



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:37 AM UTC

PDB ID : 3APY / pdb_00003apy
Title : Properties and crystal structure of methylenetetrahydrofolate reductase from
Thermus thermophilus HB8
Authors : Yamada, K.
Deposited on : 2010-10-21
Resolution : 2.80 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

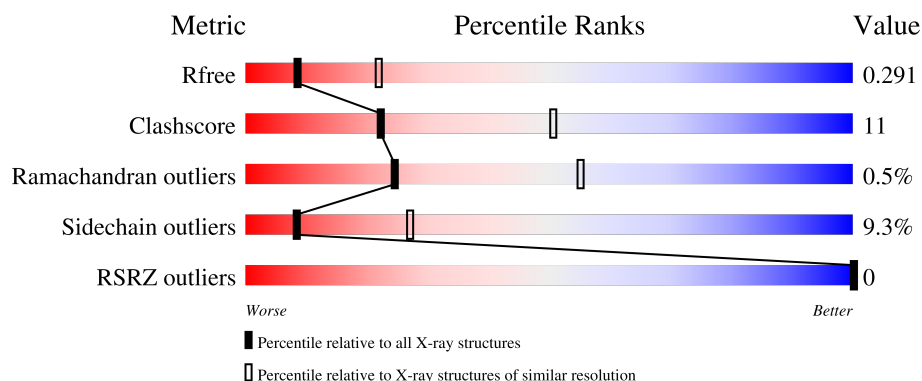
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	310	 66% 23% • 6%
1	B	310	 72% 18% • 6%
1	C	310	 70% 22% • 6%
1	D	310	 71% 18% • 6%
1	E	310	 69% 22% • 6%

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Mol	Chain	Length	Quality of chain
1	F	310	 70% 23% • 6%
1	G	310	 68% 22% • 6%
1	H	310	 71% 20% • 6%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 19755 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 Methylenetetrahydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	B	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	C	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	D	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	E	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	F	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	G	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			
1	H	291	Total	C	N	O	S	0	0	0
			2302	1469	422	406	5			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	297	ALA	-	expression tag	UNP Q5SLG6
A	298	LYS	-	expression tag	UNP Q5SLG6
A	299	LEU	-	expression tag	UNP Q5SLG6
A	300	ALA	-	expression tag	UNP Q5SLG6
A	301	ALA	-	expression tag	UNP Q5SLG6
A	302	ALA	-	expression tag	UNP Q5SLG6
A	303	LEU	-	expression tag	UNP Q5SLG6
A	304	GLU	-	expression tag	UNP Q5SLG6
A	305	HIS	-	expression tag	UNP Q5SLG6
A	306	HIS	-	expression tag	UNP Q5SLG6
A	307	HIS	-	expression tag	UNP Q5SLG6
A	308	HIS	-	expression tag	UNP Q5SLG6
A	309	HIS	-	expression tag	UNP Q5SLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
A	310	HIS	-	expression tag	UNP Q5SLG6
B	297	ALA	-	expression tag	UNP Q5SLG6
B	298	LYS	-	expression tag	UNP Q5SLG6
B	299	LEU	-	expression tag	UNP Q5SLG6
B	300	ALA	-	expression tag	UNP Q5SLG6
B	301	ALA	-	expression tag	UNP Q5SLG6
B	302	ALA	-	expression tag	UNP Q5SLG6
B	303	LEU	-	expression tag	UNP Q5SLG6
B	304	GLU	-	expression tag	UNP Q5SLG6
B	305	HIS	-	expression tag	UNP Q5SLG6
B	306	HIS	-	expression tag	UNP Q5SLG6
B	307	HIS	-	expression tag	UNP Q5SLG6
B	308	HIS	-	expression tag	UNP Q5SLG6
B	309	HIS	-	expression tag	UNP Q5SLG6
B	310	HIS	-	expression tag	UNP Q5SLG6
C	297	ALA	-	expression tag	UNP Q5SLG6
C	298	LYS	-	expression tag	UNP Q5SLG6
C	299	LEU	-	expression tag	UNP Q5SLG6
C	300	ALA	-	expression tag	UNP Q5SLG6
C	301	ALA	-	expression tag	UNP Q5SLG6
C	302	ALA	-	expression tag	UNP Q5SLG6
C	303	LEU	-	expression tag	UNP Q5SLG6
C	304	GLU	-	expression tag	UNP Q5SLG6
C	305	HIS	-	expression tag	UNP Q5SLG6
C	306	HIS	-	expression tag	UNP Q5SLG6
C	307	HIS	-	expression tag	UNP Q5SLG6
C	308	HIS	-	expression tag	UNP Q5SLG6
C	309	HIS	-	expression tag	UNP Q5SLG6
C	310	HIS	-	expression tag	UNP Q5SLG6
D	297	ALA	-	expression tag	UNP Q5SLG6
D	298	LYS	-	expression tag	UNP Q5SLG6
D	299	LEU	-	expression tag	UNP Q5SLG6
D	300	ALA	-	expression tag	UNP Q5SLG6
D	301	ALA	-	expression tag	UNP Q5SLG6
D	302	ALA	-	expression tag	UNP Q5SLG6
D	303	LEU	-	expression tag	UNP Q5SLG6
D	304	GLU	-	expression tag	UNP Q5SLG6
D	305	HIS	-	expression tag	UNP Q5SLG6
D	306	HIS	-	expression tag	UNP Q5SLG6
D	307	HIS	-	expression tag	UNP Q5SLG6
D	308	HIS	-	expression tag	UNP Q5SLG6
D	309	HIS	-	expression tag	UNP Q5SLG6

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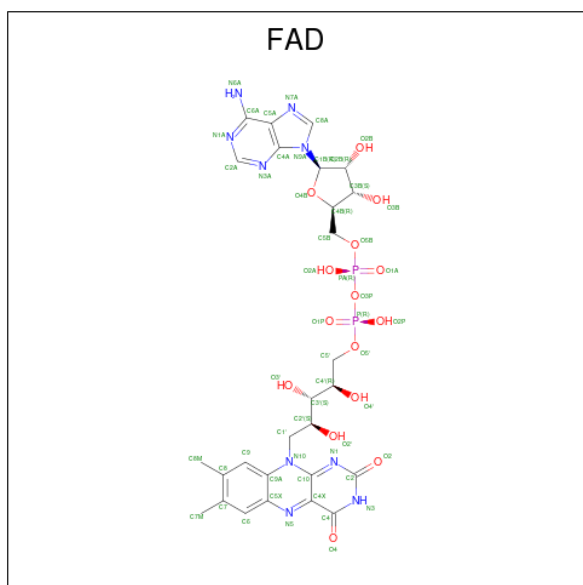
Chain	Residue	Modelled	Actual	Comment	Reference
D	310	HIS	-	expression tag	UNP Q5SLG6
E	297	ALA	-	expression tag	UNP Q5SLG6
E	298	LYS	-	expression tag	UNP Q5SLG6
E	299	LEU	-	expression tag	UNP Q5SLG6
E	300	ALA	-	expression tag	UNP Q5SLG6
E	301	ALA	-	expression tag	UNP Q5SLG6
E	302	ALA	-	expression tag	UNP Q5SLG6
E	303	LEU	-	expression tag	UNP Q5SLG6
E	304	GLU	-	expression tag	UNP Q5SLG6
E	305	HIS	-	expression tag	UNP Q5SLG6
E	306	HIS	-	expression tag	UNP Q5SLG6
E	307	HIS	-	expression tag	UNP Q5SLG6
E	308	HIS	-	expression tag	UNP Q5SLG6
E	309	HIS	-	expression tag	UNP Q5SLG6
E	310	HIS	-	expression tag	UNP Q5SLG6
F	297	ALA	-	expression tag	UNP Q5SLG6
F	298	LYS	-	expression tag	UNP Q5SLG6
F	299	LEU	-	expression tag	UNP Q5SLG6
F	300	ALA	-	expression tag	UNP Q5SLG6
F	301	ALA	-	expression tag	UNP Q5SLG6
F	302	ALA	-	expression tag	UNP Q5SLG6
F	303	LEU	-	expression tag	UNP Q5SLG6
F	304	GLU	-	expression tag	UNP Q5SLG6
F	305	HIS	-	expression tag	UNP Q5SLG6
F	306	HIS	-	expression tag	UNP Q5SLG6
F	307	HIS	-	expression tag	UNP Q5SLG6
F	308	HIS	-	expression tag	UNP Q5SLG6
F	309	HIS	-	expression tag	UNP Q5SLG6
F	310	HIS	-	expression tag	UNP Q5SLG6
G	297	ALA	-	expression tag	UNP Q5SLG6
G	298	LYS	-	expression tag	UNP Q5SLG6
G	299	LEU	-	expression tag	UNP Q5SLG6
G	300	ALA	-	expression tag	UNP Q5SLG6
G	301	ALA	-	expression tag	UNP Q5SLG6
G	302	ALA	-	expression tag	UNP Q5SLG6
G	303	LEU	-	expression tag	UNP Q5SLG6
G	304	GLU	-	expression tag	UNP Q5SLG6
G	305	HIS	-	expression tag	UNP Q5SLG6
G	306	HIS	-	expression tag	UNP Q5SLG6
G	307	HIS	-	expression tag	UNP Q5SLG6
G	308	HIS	-	expression tag	UNP Q5SLG6
G	309	HIS	-	expression tag	UNP Q5SLG6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	310	HIS	-	expression tag	UNP Q5SLG6
H	297	ALA	-	expression tag	UNP Q5SLG6
H	298	LYS	-	expression tag	UNP Q5SLG6
H	299	LEU	-	expression tag	UNP Q5SLG6
H	300	ALA	-	expression tag	UNP Q5SLG6
H	301	ALA	-	expression tag	UNP Q5SLG6
H	302	ALA	-	expression tag	UNP Q5SLG6
H	303	LEU	-	expression tag	UNP Q5SLG6
H	304	GLU	-	expression tag	UNP Q5SLG6
H	305	HIS	-	expression tag	UNP Q5SLG6
H	306	HIS	-	expression tag	UNP Q5SLG6
H	307	HIS	-	expression tag	UNP Q5SLG6
H	308	HIS	-	expression tag	UNP Q5SLG6
H	309	HIS	-	expression tag	UNP Q5SLG6
H	310	HIS	-	expression tag	UNP Q5SLG6

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	D	1	Total	Cl	0	0
			1	1		
3	E	1	Total	Cl	0	0
			1	1		
3	F	1	Total	Cl	0	0
			1	1		
3	G	1	Total	Cl	0	0
			1	1		
3	H	1	Total	Cl	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	112	Total	O	0	0
			112	112		
4	B	111	Total	O	0	0
			111	111		
4	C	117	Total	O	0	0
			117	117		
4	D	97	Total	O	0	0
			97	97		
4	E	127	Total	O	0	0
			127	127		
4	F	119	Total	O	0	0
			119	119		
4	G	111	Total	O	0	0
			111	111		

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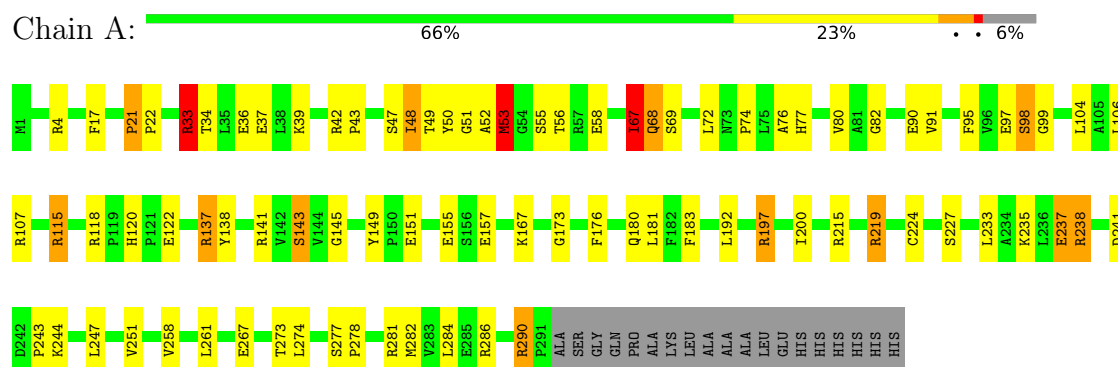
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	115	Total 115	O 115	0	0

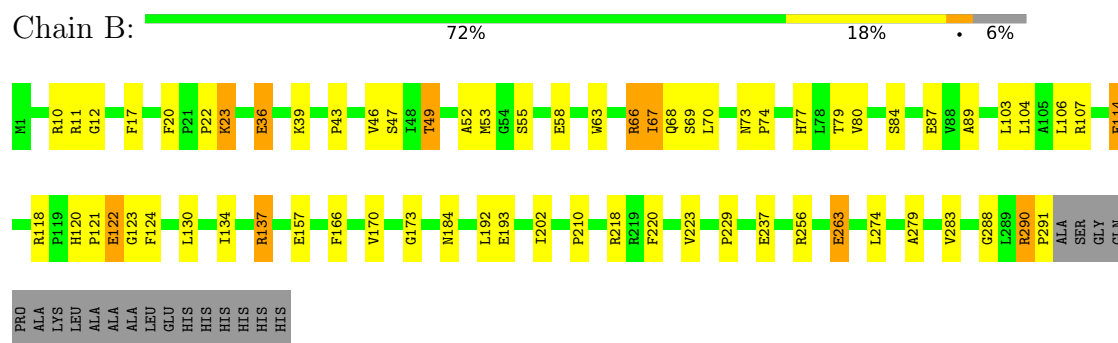
3 Residue-property plots [i](#)

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

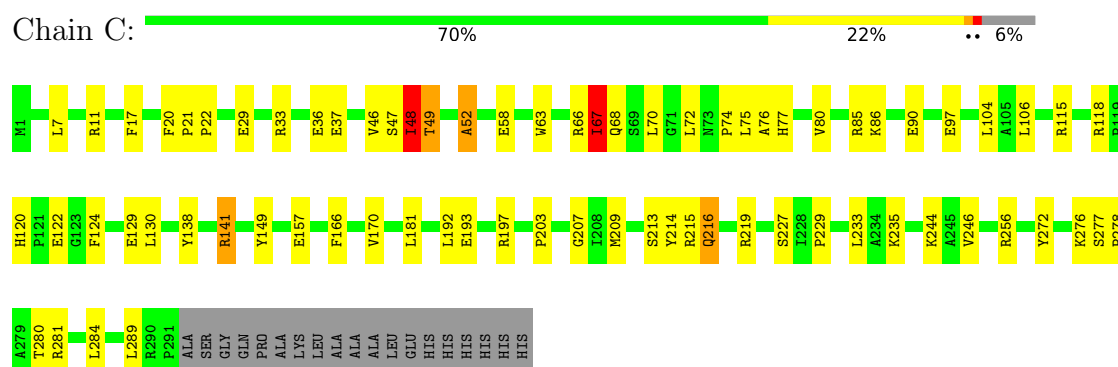
• Molecule 1: Methylenetetrahydrofolate reductase



• Molecule 1: Methylenetetrahydrofolate reductase

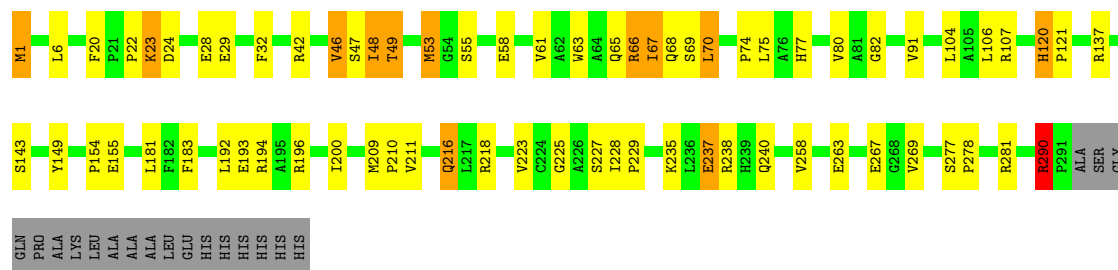


• Molecule 1: Methylenetetrahydrofolate reductase



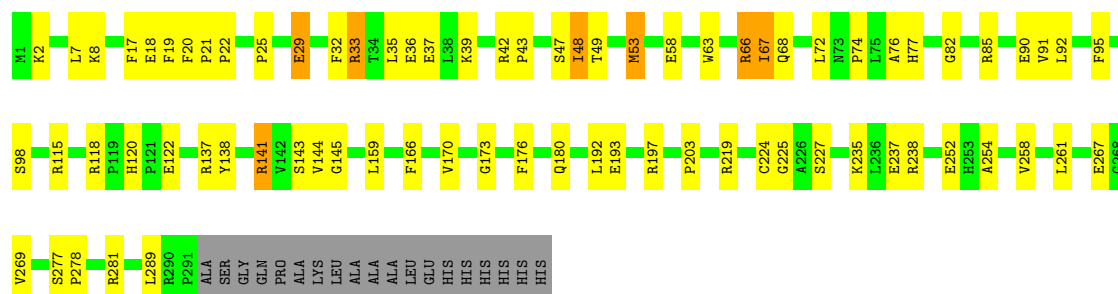
- Molecule 1: Methylenetetrahydrofolate reductase

Chain D:  71% 18% 6%



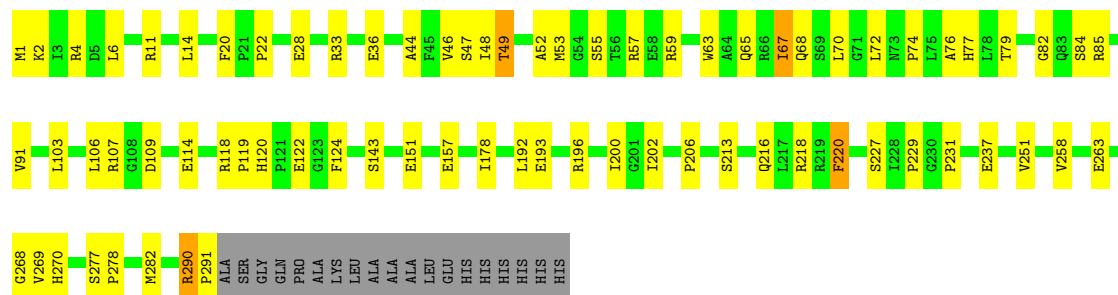
- Molecule 1: Methylenetetrahydrofolate reductase

Chain E: 69% 22% 6%

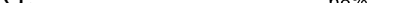


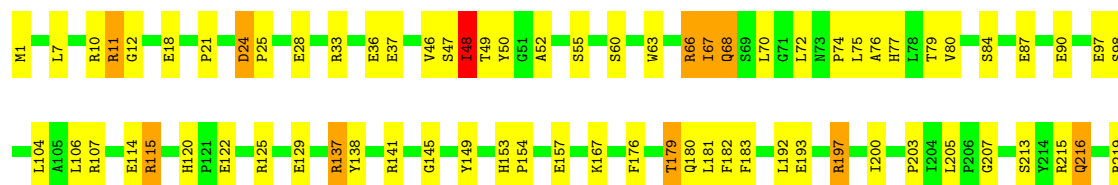
- Molecule 1: Methylenetetrahydrofolate reductase

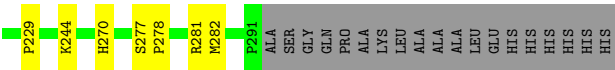
Chain F:  70% 23% • 6%



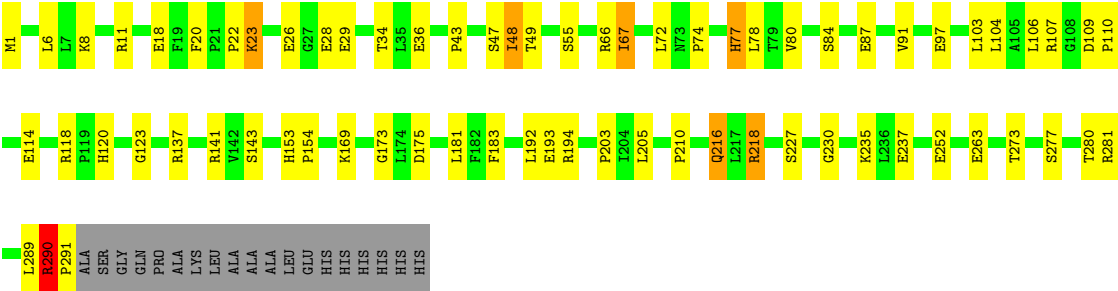
- Molecule 1: Methylenetetrahydrofolate reductase

Chain G:  68% 22% • 6%





● Molecule 1: Methylenetetrahydrofolate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	116.59Å 90.93Å 125.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.96 – 2.80 19.96 – 2.80	Depositor EDS
% Data completeness (in resolution range)	93.8 (19.96-2.80) 93.8 (19.96-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.64 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.204 , 0.306 0.203 , 0.291	Depositor DCC
R_{free} test set	3050 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	25.7	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 6.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.459 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19755	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 55.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0318e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.84	1/2357 (0.0%)	1.08	8/3185 (0.3%)
1	B	0.81	0/2357	1.02	3/3185 (0.1%)
1	C	0.82	1/2357 (0.0%)	1.04	8/3185 (0.3%)
1	D	0.83	0/2357	1.07	6/3185 (0.2%)
1	E	0.81	0/2357	1.01	2/3185 (0.1%)
1	F	0.80	0/2357	1.05	0/3185
1	G	0.84	0/2357	1.05	7/3185 (0.2%)
1	H	0.78	0/2357	1.03	4/3185 (0.1%)
All	All	0.82	2/18856 (0.0%)	1.04	38/25480 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	F	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	67	ILE	CA-CB	5.74	1.61	1.54
1	C	67	ILE	CA-CB	5.06	1.60	1.54

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	ALA	N-CA-C	-8.50	92.70	110.80
1	B	11	ARG	N-CA-C	-7.35	102.97	112.23
1	H	11	ARG	N-CA-C	-7.13	104.85	113.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	230	GLY	CA-C-N	-7.05	111.03	119.84
1	H	230	GLY	C-N-CA	-7.05	111.03	119.84
1	A	51	GLY	N-CA-C	-6.65	94.02	113.30
1	G	70	LEU	N-CA-C	-6.11	105.11	113.30
1	E	144	VAL	N-CA-C	6.07	116.61	108.11
1	C	70	LEU	N-CA-C	-6.02	105.91	113.20
1	B	229	PRO	CA-C-N	6.01	125.51	121.13
1	B	229	PRO	C-N-CA	6.01	125.51	121.13
1	G	48	ILE	N-CA-CB	5.94	118.11	110.99
1	A	33	ARG	N-CA-C	-5.89	104.77	111.07
1	C	149	TYR	CA-C-N	5.79	125.46	119.56
1	C	149	TYR	C-N-CA	5.79	125.46	119.56
1	G	21	PRO	CA-C-N	5.75	125.70	119.78
1	G	21	PRO	C-N-CA	5.75	125.70	119.78
1	A	53	MET	N-CA-C	-5.68	98.70	110.80
1	A	67	ILE	N-CA-C	5.67	116.44	110.72
1	D	120	HIS	CA-C-N	-5.62	113.10	119.28
1	D	120	HIS	C-N-CA	-5.62	113.10	119.28
1	C	48	ILE	N-CA-CB	5.61	117.72	110.99
1	A	67	ILE	CB-CA-C	5.60	119.95	112.22
1	E	33	ARG	N-CA-C	-5.36	105.34	111.07
1	G	137	ARG	N-CA-C	5.32	117.16	111.36
1	D	149	TYR	CA-C-N	5.30	125.61	119.47
1	D	149	TYR	C-N-CA	5.30	125.61	119.47
1	H	290	ARG	N-CA-C	5.28	121.48	109.81
1	C	33	ARG	N-CA-C	-5.18	105.53	111.07
1	C	46	VAL	CA-C-N	-5.15	115.46	122.93
1	C	46	VAL	C-N-CA	-5.15	115.46	122.93
1	A	21	PRO	CA-C-N	5.15	125.13	120.03
1	A	21	PRO	C-N-CA	5.15	125.13	120.03
1	G	24	ASP	CA-C-N	5.13	124.57	119.24
1	G	24	ASP	C-N-CA	5.13	124.57	119.24
1	D	290	ARG	N-CA-C	5.08	121.03	109.81
1	A	137	ARG	N-CA-C	5.04	118.69	112.54
1	D	240	GLN	N-CA-C	5.03	118.19	111.75

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	290	ARG	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2302	0	2319	71	0
1	B	2302	0	2319	46	0
1	C	2302	0	2319	47	0
1	D	2302	0	2319	47	0
1	E	2302	0	2319	55	0
1	F	2302	0	2319	46	0
1	G	2302	0	2319	66	0
1	H	2302	0	2319	41	0
2	A	53	0	31	4	0
2	B	53	0	31	2	0
2	C	53	0	31	1	0
2	D	53	0	31	1	0
2	E	53	0	31	0	0
2	F	53	0	31	2	0
2	G	53	0	31	3	0
2	H	53	0	31	1	0
3	A	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	1	0
3	H	1	0	0	0	0
4	A	112	0	0	6	0
4	B	111	0	0	5	0
4	C	117	0	0	7	0
4	D	97	0	0	4	0
4	E	127	0	0	7	0
4	F	119	0	0	5	0
4	G	111	0	0	3	0
4	H	115	0	0	8	0
All	All	19755	0	18800	399	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 11.

All (399) 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:D:67:ILE:HD11	1:D:74:PRO:HB3	1.36	1.04
1:A:290:ARG:HG2	1:A:290:ARG:HH11	1.24	0.99
1:H:290:ARG:O	1:H:290:ARG:HD3	1.61	0.99
1:A:52:ALA:HB2	1:A:56:THR:HG23	1.46	0.98
1:A:237:GLU:HG3	1:B:218:ARG:HH12	1.27	0.98
1:B:291:PRO:HG3	4:B:520:HOH:O	1.64	0.96
1:D:32:PHE:HZ	1:D:66:ARG:HD3	1.32	0.95
1:F:65:GLN:HA	1:F:68:GLN:HE21	1.32	0.92
1:A:238:ARG:HH11	1:A:238:ARG:HG2	1.36	0.91
1:D:32:PHE:CZ	1:D:66:ARG:HD3	2.07	0.89
1:A:53:MET:HG2	1:A:82:GLY:HA3	1.56	0.88
1:E:237:GLU:HG3	1:F:218:ARG:HH12	1.38	0.88
1:E:53:MET:HE3	1:E:82:GLY:HA3	1.56	0.88
1:G:120:HIS:HD2	1:G:122:GLU:H	1.22	0.87
1:A:50:TYR:O	2:A:311:FAD:N3	2.08	0.86
1:G:213:SER:HB3	1:G:216:GLN:HB3	1.58	0.86
1:H:67:ILE:HD11	1:H:74:PRO:HB3	1.56	0.86
1:E:120:HIS:HD2	1:E:122:GLU:H	1.19	0.85
1:D:67:ILE:HD11	1:D:74:PRO:CB	2.07	0.84
1:E:267:GLU:HG2	4:E:442:HOH:O	1.75	0.84
1:E:141:ARG:HH11	1:E:141:ARG:HB3	1.42	0.83
1:A:237:GLU:HG3	1:B:218:ARG:NH1	1.93	0.83
1:A:33:ARG:HG2	1:A:33:ARG:HH11	1.44	0.83
1:A:120:HIS:HD2	1:A:122:GLU:H	1.23	0.82
1:D:67:ILE:CD1	1:D:74:PRO:HB3	2.08	0.82
1:G:67:ILE:HD12	1:G:74:PRO:HB3	1.61	0.82
1:F:1:MET:HG2	1:F:6:LEU:HG	1.61	0.82
1:E:238:ARG:NH1	4:E:333:HOH:O	2.12	0.81
1:A:47:SER:HB2	1:A:77:HIS:HE1	1.44	0.81
1:B:89:ALA:HB1	1:B:137:ARG:HH22	1.45	0.81
1:A:290:ARG:HH11	1:A:290:ARG:CG	1.94	0.80
1:E:49:THR:HG22	4:E:893:HOH:O	1.82	0.79
1:A:52:ALA:HB3	1:A:55:SER:HB2	1.63	0.78
1:G:215:ARG:O	1:G:219:ARG:HG3	1.82	0.78
1:G:33:ARG:NH1	3:G:312:CL:CL	2.54	0.77
1:C:120:HIS:HD2	1:C:122:GLU:H	1.32	0.77
1:E:42:ARG:HA	4:E:965:HOH:O	1.84	0.77
1:F:52:ALA:C	1:F:53:MET:HG2	2.11	0.76
1:A:68:GLN:HG2	1:A:74:PRO:HG3	1.65	0.76
1:E:120:HIS:CD2	1:E:122:GLU:H	2.04	0.75
1:G:47:SER:HB3	1:G:77:HIS:CE1	2.22	0.74
1:H:291:PRO:HG3	4:H:340:HOH:O	1.86	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:ALA:HB1	1:B:137:ARG:NH2	2.03	0.74
1:A:120:HIS:CD2	1:A:122:GLU:H	2.05	0.73
1:A:115:ARG:HG2	4:B:984:HOH:O	1.87	0.73
1:A:47:SER:HB2	1:A:77:HIS:CE1	2.23	0.73
1:B:36:GLU:OE1	1:B:39:LYS:NZ	2.20	0.73
1:C:37:GLU:OE2	1:C:281:ARG:HD3	1.87	0.73
1:C:120:HIS:CD2	1:C:122:GLU:H	2.07	0.72
1:G:66:ARG:HG2	1:G:66:ARG:HH11	1.53	0.72
1:G:229:PRO:HA	1:H:227:SER:HB3	1.72	0.72
1:F:65:GLN:HA	1:F:68:GLN:NE2	2.05	0.71
1:B:23:LYS:HD3	4:B:323:HOH:O	1.91	0.71
1:B:73:ASN:OD1	4:B:349:HOH:O	2.08	0.71
4:A:617:HOH:O	1:B:263:GLU:HG2	1.90	0.70
1:B:63:TRP:O	1:B:67:ILE:HG23	1.91	0.70
1:F:216:GLN:O	1:F:220:PHE:HB2	1.91	0.70
1:G:104:LEU:HD23	1:G:106:LEU:HD11	1.75	0.68
1:G:149:TYR:HH	2:G:311:FAD:HO3'	1.40	0.68
1:F:63:TRP:O	1:F:67:ILE:HG23	1.94	0.68
1:H:84:SER:OG	1:H:87:GLU:HG3	1.94	0.68
1:A:238:ARG:HH11	1:A:238:ARG:CG	2.06	0.68
1:D:1:MET:HG2	1:D:6:LEU:HG	1.76	0.67
1:A:241:ASP:O	1:A:243:PRO:HD3	1.95	0.67
1:H:290:ARG:HD3	1:H:290:ARG:C	2.16	0.66
1:A:33:ARG:HH11	1:A:33:ARG:CG	2.08	0.66
1:E:32:PHE:HZ	1:E:66:ARG:HH11	1.43	0.65
1:C:48:ILE:HD12	1:C:76:ALA:HA	1.79	0.65
1:A:52:ALA:CB	1:A:55:SER:HB2	2.27	0.65
1:F:291:PRO:HG3	4:F:793:HOH:O	1.96	0.65
1:G:68:GLN:HG2	1:G:74:PRO:HG3	1.79	0.65
1:C:49:THR:HG22	4:C:378:HOH:O	1.96	0.65
1:B:120:HIS:HD2	1:B:122:GLU:H	1.42	0.64
1:E:68:GLN:HG2	1:E:74:PRO:HG3	1.78	0.64
1:A:290:ARG:HG2	1:A:290:ARG:NH1	2.00	0.64
1:A:47:SER:CB	1:A:77:HIS:HE1	2.11	0.63
1:C:256:ARG:NE	4:C:954:HOH:O	2.30	0.63
1:E:47:SER:HB3	1:E:77:HIS:CE1	2.33	0.63
1:A:238:ARG:HG2	1:A:238:ARG:NH1	2.10	0.63
1:C:235:LYS:HE3	4:C:866:HOH:O	1.98	0.62
1:A:17:PHE:CE1	1:A:43:PRO:HG3	2.33	0.62
1:A:107:ARG:HB2	2:A:311:FAD:O1A	1.98	0.62
1:H:1:MET:HG3	1:H:6:LEU:HG	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:33:ARG:NH1	1:G:281:ARG:HD2	2.15	0.62
1:F:251:VAL:HG21	1:F:282:MET:HE2	1.82	0.62
1:G:63:TRP:O	1:G:67:ILE:HG22	2.00	0.62
1:A:151:GLU:HB3	1:B:184:ASN:ND2	2.15	0.61
1:D:63:TRP:O	1:D:67:ILE:HG23	2.01	0.61
1:G:66:ARG:HH11	1:G:66:ARG:CG	2.14	0.61
1:F:2:LYS:HD3	1:F:4:ARG:NH2	2.16	0.61
1:E:141:ARG:HB3	1:E:141:ARG:NH1	2.12	0.61
1:A:34:THR:HG23	1:A:277:SER:HB3	1.82	0.61
1:D:67:ILE:HD11	1:D:74:PRO:CA	2.31	0.61
1:G:167:LYS:HB2	1:G:200:ILE:HD11	1.83	0.61
1:A:52:ALA:CB	1:A:56:THR:HG23	2.27	0.60
1:A:17:PHE:HE1	1:A:43:PRO:HG3	1.66	0.60
1:A:67:ILE:HG12	1:A:72:LEU:HB2	1.82	0.60
1:C:215:ARG:O	1:C:219:ARG:HG3	2.02	0.60
1:E:18:GLU:HA	1:E:47:SER:HB2	1.84	0.60
1:G:278:PRO:O	1:G:282:MET:HG3	2.01	0.60
1:A:167:LYS:HB2	1:A:200:ILE:HD11	1.84	0.60
1:G:179:THR:HG23	1:G:180:GLN:O	2.02	0.59
1:D:290:ARG:CD	4:D:914:HOH:O	2.51	0.59
1:E:237:GLU:HG3	1:F:218:ARG:NH1	2.14	0.59
1:C:47:SER:HB2	1:C:77:HIS:CE1	2.38	0.58
1:F:85:ARG:NE	4:F:592:HOH:O	2.36	0.58
1:G:67:ILE:CD1	1:G:74:PRO:HB3	2.32	0.58
1:A:149:TYR:OH	2:A:311:FAD:H5'2	2.04	0.58
1:G:145:GLY:HA2	1:G:176:PHE:O	2.03	0.58
1:E:37:GLU:OE1	1:E:281:ARG:HD3	2.04	0.57
1:G:120:HIS:CD2	1:G:122:GLU:H	2.13	0.57
1:H:169:LYS:NZ	4:H:452:HOH:O	2.36	0.57
1:H:290:ARG:O	1:H:290:ARG:CD	2.46	0.57
1:D:47:SER:OG	1:D:77:HIS:HE1	1.87	0.57
1:G:66:ARG:HG2	1:G:66:ARG:NH1	2.18	0.57
1:H:175:ASP:O	1:H:203:PRO:HD2	2.05	0.57
1:D:24:ASP:O	1:D:28:GLU:HG3	2.04	0.57
1:E:48:ILE:HD12	1:E:76:ALA:HA	1.86	0.57
1:G:37:GLU:OE2	1:G:281:ARG:HD3	2.04	0.57
1:E:29:GLU:HG3	4:E:580:HOH:O	2.04	0.57
1:A:282:MET:O	1:A:286:ARG:HG3	2.05	0.56
1:A:278:PRO:O	1:A:282:MET:HG3	2.06	0.56
1:A:48:ILE:HD12	1:A:76:ALA:HA	1.86	0.56
1:E:25:PRO:O	1:E:29:GLU:HG2	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:22:PRO:HG2	1:D:28:GLU:HA	1.88	0.56
1:G:115:ARG:H	1:G:115:ARG:CZ	2.19	0.56
1:E:138:TYR:O	1:E:141:ARG:HG3	2.06	0.56
1:G:67:ILE:HD13	1:G:67:ILE:C	2.31	0.56
1:B:67:ILE:C	1:B:67:ILE:HD12	2.30	0.56
1:B:130:LEU:O	1:B:134:ILE:HG13	2.06	0.56
1:H:47:SER:HB3	1:H:77:HIS:CE1	2.42	0.55
1:A:233:LEU:O	1:A:237:GLU:HB2	2.07	0.55
1:E:166:PHE:O	1:E:170:VAL:HG23	2.06	0.55
1:D:65:GLN:HA	1:D:68:GLN:HE21	1.71	0.55
1:H:22:PRO:HG2	1:H:28:GLU:HA	1.89	0.55
1:B:120:HIS:CD2	1:B:122:GLU:H	2.22	0.55
1:A:258:VAL:HA	1:A:261:LEU:HD12	1.89	0.55
1:C:229:PRO:HA	1:D:227:SER:HB3	1.89	0.54
1:E:47:SER:HB3	1:E:77:HIS:HE1	1.72	0.54
1:A:37:GLU:OE1	1:A:281:ARG:HD3	2.07	0.54
1:C:63:TRP:O	1:C:67:ILE:HG22	2.08	0.54
1:G:179:THR:HG22	1:G:207:GLY:H	1.70	0.54
1:B:10:ARG:HB2	1:B:12:GLY:O	2.07	0.54
1:E:120:HIS:HD2	1:E:122:GLU:N	1.99	0.54
1:F:106:LEU:HD13	2:F:311:FAD:C4X	2.37	0.54
1:A:95:PHE:O	1:A:98:SER:HB2	2.07	0.54
1:H:216:GLN:HG3	4:H:400:HOH:O	2.07	0.54
1:D:196:ARG:HD3	1:D:200:ILE:O	2.08	0.54
1:E:197:ARG:CZ	1:F:193:GLU:OE1	2.56	0.54
1:B:103:LEU:HD13	1:B:134:ILE:HD13	1.90	0.53
1:C:214:TYR:HE1	1:C:233:LEU:HD12	1.72	0.53
1:F:107:ARG:HB3	2:F:311:FAD:H5'2	1.90	0.53
1:D:106:LEU:HD13	2:D:311:FAD:C4X	2.37	0.53
1:A:273:THR:O	1:A:274:LEU:HB2	2.07	0.53
1:C:284:LEU:HB3	1:C:289:LEU:HD22	1.89	0.53
1:E:29:GLU:HB2	1:E:33:ARG:NH2	2.23	0.53
1:H:23:LYS:HG3	4:H:316:HOH:O	2.08	0.53
1:D:216:GLN:HG3	4:D:603:HOH:O	2.09	0.53
1:E:197:ARG:HH12	1:F:193:GLU:HB3	1.74	0.53
1:G:80:VAL:HG22	1:G:107:ARG:HA	1.91	0.53
1:A:197:ARG:NE	1:B:193:GLU:OE1	2.41	0.52
1:A:80:VAL:HG22	1:A:107:ARG:HA	1.92	0.52
1:B:120:HIS:HB3	1:B:123:GLY:HA3	1.91	0.52
1:F:20:PHE:HA	1:F:49:THR:HB	1.91	0.52
1:F:76:ALA:HB3	1:F:103:LEU:HD23	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:SER:O	4:B:930:HOH:O	2.18	0.52
1:B:106:LEU:HD13	2:B:311:FAD:C4X	2.40	0.52
1:H:106:LEU:HD13	2:H:311:FAD:C4X	2.39	0.52
1:H:210:PRO:HD2	4:H:505:HOH:O	2.09	0.52
1:F:178:ILE:HG21	1:F:270:HIS:CD2	2.45	0.52
1:G:181:LEU:HD13	1:G:183:PHE:CE2	2.45	0.51
1:E:2:LYS:HE2	1:E:141:ARG:O	2.11	0.51
1:B:52:ALA:C	1:B:53:MET:HG2	2.36	0.51
1:D:32:PHE:HE2	1:D:66:ARG:HH11	1.57	0.51
1:E:48:ILE:CD1	1:E:76:ALA:HA	2.40	0.51
1:F:85:ARG:HD3	1:F:124:PHE:CE2	2.46	0.51
1:H:109:ASP:HB3	1:H:110:PRO:HD2	1.93	0.51
1:A:267:GLU:HG2	4:A:320:HOH:O	2.11	0.51
1:D:80:VAL:HG22	1:D:107:ARG:HA	1.93	0.51
1:D:48:ILE:HD12	1:D:48:ILE:H	1.76	0.51
1:C:20:PHE:HA	1:C:49:THR:HB	1.92	0.51
1:D:277:SER:N	1:D:278:PRO:CD	2.74	0.51
1:B:80:VAL:HG22	1:B:107:ARG:HA	1.92	0.50
1:F:151:GLU:OE1	1:F:227:SER:OG	2.29	0.50
1:G:138:TYR:O	1:G:141:ARG:HG2	2.11	0.50
1:C:67:ILE:HG12	1:C:72:LEU:HB2	1.93	0.50
1:D:46:VAL:HG12	1:D:67:ILE:HD12	1.93	0.50
1:E:66:ARG:NE	4:E:326:HOH:O	2.44	0.50
1:D:181:LEU:HD12	1:D:181:LEU:C	2.36	0.50
1:D:65:GLN:HA	1:D:68:GLN:NE2	2.27	0.50
1:D:181:LEU:HD12	1:D:181:LEU:O	2.11	0.50
1:D:68:GLN:HG3	1:D:74:PRO:HG3	1.94	0.50
1:H:55:SER:HB2	4:H:334:HOH:O	2.11	0.50
1:H:218:ARG:HB3	1:H:218:ARG:HH11	1.76	0.50
1:A:181:LEU:CD1	1:A:183:PHE:CE2	2.95	0.49
1:B:89:ALA:CB	1:B:137:ARG:NH2	2.72	0.49
1:C:227:SER:HB3	1:D:229:PRO:HA	1.93	0.49
1:G:79:THR:HA	1:G:106:LEU:O	2.12	0.49
1:C:213:SER:HB3	1:C:216:GLN:HB3	1.95	0.49
1:G:18:GLU:HB2	1:G:270:HIS:NE2	2.27	0.49
1:H:104:LEU:HD23	1:H:106:LEU:HD11	1.94	0.49
1:A:180:GLN:NE2	1:A:224:CYS:HB2	2.28	0.49
1:F:79:THR:HA	1:F:106:LEU:O	2.12	0.49
1:H:290:ARG:C	1:H:290:ARG:CD	2.84	0.49
1:D:104:LEU:HD23	1:D:106:LEU:HD11	1.95	0.49
1:E:19:PHE:CE2	1:E:35:LEU:HD21	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:181:LEU:HA	1:G:207:GLY:O	2.12	0.49
1:C:67:ILE:HD13	1:C:67:ILE:C	2.38	0.49
1:H:18:GLU:O	1:H:273:THR:OG1	2.26	0.49
1:G:46:VAL:HG21	1:G:72:LEU:HD13	1.93	0.49
1:G:213:SER:CB	1:G:216:GLN:HB3	2.37	0.49
1:G:24:ASP:HB2	1:G:25:PRO:HD2	1.94	0.48
1:C:37:GLU:OE1	1:C:277:SER:OG	2.23	0.48
1:E:227:SER:HB2	1:F:227:SER:HB2	1.95	0.48
1:C:181:LEU:HA	1:C:207:GLY:O	2.12	0.48
1:C:193:GLU:HB3	4:C:318:HOH:O	2.12	0.48
1:G:179:THR:HG21	1:G:182:PHE:HE1	1.79	0.48
1:A:68:GLN:HG2	1:A:74:PRO:CG	2.38	0.48
1:A:120:HIS:HD2	1:A:122:GLU:N	2.03	0.48
1:C:67:ILE:CD1	1:C:74:PRO:HB3	2.44	0.48
1:E:254:ALA:O	1:E:258:VAL:HG23	2.13	0.48
1:C:104:LEU:HD23	1:C:106:LEU:HD11	1.95	0.48
1:H:48:ILE:HD12	1:H:48:ILE:H	1.79	0.48
1:G:18:GLU:HA	1:G:47:SER:HB2	1.95	0.48
1:G:67:ILE:HG12	1:G:72:LEU:HB2	1.95	0.48
1:A:215:ARG:O	1:A:219:ARG:HG2	2.14	0.48
1:B:79:THR:HA	1:B:106:LEU:O	2.14	0.47
1:C:246:VAL:HG21	4:C:329:HOH:O	2.14	0.47
1:D:67:ILE:HD11	1:D:74:PRO:HA	1.95	0.47
1:H:26:GLU:H	1:H:26:GLU:CD	2.22	0.47
1:F:258:VAL:HG13	1:F:269:VAL:HG21	1.96	0.47
1:A:104:LEU:HD12	1:A:145:GLY:C	2.39	0.47
1:B:80:VAL:HB	1:B:124:PHE:HB2	1.95	0.47
1:H:34:THR:HG23	1:H:277:SER:HB3	1.97	0.47
1:A:138:TYR:O	1:A:141:ARG:HG3	2.13	0.47
1:B:22:PRO:HD2	1:B:63:TRP:NE1	2.29	0.47
1:G:60:SER:HB3	4:G:467:HOH:O	2.13	0.47
1:B:84:SER:OG	1:B:87:GLU:HG3	2.15	0.47
1:A:97:GLU:HB3	4:A:704:HOH:O	2.15	0.47
1:D:53:MET:HG3	1:D:82:GLY:HA3	1.95	0.47
1:H:216:GLN:CG	4:H:400:HOH:O	2.63	0.47
1:C:17:PHE:CE2	1:C:280:THR:HG21	2.50	0.47
1:C:68:GLN:HG2	1:C:74:PRO:HG3	1.97	0.46
1:C:85:ARG:HD3	1:C:124:PHE:CE2	2.50	0.46
1:A:21:PRO:HA	1:A:22:PRO:HD3	1.75	0.46
1:B:220:PHE:CZ	1:B:274:LEU:HD11	2.50	0.46
1:A:4:ARG:HB2	1:A:143:SER:OG	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:67:ILE:HD11	1:H:74:PRO:CB	2.36	0.46
1:A:247:LEU:O	1:A:251:VAL:HG23	2.15	0.46
1:G:52:ALA:O	1:G:55:SER:HB2	2.15	0.46
1:F:85:ARG:NH1	4:F:682:HOH:O	2.48	0.46
1:E:95:PHE:O	1:E:98:SER:HB2	2.16	0.46
1:F:22:PRO:HD2	1:F:63:TRP:NE1	2.30	0.46
1:B:166:PHE:O	1:B:170:VAL:HG23	2.16	0.46
1:F:47:SER:HB3	1:F:77:HIS:CE1	2.51	0.46
1:C:120:HIS:HD2	1:C:122:GLU:N	2.07	0.46
1:B:20:PHE:HA	1:B:49:THR:HB	1.97	0.46
1:D:211:VAL:HG23	4:D:336:HOH:O	2.16	0.46
1:B:220:PHE:HZ	1:B:274:LEU:HD11	1.80	0.45
1:C:77:HIS:HD2	2:C:311:FAD:N5	2.14	0.45
1:F:22:PRO:O	1:F:59:ARG:HD3	2.15	0.45
1:F:67:ILE:HG12	1:F:74:PRO:HB3	1.98	0.45
1:A:197:ARG:HH11	1:A:197:ARG:CG	2.28	0.45
1:B:279:ALA:O	1:B:283:VAL:HG23	2.16	0.45
1:C:166:PHE:O	1:C:170:VAL:HG23	2.16	0.45
1:D:237:GLU:HG3	1:D:238:ARG:N	2.30	0.45
1:G:193:GLU:HG3	1:H:194:ARG:HG2	1.98	0.45
1:B:288:GLY:C	1:B:290:ARG:H	2.24	0.45
1:F:52:ALA:O	1:F:53:MET:HG2	2.16	0.45
1:G:277:SER:N	1:G:278:PRO:CD	2.78	0.45
1:A:145:GLY:HA2	1:A:176:PHE:O	2.15	0.45
1:E:237:GLU:CG	1:F:218:ARG:HH12	2.21	0.45
1:G:28:GLU:HG2	1:G:63:TRP:HZ2	1.81	0.45
1:E:85:ARG:HG3	1:E:122:GLU:O	2.17	0.45
1:G:179:THR:CG2	1:G:180:GLN:O	2.65	0.45
1:G:50:TYR:O	2:G:311:FAD:N3	2.43	0.45
1:H:74:PRO:HD2	4:H:601:HOH:O	2.17	0.45
1:D:290:ARG:HD3	4:D:914:HOH:O	2.17	0.45
1:B:17:PHE:CE1	1:B:43:PRO:HG3	2.52	0.45
1:B:106:LEU:HD13	2:B:311:FAD:N5	2.32	0.45
1:C:47:SER:HB3	1:C:75:LEU:HB3	1.98	0.45
1:C:86:LYS:HB2	4:C:873:HOH:O	2.17	0.45
1:D:20:PHE:HA	1:D:49:THR:HB	1.99	0.45
1:F:196:ARG:HD3	1:F:200:ILE:O	2.17	0.44
1:E:227:SER:HB3	1:F:229:PRO:HA	1.99	0.44
1:F:14:LEU:HD22	1:F:44:ALA:CB	2.48	0.44
1:G:181:LEU:CD1	1:G:183:PHE:CE2	3.01	0.44
1:A:52:ALA:HB3	1:A:55:SER:CB	2.41	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:ARG:HA	1:B:291:PRO:HD2	1.64	0.44
1:D:281:ARG:HG2	1:D:281:ARG:HH11	1.82	0.44
1:F:67:ILE:HD12	1:F:72:LEU:HB2	1.99	0.44
1:C:47:SER:HB2	1:C:77:HIS:HE1	1.82	0.44
1:C:244:LYS:O	1:C:244:LYS:HG2	2.16	0.44
1:F:120:HIS:HD2	1:F:122:GLU:H	1.65	0.44
1:A:68:GLN:HE22	1:A:99:GLY:HA3	1.83	0.44
1:C:193:GLU:HG3	1:D:194:ARG:HG2	1.98	0.44
1:C:209:MET:HG3	1:C:272:TYR:HB2	2.00	0.44
1:D:183:PHE:HB2	1:D:229:PRO:HD3	2.00	0.44
1:G:149:TYR:OH	2:G:311:FAD:O3'	2.12	0.44
1:C:124:PHE:CD2	1:C:129:GLU:HB3	2.53	0.44
1:E:47:SER:CB	1:E:77:HIS:HE1	2.31	0.44
1:E:72:LEU:HD23	1:E:72:LEU:HA	1.79	0.44
1:F:120:HIS:CD2	1:F:122:GLU:H	2.36	0.44
1:E:258:VAL:HA	1:E:261:LEU:HD12	2.00	0.44
1:G:197:ARG:HG3	1:H:193:GLU:OE1	2.18	0.44
1:B:47:SER:OG	1:B:77:HIS:HE1	2.01	0.44
1:F:277:SER:N	1:F:278:PRO:CD	2.81	0.44
1:H:289:LEU:O	1:H:290:ARG:HD3	2.18	0.44
1:C:7:LEU:HD13	1:C:203:PRO:HG2	2.00	0.43
1:G:47:SER:HB3	1:G:77:HIS:HE1	1.76	0.43
1:H:43:PRO:HG2	1:H:72:LEU:HD22	1.99	0.43
1:H:120:HIS:HB3	1:H:123:GLY:HA3	1.99	0.43
1:D:23:LYS:HG3	1:D:24:ASP:H	1.83	0.43
1:D:258:VAL:HG13	1:D:269:VAL:HG21	2.01	0.43
1:E:21:PRO:HA	1:E:22:PRO:HD3	1.76	0.43
1:G:7:LEU:HD13	1:G:203:PRO:HG2	1.99	0.43
1:A:151:GLU:OE1	1:A:227:SER:OG	2.16	0.43
1:F:2:LYS:HD3	1:F:4:ARG:HH21	1.82	0.43
1:F:109:ASP:HB3	4:F:406:HOH:O	2.17	0.43
1:B:23:LYS:HD3	1:B:23:LYS:H	1.84	0.43
1:B:67:ILE:HD11	1:B:74:PRO:N	2.32	0.43
1:B:68:GLN:HG3	1:B:74:PRO:HG3	2.00	0.43
1:E:66:ARG:CD	4:E:326:HOH:O	2.66	0.43
1:C:197:ARG:HG3	1:D:193:GLU:OE1	2.19	0.43
1:G:114:GLU:HA	1:G:115:ARG:NH1	2.34	0.43
1:B:52:ALA:O	1:B:53:MET:HG2	2.18	0.43
1:E:37:GLU:OE2	1:E:277:SER:OG	2.21	0.43
1:G:125:ARG:N	1:G:129:GLU:OE1	2.49	0.43
1:G:179:THR:HG21	1:G:182:PHE:CE1	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:277:SER:N	1:E:278:PRO:CD	2.82	0.43
1:D:209:MET:HA	1:D:210:PRO:HD3	1.84	0.43
1:E:145:GLY:HA2	1:E:176:PHE:O	2.19	0.43
1:G:63:TRP:O	1:G:67:ILE:CG2	2.67	0.42
1:C:213:SER:CB	1:C:216:GLN:HB3	2.49	0.42
1:G:84:SER:OG	1:G:87:GLU:HG3	2.18	0.42
1:C:138:TYR:O	1:C:141:ARG:HB3	2.19	0.42
1:E:17:PHE:CE1	1:E:43:PRO:HG3	2.54	0.42
1:G:48:ILE:HD12	1:G:76:ALA:HA	2.01	0.42
1:A:277:SER:N	1:A:278:PRO:CD	2.82	0.42
1:F:118:ARG:HA	1:F:119:PRO:HD2	1.84	0.42
1:E:180:GLN:NE2	1:E:224:CYS:HB2	2.34	0.42
1:E:35:LEU:O	1:E:39:LYS:HG2	2.20	0.42
1:G:244:LYS:HE3	4:G:551:HOH:O	2.20	0.42
1:A:181:LEU:HD13	1:A:183:PHE:CZ	2.55	0.42
1:D:154:PRO:HG3	1:D:225:GLY:HA3	2.01	0.42
1:E:35:LEU:HD23	1:E:35:LEU:HA	1.86	0.42
1:H:20:PHE:HA	1:H:49:THR:HB	2.01	0.42
1:H:109:ASP:HB3	1:H:110:PRO:CD	2.49	0.42
1:B:66:ARG:HA	1:B:66:ARG:HD2	1.68	0.41
1:C:21:PRO:HA	1:C:22:PRO:HD3	1.86	0.41
1:E:20:PHE:HA	1:E:49:THR:HB	2.01	0.41
1:A:53:MET:H	1:A:55:SER:H	1.68	0.41
1:G:179:THR:HB	1:G:205:LEU:O	2.20	0.41
1:A:106:LEU:HD13	2:A:311:FAD:C4X	2.50	0.41
1:A:197:ARG:CG	1:A:197:ARG:NH1	2.83	0.41
1:H:47:SER:OG	1:H:77:HIS:HE1	2.03	0.41
1:H:181:LEU:HD13	1:H:183:PHE:CE2	2.54	0.41
1:A:215:ARG:NH2	4:A:546:HOH:O	2.52	0.41
1:B:210:PRO:HB2	1:B:279:ALA:HB1	2.02	0.41
1:H:80:VAL:HG22	1:H:107:ARG:HA	2.01	0.41
1:A:115:ARG:HG3	4:A:981:HOH:O	2.20	0.41
1:C:277:SER:N	1:C:278:PRO:CD	2.83	0.41
1:E:141:ARG:NH1	1:E:141:ARG:CB	2.83	0.41
1:F:14:LEU:HD22	1:F:44:ALA:HB2	2.01	0.41
1:H:280:THR:O	1:H:281:ARG:C	2.62	0.41
1:D:70:LEU:HD12	1:D:70:LEU:HA	1.94	0.41
1:E:7:LEU:HD13	1:E:203:PRO:HG2	2.03	0.41
1:F:82:GLY:HA2	4:F:753:HOH:O	2.21	0.41
1:C:215:ARG:HE	1:C:219:ARG:HD3	1.86	0.41
1:D:48:ILE:HD12	1:D:75:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:47:SER:CB	1:G:77:HIS:HE1	2.33	0.41
1:A:33:ARG:CG	1:A:33:ARG:NH1	2.76	0.41
1:C:63:TRP:O	1:C:67:ILE:CG2	2.69	0.41
1:E:258:VAL:HG13	1:E:269:VAL:HG21	2.03	0.41
1:A:39:LYS:NZ	4:A:595:HOH:O	2.53	0.41
1:A:52:ALA:HB2	1:A:56:THR:CG2	2.33	0.41
1:B:104:LEU:HD23	1:B:106:LEU:HD11	2.03	0.41
1:C:80:VAL:CG1	1:C:130:LEU:HB2	2.51	0.41
1:G:1:MET:N	4:G:544:HOH:O	2.53	0.41
1:G:10:ARG:HB2	1:G:12:GLY:O	2.21	0.41
1:G:11:ARG:HG2	1:G:11:ARG:HH11	1.86	0.41
1:G:47:SER:HA	1:G:75:LEU:O	2.21	0.41
1:H:289:LEU:O	1:H:290:ARG:O	2.39	0.41
1:B:120:HIS:CG	1:B:121:PRO:HD2	2.56	0.41
1:F:206:PRO:HD2	1:F:268:GLY:O	2.21	0.41
1:G:114:GLU:HG2	1:G:115:ARG:HH12	1.85	0.41
1:H:153:HIS:HA	1:H:154:PRO:HD3	1.93	0.41
1:A:284:LEU:HD23	1:A:284:LEU:HA	1.88	0.40
1:C:118:ARG:NH1	4:C:862:HOH:O	2.54	0.40
1:D:277:SER:H	1:D:278:PRO:CD	2.34	0.40
1:F:22:PRO:HB2	1:F:28:GLU:HG2	2.03	0.40
1:G:153:HIS:HA	1:G:154:PRO:HD2	1.87	0.40
1:E:63:TRP:O	1:E:67:ILE:HG23	2.21	0.40
1:E:225:GLY:HA2	1:F:231:PRO:HD3	2.03	0.40
1:D:120:HIS:O	1:D:121:PRO:C	2.63	0.40
1:A:107:ARG:HH22	1:A:155:GLU:CD	2.30	0.40
1:D:277:SER:N	1:D:278:PRO:HD2	2.36	0.40
1:G:24:ASP:HB2	1:G:25:PRO:CD	2.51	0.40
1:G:215:ARG:HE	1:G:219:ARG:HD3	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	289/310 (93%)	280 (97%)	8 (3%)	1 (0%)	36	66
1	B	289/310 (93%)	278 (96%)	8 (3%)	3 (1%)	12	38
1	C	289/310 (93%)	275 (95%)	13 (4%)	1 (0%)	36	66
1	D	289/310 (93%)	274 (95%)	13 (4%)	2 (1%)	18	47
1	E	289/310 (93%)	277 (96%)	11 (4%)	1 (0%)	36	66
1	F	289/310 (93%)	275 (95%)	12 (4%)	2 (1%)	18	47
1	G	289/310 (93%)	272 (94%)	17 (6%)	0	100	100
1	H	289/310 (93%)	279 (96%)	8 (3%)	2 (1%)	18	47
All	All	2312/2480 (93%)	2210 (96%)	90 (4%)	12 (0%)	24	55

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	290	ARG
1	C	52	ALA
1	D	290	ARG
1	H	290	ARG
1	B	114	GLU
1	F	290	ARG
1	A	173	GLY
1	F	57	ARG
1	D	61	VAL
1	B	173	GLY
1	E	173	GLY
1	H	173	GLY

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	236/249 (95%)	210 (89%)	26 (11%)	6	20
1	B	236/249 (95%)	216 (92%)	20 (8%)	10	31
1	C	236/249 (95%)	220 (93%)	16 (7%)	14	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	236/249 (95%)	208 (88%)	28 (12%)	5	17
1	E	236/249 (95%)	213 (90%)	23 (10%)	8	25
1	F	236/249 (95%)	216 (92%)	20 (8%)	10	31
1	G	236/249 (95%)	219 (93%)	17 (7%)	13	39
1	H	236/249 (95%)	210 (89%)	26 (11%)	6	20
All	All	1888/1992 (95%)	1712 (91%)	176 (9%)	8	27

All (176) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	ARG
1	A	36	GLU
1	A	42	ARG
1	A	48	ILE
1	A	49	THR
1	A	53	MET
1	A	58	GLU
1	A	67	ILE
1	A	68	GLN
1	A	69	SER
1	A	90	GLU
1	A	91	VAL
1	A	98	SER
1	A	115	ARG
1	A	118	ARG
1	A	137	ARG
1	A	143	SER
1	A	157	GLU
1	A	192	LEU
1	A	197	ARG
1	A	219	ARG
1	A	235	LYS
1	A	237	GLU
1	A	238	ARG
1	A	244	LYS
1	A	290	ARG
1	B	23	LYS
1	B	36	GLU
1	B	46	VAL
1	B	49	THR

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Mol	Chain	Res	Type
1	B	55	SER
1	B	58	GLU
1	B	66	ARG
1	B	67	ILE
1	B	70	LEU
1	B	114	GLU
1	B	118	ARG
1	B	122	GLU
1	B	137	ARG
1	B	157	GLU
1	B	192	LEU
1	B	202	ILE
1	B	223	VAL
1	B	237	GLU
1	B	256	ARG
1	B	263	GLU
1	C	11	ARG
1	C	29	GLU
1	C	36	GLU
1	C	48	ILE
1	C	49	THR
1	C	58	GLU
1	C	66	ARG
1	C	67	ILE
1	C	90	GLU
1	C	97	GLU
1	C	115	ARG
1	C	141	ARG
1	C	157	GLU
1	C	192	LEU
1	C	216	GLN
1	C	276	LYS
1	D	1	MET
1	D	23	LYS
1	D	29	GLU
1	D	42	ARG
1	D	46	VAL
1	D	48	ILE
1	D	49	THR
1	D	53	MET
1	D	55	SER
1	D	58	GLU

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Mol	Chain	Res	Type
1	D	66	ARG
1	D	67	ILE
1	D	69	SER
1	D	70	LEU
1	D	91	VAL
1	D	137	ARG
1	D	143	SER
1	D	155	GLU
1	D	192	LEU
1	D	216	GLN
1	D	218	ARG
1	D	223	VAL
1	D	228	ILE
1	D	235	LYS
1	D	237	GLU
1	D	263	GLU
1	D	267	GLU
1	D	290	ARG
1	E	8	LYS
1	E	29	GLU
1	E	36	GLU
1	E	48	ILE
1	E	53	MET
1	E	58	GLU
1	E	66	ARG
1	E	67	ILE
1	E	90	GLU
1	E	91	VAL
1	E	92	LEU
1	E	115	ARG
1	E	118	ARG
1	E	137	ARG
1	E	141	ARG
1	E	143	SER
1	E	159	LEU
1	E	192	LEU
1	E	193	GLU
1	E	219	ARG
1	E	235	LYS
1	E	252	GLU
1	E	289	LEU
1	F	11	ARG

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Mol	Chain	Res	Type
1	F	33	ARG
1	F	36	GLU
1	F	46	VAL
1	F	48	ILE
1	F	49	THR
1	F	55	SER
1	F	67	ILE
1	F	70	LEU
1	F	84	SER
1	F	91	VAL
1	F	114	GLU
1	F	143	SER
1	F	157	GLU
1	F	192	LEU
1	F	202	ILE
1	F	213	SER
1	F	220	PHE
1	F	237	GLU
1	F	263	GLU
1	G	11	ARG
1	G	36	GLU
1	G	48	ILE
1	G	49	THR
1	G	66	ARG
1	G	67	ILE
1	G	68	GLN
1	G	90	GLU
1	G	97	GLU
1	G	98	SER
1	G	115	ARG
1	G	137	ARG
1	G	157	GLU
1	G	179	THR
1	G	192	LEU
1	G	197	ARG
1	G	216	GLN
1	H	8	LYS
1	H	23	LYS
1	H	29	GLU
1	H	36	GLU
1	H	48	ILE
1	H	66	ARG

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Mol	Chain	Res	Type
1	H	67	ILE
1	H	77	HIS
1	H	78	LEU
1	H	91	VAL
1	H	97	GLU
1	H	103	LEU
1	H	114	GLU
1	H	118	ARG
1	H	137	ARG
1	H	141	ARG
1	H	143	SER
1	H	192	LEU
1	H	205	LEU
1	H	216	GLN
1	H	218	ARG
1	H	235	LYS
1	H	237	GLU
1	H	252	GLU
1	H	263	GLU
1	H	290	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	68	GLN
1	A	73	ASN
1	A	77	HIS
1	A	120	HIS
1	A	216	GLN
1	A	240	GLN
1	B	65	GLN
1	B	68	GLN
1	B	77	HIS
1	B	120	HIS
1	B	165	HIS
1	B	239	HIS
1	C	65	GLN
1	C	73	ASN
1	C	77	HIS
1	C	120	HIS
1	C	216	GLN

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Mol	Chain	Res	Type
1	C	239	HIS
1	D	65	GLN
1	D	68	GLN
1	D	77	HIS
1	D	83	GLN
1	D	165	HIS
1	D	216	GLN
1	D	239	HIS
1	E	65	GLN
1	E	68	GLN
1	E	77	HIS
1	E	120	HIS
1	F	65	GLN
1	F	68	GLN
1	F	77	HIS
1	F	120	HIS
1	F	239	HIS
1	G	65	GLN
1	G	73	ASN
1	G	77	HIS
1	G	120	HIS
1	H	65	GLN
1	H	68	GLN
1	H	77	HIS
1	H	120	HIS
1	H	239	HIS
1	H	240	GLN

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 14 ligands modelled in this entry, 6 are monoatomic - leaving 8 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
2	FAD	D	311	-	58,58,58	1.09	5 (8%)	85,89,89	1.74	18 (21%)
2	FAD	G	311	-	58,58,58	1.23	7 (12%)	85,89,89	1.64	17 (20%)
2	FAD	A	311	-	58,58,58	1.21	6 (10%)	85,89,89	1.51	16 (18%)
2	FAD	B	311	-	58,58,58	1.20	4 (6%)	85,89,89	1.64	18 (21%)
2	FAD	E	311	-	58,58,58	1.15	6 (10%)	85,89,89	1.64	17 (20%)
2	FAD	F	311	-	58,58,58	1.23	5 (8%)	85,89,89	1.82	20 (23%)
2	FAD	H	311	-	58,58,58	1.15	5 (8%)	85,89,89	1.53	14 (16%)
2	FAD	C	311	-	58,58,58	1.07	3 (5%)	85,89,89	1.69	18 (21%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	D	311	-	-	6/34/50/50	0/6/6/6
2	FAD	G	311	-	-	10/34/50/50	0/6/6/6
2	FAD	A	311	-	-	10/34/50/50	0/6/6/6
2	FAD	B	311	-	-	6/34/50/50	0/6/6/6
2	FAD	E	311	-	-	0/34/50/50	0/6/6/6
2	FAD	F	311	-	-	10/34/50/50	0/6/6/6
2	FAD	H	311	-	-	4/34/50/50	0/6/6/6
2	FAD	C	311	-	-	5/34/50/50	0/6/6/6

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	311	FAD	C4X-N5	4.12	1.39	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	311	FAD	C4X-N5	3.91	1.39	1.30
2	B	311	FAD	C4X-N5	3.90	1.39	1.30
2	C	311	FAD	C4X-N5	3.63	1.38	1.30
2	D	311	FAD	C4X-N5	3.54	1.38	1.30
2	F	311	FAD	C2A-N1A	3.42	1.40	1.33
2	A	311	FAD	C2A-N1A	3.39	1.39	1.33
2	F	311	FAD	C5'-C4'	3.33	1.56	1.51
2	B	311	FAD	C2A-N1A	3.28	1.39	1.33
2	F	311	FAD	C4X-N5	3.27	1.37	1.30
2	H	311	FAD	C4X-N5	3.19	1.37	1.30
2	E	311	FAD	C4X-N5	3.09	1.37	1.30
2	C	311	FAD	C10-N1	3.09	1.39	1.33
2	A	311	FAD	C2A-N3A	3.06	1.39	1.33
2	F	311	FAD	C2A-N3A	3.06	1.39	1.33
2	G	311	FAD	C8A-N7A	2.65	1.36	1.31
2	E	311	FAD	C10-N1	2.65	1.38	1.33
2	E	311	FAD	C2A-N1A	2.64	1.38	1.33
2	D	311	FAD	C2A-N1A	2.63	1.38	1.33
2	H	311	FAD	C8A-N7A	2.62	1.36	1.31
2	H	311	FAD	C2A-N1A	2.61	1.38	1.33
2	E	311	FAD	C2A-N3A	2.60	1.38	1.33
2	F	311	FAD	C8A-N7A	2.53	1.36	1.31
2	A	311	FAD	C8A-N7A	2.52	1.36	1.31
2	B	311	FAD	C8A-N7A	2.51	1.36	1.31
2	D	311	FAD	C10-N1	2.51	1.38	1.33
2	D	311	FAD	C8A-N7A	2.48	1.36	1.31
2	D	311	FAD	C2A-N3A	2.46	1.38	1.33
2	G	311	FAD	C2A-N1A	2.43	1.38	1.33
2	B	311	FAD	C2A-N3A	2.42	1.38	1.33
2	G	311	FAD	P-O3P	2.35	1.62	1.59
2	H	311	FAD	C2A-N3A	2.33	1.38	1.33
2	C	311	FAD	C8A-N7A	2.30	1.36	1.31
2	E	311	FAD	C4X-C10	-2.22	1.37	1.44
2	E	311	FAD	C8A-N7A	2.20	1.35	1.31
2	G	311	FAD	C2A-N3A	2.18	1.37	1.33
2	A	311	FAD	C5A-N7A	-2.14	1.35	1.39
2	H	311	FAD	C10-N1	2.12	1.37	1.33
2	G	311	FAD	O2'-C2'	-2.04	1.39	1.43
2	G	311	FAD	C4X-C10	-2.03	1.38	1.44
2	A	311	FAD	O2'-C2'	-2.00	1.39	1.43

All (138) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	311	FAD	N3A-C2A-N1A	-6.83	118.25	128.58
2	H	311	FAD	N3A-C2A-N1A	-6.33	118.99	128.58
2	G	311	FAD	N3A-C2A-N1A	-6.00	119.50	128.58
2	C	311	FAD	N3A-C2A-N1A	-5.79	119.82	128.58
2	D	311	FAD	C5A-C4A-N3A	-5.71	118.86	126.72
2	D	311	FAD	N3A-C2A-N1A	-5.41	120.39	128.58
2	E	311	FAD	N3A-C2A-N1A	-5.21	120.69	128.58
2	F	311	FAD	N3A-C2A-N1A	-5.12	120.84	128.58
2	A	311	FAD	N3A-C2A-N1A	-4.67	121.52	128.58
2	C	311	FAD	C5A-C4A-N3A	-4.53	120.48	126.72
2	C	311	FAD	N9A-C8A-N7A	-4.38	107.72	113.94
2	F	311	FAD	C5A-C4A-N3A	-4.34	120.75	126.72
2	A	311	FAD	C5A-C4A-N3A	-4.03	121.17	126.72
2	F	311	FAD	O4'-C4'-C5'	3.93	118.65	109.99
2	C	311	FAD	C2A-N3A-C4A	3.92	121.40	111.83
2	H	311	FAD	C5A-C4A-N3A	-3.90	121.34	126.72
2	D	311	FAD	C4-N3-C2	-3.89	118.74	125.64
2	G	311	FAD	C4X-C10-N10	3.88	122.04	116.48
2	F	311	FAD	N9A-C8A-N7A	-3.85	108.48	113.94
2	E	311	FAD	N9A-C8A-N7A	-3.83	108.50	113.94
2	D	311	FAD	C2A-N3A-C4A	3.83	121.17	111.83
2	E	311	FAD	C5A-C4A-N3A	-3.77	121.53	126.72
2	D	311	FAD	N9A-C8A-N7A	-3.74	108.63	113.94
2	A	311	FAD	C4X-C10-N10	3.74	121.83	116.48
2	C	311	FAD	C5A-N7A-C8A	3.70	109.27	103.45
2	D	311	FAD	C5A-N7A-C8A	3.70	109.27	103.45
2	D	311	FAD	N3A-C4A-N9A	3.58	133.25	127.17
2	F	311	FAD	C4-N3-C2	-3.52	119.39	125.64
2	F	311	FAD	C5A-N7A-C8A	3.49	108.94	103.45
2	B	311	FAD	N9A-C8A-N7A	-3.49	108.98	113.94
2	H	311	FAD	C2A-N3A-C4A	3.42	120.19	111.83
2	G	311	FAD	C5A-C4A-N3A	-3.42	122.01	126.72
2	B	311	FAD	C2A-N3A-C4A	3.39	120.10	111.83
2	E	311	FAD	C9A-C5X-N5	-3.37	118.87	122.45
2	H	311	FAD	C4-N3-C2	-3.37	119.66	125.64
2	G	311	FAD	O2P-P-O3P	3.36	116.36	107.27
2	F	311	FAD	O3P-P-O1P	-3.35	100.64	110.70
2	B	311	FAD	C5A-C4A-N3A	-3.34	122.11	126.72
2	G	311	FAD	C4-N3-C2	-3.29	119.80	125.64
2	C	311	FAD	O4B-C1B-N9A	-3.23	101.88	108.09
2	G	311	FAD	C2A-N3A-C4A	3.22	119.70	111.83
2	F	311	FAD	O2-C2-N1	-3.22	116.46	121.80
2	F	311	FAD	C2A-N3A-C4A	3.15	119.53	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	311	FAD	C5A-N7A-C8A	3.14	108.39	103.45
2	F	311	FAD	C4A-C5A-N7A	-3.13	107.00	110.58
2	E	311	FAD	C4-N3-C2	-3.12	120.09	125.64
2	B	311	FAD	C4X-C10-N10	3.09	120.91	116.48
2	H	311	FAD	N9A-C8A-N7A	-3.09	109.55	113.94
2	E	311	FAD	C2A-N3A-C4A	3.08	119.36	111.83
2	G	311	FAD	C4X-C4-N3	3.06	121.05	113.25
2	G	311	FAD	C9A-C5X-N5	-3.03	119.24	122.45
2	F	311	FAD	C4X-C10-N10	3.01	120.78	116.48
2	D	311	FAD	C10-C4X-N5	-3.00	118.67	124.81
2	B	311	FAD	C4-N3-C2	-3.00	120.32	125.64
2	D	311	FAD	C4X-C10-N10	2.99	120.77	116.48
2	B	311	FAD	O2-C2-N1	-2.98	116.84	121.80
2	A	311	FAD	C2A-N3A-C4A	2.95	119.04	111.83
2	F	311	FAD	C10-C4X-N5	-2.95	118.79	124.81
2	H	311	FAD	C4X-C10-N10	2.94	120.69	116.48
2	B	311	FAD	C5A-N7A-C8A	2.93	108.06	103.45
2	D	311	FAD	C4A-C5A-N7A	-2.90	107.27	110.58
2	C	311	FAD	N3A-C4A-N9A	2.89	132.09	127.17
2	C	311	FAD	C9A-C5X-N5	-2.89	119.39	122.45
2	E	311	FAD	C4X-C4-N3	2.86	120.53	113.25
2	G	311	FAD	C10-C4X-N5	-2.86	118.97	124.81
2	A	311	FAD	C4-N3-C2	-2.86	120.56	125.64
2	G	311	FAD	C4-C4X-N5	2.85	122.14	118.21
2	A	311	FAD	C4X-C4-N3	2.81	120.41	113.25
2	F	311	FAD	O5'-C5'-C4'	2.81	116.86	109.36
2	C	311	FAD	C4A-C5A-N7A	-2.79	107.40	110.58
2	B	311	FAD	O4-C4-C4X	-2.76	119.24	126.53
2	H	311	FAD	O4-C4-C4X	-2.75	119.27	126.53
2	D	311	FAD	C4X-C4-N3	2.75	120.24	113.25
2	E	311	FAD	O3P-PA-O1A	-2.74	102.46	110.70
2	B	311	FAD	C10-C4X-N5	-2.71	119.28	124.81
2	A	311	FAD	O2P-P-O3P	-2.70	99.97	107.27
2	E	311	FAD	O4-C4-C4X	-2.68	119.47	126.53
2	E	311	FAD	C10-C4X-N5	-2.67	119.36	124.81
2	G	311	FAD	C5X-C9A-N10	2.67	120.38	117.97
2	E	311	FAD	C4X-C10-N10	2.67	120.30	116.48
2	C	311	FAD	C4X-C4-N3	2.67	120.04	113.25
2	G	311	FAD	C10-N1-C2	2.66	122.62	116.85
2	C	311	FAD	C4X-C10-N10	2.64	120.27	116.48
2	C	311	FAD	C10-C4X-N5	-2.64	119.41	124.81
2	B	311	FAD	O2P-P-O3P	2.62	114.34	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	311	FAD	O3'-C3'-C2'	2.59	114.83	108.93
2	F	311	FAD	C4'-C3'-C2'	-2.59	109.26	113.57
2	A	311	FAD	N3A-C4A-N9A	2.58	131.56	127.17
2	B	311	FAD	O2-C2-N3	2.56	123.49	118.58
2	C	311	FAD	C4-N3-C2	-2.54	121.13	125.64
2	F	311	FAD	N3A-C4A-N9A	2.52	131.46	127.17
2	H	311	FAD	C4X-C4-N3	2.52	119.68	113.25
2	D	311	FAD	O4'-C4'-C3'	-2.52	103.35	109.25
2	G	311	FAD	C4X-C10-N1	-2.50	118.45	124.59
2	F	311	FAD	O2P-P-O3P	2.49	114.01	107.27
2	A	311	FAD	C10-C4X-N5	-2.49	119.72	124.81
2	H	311	FAD	C5A-N7A-C8A	2.46	107.32	103.45
2	H	311	FAD	C10-C4X-N5	-2.45	119.81	124.81
2	A	311	FAD	C4-C4X-N5	2.45	121.59	118.21
2	E	311	FAD	N3A-C4A-N9A	2.44	131.32	127.17
2	H	311	FAD	O2P-P-O3P	2.37	113.68	107.27
2	B	311	FAD	C2A-N1A-C6A	2.36	122.61	118.73
2	F	311	FAD	O4-C4-C4X	-2.36	120.31	126.53
2	B	311	FAD	C4X-C4-N3	2.32	119.16	113.25
2	F	311	FAD	C4-C4X-C10	2.31	120.89	116.93
2	C	311	FAD	C4A-N9A-C8A	2.30	108.15	105.74
2	C	311	FAD	C2B-C1B-N9A	2.30	119.02	113.30
2	D	311	FAD	C4X-C10-N1	-2.29	118.98	124.59
2	H	311	FAD	N3A-C4A-N9A	2.27	131.02	127.17
2	E	311	FAD	C5X-N5-C4X	2.27	121.75	118.09
2	C	311	FAD	C4X-C10-N1	-2.25	119.08	124.59
2	A	311	FAD	C2B-C1B-N9A	2.24	118.88	113.30
2	H	311	FAD	O2-C2-N1	-2.23	118.10	121.80
2	E	311	FAD	C10-N1-C2	2.20	121.62	116.85
2	G	311	FAD	O4-C4-N3	-2.17	116.04	120.11
2	G	311	FAD	O4B-C1B-C2B	-2.16	101.98	106.62
2	B	311	FAD	C4-C4X-N5	2.16	121.19	118.21
2	D	311	FAD	O2-C2-N1	-2.15	118.22	121.80
2	C	311	FAD	O3P-PA-O1A	-2.14	104.26	110.70
2	A	311	FAD	C9A-C5X-N5	-2.14	120.19	122.45
2	A	311	FAD	C10-N1-C2	2.13	121.47	116.85
2	A	311	FAD	N9A-C8A-N7A	-2.13	110.91	113.94
2	E	311	FAD	C4A-C5A-N7A	-2.13	108.15	110.58
2	D	311	FAD	C6A-C5A-C4A	2.11	120.06	117.18
2	B	311	FAD	C4A-N9A-C1B	-2.11	121.69	126.63
2	D	311	FAD	O4-C4-C4X	-2.11	120.97	126.53
2	F	311	FAD	C4X-C4-N3	2.11	118.61	113.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	311	FAD	C9A-C5X-N5	-2.08	120.25	122.45
2	G	311	FAD	N9A-C8A-N7A	-2.07	111.00	113.94
2	A	311	FAD	O5B-PA-O1A	2.06	117.12	108.94
2	B	311	FAD	C4A-C5A-N7A	-2.06	108.23	110.58
2	B	311	FAD	N3A-C4A-N9A	2.04	130.64	127.17
2	D	311	FAD	C9A-C5X-N5	-2.04	120.29	122.45
2	C	311	FAD	C10-N1-C2	2.03	121.24	116.85
2	G	311	FAD	C2A-N1A-C6A	2.02	122.06	118.73
2	E	311	FAD	C4X-C10-N1	-2.02	119.64	124.59
2	A	311	FAD	O4B-C1B-N9A	-2.02	104.21	108.09
2	D	311	FAD	C4-C4X-C10	2.01	120.38	116.93

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	311	FAD	C5B-O5B-PA-O1A
2	A	311	FAD	C5B-O5B-PA-O2A
2	A	311	FAD	C5B-O5B-PA-O3P
2	A	311	FAD	C3'-C4'-C5'-O5'
2	A	311	FAD	O4'-C4'-C5'-O5'
2	A	311	FAD	C5'-O5'-P-O1P
2	A	311	FAD	C5'-O5'-P-O2P
2	A	311	FAD	C5'-O5'-P-O3P
2	D	311	FAD	C5B-O5B-PA-O1A
2	D	311	FAD	C5B-O5B-PA-O3P
2	F	311	FAD	O4'-C4'-C5'-O5'
2	F	311	FAD	C5'-O5'-P-O1P
2	H	311	FAD	C5B-O5B-PA-O1A
2	F	311	FAD	C3'-C4'-C5'-O5'
2	C	311	FAD	O3'-C3'-C4'-C5'
2	G	311	FAD	P-O3P-PA-O1A
2	C	311	FAD	C2'-C3'-C4'-C5'
2	F	311	FAD	C2'-C3'-C4'-C5'
2	G	311	FAD	C2'-C3'-C4'-C5'
2	G	311	FAD	C2'-C3'-C4'-O4'
2	A	311	FAD	PA-O3P-P-O5'
2	F	311	FAD	P-O3P-PA-O5B
2	G	311	FAD	O2'-C2'-C3'-C4'
2	D	311	FAD	C1'-C2'-C3'-O3'
2	F	311	FAD	C1'-C2'-C3'-O3'
2	G	311	FAD	C1'-C2'-C3'-O3'

Continued on next page...

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Mol	Chain	Res	Type	Atoms
2	B	311	FAD	C3B-C4B-C5B-O5B
2	B	311	FAD	P-O3P-PA-O2A
2	D	311	FAD	P-O3P-PA-O2A
2	F	311	FAD	P-O3P-PA-O1A
2	C	311	FAD	O3'-C3'-C4'-O4'
2	F	311	FAD	C5'-O5'-P-O2P
2	F	311	FAD	C5'-O5'-P-O3P
2	H	311	FAD	C5B-O5B-PA-O3P
2	C	311	FAD	C2'-C3'-C4'-O4'
2	C	311	FAD	P-O3P-PA-O1A
2	H	311	FAD	P-O3P-PA-O2A
2	B	311	FAD	C2B-C1B-N9A-C8A
2	G	311	FAD	O3'-C3'-C4'-C5'
2	G	311	FAD	O3'-C3'-C4'-O4'
2	B	311	FAD	O4B-C4B-C5B-O5B
2	F	311	FAD	O3'-C3'-C4'-C5'
2	B	311	FAD	O4B-C1B-N9A-C8A
2	D	311	FAD	C2'-C3'-C4'-O4'
2	G	311	FAD	C1'-C2'-C3'-C4'
2	A	311	FAD	C2B-C1B-N9A-C8A
2	B	311	FAD	P-O3P-PA-O1A
2	G	311	FAD	P-O3P-PA-O2A
2	H	311	FAD	P-O3P-PA-O1A
2	G	311	FAD	C2B-C1B-N9A-C8A
2	D	311	FAD	P-O3P-PA-O1A

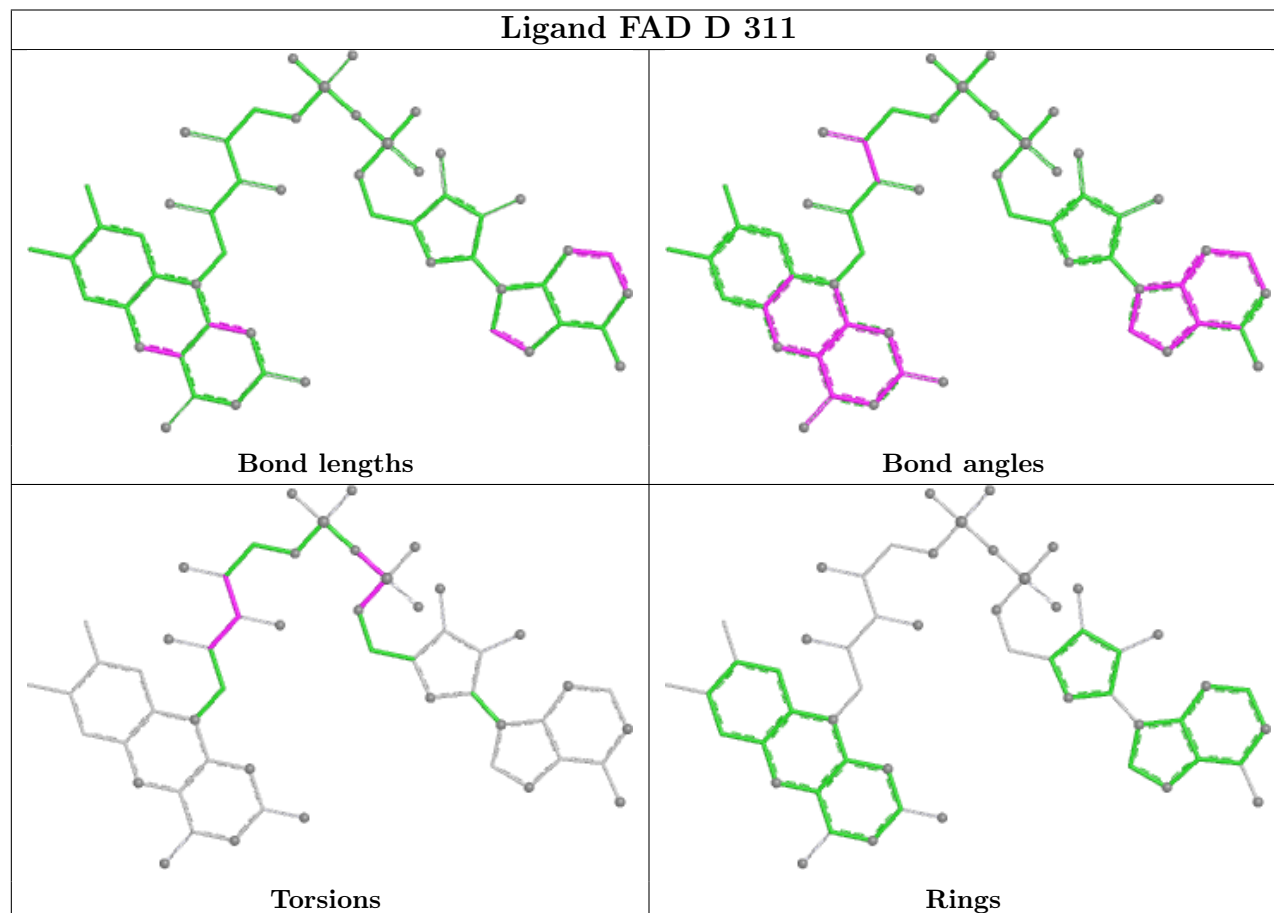
There are no ring outliers.

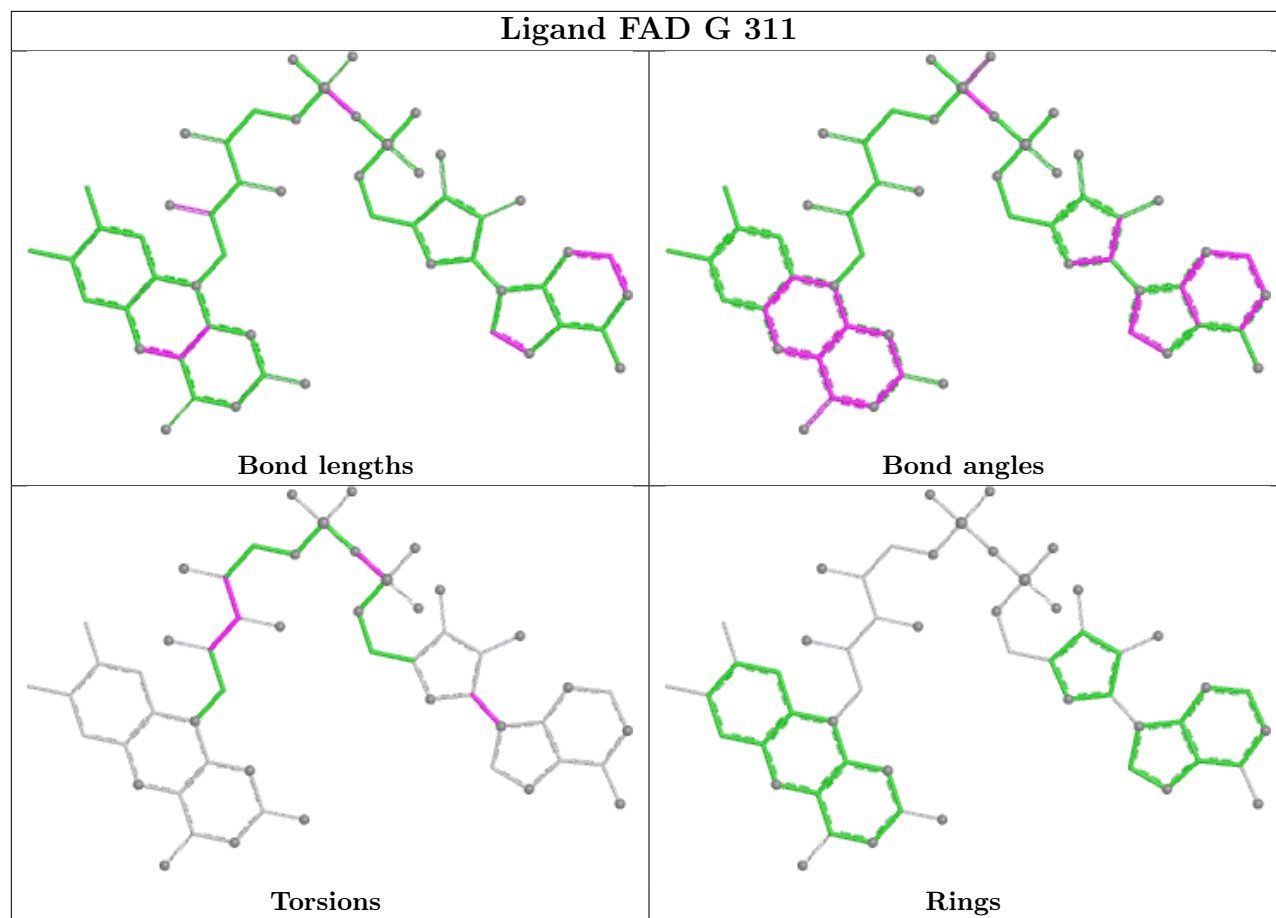
7 monomers are involved in 14 short contacts:

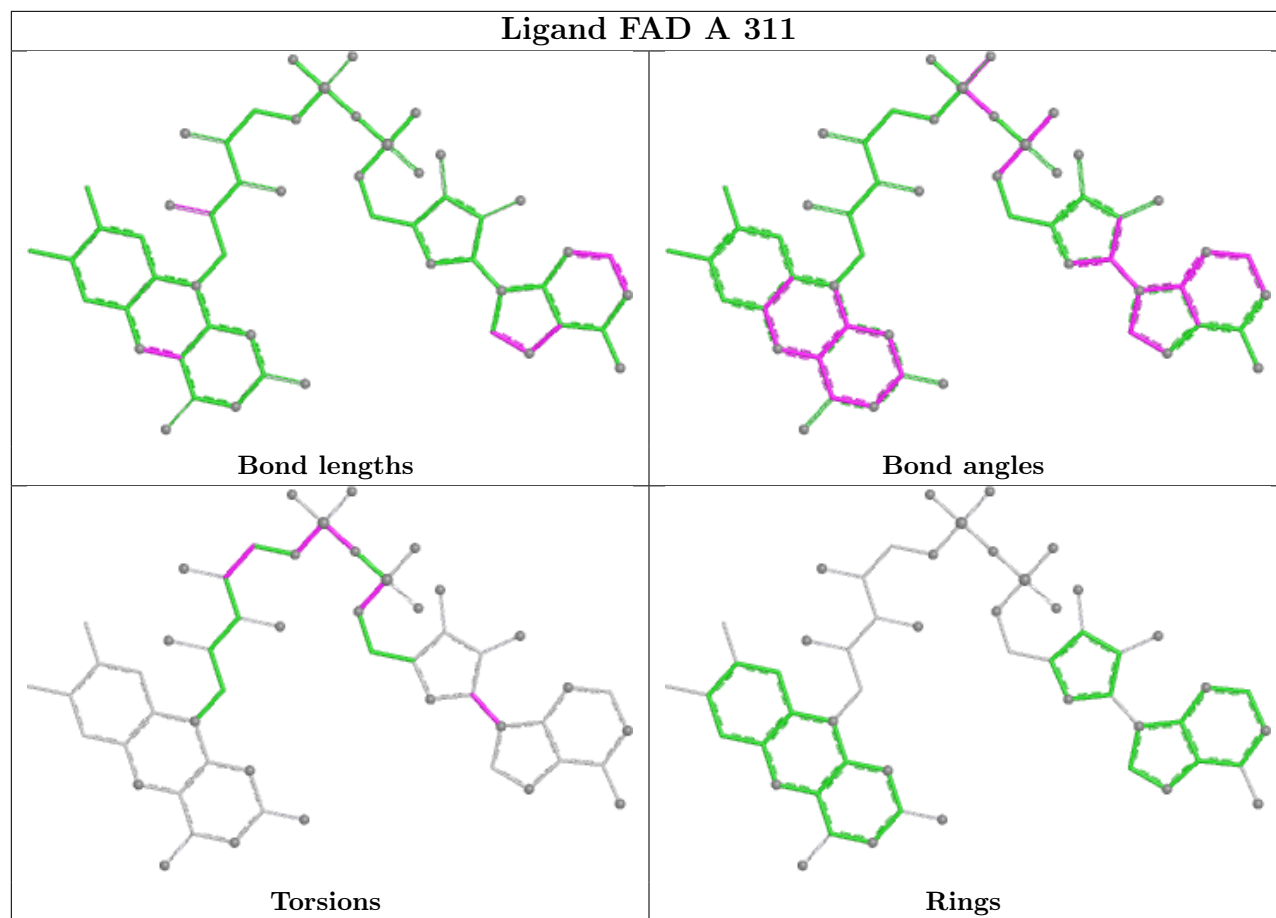
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	311	FAD	1	0
2	G	311	FAD	3	0
2	A	311	FAD	4	0
2	B	311	FAD	2	0
2	F	311	FAD	2	0
2	H	311	FAD	1	0
2	C	311	FAD	1	0

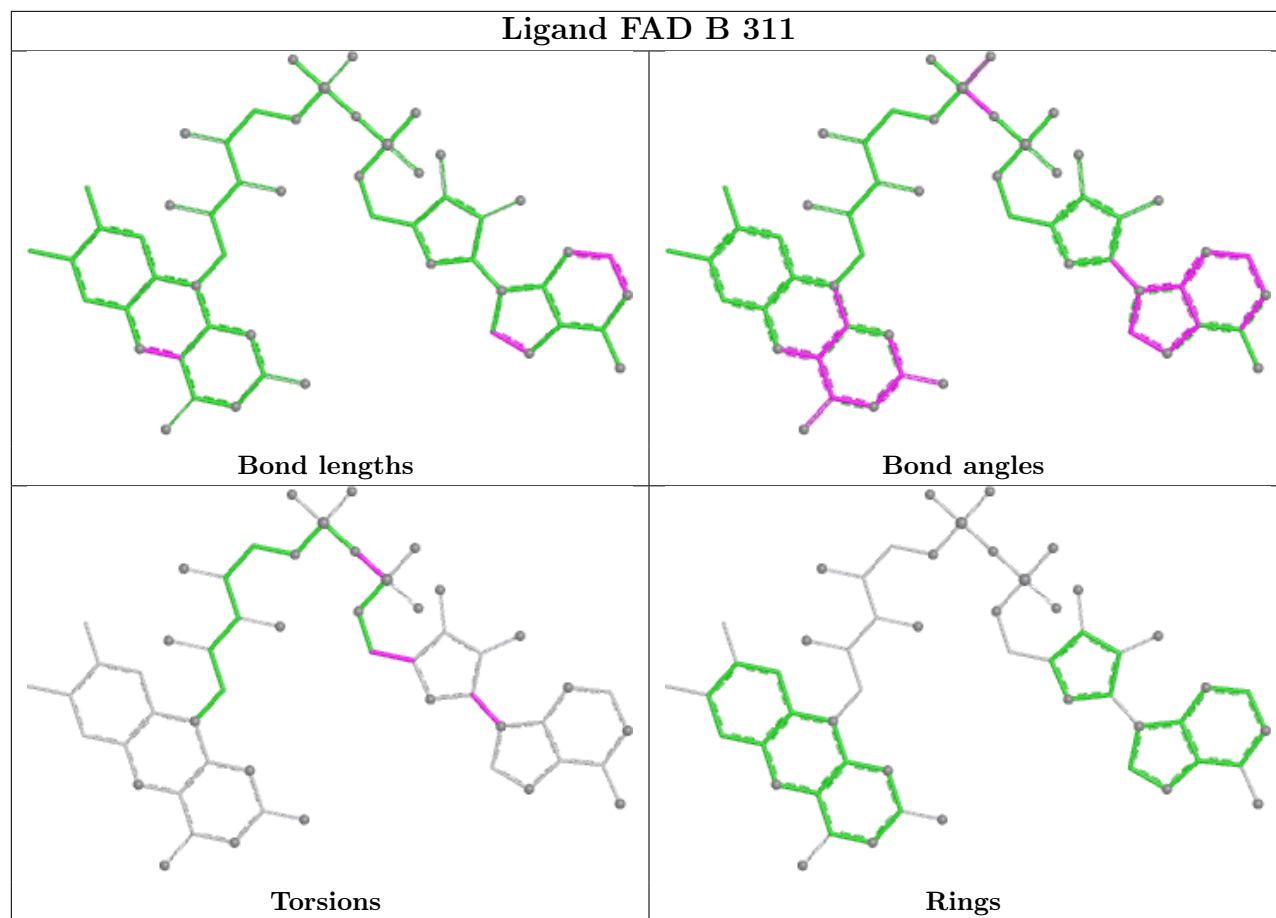
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is

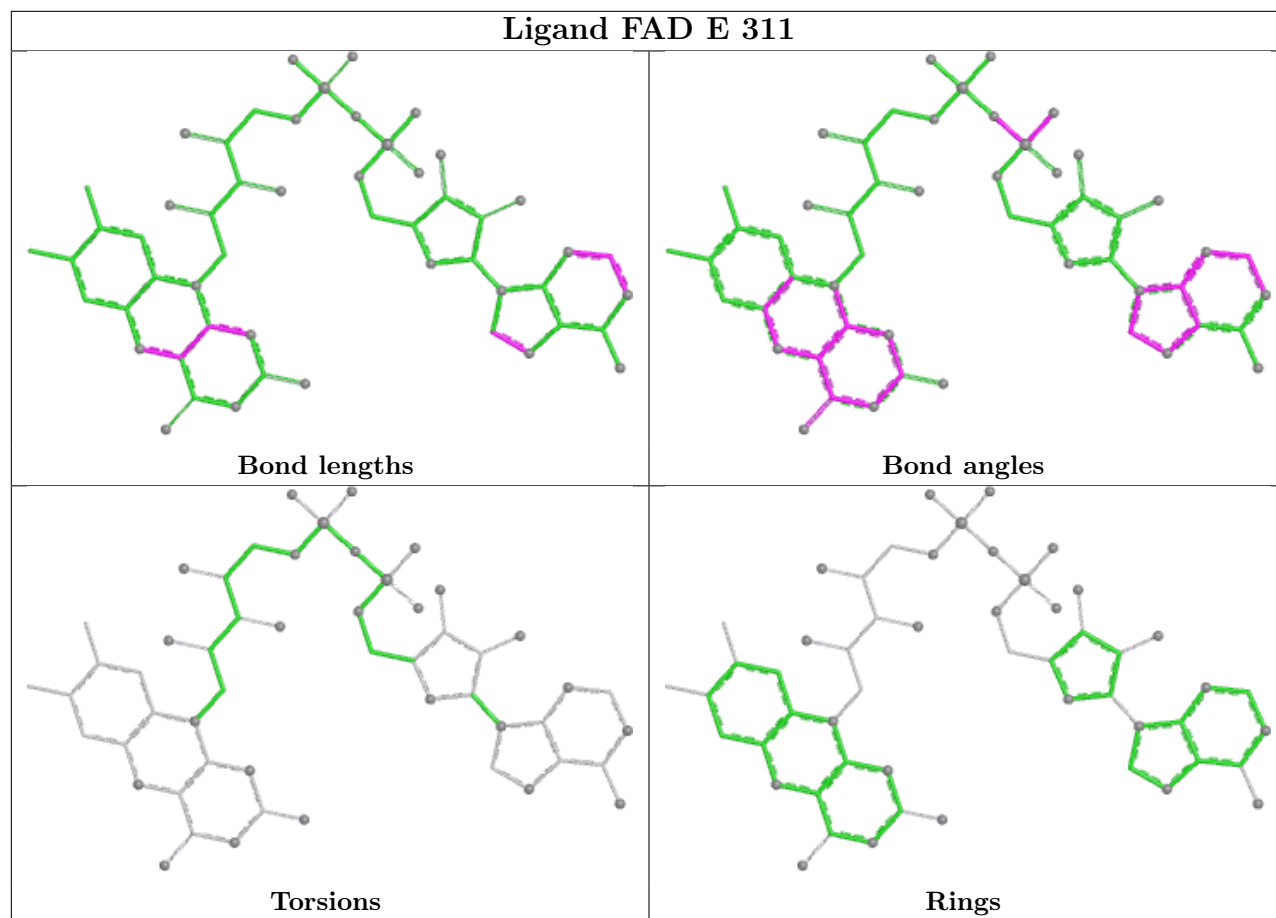
within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



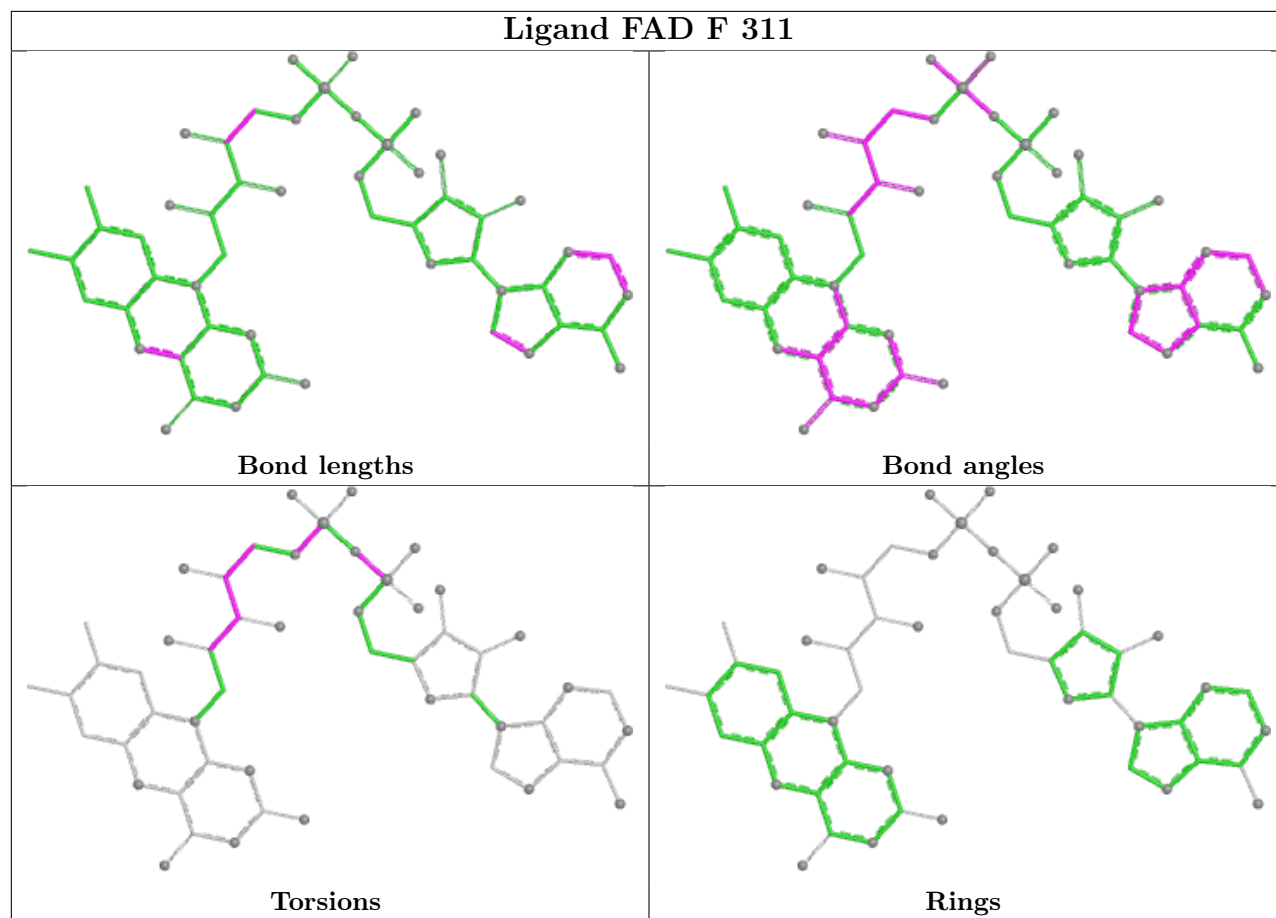


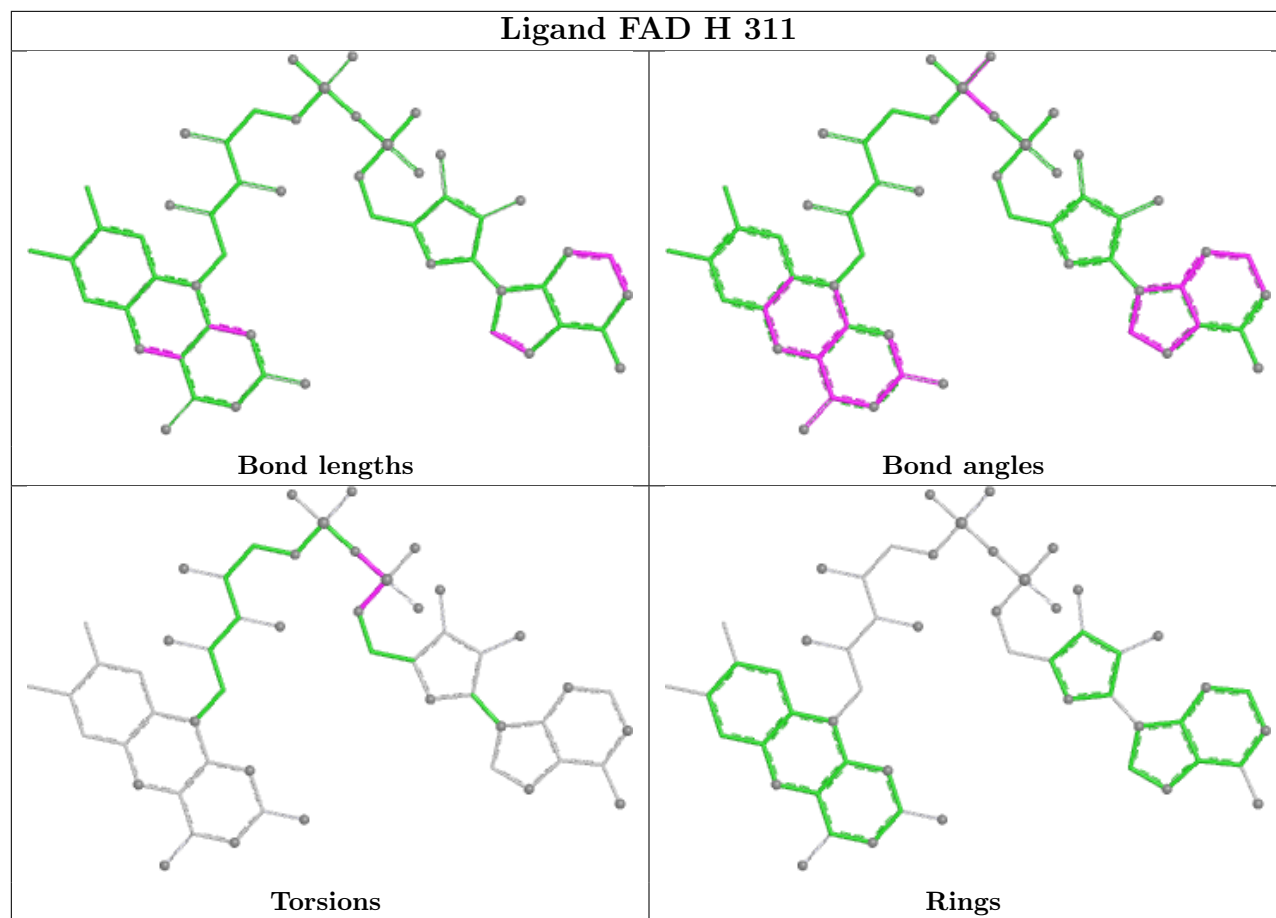


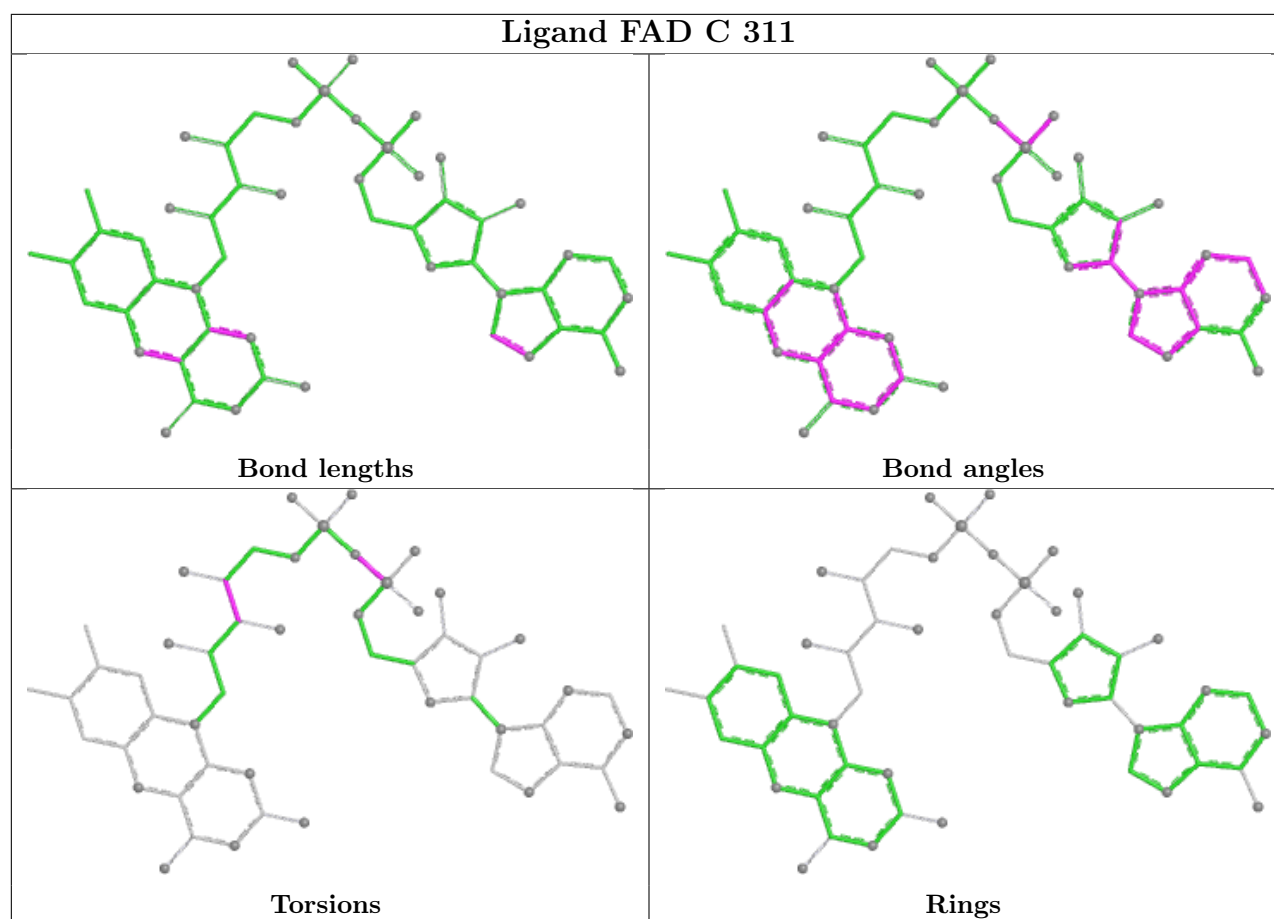




Ligand FAD F 311







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	291/310 (93%)	-1.81	0 100 100	6, 17, 33, 45	1 (0%)
1	B	291/310 (93%)	-1.79	0 100 100	5, 18, 34, 46	1 (0%)
1	C	291/310 (93%)	-1.80	0 100 100	5, 18, 34, 56	1 (0%)
1	D	291/310 (93%)	-1.79	0 100 100	5, 18, 33, 46	1 (0%)
1	E	291/310 (93%)	-1.81	0 100 100	7, 18, 36, 52	1 (0%)
1	F	291/310 (93%)	-1.79	0 100 100	7, 18, 35, 47	1 (0%)
1	G	291/310 (93%)	-1.80	0 100 100	6, 17, 33, 50	1 (0%)
1	H	291/310 (93%)	-1.80	0 100 100	5, 18, 35, 47	1 (0%)
All	All	2328/2480 (93%)	-1.80	0 100 100	5, 18, 34, 56	8 (0%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

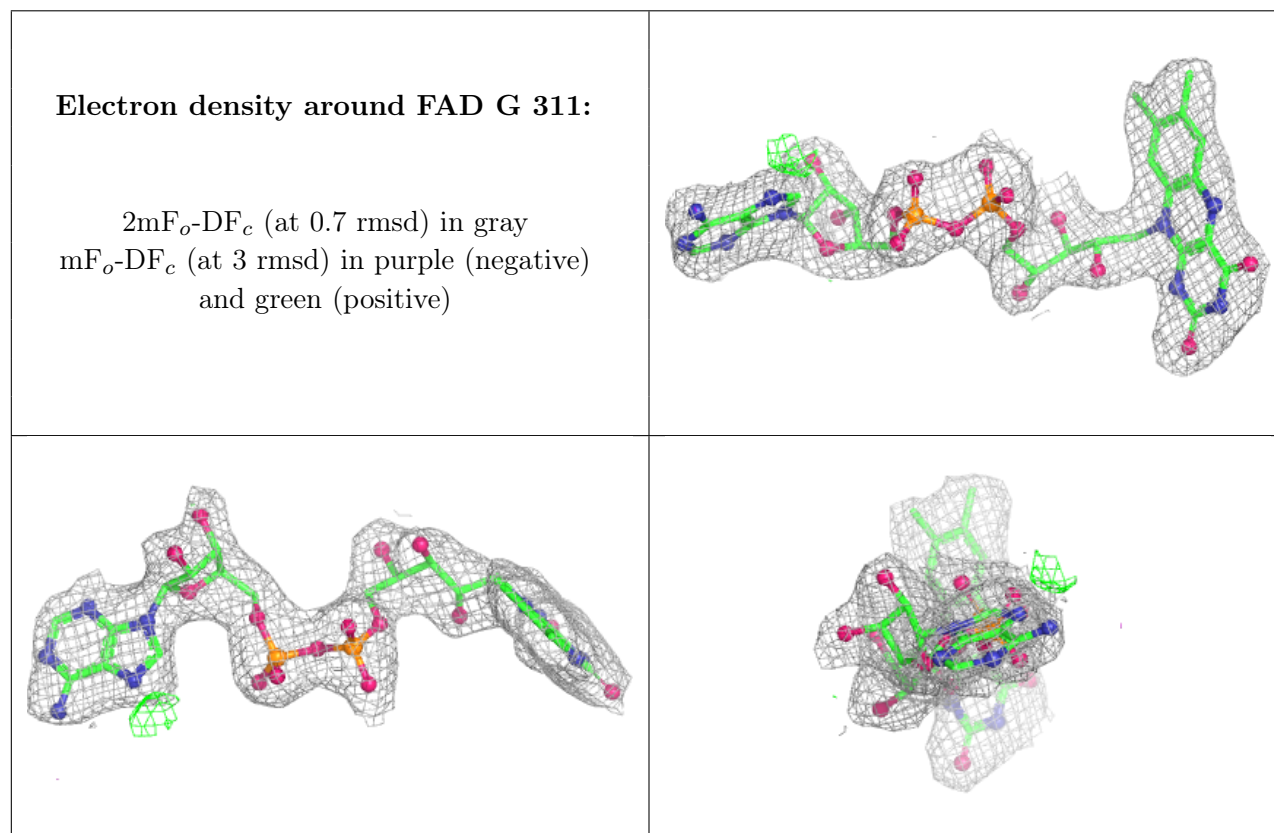
There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

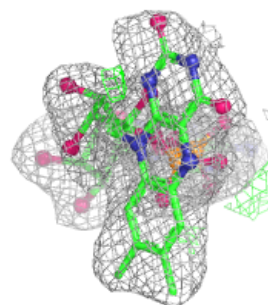
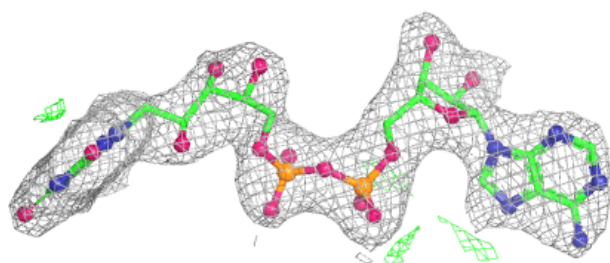
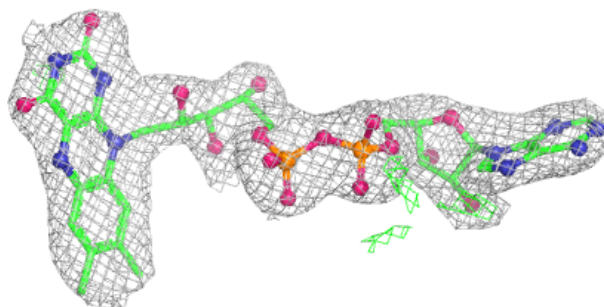
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CL	F	312	1/1	0.97	0.04	41,41,41,41	0
3	CL	E	312	1/1	0.98	0.05	37,37,37,37	0
3	CL	D	312	1/1	0.98	0.07	40,40,40,40	0
3	CL	A	312	1/1	0.99	0.03	31,31,31,31	0
3	CL	G	312	1/1	0.99	0.03	38,38,38,38	0
3	CL	H	312	1/1	0.99	0.06	31,31,31,31	0
2	FAD	G	311	53/53	1.00	0.02	4,9,13,13	0
2	FAD	H	311	53/53	1.00	0.02	11,15,18,19	0
2	FAD	A	311	53/53	1.00	0.02	9,11,14,14	0
2	FAD	B	311	53/53	1.00	0.02	6,14,18,19	0
2	FAD	C	311	53/53	1.00	0.02	5,10,15,15	0
2	FAD	D	311	53/53	1.00	0.02	7,12,18,19	0
2	FAD	E	311	53/53	1.00	0.02	6,9,13,14	0
2	FAD	F	311	53/53	1.00	0.02	2,16,19,22	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.

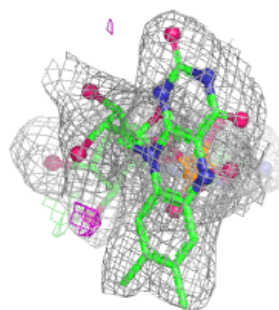
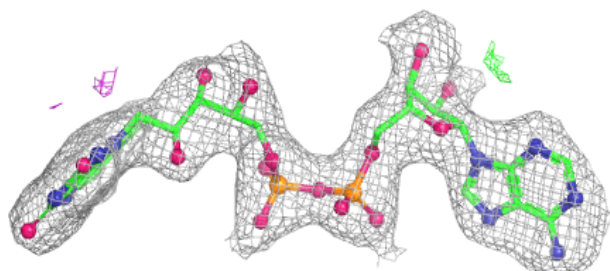
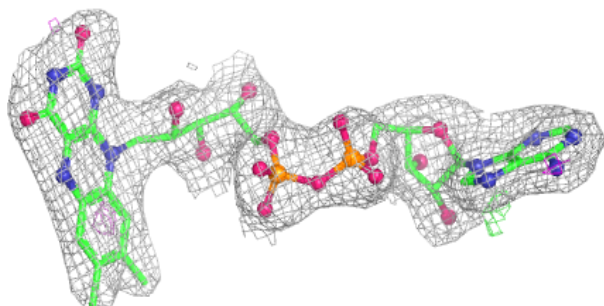


Electron density around FAD H 311:

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

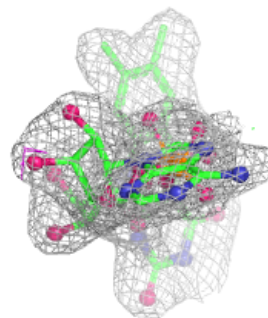
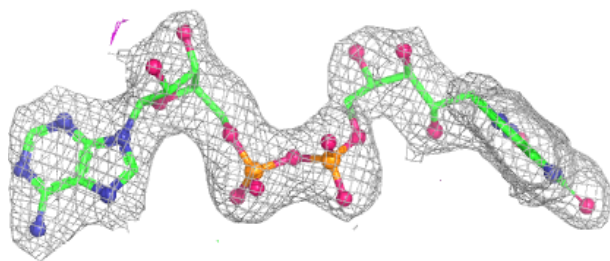
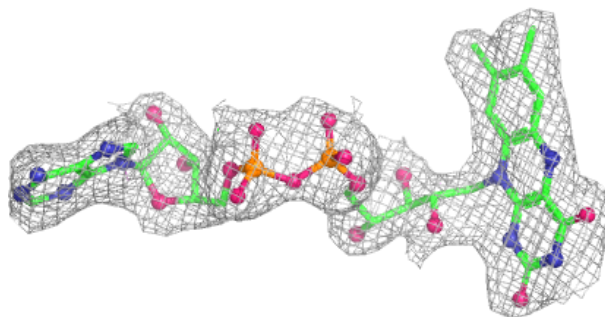
**Electron density around FAD A 311:**

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

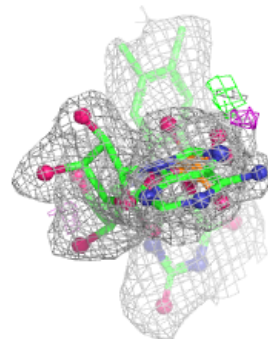
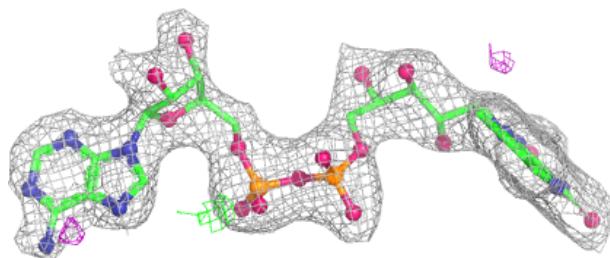
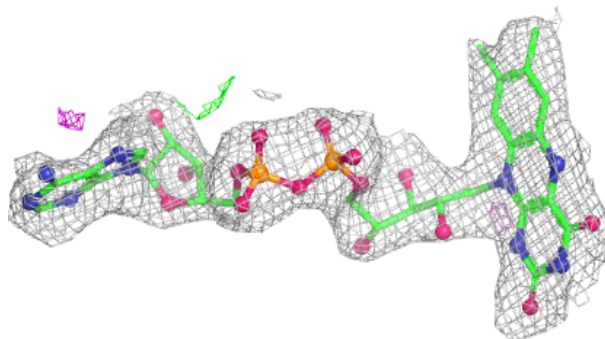


Electron density around FAD B 311:

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

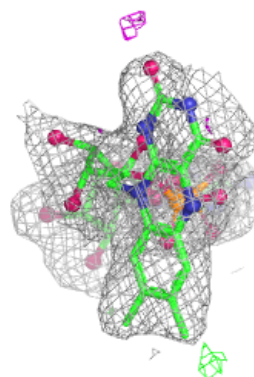
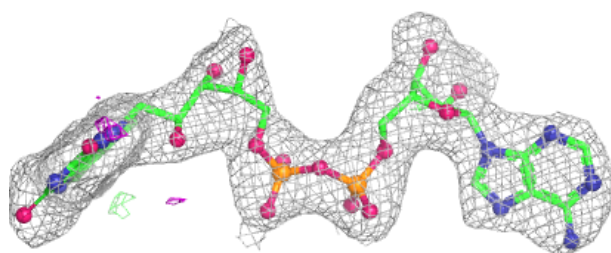
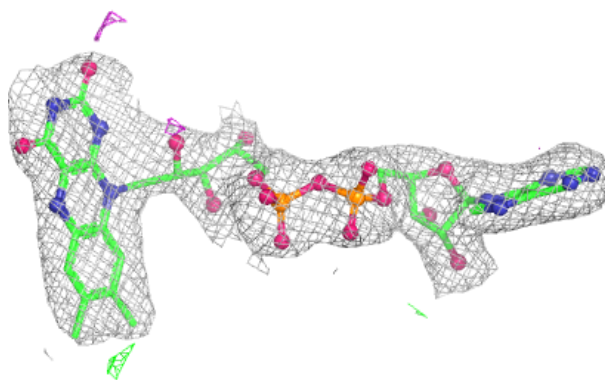
**Electron density around FAD C 311:**

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

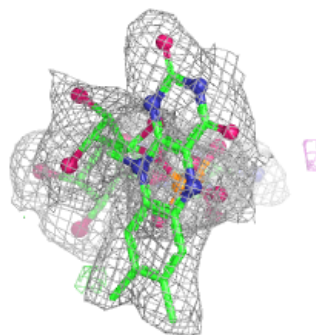
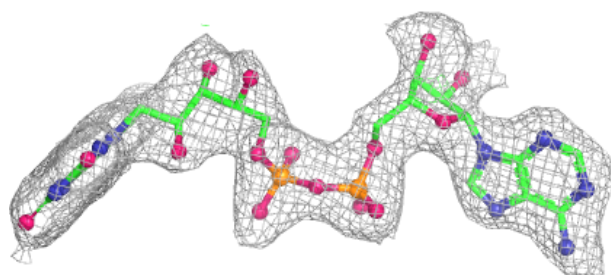
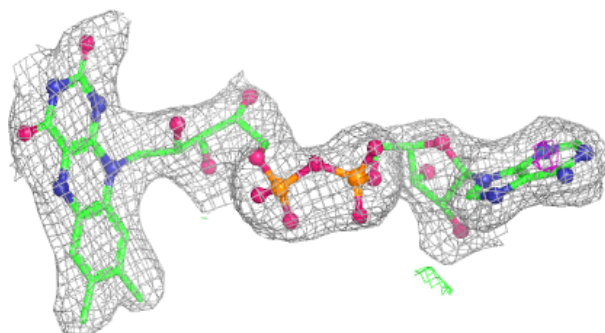


Electron density around FAD D 311:

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

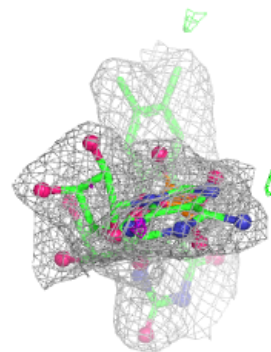
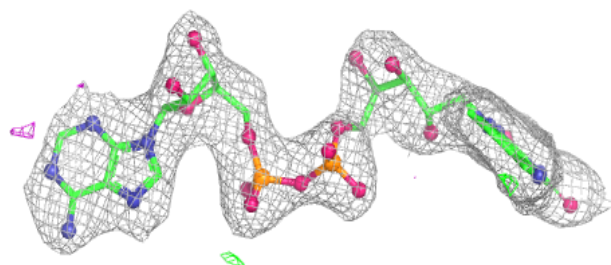
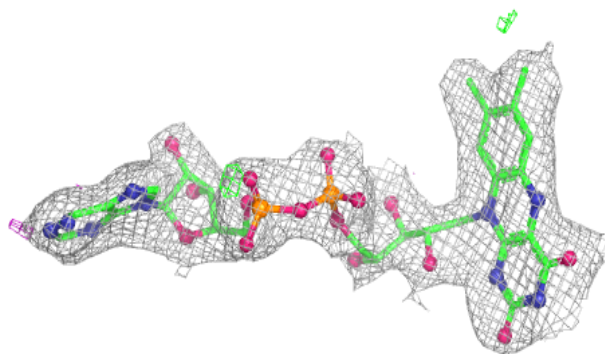
**Electron density around FAD E 311:**

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



Electron density around FAD F 311:

$2mF_o - DF_c$ (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.