A-PA-O3A	2.82	114.90	107.27
3	D	601	FAD	C4X-C4-N3	2.80	120.38	113.25
3	C	600	FAD	O4-C4-C4X	-2.78	119.19	126.53
3	A	602	FAD	O4-C4-C4X	-2.75	119.26	126.53
3	A	602	FAD	C1'-N10-C9A	-2.75	115.28	120.63
3	G	607	FAD	C4X-C10-N10	2.74	120.41	116.48
3	B	604	FAD	O2'-C2'-C3'	2.73	115.64	109.25
3	D	601	FAD	C5'-C4'-C3'	-2.72	107.09	112.22
3	B	604	FAD	O5B-C5B-C4B	2.70	118.19	108.99
3	C	600	FAD	C4X-C10-N10	2.69	120.34	116.48
3	B	604	FAD	C1'-N10-C9A	-2.69	115.40	120.63
3	E	606	FAD	O2'-C2'-C3'	2.67	115.49	109.25
3	E	606	FAD	C3B-C2B-C1B	2.65	106.48	101.46
3	B	604	FAD	C4-C4X-N5	2.65	121.87	118.21
3	D	601	FAD	C4A-N9A-C1B	-2.65	120.44	126.63
2	C	800	UDP	O4-C4-C5	-2.64	120.61	125.16
2	G	800	UDP	O2A-PA-O3A	2.61	114.32	107.27
3	H	605	FAD	O2-C2-N1	-2.60	117.49	121.80
3	F	603	FAD	O5B-C5B-C4B	2.59	117.81	108.99
2	H	800	UDP	O2A-PA-O3A	2.58	114.25	107.27
3	C	600	FAD	C10-C4X-N5	-2.57	119.56	124.81
2	A	800	UDP	O4-C4-C5	-2.57	120.73	125.16
3	G	607	FAD	O5B-C5B-C4B	2.56	117.72	108.99
2	A	800	UDP	O2-C2-N1	-2.56	119.46	122.80
2	C	800	UDP	O2A-PA-O3A	2.56	114.19	107.27
2	D	800	UDP	O4-C4-C5	-2.54	120.78	125.16
3	F	603	FAD	C4X-C10-N10	2.54	120.11	116.48
2	D	800	UDP	O2A-PA-O3A	2.53	114.12	107.27
2	E	800	UDP	O4-C4-C5	-2.53	120.80	125.16
2	F	800	UDP	O4-C4-C5	-2.53	120.80	125.16
3	E	606	FAD	C4A-N9A-C1B	-2.52	120.73	126.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	601	FAD	C4'-C3'-C2'	2.52	117.77	113.57
3	H	605	FAD	O2'-C2'-C1'	2.50	120.47	110.20
2	G	800	UDP	O4-C4-C5	-2.50	120.86	125.16
3	A	602	FAD	O5'-C5'-C4'	2.48	115.99	109.36
3	F	603	FAD	C10-C4X-N5	-2.47	119.77	124.81
3	F	603	FAD	C3B-C2B-C1B	2.47	106.14	101.46
2	F	800	UDP	O2A-PA-O3A	2.47	113.94	107.27
3	G	607	FAD	C10-C4X-N5	-2.45	119.80	124.81
3	A	602	FAD	C10-C4X-N5	-2.44	119.83	124.81
3	F	603	FAD	C4'-C3'-C2'	2.44	117.63	113.57
2	A	800	UDP	O2A-PA-O3A	2.43	113.85	107.27
3	D	601	FAD	C10-C4X-N5	-2.43	119.84	124.81
2	H	800	UDP	O4-C4-C5	-2.41	121.00	125.16
3	E	606	FAD	C4X-C10-N10	2.41	119.93	116.48
3	A	602	FAD	C4X-C10-N10	2.36	119.86	116.48
3	D	601	FAD	C3B-C2B-C1B	2.35	105.92	101.46
2	B	800	UDP	O4-C4-C5	-2.34	121.13	125.16
3	C	600	FAD	O5'-C5'-C4'	2.31	115.53	109.36
3	E	606	FAD	C10-C4X-N5	-2.29	120.13	124.81
2	E	800	UDP	O2-C2-N1	-2.28	119.82	122.80
3	C	600	FAD	C3B-C2B-C1B	2.27	105.76	101.46
3	B	604	FAD	C10-C4X-N5	-2.25	120.21	124.81
2	H	800	UDP	O2-C2-N1	-2.21	119.92	122.80
3	F	603	FAD	O2-C2-N1	-2.20	118.15	121.80
3	E	606	FAD	O2-C2-N1	-2.19	118.16	121.80
3	C	600	FAD	O2'-C2'-C1'	2.18	119.15	110.20
3	D	601	FAD	O2'-C2'-C1'	2.17	119.13	110.20
3	H	605	FAD	C5'-C4'-C3'	-2.15	108.16	112.22
3	D	601	FAD	O2-C2-N1	-2.14	118.25	121.80
3	C	600	FAD	O5B-C5B-C4B	2.13	116.25	108.99
2	C	800	UDP	C6-N1-C2	-2.13	118.41	121.00
3	E	606	FAD	O2'-C2'-C1'	2.12	118.91	110.20
3	F	603	FAD	O2'-C2'-C1'	2.11	118.88	110.20
3	A	602	FAD	O2-C2-N1	-2.11	118.30	121.80
3	C	600	FAD	O2-C2-N1	-2.11	118.30	121.80
3	B	604	FAD	C4X-C10-N10	2.10	119.49	116.48
3	G	607	FAD	O2'-C2'-C1'	2.08	118.76	110.20
2	D	800	UDP	O2-C2-N1	-2.08	120.09	122.80
3	C	600	FAD	O4B-C4B-C3B	2.07	109.26	105.15
3	A	602	FAD	C3B-C2B-C1B	2.06	105.37	101.46
3	G	607	FAD	O2-C2-N1	-2.06	118.37	121.80
3	B	604	FAD	C4X-C10-N1	-2.06	119.54	124.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	607	FAD	O4'-C4'-C5'	-2.05	105.46	109.99
3	A	602	FAD	C9-C9A-N10	-2.05	119.10	121.85
3	A	602	FAD	C4X-C10-N1	-2.04	119.58	124.59
3	H	605	FAD	C4'-C3'-C2'	2.03	116.95	113.57
2	C	800	UDP	O2-C2-N1	-2.00	120.19	122.80
3	A	602	FAD	C4'-C3'-C2'	2.00	116.90	113.57

There are no chirality outliers.

All (112) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	800	UDP	C5'-O5'-PA-O1A
2	A	800	UDP	C5'-O5'-PA-O3A
2	A	800	UDP	PA-O3A-PB-O2B
2	A	800	UDP	PA-O3A-PB-O3B
2	B	800	UDP	O4'-C4'-C5'-O5'
2	B	800	UDP	PA-O3A-PB-O2B
2	B	800	UDP	PA-O3A-PB-O3B
2	C	800	UDP	PA-O3A-PB-O2B
2	C	800	UDP	PA-O3A-PB-O3B
2	E	800	UDP	C5'-O5'-PA-O2A
2	F	800	UDP	C5'-O5'-PA-O1A
2	F	800	UDP	PA-O3A-PB-O3B
2	G	800	UDP	C5'-O5'-PA-O1A
2	G	800	UDP	PA-O3A-PB-O3B
2	H	800	UDP	C5'-O5'-PA-O1A
2	H	800	UDP	PA-O3A-PB-O2B
2	H	800	UDP	PA-O3A-PB-O3B
3	A	602	FAD	C5B-O5B-PA-O1A
3	A	602	FAD	C5B-O5B-PA-O2A
3	A	602	FAD	C5B-O5B-PA-O3P
3	B	604	FAD	C5B-O5B-PA-O1A
3	B	604	FAD	C5B-O5B-PA-O2A
3	B	604	FAD	C5B-O5B-PA-O3P
3	C	600	FAD	C5B-O5B-PA-O1A
3	C	600	FAD	C5B-O5B-PA-O2A
3	C	600	FAD	C5B-O5B-PA-O3P
3	C	600	FAD	P-O3P-PA-O5B
3	D	601	FAD	C5B-O5B-PA-O1A
3	D	601	FAD	C5B-O5B-PA-O2A
3	D	601	FAD	C5B-O5B-PA-O3P
3	D	601	FAD	C5'-O5'-P-O2P

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Mol	Chain	Res	Type	Atoms
3	E	606	FAD	C5B-O5B-PA-O1A
3	E	606	FAD	C5B-O5B-PA-O2A
3	E	606	FAD	C5B-O5B-PA-O3P
3	F	603	FAD	C5B-O5B-PA-O1A
3	F	603	FAD	C5B-O5B-PA-O2A
3	F	603	FAD	C5B-O5B-PA-O3P
3	G	607	FAD	C5B-O5B-PA-O2A
3	G	607	FAD	P-O3P-PA-O5B
3	H	605	FAD	C5B-O5B-PA-O2A
3	H	605	FAD	C5B-O5B-PA-O3P
3	H	605	FAD	N10-C1'-C2'-O2'
3	B	604	FAD	O2'-C2'-C3'-C4'
3	G	607	FAD	O2'-C2'-C3'-C4'
2	B	800	UDP	C3'-C4'-C5'-O5'
2	E	800	UDP	O4'-C4'-C5'-O5'
3	C	600	FAD	O4B-C4B-C5B-O5B
3	C	600	FAD	C3B-C4B-C5B-O5B
3	B	604	FAD	O4B-C4B-C5B-O5B
3	D	601	FAD	O4B-C4B-C5B-O5B
3	F	603	FAD	O4B-C4B-C5B-O5B
3	H	605	FAD	O4B-C4B-C5B-O5B
3	F	603	FAD	O2'-C2'-C3'-C4'
2	D	800	UDP	C3'-C4'-C5'-O5'
3	A	602	FAD	C3B-C4B-C5B-O5B
3	H	605	FAD	C3B-C4B-C5B-O5B
3	B	604	FAD	O2'-C2'-C3'-O3'
3	A	602	FAD	O2'-C2'-C3'-C4'
3	G	607	FAD	O4'-C4'-C5'-O5'
3	A	602	FAD	O4B-C4B-C5B-O5B
3	B	604	FAD	C3B-C4B-C5B-O5B
3	D	601	FAD	C3B-C4B-C5B-O5B
3	E	606	FAD	O4B-C4B-C5B-O5B
3	F	603	FAD	C3B-C4B-C5B-O5B
2	D	800	UDP	O4'-C4'-C5'-O5'
3	E	606	FAD	C3B-C4B-C5B-O5B
3	G	607	FAD	O4B-C4B-C5B-O5B
3	G	607	FAD	C3'-C4'-C5'-O5'
3	G	607	FAD	O2'-C2'-C3'-O3'
2	E	800	UDP	C3'-C4'-C5'-O5'
3	G	607	FAD	C3B-C4B-C5B-O5B
3	A	602	FAD	P-O3P-PA-O5B
3	B	604	FAD	P-O3P-PA-O5B

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Mol	Chain	Res	Type	Atoms
3	D	601	FAD	P-O3P-PA-O5B
3	E	606	FAD	P-O3P-PA-O5B
3	F	603	FAD	P-O3P-PA-O5B
3	H	605	FAD	P-O3P-PA-O5B
3	D	601	FAD	C3'-C4'-C5'-O5'
3	E	606	FAD	PA-O3P-P-O1P
3	H	605	FAD	PA-O3P-P-O1P
3	D	601	FAD	O4'-C4'-C5'-O5'
3	A	602	FAD	O2'-C2'-C3'-O3'
3	F	603	FAD	O2'-C2'-C3'-O3'
3	D	601	FAD	N10-C1'-C2'-O2'
2	C	800	UDP	C5'-O5'-PA-O1A
2	E	800	UDP	C5'-O5'-PA-O1A
2	E	800	UDP	C5'-O5'-PA-O3A
2	G	800	UDP	C5'-O5'-PA-O3A
3	D	601	FAD	C5'-O5'-P-O1P
3	D	601	FAD	C5'-O5'-P-O3P
3	F	603	FAD	C5'-O5'-P-O2P
3	F	603	FAD	C5'-O5'-P-O3P
3	G	607	FAD	C5B-O5B-PA-O1A
3	G	607	FAD	C5B-O5B-PA-O3P
3	H	605	FAD	C5B-O5B-PA-O1A
3	C	600	FAD	O2'-C2'-C3'-C4'
3	F	603	FAD	PA-O3P-P-O1P
2	D	800	UDP	PB-O3A-PA-O1A
3	B	604	FAD	PA-O3P-P-O2P
3	H	605	FAD	C1'-C2'-C3'-O3'
2	F	800	UDP	PA-O3A-PB-O1B
2	G	800	UDP	PA-O3A-PB-O1B
2	D	800	UDP	PA-O3A-PB-O2B
2	E	800	UDP	PA-O3A-PB-O2B
2	F	800	UDP	PA-O3A-PB-O2B
2	G	800	UDP	PA-O3A-PB-O2B
3	D	601	FAD	PA-O3P-P-O1P
3	G	607	FAD	C2'-C3'-C4'-C5'
3	A	602	FAD	PA-O3P-P-O2P
3	D	601	FAD	PA-O3P-P-O2P
3	E	606	FAD	PA-O3P-P-O2P
3	H	605	FAD	PA-O3P-P-O2P

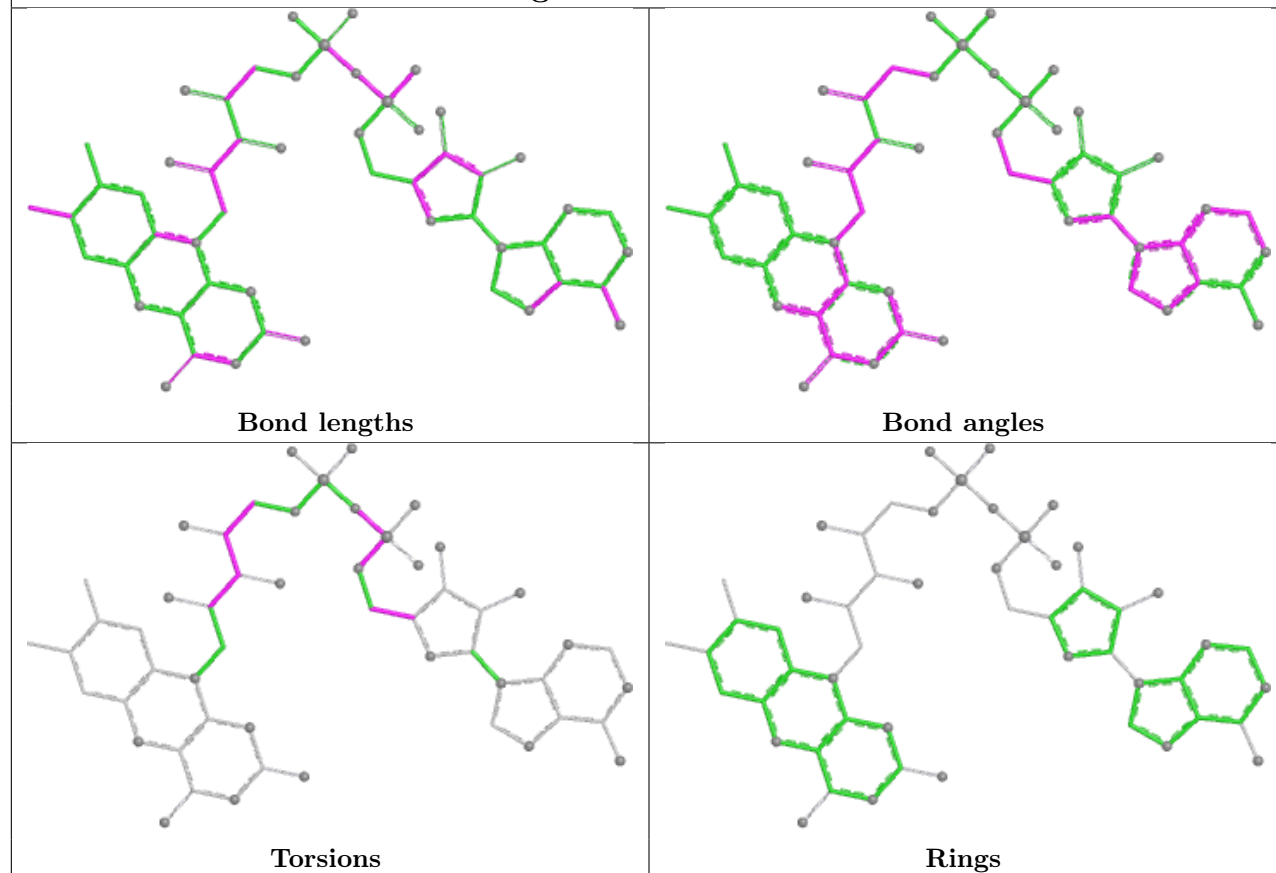
There are no ring outliers.

16 monomers are involved in 32 short contacts:

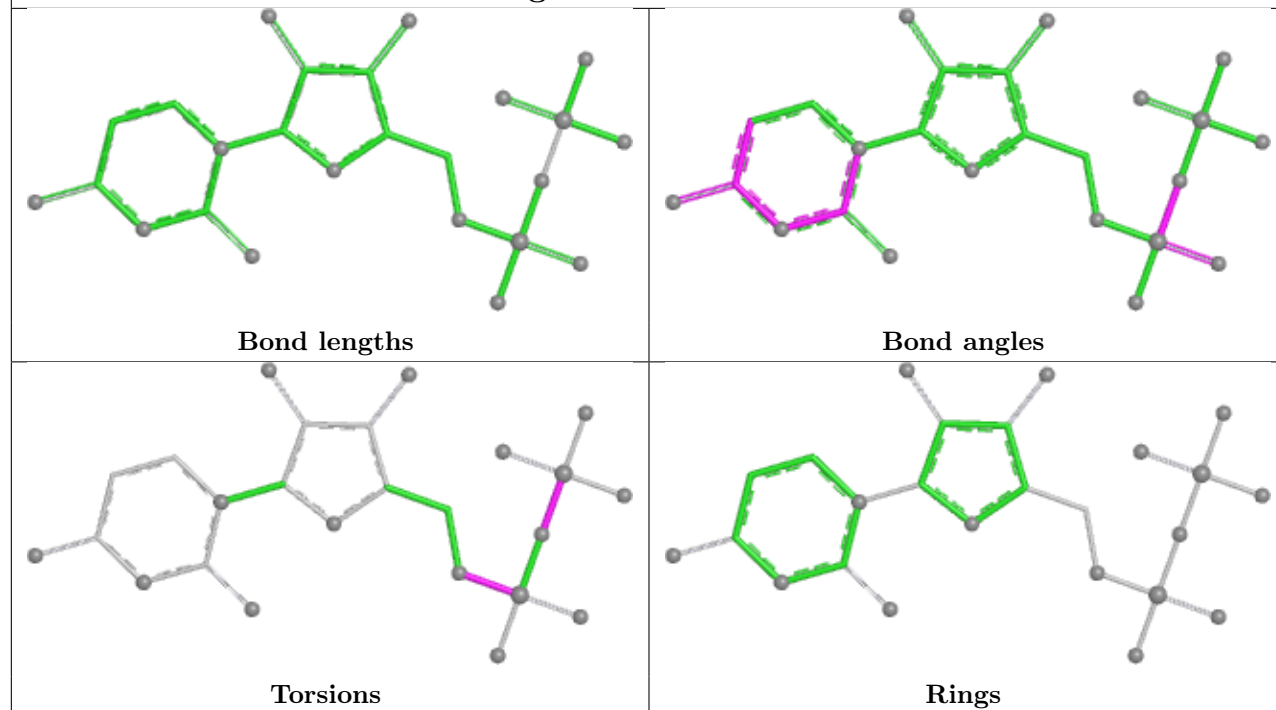
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	607	FAD	2	0
2	G	800	UDP	3	0
3	F	603	FAD	1	0
2	F	800	UDP	2	0
2	H	800	UDP	2	0
3	B	604	FAD	1	0
2	C	800	UDP	4	0
3	A	602	FAD	1	0
3	C	600	FAD	1	0
2	D	800	UDP	3	0
3	D	601	FAD	1	0
3	E	606	FAD	1	0
2	B	800	UDP	2	0
2	A	800	UDP	3	0
2	E	800	UDP	2	0
3	H	605	FAD	3	0

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.

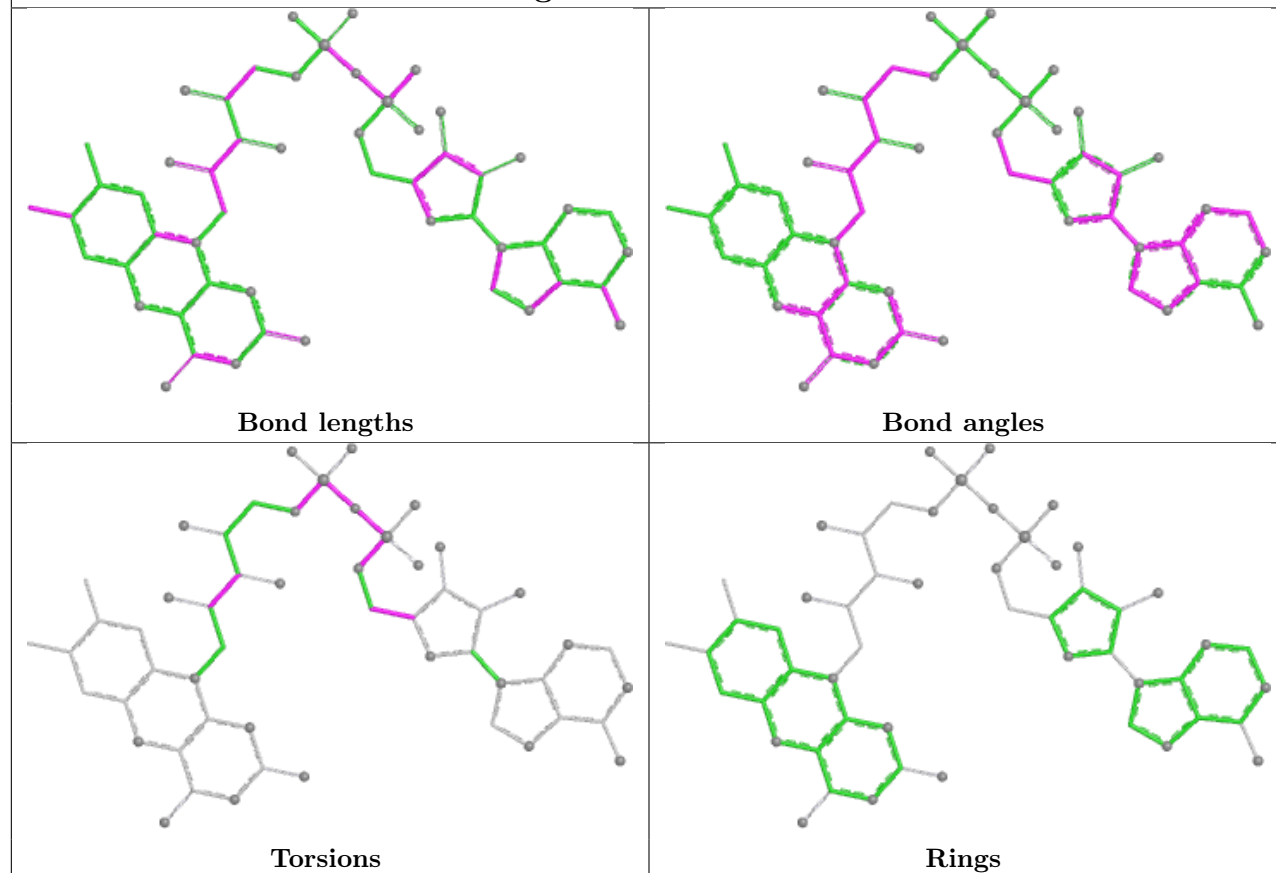
Ligand FAD G 607



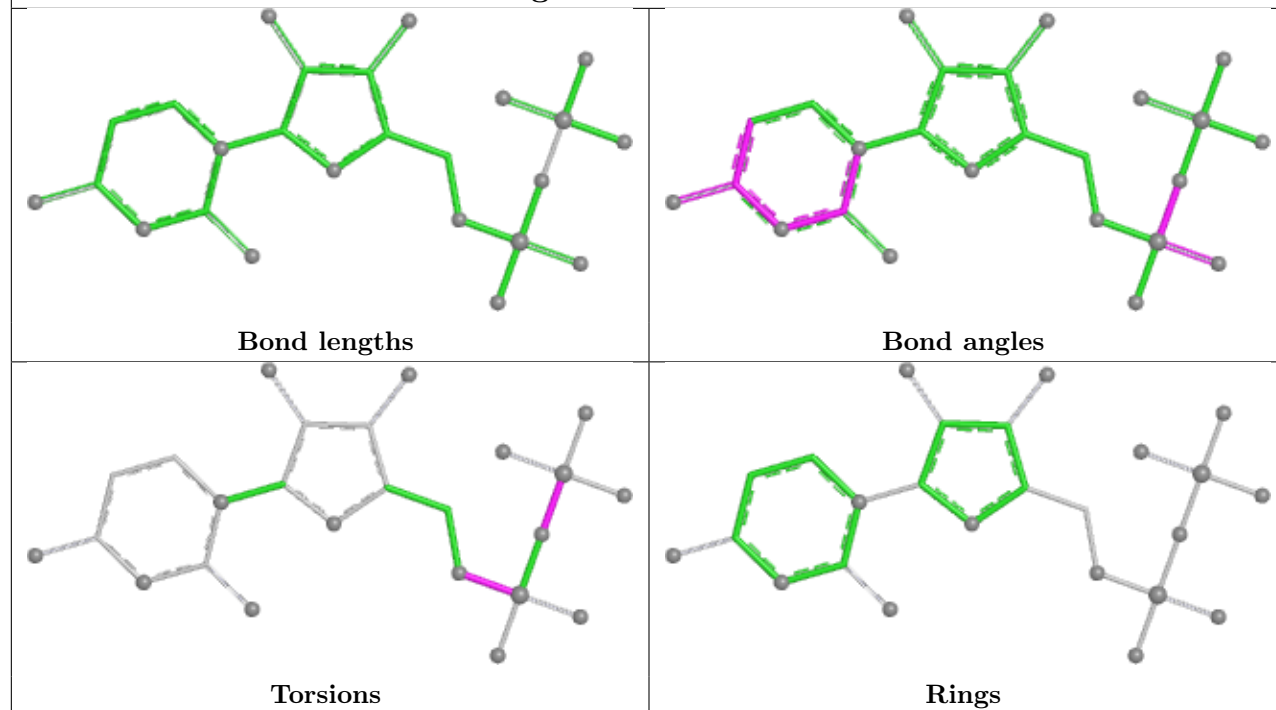
Ligand UDP G 800

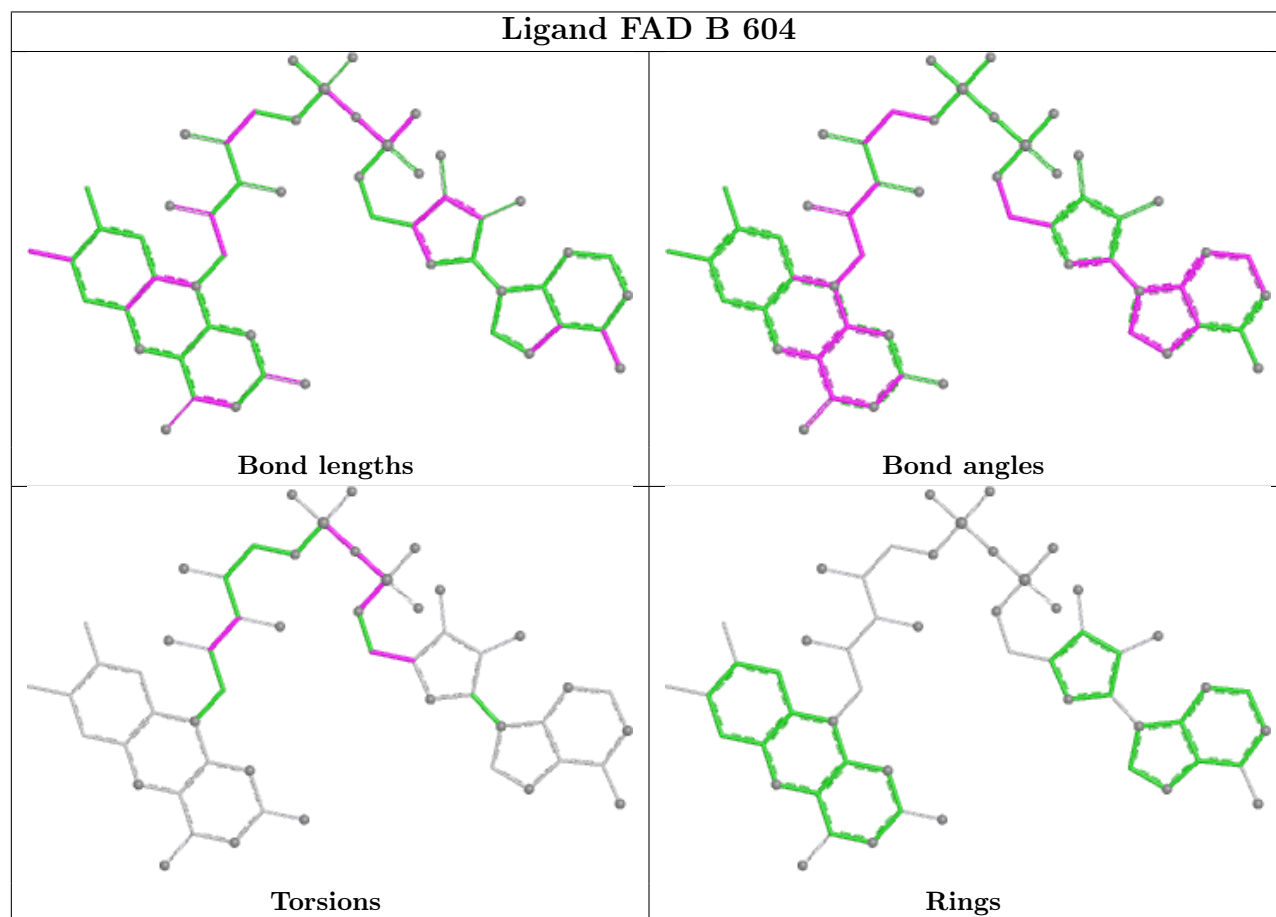
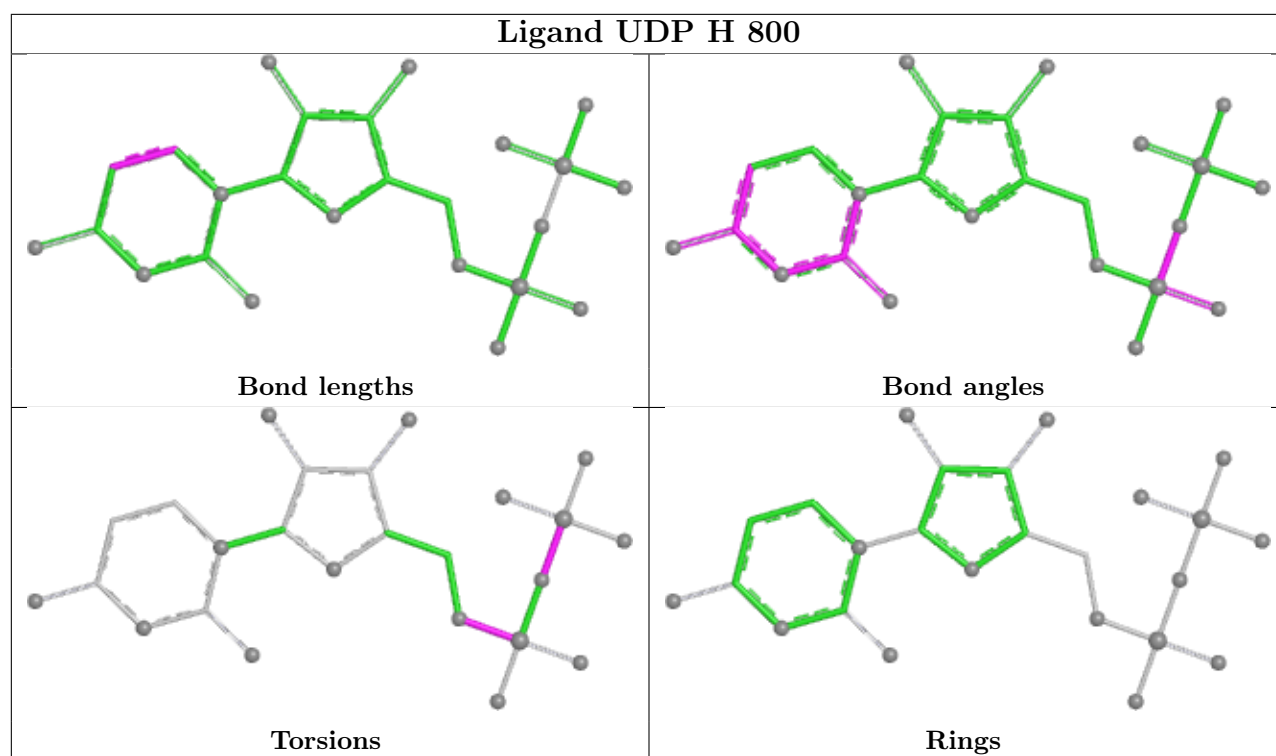


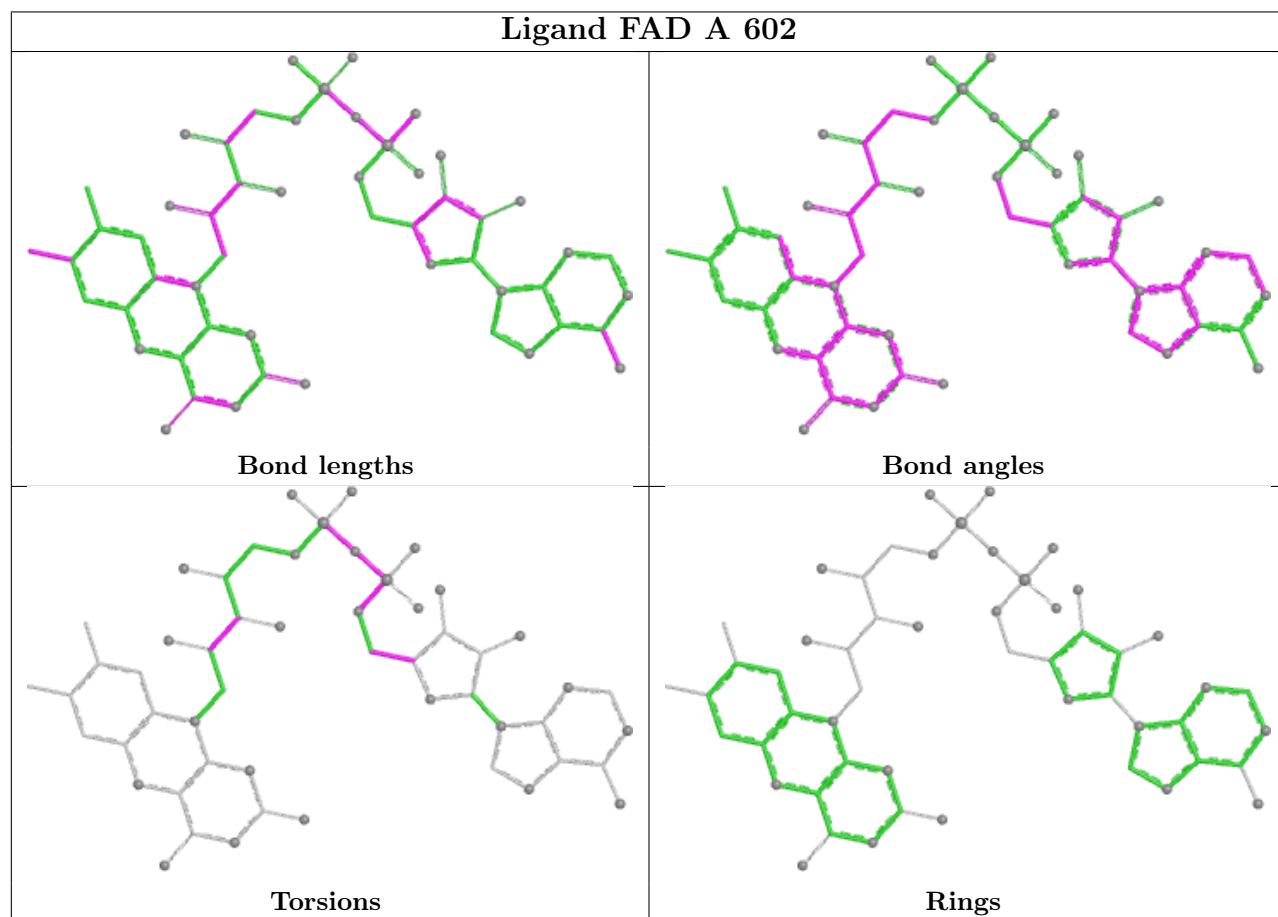
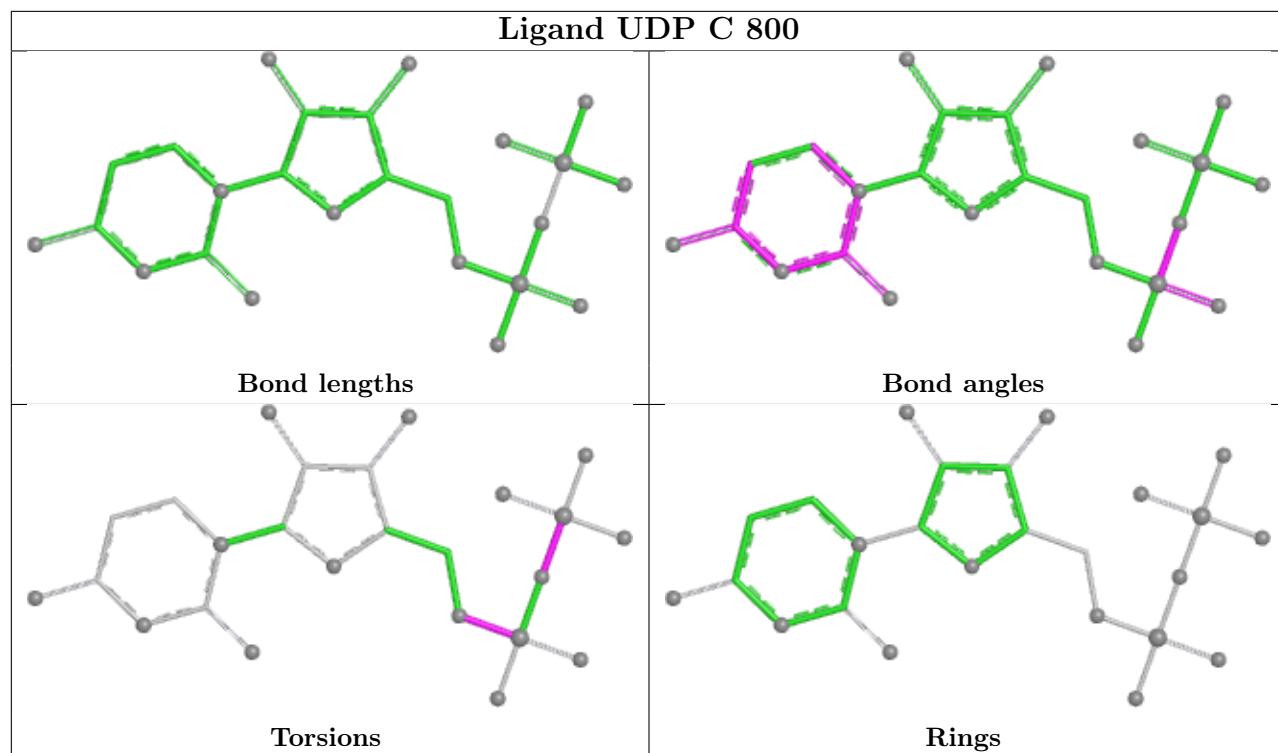
Ligand FAD F 603



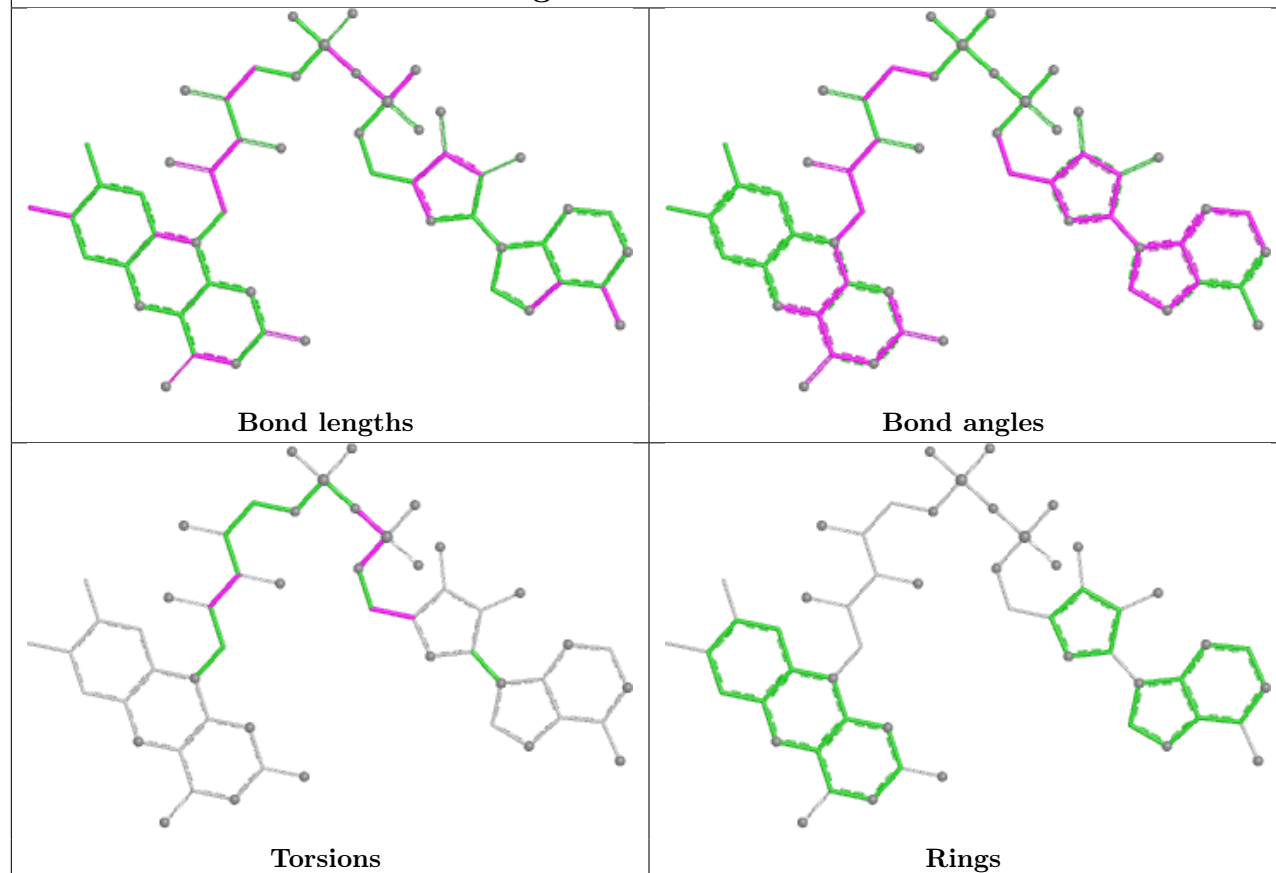
Ligand UDP F 800



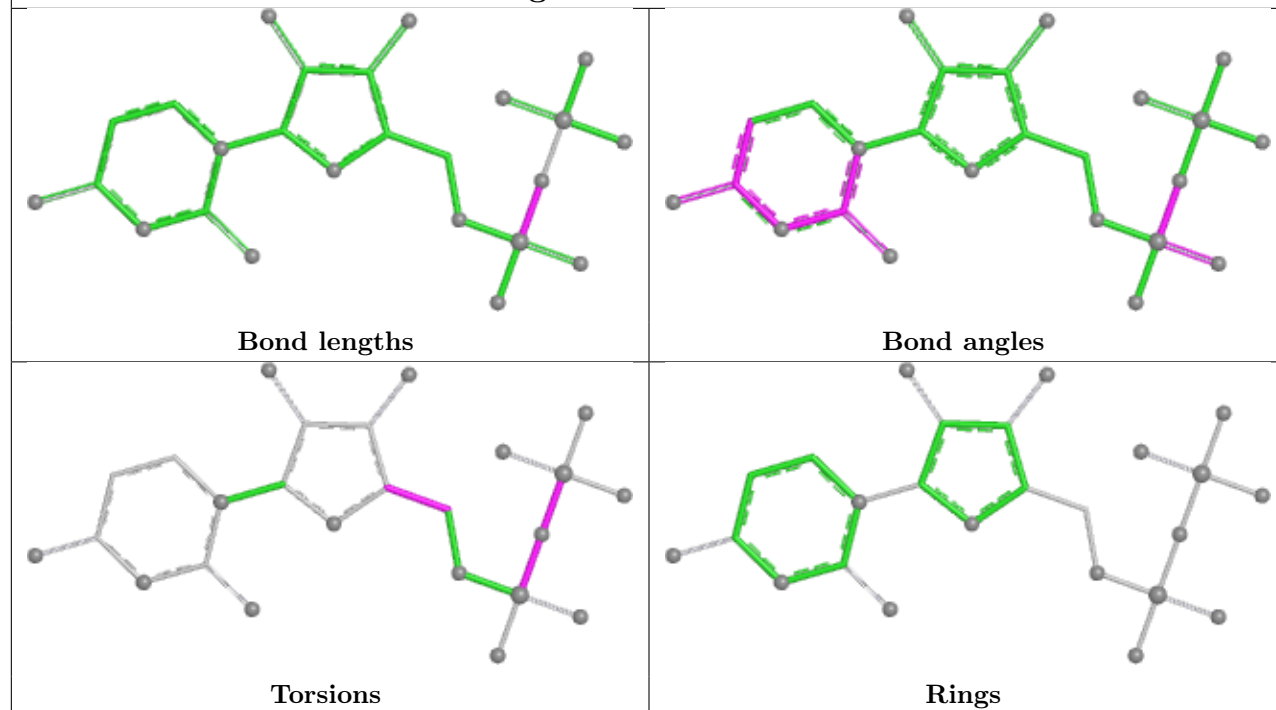


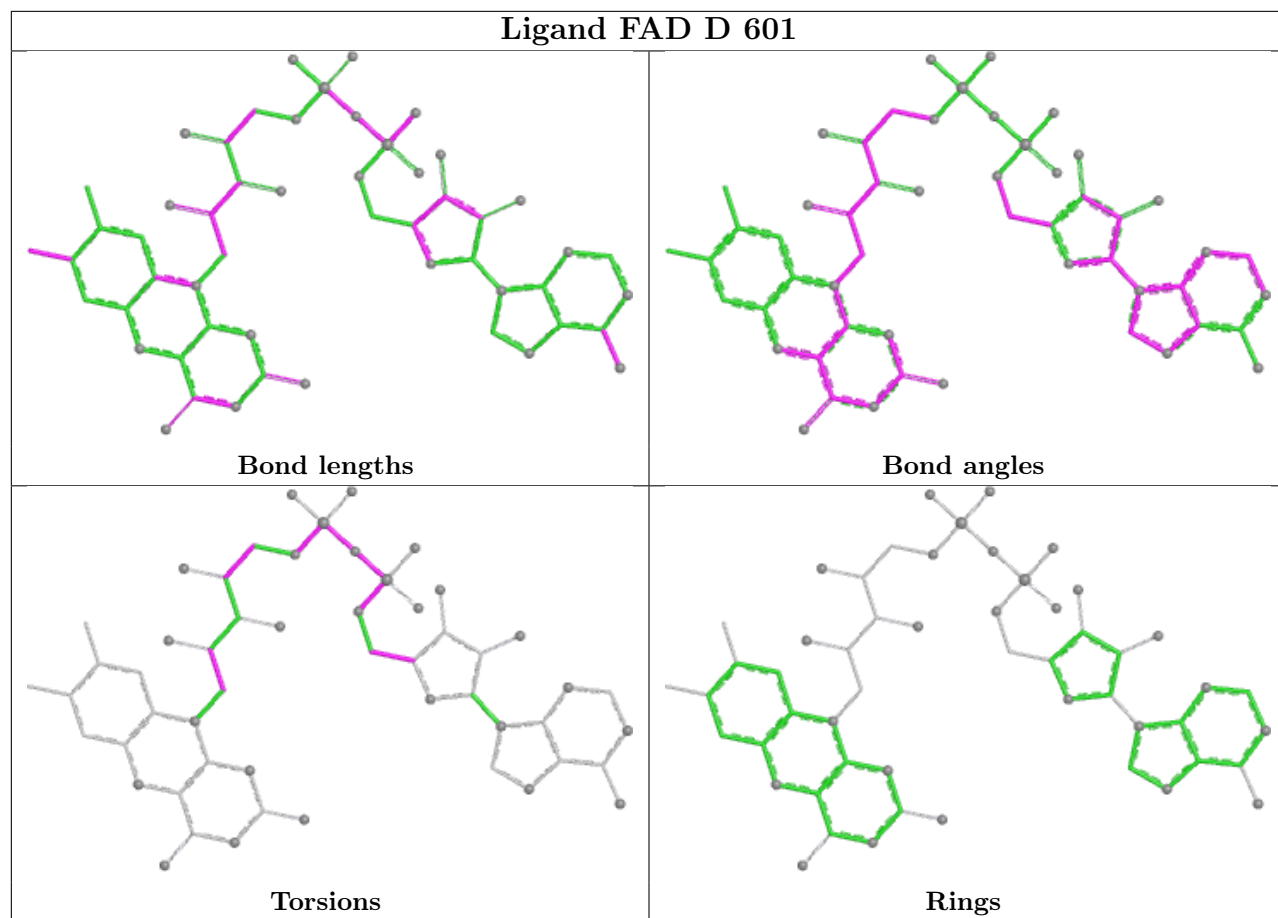


Ligand FAD C 600

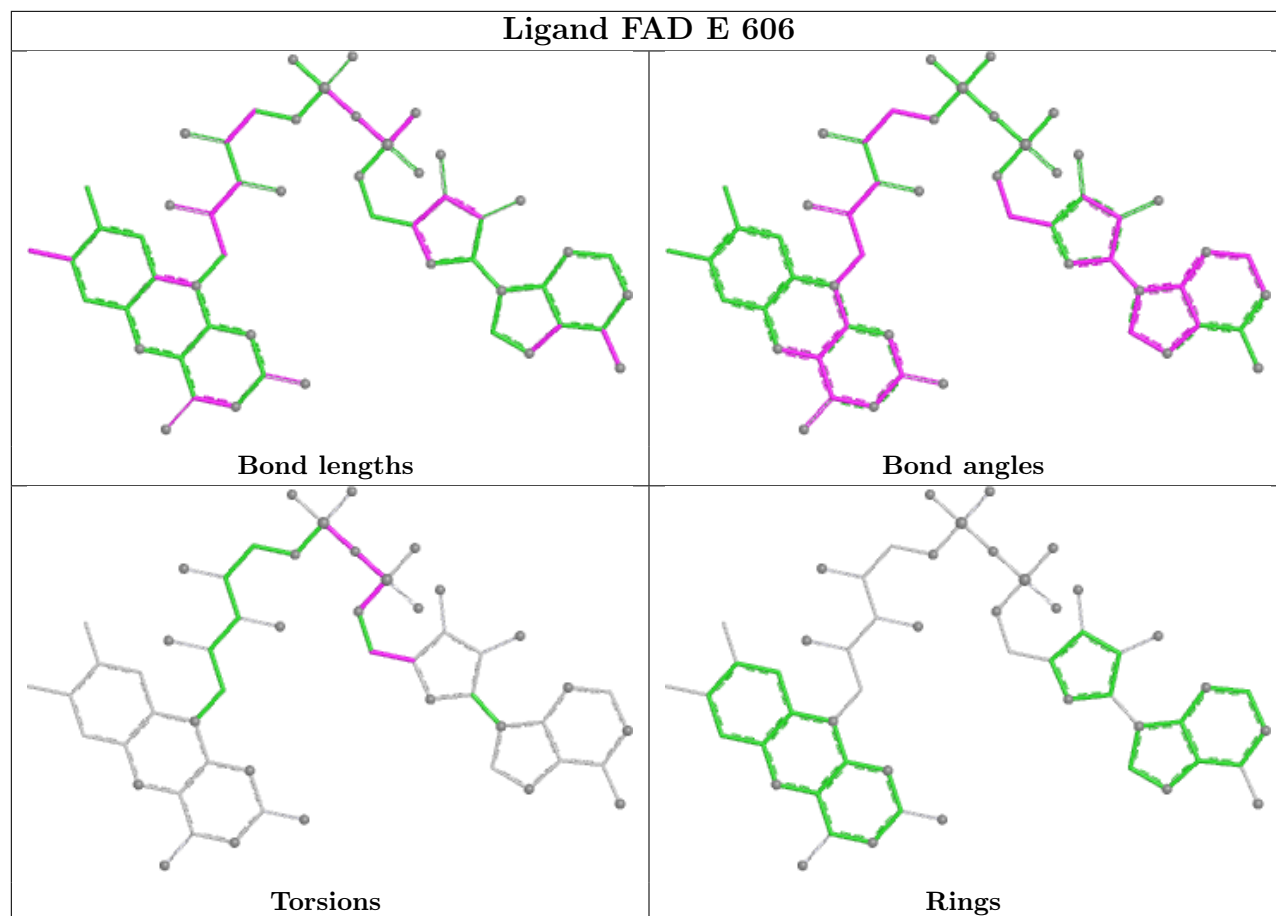


Ligand UDP D 800

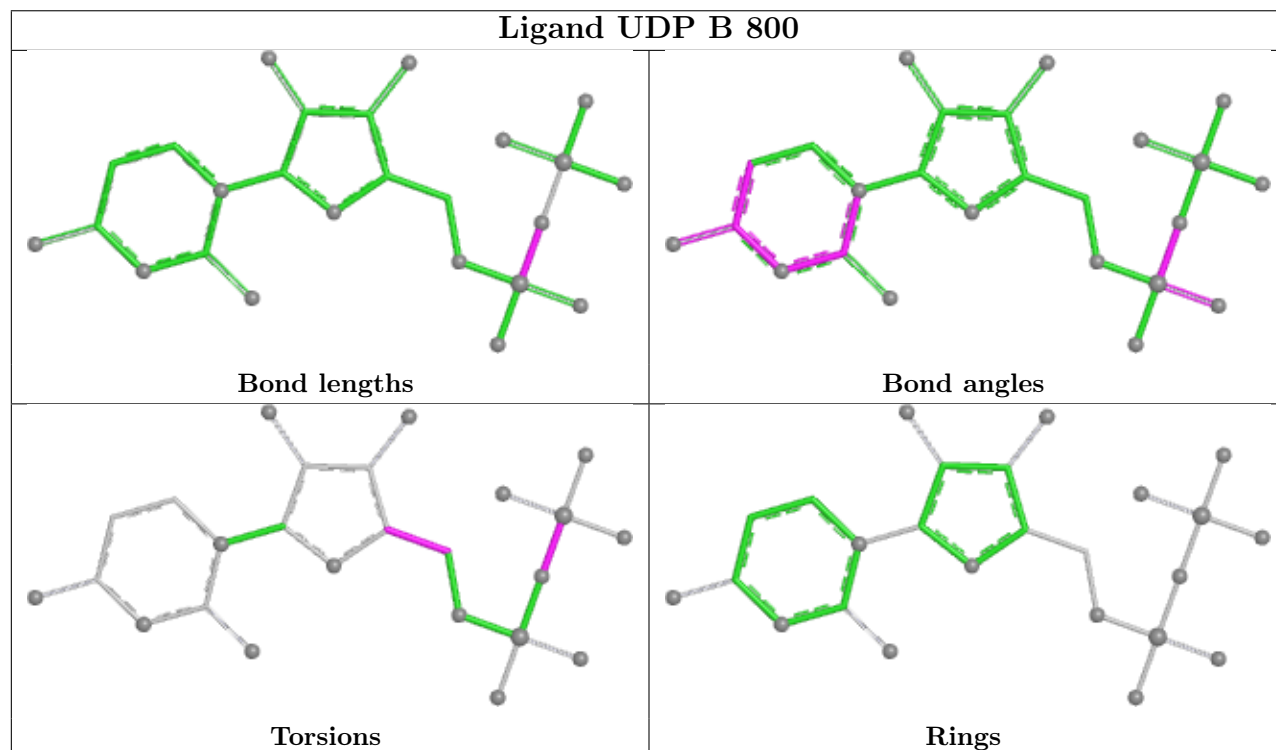


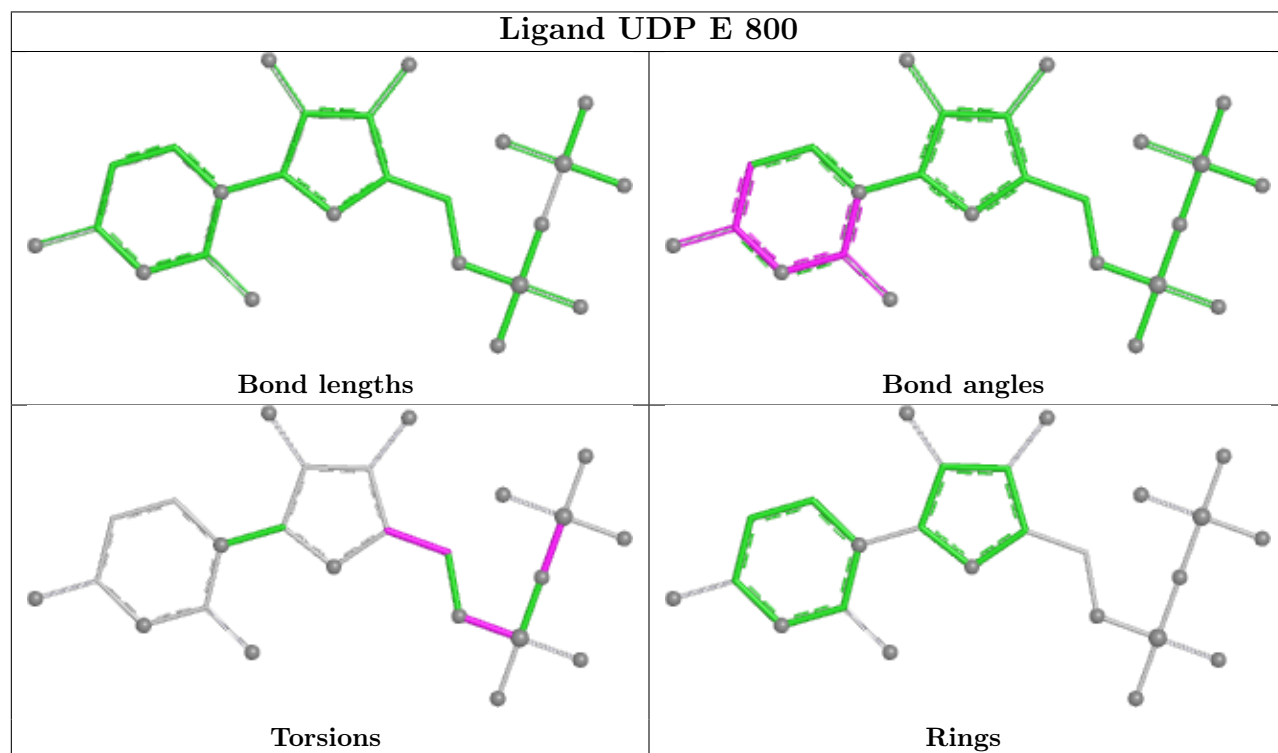
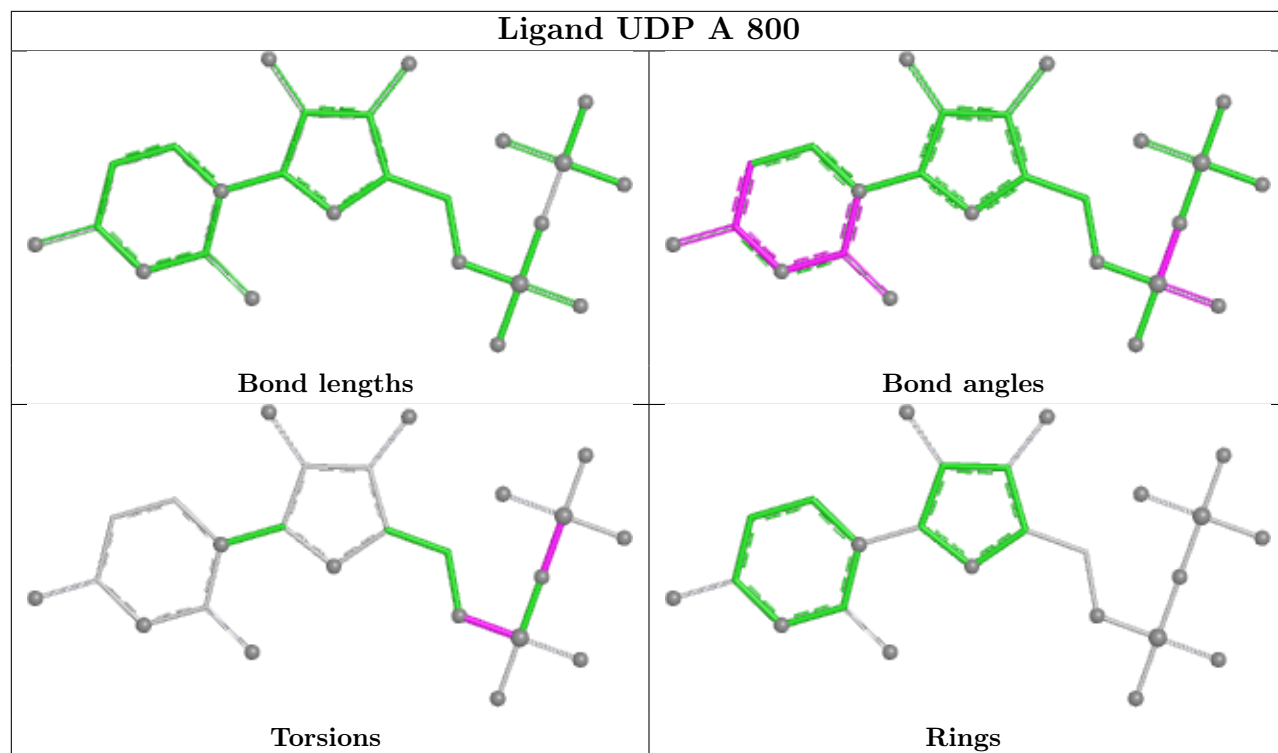


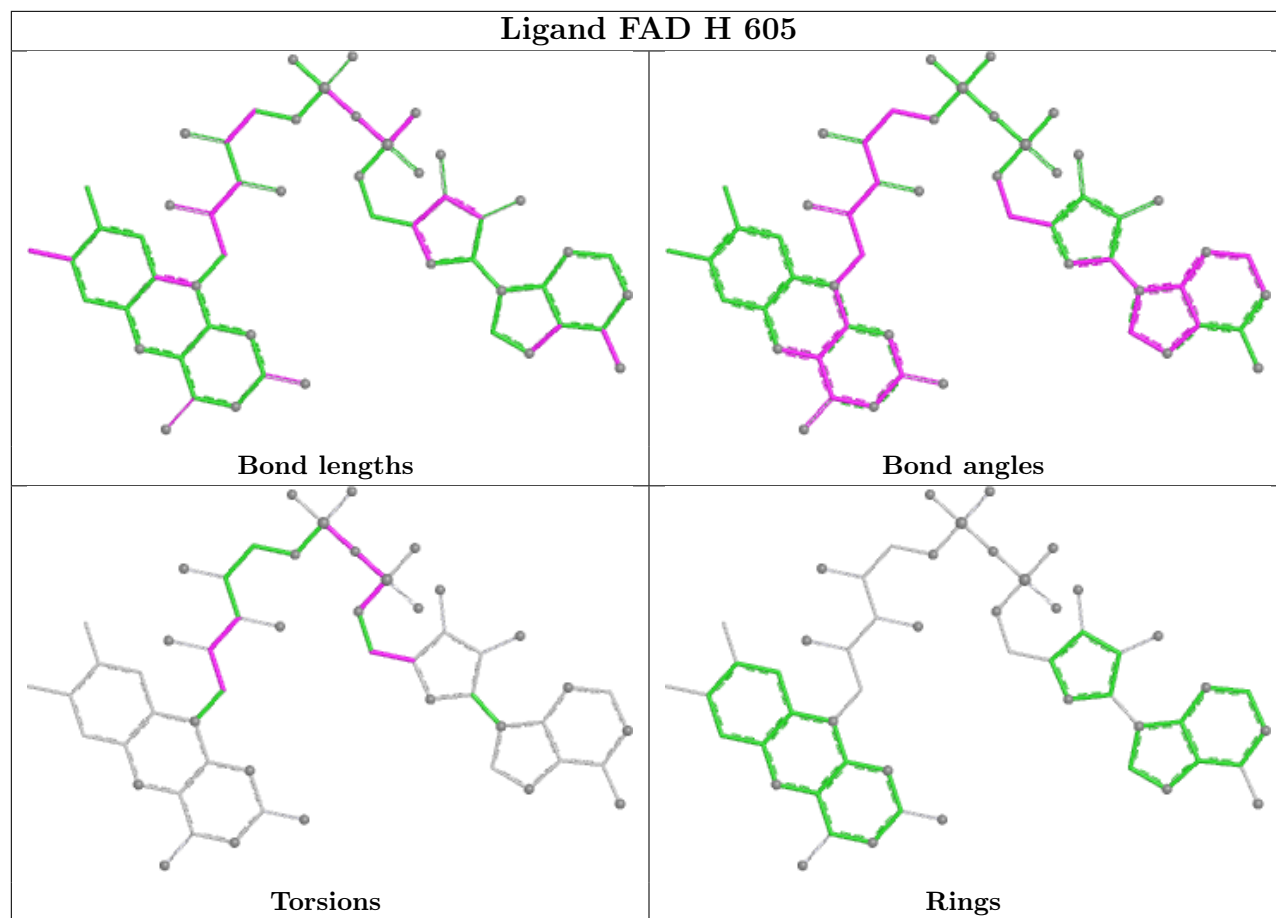
Ligand FAD E 606



Ligand UDP B 800







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	510/510 (100%)	-0.40	14 (2%) 56 51	21, 41, 77, 114	0
1	B	510/510 (100%)	-0.42	10 (1%) 65 61	20, 40, 77, 114	1 (0%)
1	C	510/510 (100%)	-0.36	11 (2%) 62 58	22, 41, 77, 118	1 (0%)
1	D	510/510 (100%)	-0.32	14 (2%) 56 51	23, 43, 79, 119	0
1	E	510/510 (100%)	-0.36	8 (1%) 70 67	20, 43, 77, 118	1 (0%)
1	F	510/510 (100%)	-0.28	9 (1%) 67 64	22, 44, 79, 119	0
1	G	510/510 (100%)	-0.26	12 (2%) 59 55	25, 44, 78, 116	0
1	H	510/510 (100%)	-0.13	13 (2%) 58 54	24, 45, 80, 117	1 (0%)
All	All	4080/4080 (100%)	-0.32	91 (2%) 62 58	20, 43, 79, 119	4 (0%)

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	208	ALA	5.1
1	G	208	ALA	5.0
1	F	208	ALA	5.0
1	H	208	ALA	4.9
1	B	208	ALA	4.5
1	H	511	LEU	4.4
1	D	208	ALA	4.4
1	D	204	TRP	4.2
1	A	208	ALA	4.2
1	H	211[A]	ARG	3.9
1	C	204	TRP	3.7
1	E	208	ALA	3.7
1	H	2	THR	3.6
1	C	206	PRO	3.6
1	G	511	LEU	3.6
1	B	2	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	511	LEU	3.5
1	E	511	LEU	3.5
1	D	205	GLY	3.4
1	D	511	LEU	3.4
1	H	204	TRP	3.3
1	B	204	TRP	3.3
1	D	206	PRO	3.2
1	A	62	GLY	3.2
1	F	206	PRO	3.2
1	H	180	GLY	3.2
1	F	65	ILE	3.1
1	E	204	TRP	3.0
1	H	62	GLY	3.0
1	F	201	ALA	3.0
1	C	511	LEU	2.9
1	H	209	THR	2.9
1	H	206	PRO	2.9
1	A	64	VAL	2.8
1	D	64	VAL	2.8
1	H	64	VAL	2.8
1	A	65	ILE	2.8
1	G	62	GLY	2.8
1	E	64	VAL	2.8
1	H	201	ALA	2.7
1	A	180	GLY	2.7
1	A	207	ASN	2.7
1	A	184	ALA	2.7
1	D	180	GLY	2.7
1	D	62	GLY	2.6
1	B	66[A]	PHE	2.6
1	D	2	THR	2.6
1	E	2	THR	2.6
1	A	2	THR	2.5
1	C	83	GLU	2.5
1	G	2	THR	2.5
1	F	511	LEU	2.5
1	A	205	GLY	2.5
1	G	202	GLY	2.5
1	E	358	GLN	2.4
1	E	201	ALA	2.4
1	C	202	GLY	2.4
1	E	206	PRO	2.4

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Mol	Chain	Res	Type	RSRZ
1	H	244	LYS	2.4
1	G	64	VAL	2.4
1	G	65	ILE	2.4
1	A	183	VAL	2.4
1	B	356	ARG	2.4
1	C	209	THR	2.3
1	C	507	SER	2.3
1	A	204	TRP	2.3
1	B	201	ALA	2.3
1	D	184	ALA	2.3
1	F	2	THR	2.3
1	G	204	TRP	2.3
1	G	205	GLY	2.3
1	B	511	LEU	2.3
1	F	207	ASN	2.3
1	C	180	GLY	2.2
1	B	62	GLY	2.2
1	A	201	ALA	2.2
1	G	356	ARG	2.2
1	D	281	GLN	2.2
1	A	199	LYS	2.2
1	D	209	THR	2.1
1	G	358	GLN	2.1
1	G	201	ALA	2.1
1	D	198	GLY	2.1
1	C	207	ASN	2.1
1	D	207	ASN	2.1
1	C	183	VAL	2.1
1	B	359	SER	2.0
1	F	204	TRP	2.0
1	B	202	GLY	2.0
1	F	202	GLY	2.0
1	H	65	ILE	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 oligosaccharides in this entry.

6.4 Ligands ⓘ

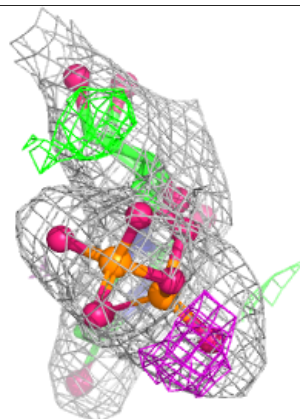
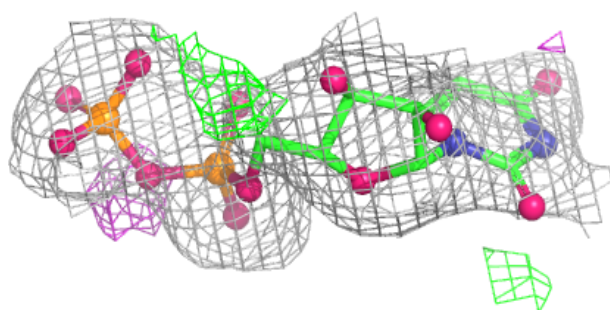
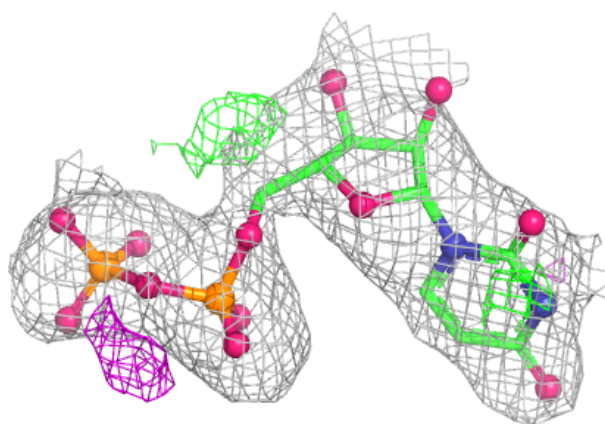
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CL	F	512	1/1	0.76	0.20	81,81,81,81	0
4	CL	E	1	1/1	0.83	0.17	75,75,75,75	0
4	CL	H	1	1/1	0.85	0.12	66,66,66,66	0
4	CL	F	1	1/1	0.86	0.16	85,85,85,85	0
2	UDP	D	800	25/25	0.90	0.12	41,86,121,130	0
4	CL	A	512	1/1	0.90	0.24	79,79,79,79	0
2	UDP	G	800	25/25	0.92	0.09	45,69,97,117	0
2	UDP	F	800	25/25	0.92	0.09	47,68,104,116	0
2	UDP	C	800	25/25	0.93	0.08	43,61,94,104	0
4	CL	C	1	1/1	0.94	0.08	58,58,58,58	0
2	UDP	H	800	25/25	0.94	0.09	50,80,110,113	0
4	CL	A	1	1/1	0.94	0.09	62,62,62,62	0
2	UDP	A	800	25/25	0.94	0.09	53,74,109,122	0
4	CL	B	1	1/1	0.94	0.10	63,63,63,63	0
3	FAD	H	605	53/53	0.95	0.08	27,44,64,69	0
2	UDP	E	800	25/25	0.95	0.08	32,63,96,107	0
2	UDP	B	800	25/25	0.95	0.07	30,50,91,101	0
3	FAD	D	601	53/53	0.97	0.07	23,37,50,57	0
3	FAD	F	603	53/53	0.97	0.07	19,39,52,67	0
3	FAD	G	607	53/53	0.97	0.06	22,39,53,62	0
3	FAD	C	600	53/53	0.97	0.07	11,33,48,54	0
3	FAD	E	606	53/53	0.98	0.06	19,31,47,59	0
3	FAD	A	602	53/53	0.98	0.06	16,33,53,70	0
3	FAD	B	604	53/53	0.98	0.06	18,35,47,51	0

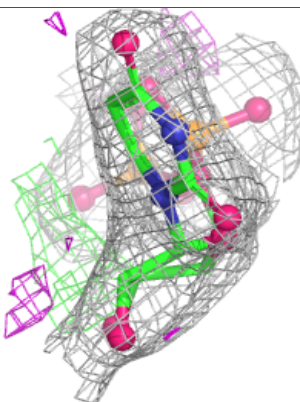
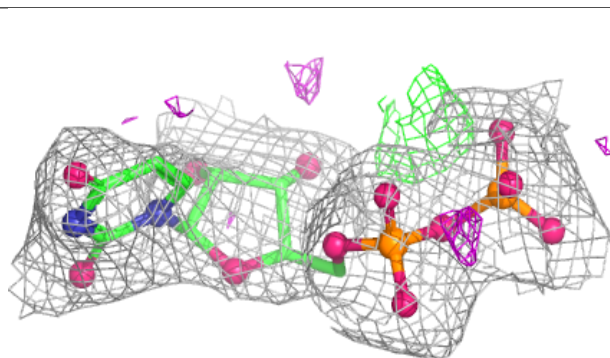
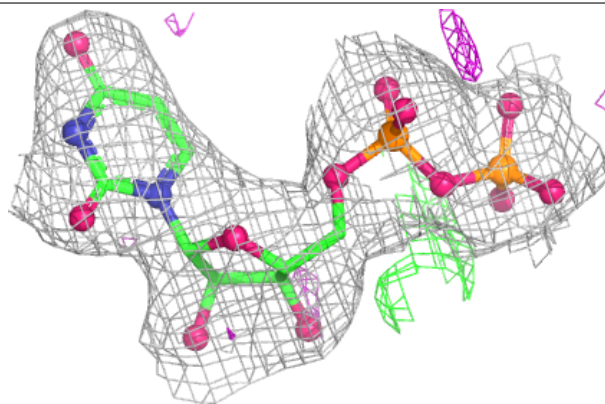
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UDP D 800:

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

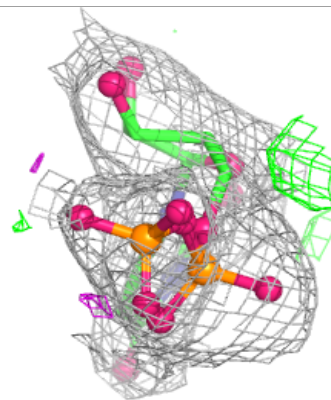
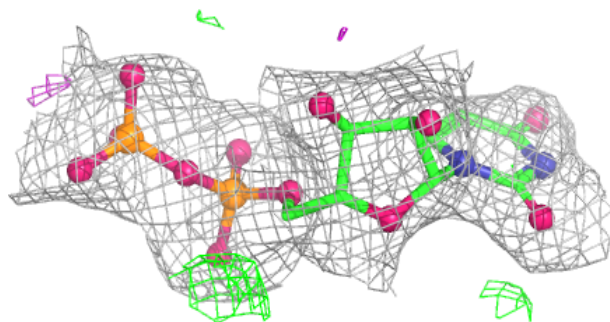
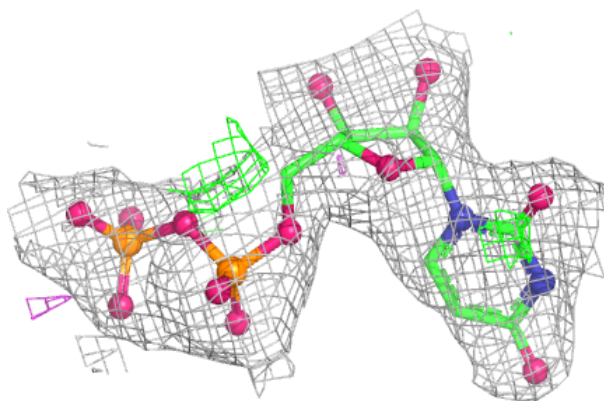
**Electron density around UDP G 800:**

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

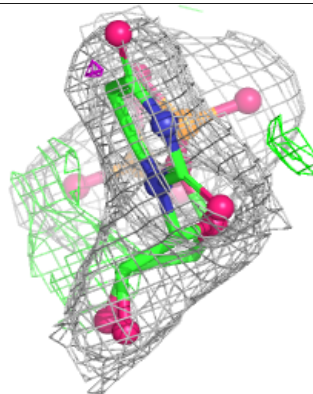
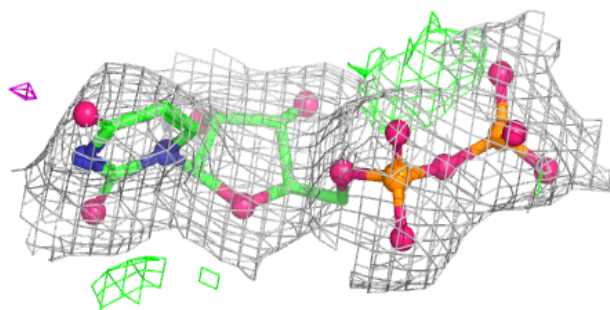
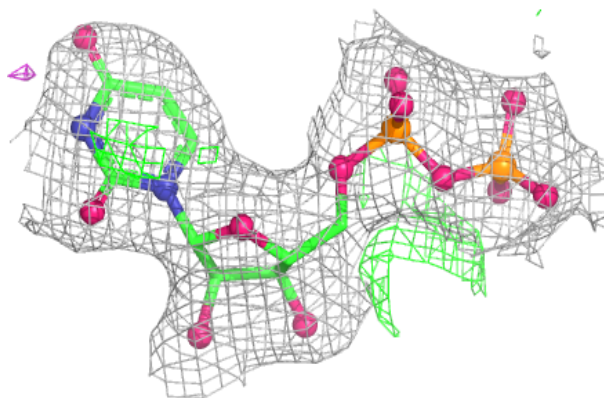


Electron density around UDP F 800:

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

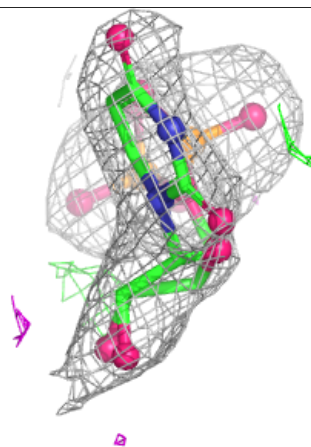
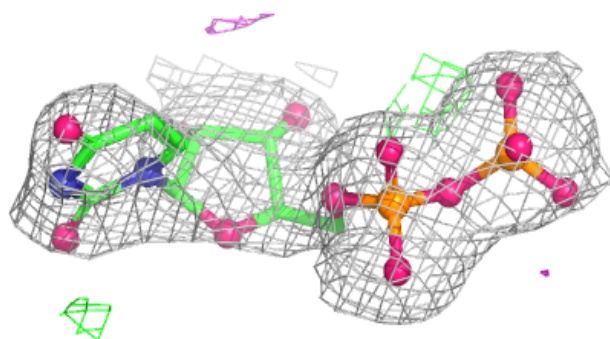
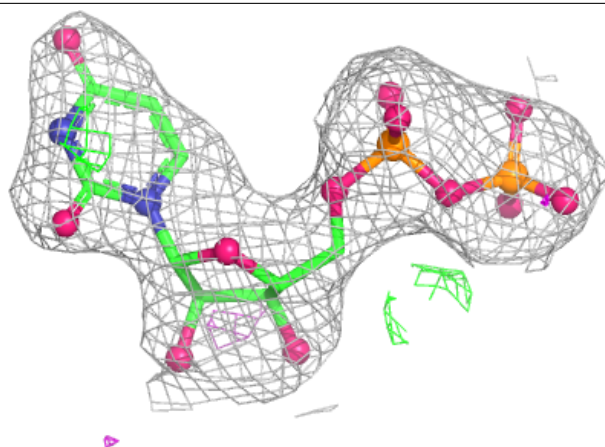
**Electron density around UDP C 800:**

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



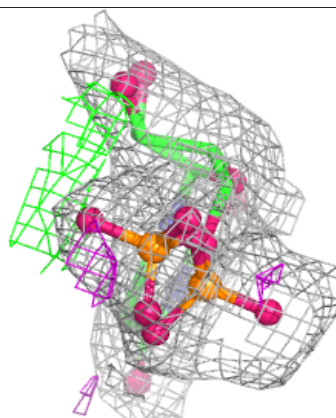
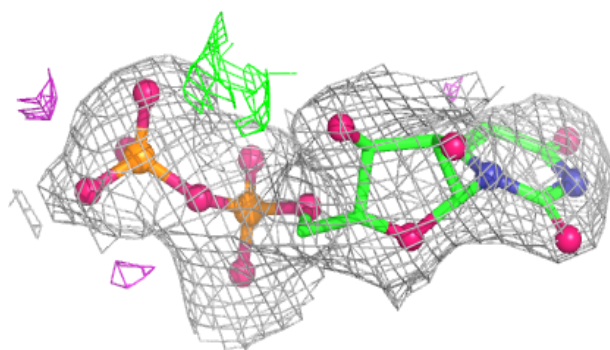
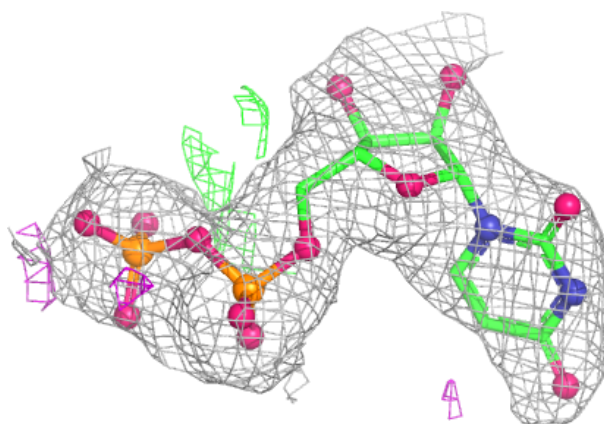
Electron density around UDP H 800:

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

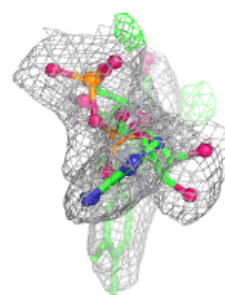
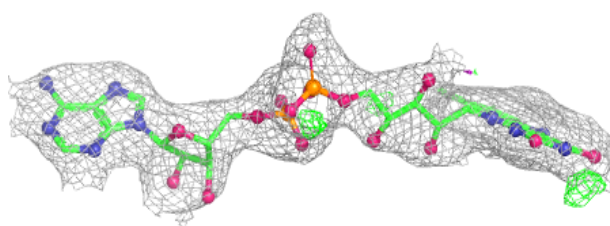
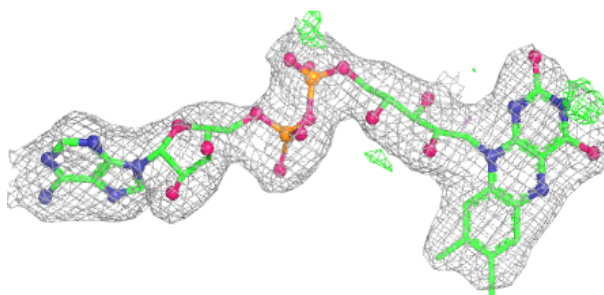


Electron density around UDP A 800:

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

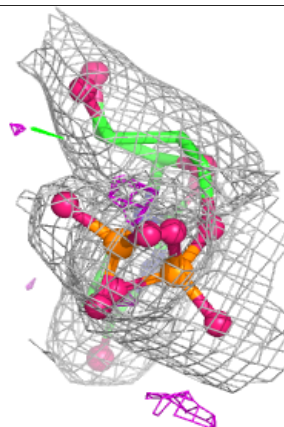
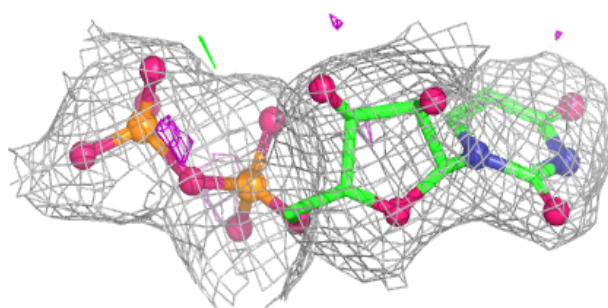
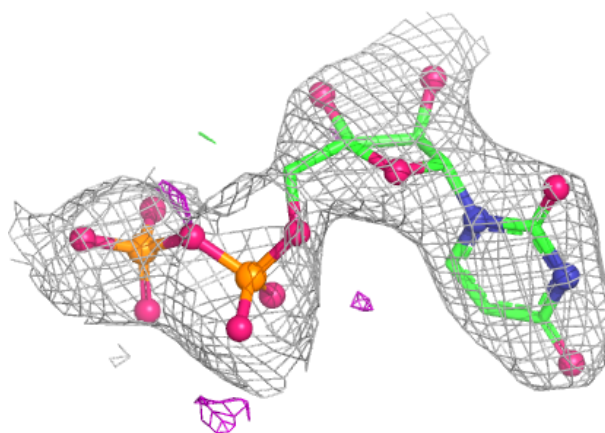
**Electron density around FAD H 605:**

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

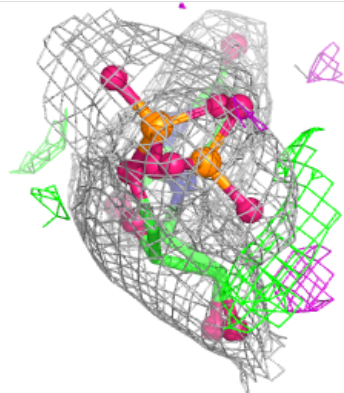
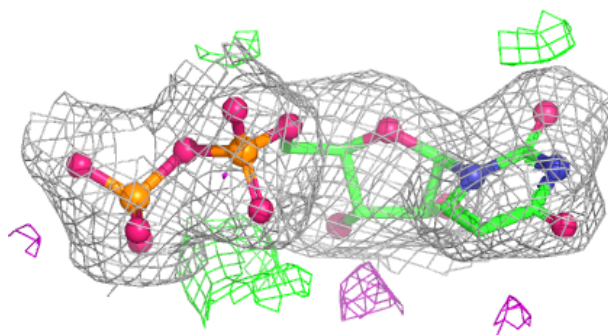
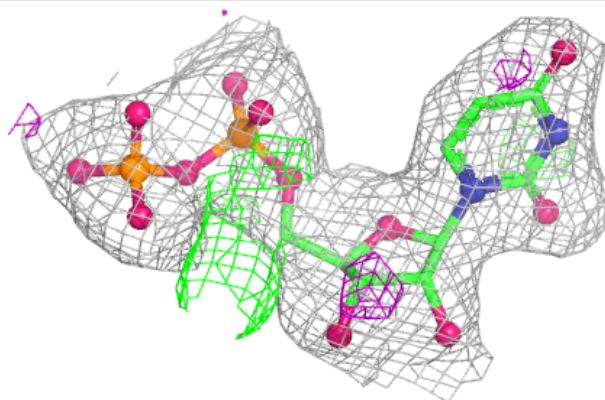


Electron density around UDP E 800:

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

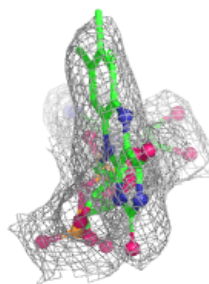
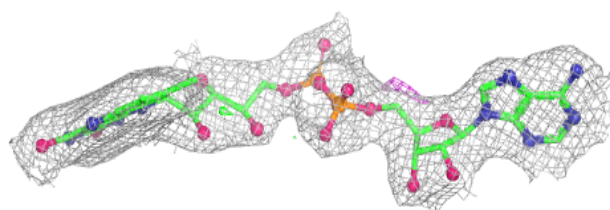
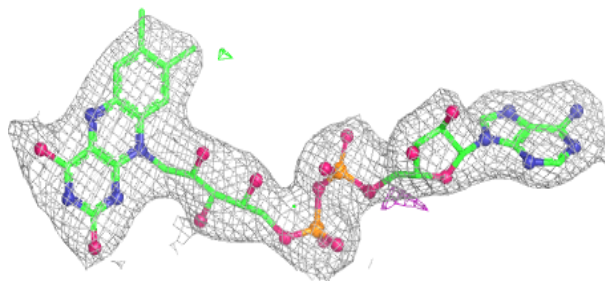
**Electron density around UDP B 800:**

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

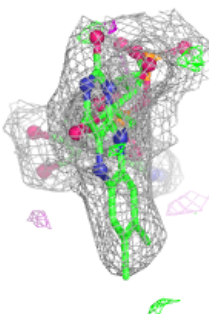
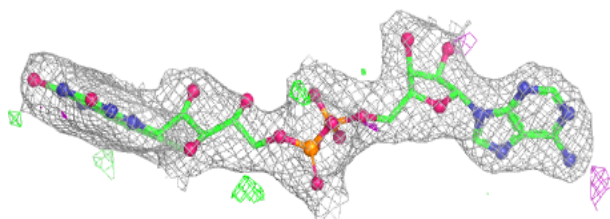
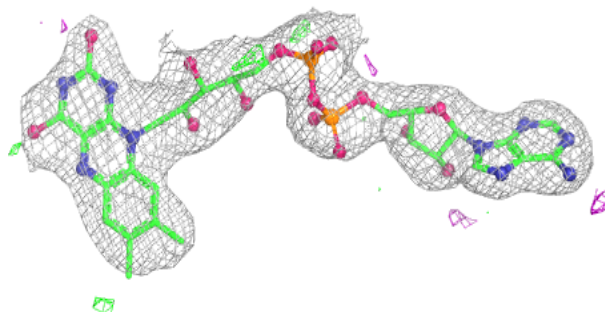


Electron density around FAD D 601:

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

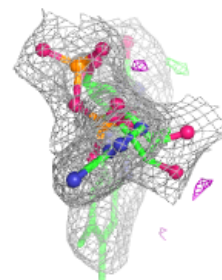
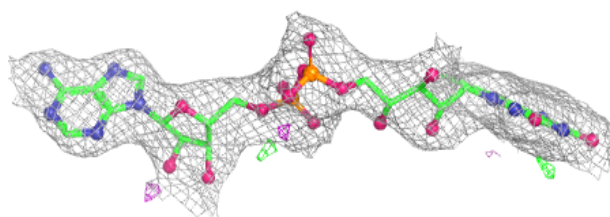
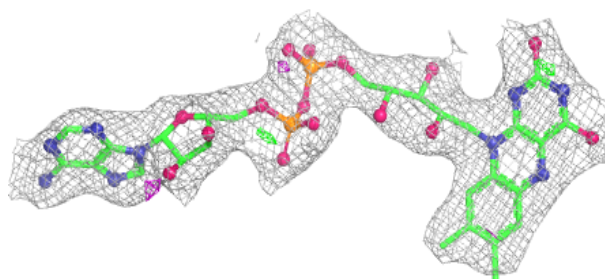
**Electron density around FAD F 603:**

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

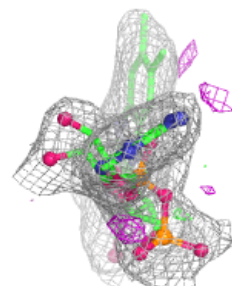
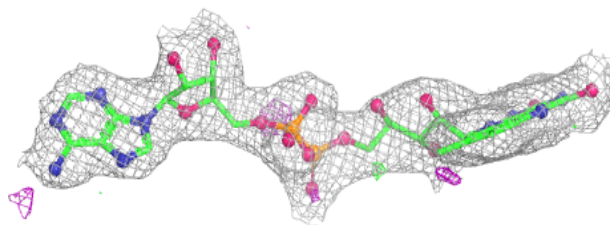
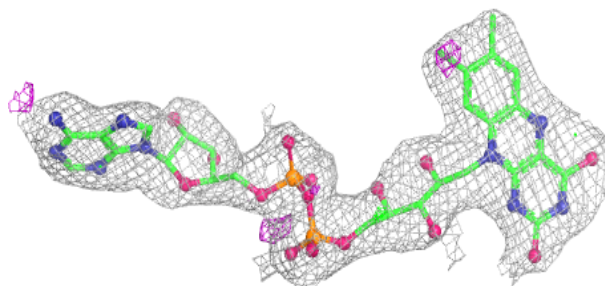


Electron density around FAD G 607:

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

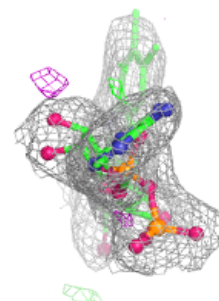
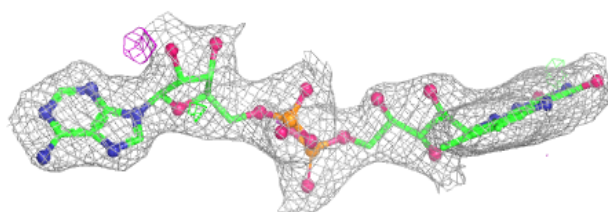
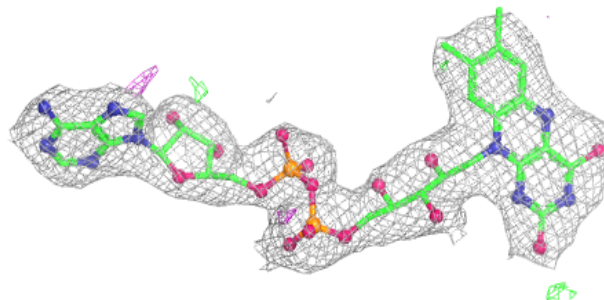
**Electron density around FAD C 600:**

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

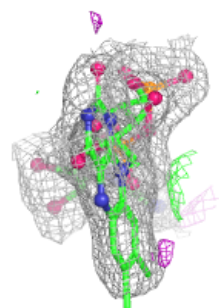
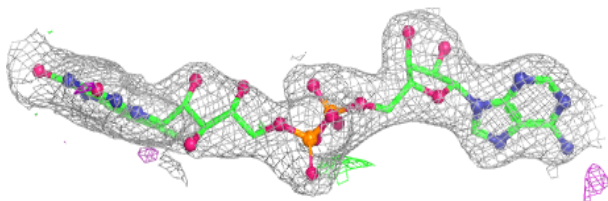
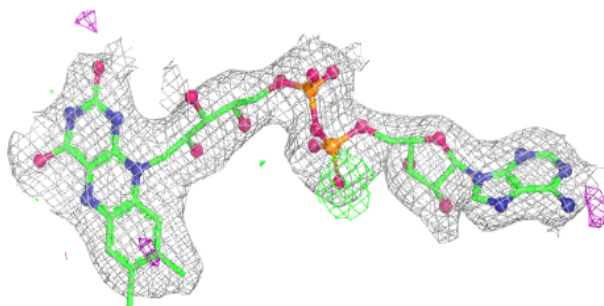


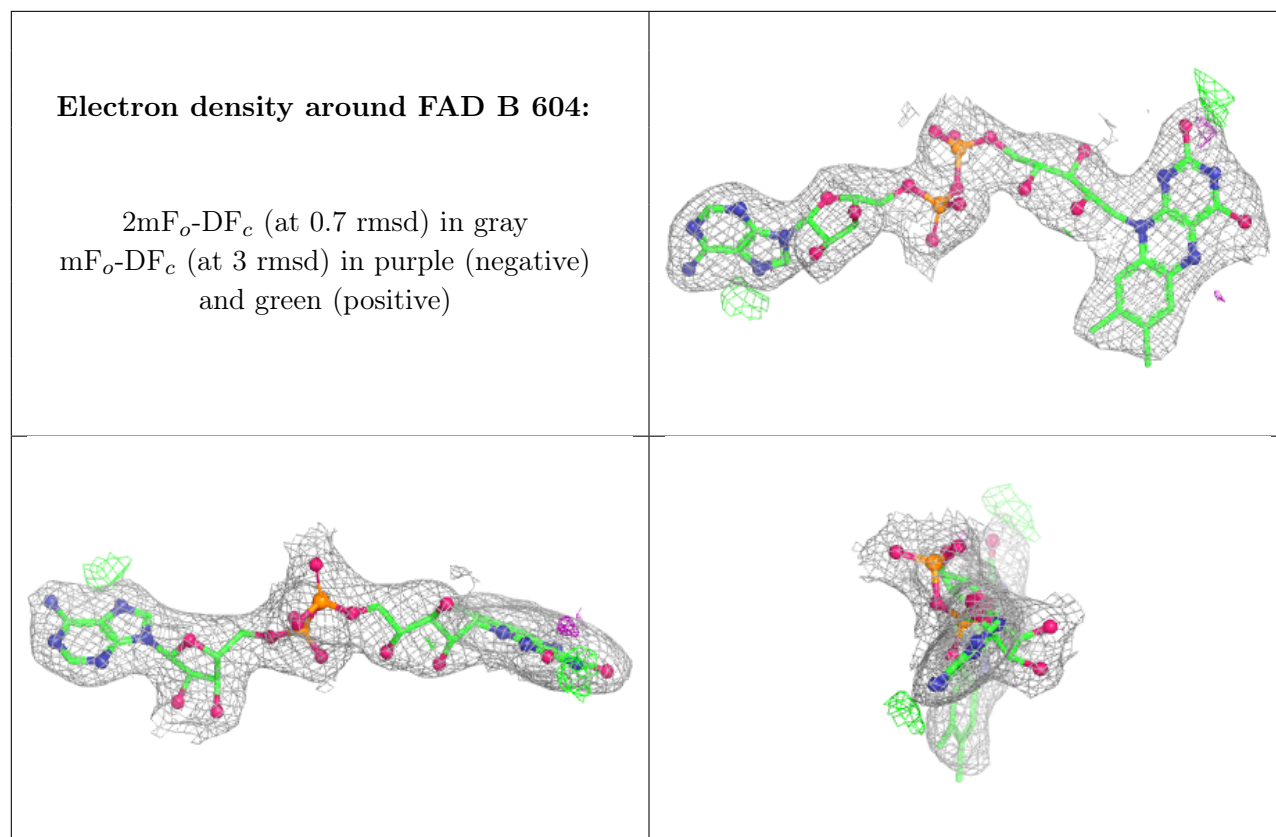
Electron density around FAD E 606:

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

**Electron density around FAD A 602:**

$2mF_o-DF_c$ (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.