Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 01:40 pm BST

PDB ID	:	1QLT
Title	:	STRUCTURE OF THE H422A MUTANT OF THE FLAVOENZYME VANILLYL-ALCOHOL OXIDASE
Authors	:	Mattevi, A.; Fraaije, M.
Deposited on	:	1999-09-16
Resolution	:	2.20 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

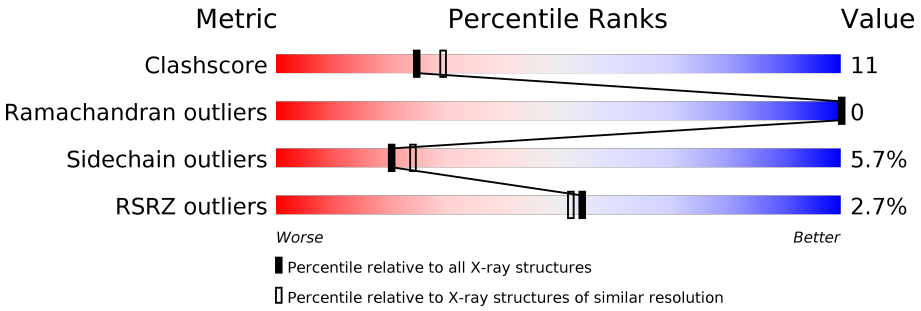
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

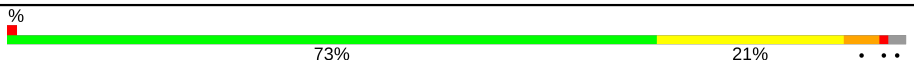
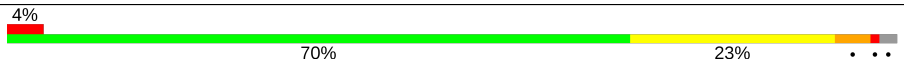
The reported resolution of this entry is 2.20 Å.

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))
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	560	
1	B	560	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9178 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.

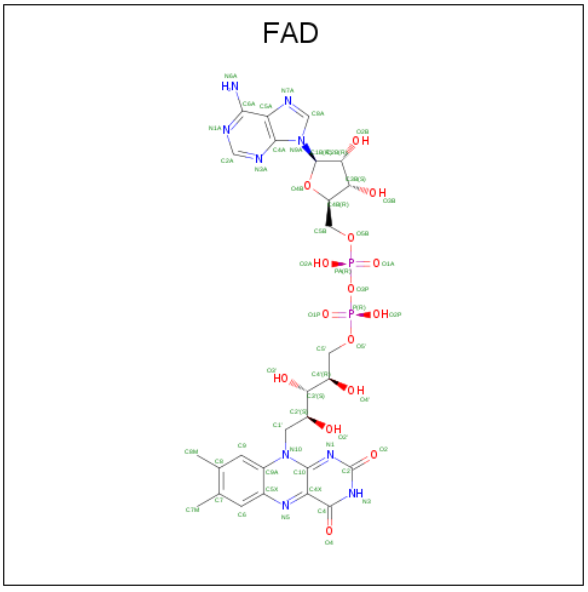
- Molecule 1 is a protein called VANILLYL-ALCOHOL OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	550	Total	C	N	O	S	68	0	0
			4346	2790	742	790	24			
1	B	550	Total	C	N	O	S	68	0	0
			4346	2790	742	790	24			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	422	ALA	HIS	engineered mutation	? ?
B	422	ALA	HIS	engineered mutation	? ?

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



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

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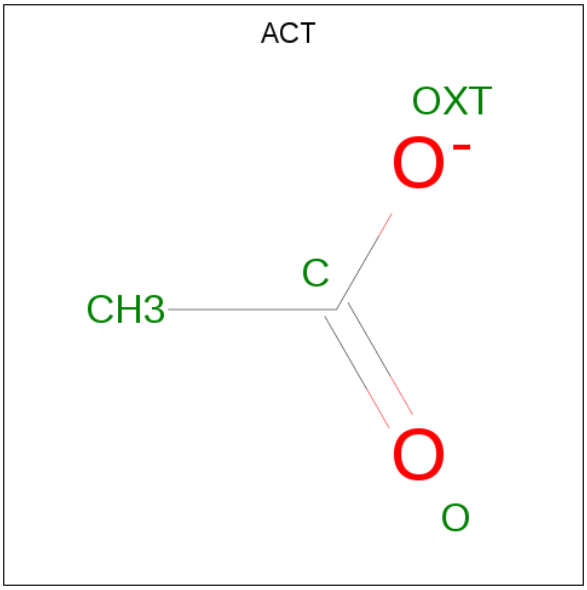
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is water.

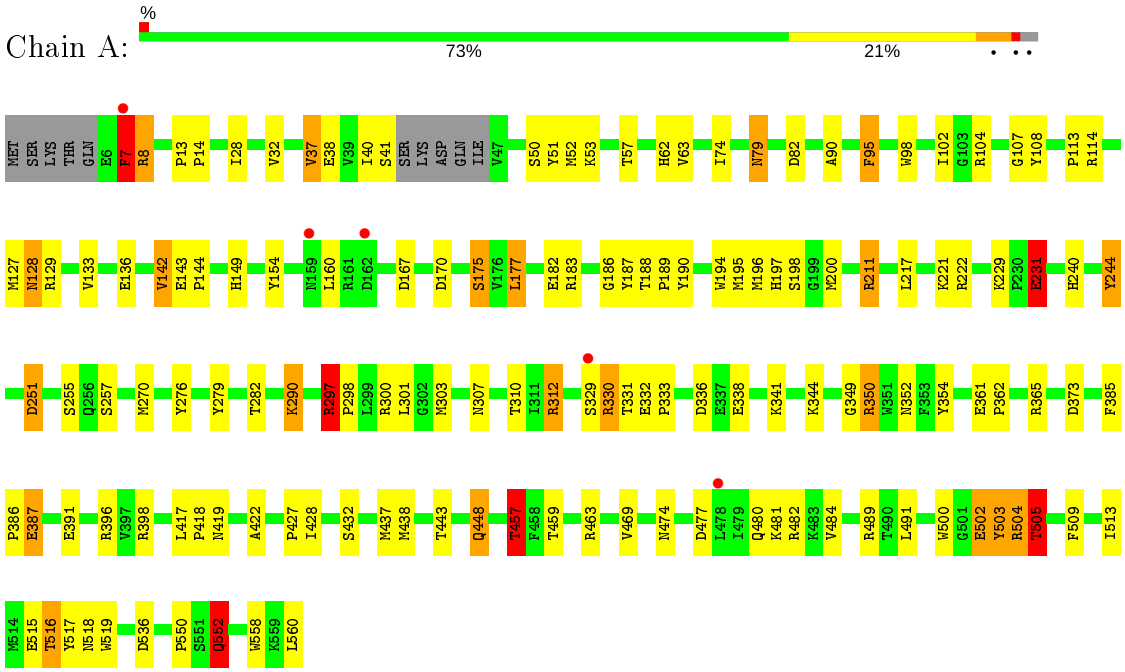
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	207	Total	O	0	0
			207	207		
4	B	165	Total	O	0	0
			165	165		

3 Residue-property plots (i)

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

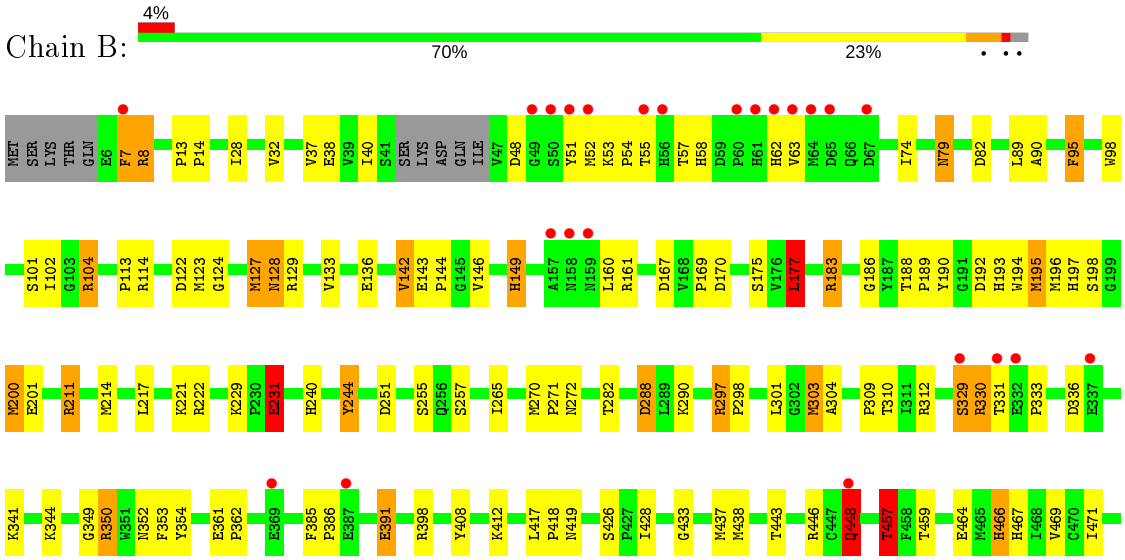
- Molecule 1: VANILLYL-ALCOHOL OXIDASE

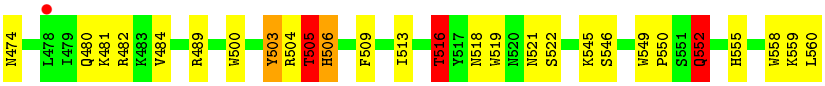
Chain A:



- Molecule 1: VANILLYL-ALCOHOL OXIDASE

Chain B:




N474
L478
I479
Q480
K481
R482
K483
V484
R489
W500
Y503
R504
T505
H506
F509
I513
T516
Y517
N518
W519
N520
N521
S522
K545
S546
W549
P550
S551
Q552
H555
W558
K559
L560

4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α, β, γ	129.66Å 129.66Å 132.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 34.70 – 2.20	Depositor EDS
% Data completeness (in resolution range)	93.1 (30.00-2.20) 93.0 (34.70-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	2.78 (at 2.20Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.210 , 0.264 0.204 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	27.4	Xtriage
Anisotropy	0.601	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 48.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.005 for l,-k,h 0.017 for -l,-k,-h 0.022 for -h,-l,-k 0.006 for -h,l,k 0.034 for -h,k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	9178	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, ACT

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	0.91	9/4464 (0.2%)	1.78	78/6067 (1.3%)
1	B	0.86	8/4464 (0.2%)	1.70	74/6067 (1.2%)
All	All	0.89	17/8928 (0.2%)	1.74	152/12134 (1.3%)

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	A	0	1
1	B	0	1
All	All	0	2

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	52	MET	CB-CG	29.10	2.44	1.51
1	A	52	MET	CB-CG	29.09	2.44	1.51
1	A	53	LYS	CB-CG	14.72	1.92	1.52
1	A	350	ARG	CZ-NH1	-12.21	1.17	1.33
1	B	350	ARG	CZ-NH1	-10.75	1.19	1.33
1	A	350	ARG	CZ-NH2	8.38	1.44	1.33
1	A	7	PHE	CB-CG	-8.13	1.37	1.51
1	B	221	LYS	CG-CD	-8.12	1.24	1.52
1	B	53	LYS	CB-CG	7.40	1.72	1.52
1	B	7	PHE	CB-CG	-7.25	1.39	1.51
1	A	221	LYS	CG-CD	-6.86	1.29	1.52
1	B	161	ARG	CB-CG	-6.65	1.34	1.52
1	A	341	LYS	CB-CG	-6.53	1.34	1.52
1	A	344	LYS	CB-CG	-6.04	1.36	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	341	LYS	CB-CG	-5.66	1.37	1.52
1	A	365	ARG	CD-NE	-5.18	1.37	1.46
1	B	344	LYS	CB-CG	-5.01	1.39	1.52

All (152) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	350	ARG	NE-CZ-NH1	-43.73	98.43	120.30
1	B	350	ARG	NE-CZ-NH1	-40.19	100.20	120.30
1	A	211	ARG	CD-NE-CZ	35.88	173.83	123.60
1	B	211	ARG	CD-NE-CZ	34.67	172.14	123.60
1	A	504	ARG	CD-NE-CZ	20.60	152.44	123.60
1	B	504	ARG	CD-NE-CZ	20.44	152.21	123.60
1	B	52	MET	CA-CB-CG	-18.32	82.16	113.30
1	A	52	MET	CA-CB-CG	-18.09	82.55	113.30
1	A	448	GLN	OE1-CD-NE2	-17.91	80.70	121.90
1	A	350	ARG	NE-CZ-NH2	15.34	127.97	120.30
1	A	448	GLN	CG-CD-NE2	14.24	150.88	116.70
1	A	53	LYS	CB-CG-CD	-13.78	75.77	111.60
1	A	504	ARG	NE-CZ-NH2	-13.72	113.44	120.30
1	A	37	VAL	CA-C-N	13.43	146.74	117.20
1	A	114	ARG	NE-CZ-NH2	-12.55	114.03	120.30
1	B	53	LYS	CB-CG-CD	-12.39	79.38	111.60
1	A	37	VAL	CA-C-O	-11.60	95.75	120.10
1	A	398	ARG	NE-CZ-NH1	11.57	126.08	120.30
1	B	446	ARG	NE-CZ-NH1	11.54	126.07	120.30
1	A	177	LEU	CA-CB-CG	11.51	141.78	115.30
1	B	448	GLN	OE1-CD-NE2	-11.37	95.75	121.90
1	A	504	ARG	NE-CZ-NH1	11.10	125.85	120.30
1	B	211	ARG	CA-CB-CG	10.87	137.31	113.40
1	B	504	ARG	NE-CZ-NH1	10.65	125.63	120.30
1	A	211	ARG	CA-CB-CG	10.55	136.60	113.40
1	A	297	ARG	NE-CZ-NH2	-10.48	115.06	120.30
1	B	504	ARG	NE-CZ-NH2	-10.47	115.06	120.30
1	A	341	LYS	CA-CB-CG	10.33	136.13	113.40
1	A	53	LYS	N-CA-CB	10.06	128.72	110.60
1	A	312	ARG	NE-CZ-NH1	-10.06	115.27	120.30
1	A	231	GLU	OE1-CD-OE2	-9.93	111.39	123.30
1	B	552	GLN	CA-CB-CG	9.60	134.51	113.40
1	B	448	GLN	CG-CD-NE2	9.56	139.66	116.70
1	A	196	MET	CG-SD-CE	9.52	115.43	100.20
1	B	350	ARG	NE-CZ-NH2	9.34	124.97	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	211	ARG	CB-CG-CD	9.29	135.76	111.60
1	B	341	LYS	CA-CB-CG	9.12	133.46	113.40
1	B	244	TYR	CB-CG-CD1	9.05	126.43	121.00
1	B	177	LEU	CA-CB-CG	9.01	136.01	115.30
1	B	244	TYR	CB-CG-CD2	-8.74	115.75	121.00
1	B	127	MET	C-N-CA	8.69	143.43	121.70
1	A	297	ARG	CD-NE-CZ	8.69	135.76	123.60
1	A	552	GLN	CA-CB-CG	8.52	132.13	113.40
1	B	221	LYS	CB-CG-CD	8.45	133.58	111.60
1	A	221	LYS	CB-CG-CD	8.32	133.25	111.60
1	A	129	ARG	NE-CZ-NH1	8.12	124.36	120.30
1	B	128	ASN	CB-CA-C	-8.10	94.21	110.40
1	A	251	ASP	CB-CG-OD1	8.04	125.53	118.30
1	A	128	ASN	CB-CA-C	-7.91	94.57	110.40
1	A	489	ARG	NE-CZ-NH1	7.84	124.22	120.30
1	A	536	ASP	CB-CG-OD1	7.79	125.31	118.30
1	B	211	ARG	CB-CG-CD	7.76	131.78	111.60
1	A	128	ASN	N-CA-CB	-7.74	96.67	110.60
1	B	231	GLU	OE1-CD-OE2	-7.72	114.04	123.30
1	B	482	ARG	NE-CZ-NH2	-7.59	116.50	120.30
1	A	504	ARG	CG-CD-NE	7.54	127.64	111.80
1	B	504	ARG	CG-CD-NE	7.51	127.56	111.80
1	A	463	ARG	NE-CZ-NH2	-7.50	116.55	120.30
1	A	127	MET	C-N-CA	7.46	140.34	121.70
1	A	129	ARG	CD-NE-CZ	7.44	134.01	123.60
1	B	297	ARG	NE-CZ-NH2	-7.17	116.71	120.30
1	B	129	ARG	CD-NE-CZ	7.00	133.40	123.60
1	A	330	ARG	NE-CZ-NH2	6.97	123.79	120.30
1	B	196	MET	CG-SD-CE	6.95	111.31	100.20
1	B	38	GLU	OE1-CD-OE2	-6.81	115.13	123.30
1	B	505	THR	CB-CA-C	-6.80	93.23	111.60
1	B	457	THR	CB-CA-C	-6.77	93.32	111.60
1	A	457	THR	CB-CA-C	-6.75	93.36	111.60
1	A	7	PHE	CA-CB-CG	6.70	129.99	113.90
1	B	142	VAL	N-CA-CB	-6.63	96.92	111.50
1	B	297	ARG	CG-CD-NE	-6.61	97.91	111.80
1	B	211	ARG	NE-CZ-NH1	-6.61	117.00	120.30
1	B	122	ASP	CB-CG-OD2	6.58	124.22	118.30
1	A	279	TYR	CB-CG-CD2	6.40	124.84	121.00
1	A	482	ARG	CD-NE-CZ	6.40	132.56	123.60
1	A	142	VAL	N-CA-CB	-6.39	97.44	111.50
1	B	104	ARG	CD-NE-CZ	6.34	132.48	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	128	ASN	N-CA-CB	-6.29	99.28	110.60
1	B	222	ARG	NE-CZ-NH1	-6.26	117.17	120.30
1	B	231	GLU	CG-CD-OE2	6.19	130.68	118.30
1	B	560	LEU	CA-C-O	-6.16	107.16	120.10
1	B	183	ARG	CD-NE-CZ	6.06	132.09	123.60
1	B	350	ARG	NH1-CZ-NH2	6.05	126.06	119.40
1	A	128	ASN	N-CA-C	6.04	127.31	111.00
1	A	505	THR	CB-CA-C	-6.04	95.30	111.60
1	A	560	LEU	CA-C-O	-6.00	107.49	120.10
1	A	457	THR	N-CA-CB	5.98	121.67	110.30
1	B	398	ARG	CD-NE-CZ	5.96	131.95	123.60
1	B	104	ARG	NE-CZ-NH2	-5.96	117.32	120.30
1	B	211	ARG	N-CA-CB	5.92	121.25	110.60
1	A	398	ARG	NH1-CZ-NH2	-5.90	112.91	119.40
1	B	7	PHE	CA-CB-CG	5.90	128.06	113.90
1	B	426	SER	N-CA-CB	-5.90	101.65	110.50
1	A	170	ASP	CB-CG-OD1	5.89	123.60	118.30
1	A	300	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	A	279	TYR	CB-CG-CD1	-5.81	117.51	121.00
1	A	297	ARG	CG-CD-NE	-5.79	99.65	111.80
1	A	515	GLU	OE1-CD-OE2	-5.69	116.47	123.30
1	A	8	ARG	CB-CG-CD	5.68	126.36	111.60
1	A	396	ARG	CD-NE-CZ	5.68	131.55	123.60
1	A	183	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	A	154	TYR	CB-CG-CD2	-5.65	117.61	121.00
1	A	136	GLU	CA-CB-CG	5.64	125.80	113.40
1	A	312	ARG	NE-CZ-NH2	5.61	123.10	120.30
1	A	244	TYR	CB-CG-CD2	-5.60	117.64	121.00
1	B	104	ARG	NE-CZ-NH1	5.59	123.09	120.30
1	B	7	PHE	N-CA-C	5.59	126.08	111.00
1	B	128	ASN	N-CA-C	5.59	126.08	111.00
1	A	222	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	A	190	TYR	CB-CG-CD2	-5.57	117.66	121.00
1	B	214	MET	CG-SD-CE	5.56	109.10	100.20
1	A	244	TYR	CB-CG-CD1	5.55	124.33	121.00
1	B	192	ASP	CB-CG-OD1	5.55	123.29	118.30
1	A	129	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	A	183	ARG	NE-CZ-NH1	-5.51	117.55	120.30
1	A	505	THR	N-CA-CB	5.51	120.76	110.30
1	B	255	SER	N-CA-CB	5.48	118.72	110.50
1	A	463	ARG	NE-CZ-NH1	5.47	123.03	120.30
1	B	8	ARG	NE-CZ-NH2	5.46	123.03	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	489	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	B	288	ASP	CB-CG-OD1	5.40	123.16	118.30
1	B	457	THR	N-CA-CB	5.39	120.54	110.30
1	B	170	ASP	CB-CG-OD1	5.39	123.15	118.30
1	B	114	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	B	136	GLU	CA-CB-CG	5.38	125.24	113.40
1	B	37	VAL	CA-C-O	-5.38	108.81	120.10
1	A	177	LEU	CB-CG-CD1	5.38	120.14	111.00
1	B	195	MET	CG-SD-CE	-5.37	91.61	100.20
1	B	146	VAL	CA-CB-CG2	-5.36	102.86	110.90
1	B	552	GLN	CB-CG-CD	5.35	125.50	111.60
1	B	505	THR	N-CA-CB	5.34	120.45	110.30
1	A	552	GLN	CB-CG-CD	5.33	125.45	111.60
1	A	276	TYR	CB-CG-CD2	5.31	124.19	121.00
1	B	190	TYR	CB-CG-CD1	5.24	124.14	121.00
1	B	489	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	B	516	THR	O-C-N	-5.21	114.37	122.70
1	A	182	GLU	OE1-CD-OE2	5.17	129.50	123.30
1	B	303	MET	O-C-N	-5.16	114.44	122.70
1	A	7	PHE	N-CA-C	5.13	124.86	111.00
1	A	8	ARG	CA-CB-CG	5.13	124.69	113.40
1	B	330	ARG	N-CA-CB	-5.12	101.38	110.60
1	B	48	ASP	CB-CG-OD1	5.10	122.89	118.30
1	B	177	LEU	CB-CG-CD1	5.08	119.64	111.00
1	A	509	PHE	CB-CG-CD1	-5.08	117.25	120.80
1	A	183	ARG	NH1-CZ-NH2	5.07	124.97	119.40
1	A	290	LYS	CB-CG-CD	5.06	124.76	111.60
1	A	255	SER	N-CA-CB	5.06	118.09	110.50
1	B	506	HIS	N-CA-CB	-5.06	101.49	110.60
1	B	8	ARG	CB-CG-CD	5.06	124.75	111.60
1	B	190	TYR	CB-CG-CD2	-5.02	117.99	121.00
1	A	244	TYR	C-N-CA	-5.01	111.77	122.30
1	B	201	GLU	OE1-CD-OE2	5.01	129.32	123.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	350	ARG	Sidechain
1	B	350	ARG	Sidechain

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4346	0	4288	83	4
1	B	4346	0	4287	100	4
2	A	53	0	31	5	0
2	B	53	0	31	8	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	207	0	0	5	0
4	B	165	0	0	6	0
All	All	9178	0	8643	184	4

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 (184) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:600:FAD:H8A	2:B:600:FAD:C5B	1.66	1.25
2:A:600:FAD:C8A	2:A:600:FAD:H51A	1.69	1.20
2:A:600:FAD:C5B	2:A:600:FAD:H8A	1.73	1.19
2:B:600:FAD:C8A	2:B:600:FAD:H51A	1.73	1.18
1:B:349:GLY:H	1:B:352:ASN:HD21	1.09	0.99
1:A:349:GLY:H	1:A:352:ASN:HD21	0.99	0.97
1:A:550:PRO:HB2	1:A:552:GLN:NE2	1.81	0.95
1:B:550:PRO:HB2	1:B:552:GLN:HE22	1.35	0.91
1:B:550:PRO:HB2	1:B:552:GLN:NE2	1.89	0.88
1:B:57:THR:HG22	1:B:74:ILE:HD11	1.56	0.86
1:A:57:THR:HG22	1:A:74:ILE:HD11	1.59	0.85
2:B:600:FAD:H51A	2:B:600:FAD:H8A	0.84	0.80
1:A:550:PRO:HB2	1:A:552:GLN:HE22	1.46	0.79
1:A:516:THR:HG21	4:A:2173:HOH:O	1.86	0.76
2:B:600:FAD:O2'	2:B:600:FAD:H9	1.90	0.71
1:B:312:ARG:HG2	1:B:457:THR:HG23	1.74	0.70
1:A:167:ASP:OD1	1:A:186:GLY:HA3	1.94	0.68
1:A:428:ILE:HD11	1:A:503:TYR:HB3	1.76	0.66
2:B:600:FAD:C8A	2:B:600:FAD:C5B	2.51	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:GLY:N	1:A:352:ASN:HD21	1.83	0.65
1:A:312:ARG:HG2	1:A:457:THR:HG23	1.78	0.65
1:B:79:ASN:ND2	1:B:82:ASP:H	1.95	0.64
2:A:600:FAD:H8A	2:A:600:FAD:H51A	0.78	0.64
1:A:519:TRP:CZ3	1:B:211:ARG:HG2	2.32	0.64
1:B:79:ASN:HD22	1:B:82:ASP:H	1.44	0.64
1:B:194:TRP:O	1:B:197:HIS:HD2	1.81	0.64
1:B:310:THR:HG22	1:B:459:THR:HG22	1.80	0.63
1:B:480:GLN:O	1:B:484:VAL:HG23	1.99	0.63
1:B:102:ILE:HG12	1:B:175:SER:HB2	1.80	0.63
1:B:505:THR:HG23	4:B:2133:HOH:O	1.99	0.63
1:A:79:ASN:ND2	1:A:82:ASP:H	1.97	0.63
1:B:167:ASP:OD1	1:B:186:GLY:HA3	1.98	0.63
1:A:188:THR:HB	1:A:189:PRO:CD	2.30	0.62
1:B:297:ARG:HB3	1:B:298:PRO:CD	2.29	0.62
1:B:188:THR:HB	1:B:189:PRO:CD	2.30	0.61
1:A:194:TRP:O	1:A:197:HIS:HD2	1.83	0.60
1:B:51:TYR:CE1	1:B:104:ARG:HD3	2.36	0.60
1:B:304:ALA:HB2	4:B:2103:HOH:O	2.01	0.60
1:A:505:THR:HG21	1:A:513:ILE:HD12	1.82	0.60
1:B:189:PRO:HB2	1:B:270:MET:HE3	1.82	0.60
1:B:516:THR:HG21	4:B:2131:HOH:O	2.01	0.60
1:A:310:THR:HG22	1:A:459:THR:HG22	1.83	0.60
1:B:552:GLN:NE2	1:B:552:GLN:H	2.00	0.60
1:A:513:ILE:O	1:A:516:THR:HB	2.02	0.59
1:A:211:ARG:HG2	1:B:519:TRP:CZ3	2.37	0.59
1:B:282:THR:HG22	1:B:352:ASN:HD22	1.67	0.59
1:A:244:TYR:OH	1:B:195:MET:HG2	2.04	0.58
1:B:438:MET:HG2	1:B:500:TRP:HH2	1.67	0.58
1:B:505:THR:HG21	1:B:513:ILE:HD12	1.85	0.58
1:A:297:ARG:HB3	1:A:298:PRO:CD	2.33	0.58
1:A:349:GLY:H	1:A:352:ASN:ND2	1.84	0.58
1:A:282:THR:HG22	1:A:352:ASN:HD22	1.69	0.57
1:B:349:GLY:H	1:B:352:ASN:ND2	1.92	0.57
1:B:133:VAL:HG11	1:B:160:LEU:HD13	1.86	0.57
1:B:51:TYR:CZ	1:B:104:ARG:HD3	2.40	0.57
1:B:229:LYS:HD2	1:B:231:GLU:OE2	2.04	0.56
2:B:600:FAD:O2'	2:B:600:FAD:C9	2.52	0.56
1:A:333:PRO:HB2	1:A:448:GLN:NE2	2.20	0.56
1:A:519:TRP:CE3	1:B:211:ARG:HG2	2.42	0.55
1:A:385:PHE:HB3	1:A:386:PRO:HD2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:438:MET:HG2	1:A:500:TRP:HH2	1.71	0.55
1:A:98:TRP:CD2	1:A:113:PRO:HA	2.42	0.55
1:A:189:PRO:HG2	1:A:270:MET:HE1	1.89	0.55
1:B:282:THR:HG22	1:B:352:ASN:ND2	2.22	0.55
1:B:149:HIS:HD1	1:B:408:TYR:HH	1.53	0.54
1:B:513:ILE:O	1:B:516:THR:HB	2.06	0.54
1:B:428:ILE:HD11	1:B:503:TYR:HB3	1.88	0.54
1:A:417:LEU:HB3	1:A:418:PRO:HD2	1.88	0.54
1:A:217:LEU:CD2	1:B:516:THR:HG23	2.37	0.53
1:B:443:THR:HG21	1:B:469:VAL:HG21	1.91	0.53
1:A:197:HIS:HE1	1:A:251:ASP:OD2	1.92	0.53
1:B:310:THR:CG2	1:B:459:THR:HG22	2.39	0.53
1:A:282:THR:HG22	1:A:352:ASN:ND2	2.23	0.52
1:B:40:ILE:HD11	1:B:74:ILE:CD1	2.39	0.52
1:A:51:TYR:CZ	1:A:104:ARG:HD3	2.44	0.52
1:B:189:PRO:HG2	1:B:270:MET:HE1	1.91	0.52
1:B:385:PHE:HB3	1:B:386:PRO:HD2	1.91	0.52
1:B:419:ASN:HB2	1:B:474:ASN:OD1	2.09	0.52
1:B:90:ALA:HB1	1:B:95:PHE:O	2.09	0.52
1:B:457:THR:HG21	4:B:2104:HOH:O	2.09	0.51
2:A:600:FAD:C8A	2:A:600:FAD:C5B	2.57	0.51
1:B:13:PRO:HG3	1:B:95:PHE:CE1	2.44	0.51
1:A:505:THR:HG23	4:A:2174:HOH:O	2.11	0.51
1:A:40:ILE:HD11	1:A:74:ILE:CD1	2.41	0.51
1:B:505:THR:CG2	1:B:513:ILE:HD12	2.40	0.51
1:B:102:ILE:HB	2:B:600:FAD:O2P	2.11	0.51
1:A:79:ASN:HD22	1:A:82:ASP:H	1.59	0.51
1:B:333:PRO:HB2	1:B:448:GLN:OE1	2.11	0.50
1:B:464:GLU:OE1	1:B:466:HIS:ND1	2.40	0.50
1:A:505:THR:CG2	1:A:513:ILE:HD12	2.41	0.50
1:A:550:PRO:HB2	1:A:552:GLN:HE21	1.68	0.50
1:A:14:PRO:HG3	1:A:558:TRP:CZ2	2.47	0.50
1:A:211:ARG:HG2	1:B:519:TRP:CE3	2.47	0.50
1:B:98:TRP:CD2	1:B:113:PRO:HA	2.46	0.50
1:A:257:SER:HA	4:A:2107:HOH:O	2.12	0.50
1:A:516:THR:HG23	1:B:217:LEU:CD2	2.41	0.50
1:B:297:ARG:HB3	1:B:298:PRO:HD3	1.95	0.49
1:A:189:PRO:HB2	1:A:270:MET:HE3	1.94	0.49
1:A:552:GLN:H	1:A:552:GLN:NE2	2.09	0.49
1:A:51:TYR:CE1	1:A:104:ARG:HD3	2.47	0.49
1:A:361:GLU:N	1:A:362:PRO:HD2	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:600:FAD:H9	2:A:600:FAD:O2'	2.12	0.49
1:B:312:ARG:HD3	1:B:354:TYR:CE1	2.48	0.49
1:B:361:GLU:N	1:B:362:PRO:HD2	2.27	0.49
1:A:62:HIS:O	1:A:481:LYS:HE3	2.13	0.49
1:B:198:SER:O	1:B:240:HIS:HD2	1.96	0.49
1:A:90:ALA:HB1	1:A:95:PHE:O	2.13	0.48
1:B:505:THR:HG21	1:B:509:PHE:HB2	1.95	0.48
1:A:188:THR:CB	1:A:189:PRO:CD	2.91	0.48
1:A:480:GLN:O	1:A:484:VAL:HG23	2.13	0.48
1:A:312:ARG:HD3	1:A:354:TYR:CE1	2.47	0.48
1:B:349:GLY:N	1:B:352:ASN:HD21	1.93	0.48
1:B:28:ILE:O	1:B:32:VAL:HG22	2.14	0.48
1:A:297:ARG:HB3	1:A:298:PRO:HD3	1.95	0.47
1:B:188:THR:CB	1:B:189:PRO:CD	2.91	0.47
1:A:102:ILE:HG12	1:A:175:SER:HB2	1.95	0.47
1:A:195:MET:HG2	1:B:244:TYR:OH	2.15	0.47
1:B:197:HIS:HE1	1:B:251:ASP:OD2	1.97	0.47
1:A:330:ARG:NH2	1:A:333:PRO:O	2.47	0.47
1:A:195:MET:SD	1:B:195:MET:HE1	2.54	0.47
1:A:290:LYS:HB2	1:A:437:MET:CG	2.44	0.47
1:B:471:ILE:HG21	1:B:484:VAL:HG13	1.96	0.47
1:A:516:THR:HG22	1:A:517:TYR:CD1	2.50	0.47
1:B:505:THR:HG22	1:B:506:HIS:H	1.80	0.47
1:A:310:THR:CG2	1:A:459:THR:HG22	2.44	0.46
1:A:387:GLU:HG2	4:A:2136:HOH:O	2.16	0.46
1:B:167:ASP:OD1	1:B:193:HIS:NE2	2.49	0.46
1:B:309:PRO:HB2	1:B:353:PHE:CE1	2.51	0.46
1:A:443:THR:HG21	1:A:469:VAL:HG21	1.98	0.45
1:A:419:ASN:HB2	1:A:474:ASN:OD1	2.17	0.45
1:B:271:PRO:O	1:B:272:ASN:C	2.54	0.45
1:A:133:VAL:HG11	1:A:160:LEU:HD13	1.97	0.45
1:A:198:SER:O	1:A:240:HIS:HD2	2.00	0.45
1:A:7:PHE:HB3	4:A:2007:HOH:O	2.16	0.45
1:A:107:GLY:HA2	1:A:422:ALA:O	2.17	0.45
1:A:143:GLU:HB3	1:A:144:PRO:HD2	1.98	0.45
1:A:28:ILE:O	1:A:32:VAL:HG22	2.16	0.45
1:A:330:ARG:NH1	1:A:338:GLU:OE1	2.50	0.45
1:A:229:LYS:HD2	1:A:231:GLU:OE2	2.18	0.44
1:A:427:PRO:HA	1:A:502:GLU:HA	2.00	0.44
1:B:14:PRO:HG3	1:B:558:TRP:CZ2	2.52	0.44
1:A:187:TYR:O	1:A:307:ASN:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:ILE:HG12	1:B:89:LEU:HD12	2.00	0.44
1:B:330:ARG:NH2	1:B:333:PRO:O	2.48	0.44
1:A:13:PRO:HG3	1:A:95:PHE:CE1	2.53	0.43
1:B:62:HIS:O	1:B:481:LYS:HE3	2.18	0.43
1:B:188:THR:HB	1:B:189:PRO:HD2	1.99	0.43
1:A:188:THR:HB	1:A:189:PRO:HD3	2.00	0.43
1:B:290:LYS:HB2	1:B:437:MET:HG3	2.01	0.43
1:B:51:TYR:O	1:B:54:PRO:HD3	2.19	0.43
1:A:108:TYR:CZ	1:A:504:ARG:HG2	2.54	0.42
1:B:123:MET:O	1:B:127:MET:HB2	2.19	0.42
1:B:257:SER:HA	4:B:2154:HOH:O	2.18	0.42
1:B:518:ASN:O	1:B:519:TRP:C	2.57	0.42
1:B:433:GLY:O	1:B:437:MET:HG2	2.18	0.42
1:B:549:TRP:CH2	1:B:558:TRP:HB3	2.55	0.42
1:B:102:ILE:CG1	1:B:175:SER:HB2	2.49	0.42
1:B:417:LEU:HB3	1:B:418:PRO:HD2	2.02	0.42
1:B:385:PHE:HB3	1:B:386:PRO:CD	2.50	0.42
1:B:101:SER:O	1:B:124:GLY:HA3	2.20	0.42
1:B:194:TRP:O	1:B:197:HIS:CD2	2.65	0.42
1:B:169:PRO:HB3	2:B:600:FAD:N3	2.34	0.42
1:A:195:MET:HE1	1:B:195:MET:SD	2.60	0.42
1:A:38:GLU:N	1:A:74:ILE:O	2.50	0.41
1:B:144:PRO:HA	1:B:177:LEU:HG	2.01	0.41
1:B:183:ARG:HG2	1:B:194:TRP:HB2	2.02	0.41
1:B:312:ARG:CG	1:B:457:THR:HG23	2.48	0.41
1:B:55:THR:HG21	1:B:58:HIS:CE1	2.55	0.41
1:A:477:ASP:O	1:A:481:LYS:HG3	2.20	0.41
1:B:143:GLU:HB3	1:B:144:PRO:HD2	2.02	0.41
1:B:333:PRO:CB	1:B:448:GLN:OE1	2.68	0.41
1:A:312:ARG:HD3	1:A:354:TYR:CD1	2.56	0.41
1:A:491:LEU:HD23	1:A:491:LEU:HA	1.85	0.41
1:A:332:GLU:HB3	1:A:333:PRO:CD	2.51	0.41
1:B:79:ASN:C	1:B:79:ASN:HD22	2.24	0.41
1:A:143:GLU:HB3	1:A:144:PRO:CD	2.52	0.40
1:B:545:LYS:HD2	4:B:2155:HOH:O	2.20	0.40
1:B:545:LYS:O	1:B:546:SER:HB2	2.21	0.40
1:B:521:ASN:O	1:B:522:SER:HB2	2.21	0.40
1:B:555:HIS:HB3	1:B:559:LYS:HE3	2.04	0.40
1:B:200:MET:HE3	1:B:265:ILE:HD12	2.03	0.40
1:A:518:ASN:O	1:A:519:TRP:C	2.58	0.40
1:B:231:GLU:HG3	1:B:231:GLU:H	1.31	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:TRP:CG	1:A:113:PRO:HA	2.57	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:VAL:O	1:B:391:GLU:OE2[6_655]	1.81	0.39
1:A:50:SER:OG	1:B:412:LYS:NZ[6_655]	1.94	0.26
1:A:419:ASN:OD1	1:B:329:SER:OG[6_655]	2.16	0.04
1:A:373:ASP:OD2	1:B:288:ASP:OD1[4_665]	2.19	0.01

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	546/560 (98%)	523 (96%)	23 (4%)	0	100	100
1	B	546/560 (98%)	523 (96%)	23 (4%)	0	100	100
All	All	1092/1120 (98%)	1046 (96%)	46 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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	469/480 (98%)	441 (94%)	28 (6%)	19 22
1	B	469/480 (98%)	444 (95%)	25 (5%)	22 27
All	All	938/960 (98%)	885 (94%)	53 (6%)	20 24

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	PHE
1	A	8	ARG
1	A	41	SER
1	A	63	VAL
1	A	79	ASN
1	A	95	PHE
1	A	128	ASN
1	A	142	VAL
1	A	149	HIS
1	A	175	SER
1	A	177	LEU
1	A	200	MET
1	A	231	GLU
1	A	297	ARG
1	A	301	LEU
1	A	303	MET
1	A	329	SER
1	A	331	THR
1	A	336	ASP
1	A	387	GLU
1	A	391	GLU
1	A	432	SER
1	A	457	THR
1	A	502	GLU
1	A	503	TYR
1	A	505	THR
1	A	516	THR
1	A	552	GLN
1	B	7	PHE
1	B	8	ARG
1	B	63	VAL
1	B	79	ASN
1	B	95	PHE
1	B	128	ASN
1	B	142	VAL

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Mol	Chain	Res	Type
1	B	149	HIS
1	B	177	LEU
1	B	200	MET
1	B	231	GLU
1	B	301	LEU
1	B	303	MET
1	B	329	SER
1	B	331	THR
1	B	336	ASP
1	B	391	GLU
1	B	448	GLN
1	B	457	THR
1	B	466	HIS
1	B	467	HIS
1	B	503	TYR
1	B	505	THR
1	B	516	THR
1	B	552	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	128	ASN
1	A	158	ASN
1	A	197	HIS
1	A	240	HIS
1	A	352	ASN
1	A	439	GLN
1	A	467	HIS
1	A	485	GLN
1	A	520	ASN
1	A	552	GLN
1	A	555	HIS
1	B	79	ASN
1	B	128	ASN
1	B	197	HIS
1	B	240	HIS
1	B	352	ASN
1	B	439	GLN
1	B	467	HIS
1	B	520	ASN

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Mol	Chain	Res	Type
1	B	552	GLN
1	B	555	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	ACT	A	601	-	1,3,3	3.69	1 (100%)	0,3,3	0.00	-
2	FAD	B	600	-	51,58,58	2.06	12 (23%)	60,89,89	2.06	17 (28%)
2	FAD	A	600	-	51,58,58	1.95	10 (19%)	60,89,89	2.14	15 (25%)
3	ACT	B	601	-	1,3,3	3.75	1 (100%)	0,3,3	0.00	-

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	B	600	-	-	10/30/50/50	0/6/6/6
2	FAD	A	600	-	-	7/30/50/50	0/6/6/6

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	C4X-C10	7.35	1.46	1.38
2	B	600	FAD	C4X-C10	6.81	1.45	1.38
2	B	600	FAD	O4B-C1B	5.94	1.49	1.41
2	B	600	FAD	O5'-C5'	4.47	1.62	1.44
2	A	600	FAD	PA-O2A	-4.46	1.34	1.55
2	A	600	FAD	O4B-C1B	4.35	1.47	1.41
2	B	600	FAD	PA-O2A	-4.12	1.36	1.55
3	B	601	ACT	CH3-C	3.75	1.53	1.48
3	A	601	ACT	CH3-C	3.69	1.53	1.48
2	A	600	FAD	P-O2P	-3.58	1.38	1.55
2	B	600	FAD	C2B-C1B	-3.57	1.48	1.53
2	A	600	FAD	O5'-C5'	3.39	1.57	1.44
2	A	600	FAD	C2-N1	-3.23	1.31	1.38
2	B	600	FAD	C4-N3	3.18	1.38	1.33
2	B	600	FAD	P-O2P	-3.17	1.40	1.55
2	A	600	FAD	PA-O5B	-2.89	1.47	1.59
2	B	600	FAD	C2-N3	2.79	1.43	1.38
2	A	600	FAD	C2B-C1B	-2.64	1.49	1.53
2	A	600	FAD	C2-N3	2.62	1.43	1.38
2	B	600	FAD	O4'-C4'	2.56	1.48	1.43
2	B	600	FAD	C9A-N10	2.54	1.41	1.38
2	A	600	FAD	C9A-N10	2.52	1.41	1.38
2	B	600	FAD	C2-N1	-2.40	1.33	1.38
2	B	600	FAD	C2A-N3A	2.23	1.35	1.32

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	600	FAD	C4-N3-C2	7.47	121.45	115.14
2	B	600	FAD	C4-N3-C2	5.97	120.19	115.14
2	A	600	FAD	P-O3P-PA	5.74	152.51	132.83
2	B	600	FAD	P-O3P-PA	4.95	149.79	132.83
2	A	600	FAD	C4X-C4-N3	-4.82	116.83	123.43
2	B	600	FAD	C4X-C4-N3	-4.66	117.05	123.43
2	B	600	FAD	C1B-N9A-C4A	-4.47	118.79	126.64
2	B	600	FAD	O5B-PA-O1A	-4.33	92.17	109.07
2	A	600	FAD	O5B-PA-O1A	-4.05	93.23	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	600	FAD	C1B-N9A-C4A	-3.82	119.93	126.64
2	A	600	FAD	C2A-N1A-C6A	3.70	125.08	118.75
2	B	600	FAD	N6A-C6A-N1A	3.46	125.75	118.57
2	A	600	FAD	C4A-C5A-N7A	3.30	112.84	109.40
2	A	600	FAD	O4B-C4B-C5B	-3.28	98.58	109.37
2	A	600	FAD	N3A-C2A-N1A	-3.23	123.63	128.68
2	B	600	FAD	C4A-C5A-N7A	3.21	112.74	109.40
2	B	600	FAD	C5X-C9A-N10	-2.83	115.66	117.72
2	B	600	FAD	C5A-C6A-N1A	-2.77	114.06	120.35
2	B	600	FAD	O4B-C4B-C5B	-2.76	100.28	109.37
2	B	600	FAD	O3B-C3B-C4B	2.72	118.93	111.05
2	A	600	FAD	O2P-P-O1P	2.55	124.83	112.24
2	B	600	FAD	C7-C6-C5X	2.47	124.71	121.22
2	B	600	FAD	C9A-C5X-N5	2.46	126.21	122.36
2	B	600	FAD	C2A-N1A-C6A	2.45	122.94	118.75
2	A	600	FAD	O5'-P-O1P	-2.31	100.04	109.07
2	B	600	FAD	O3'-C3'-C2'	-2.31	103.24	108.81
2	A	600	FAD	C5A-C6A-N1A	-2.26	115.23	120.35
2	A	600	FAD	PA-O5B-C5B	2.21	134.63	121.68
2	A	600	FAD	O2A-PA-O5B	2.15	117.72	107.75
2	A	600	FAD	C9-C8-C7	-2.12	116.33	119.91
2	B	600	FAD	O5'-P-O1P	-2.02	101.17	109.07
2	B	600	FAD	C6-C5X-N5	-2.01	116.84	119.05

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	600	FAD	C3'-C4'-C5'-O5'
2	B	600	FAD	C5'-O5'-P-O1P
2	A	600	FAD	C5'-O5'-P-O1P
2	B	600	FAD	O4'-C4'-C5'-O5'
2	A	600	FAD	C3'-C4'-C5'-O5'
2	A	600	FAD	C4'-C5'-O5'-P
2	B	600	FAD	C5'-O5'-P-O3P
2	A	600	FAD	C5'-O5'-P-O3P
2	B	600	FAD	O4B-C4B-C5B-O5B
2	B	600	FAD	C5B-O5B-PA-O1A
2	B	600	FAD	C5'-O5'-P-O2P
2	A	600	FAD	C5'-O5'-P-O2P
2	A	600	FAD	O4'-C4'-C5'-O5'
2	B	600	FAD	C4'-C5'-O5'-P

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Mol	Chain	Res	Type	Atoms
2	B	600	FAD	C3B-C4B-C5B-O5B
2	B	600	FAD	P-O3P-PA-O2A
2	A	600	FAD	O4B-C4B-C5B-O5B

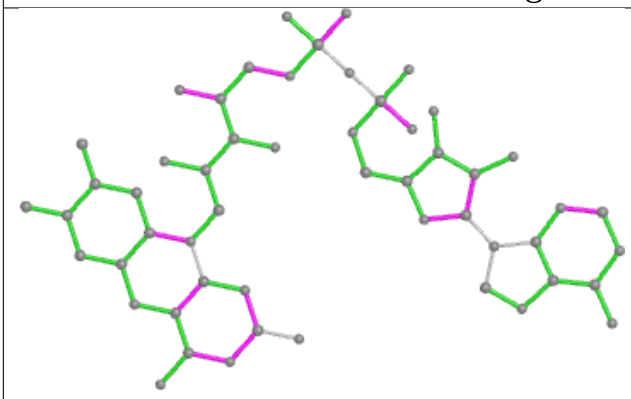
There are no ring outliers.

2 monomers are involved in 13 short contacts:

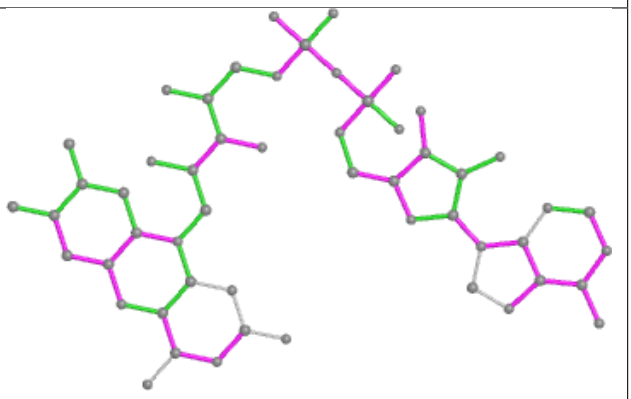
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	600	FAD	8	0
2	A	600	FAD	5	0

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 then 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.

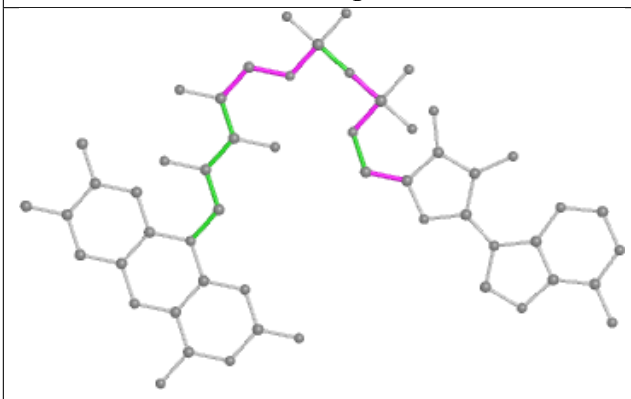
Ligand FAD B 600



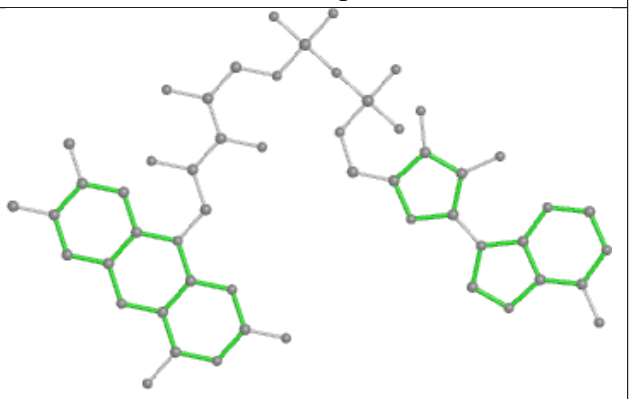
Bond lengths



Bond angles

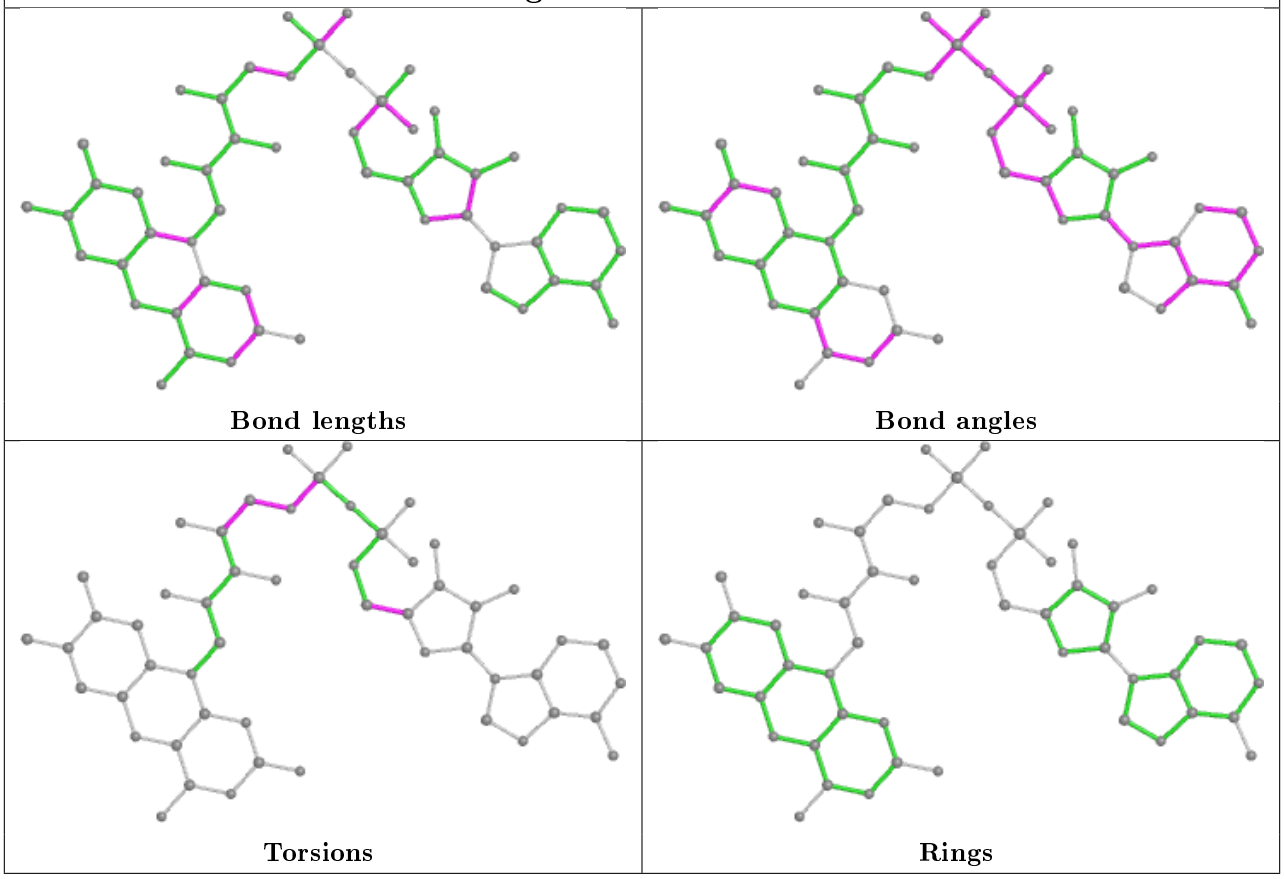


Torsions



Rings

Ligand FAD A 600



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	550/560 (98%)	-0.25	5 (0%) 84 83	18, 30, 53, 74	15 (2%)
1	B	550/560 (98%)	-0.05	25 (4%) 33 32	18, 30, 53, 74	15 (2%)
All	All	1100/1120 (98%)	-0.15	30 (2%) 54 52	18, 30, 54, 74	30 (2%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	51	TYR	5.1
1	A	329	SER	4.6
1	B	159	ASN	4.5
1	B	329	SER	4.2
1	B	49	GLY	3.9
1	A	159	ASN	3.7
1	B	52	MET	3.7
1	B	63	VAL	3.4
1	B	387	GLU	3.1
1	B	50	SER	3.0
1	B	55	THR	3.0
1	B	64	MET	2.9
1	B	478	LEU	2.9
1	B	332	GLU	2.8
1	B	62	HIS	2.8
1	B	337	GLU	2.7
1	B	157	ALA	2.6
1	B	56	HIS	2.5
1	B	158	ASN	2.4
1	B	60	PRO	2.4
1	B	7	PHE	2.4
1	B	448	GLN	2.3
1	B	369	GLU	2.2
1	B	65	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	7	PHE	2.2
1	B	331	THR	2.1
1	B	67	ASP	2.1
1	B	61	HIS	2.1
1	A	478	LEU	2.0
1	A	162	ASP	2.0

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 carbohydrates in this entry.

6.4 Ligands (i)

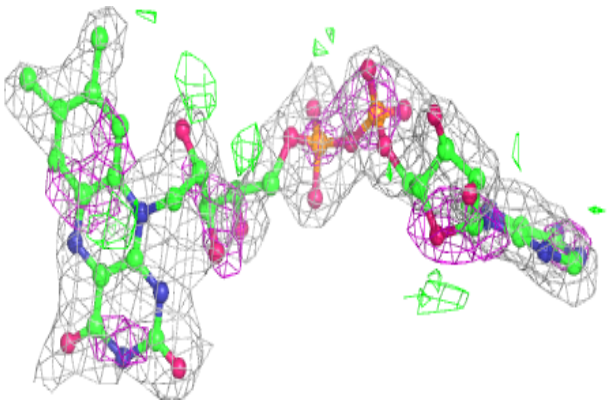
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

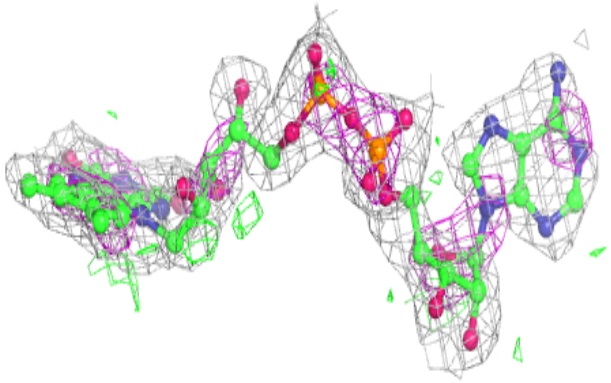
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FAD	B	600	53/53	0.87	0.21	34,37,39,40	0
3	ACT	A	601	4/4	0.92	0.17	34,35,35,39	0
2	FAD	A	600	53/53	0.94	0.16	33,36,39,40	0
3	ACT	B	601	4/4	0.95	0.12	34,35,36,39	0

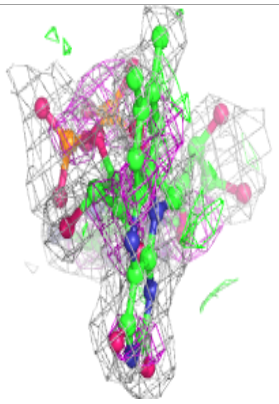
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD B 600:

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

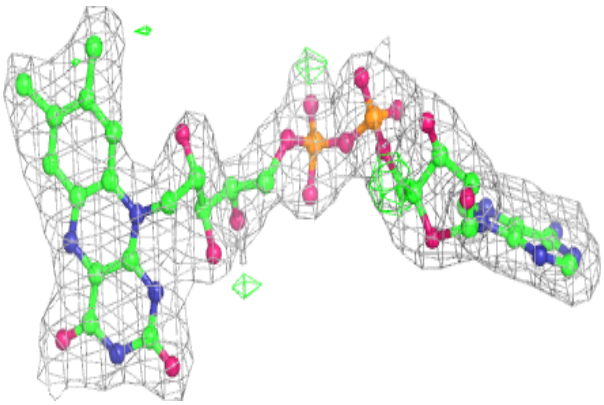


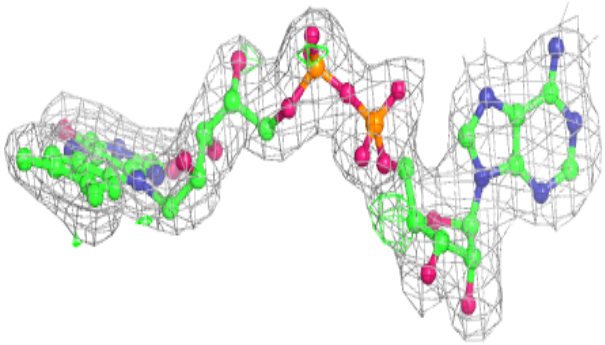


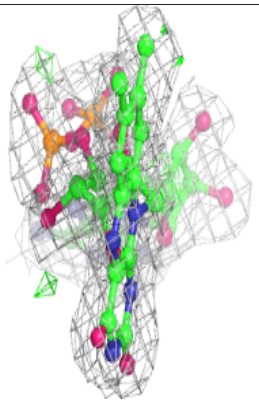


Electron density around FAD A 600:

$2mF_o$-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.