



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:36 PM UTC

PDB ID : 4NWZ / pdb_00004nwz
Title : Structure of bacterial type II NADH dehydrogenase from *Caldalkalibacillus thermarum* at 2.5Å resolution
Authors : Nakatani, Y.; Heikal, A.; Lott, J.S.; Sazanov, L.A.; Baker, E.N.; Cook, G.M.
Deposited on : 2013-12-07
Resolution : 2.50 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

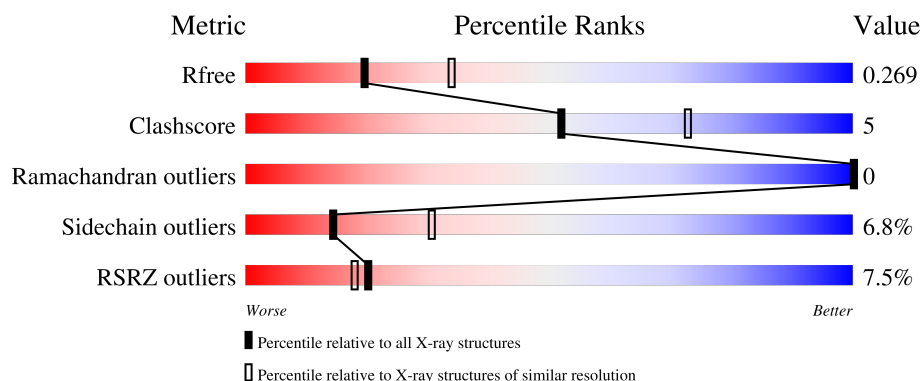
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	405	0% 82% 14% ..
1	B	405	5% 82% 13% ..
1	C	405	2% 83% 12% ..
1	D	405	20% 78% 12% . 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12454 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 FAD-dependent pyridine nucleotide-disulfide oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	84	1	0
			3049	1944	529	567	9			
1	B	395	Total	C	N	O	S	40	2	0
			3055	1948	529	569	9			
1	C	394	Total	C	N	O	S	80	0	0
			3039	1938	528	564	9			
1	D	373	Total	C	N	O	S	106	1	0
			2886	1846	493	538	9			

There are 24 discrepancies between the modelled and reference sequences:

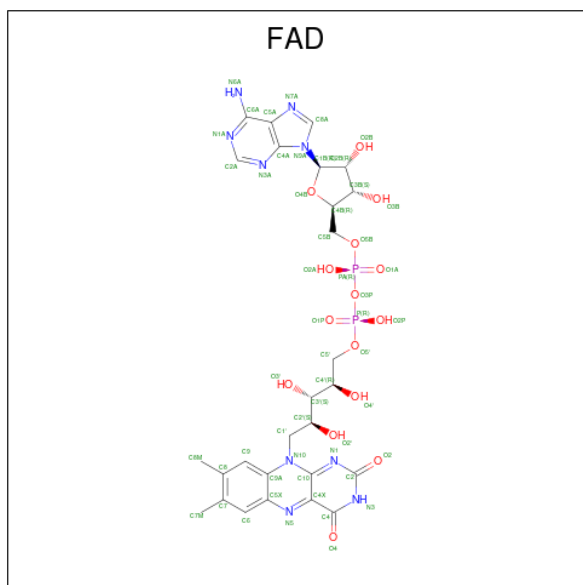
Chain	Residue	Modelled	Actual	Comment	Reference
A	400	HIS	-	expression tag	UNP F5L3B8
A	401	HIS	-	expression tag	UNP F5L3B8
A	402	HIS	-	expression tag	UNP F5L3B8
A	403	HIS	-	expression tag	UNP F5L3B8
A	404	HIS	-	expression tag	UNP F5L3B8
A	405	HIS	-	expression tag	UNP F5L3B8
B	400	HIS	-	expression tag	UNP F5L3B8
B	401	HIS	-	expression tag	UNP F5L3B8
B	402	HIS	-	expression tag	UNP F5L3B8
B	403	HIS	-	expression tag	UNP F5L3B8
B	404	HIS	-	expression tag	UNP F5L3B8
B	405	HIS	-	expression tag	UNP F5L3B8
C	400	HIS	-	expression tag	UNP F5L3B8
C	401	HIS	-	expression tag	UNP F5L3B8
C	402	HIS	-	expression tag	UNP F5L3B8
C	403	HIS	-	expression tag	UNP F5L3B8
C	404	HIS	-	expression tag	UNP F5L3B8
C	405	HIS	-	expression tag	UNP F5L3B8
D	400	HIS	-	expression tag	UNP F5L3B8
D	401	HIS	-	expression tag	UNP F5L3B8
D	402	HIS	-	expression tag	UNP F5L3B8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	403	HIS	-	expression tag	UNP F5L3B8
D	404	HIS	-	expression tag	UNP F5L3B8
D	405	HIS	-	expression tag	UNP F5L3B8

- 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	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

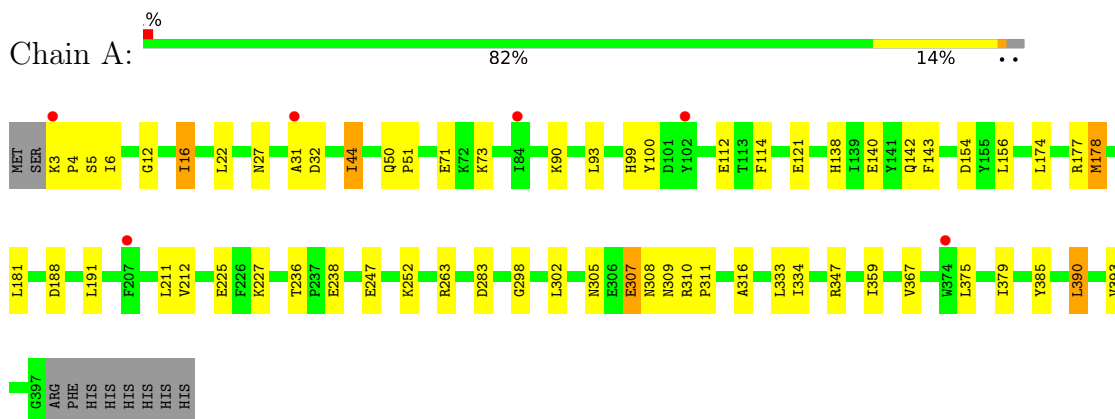
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	65	Total	O	0	0
			65	65		
3	B	77	Total	O	0	0
			77	77		
3	C	53	Total	O	0	0
			53	53		
3	D	18	Total	O	0	0
			18	18		

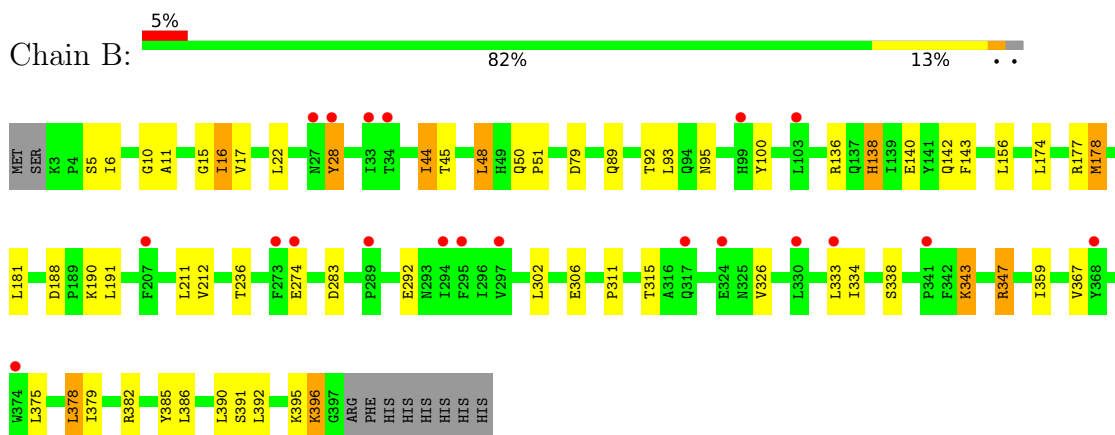
3 Residue-property plots

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

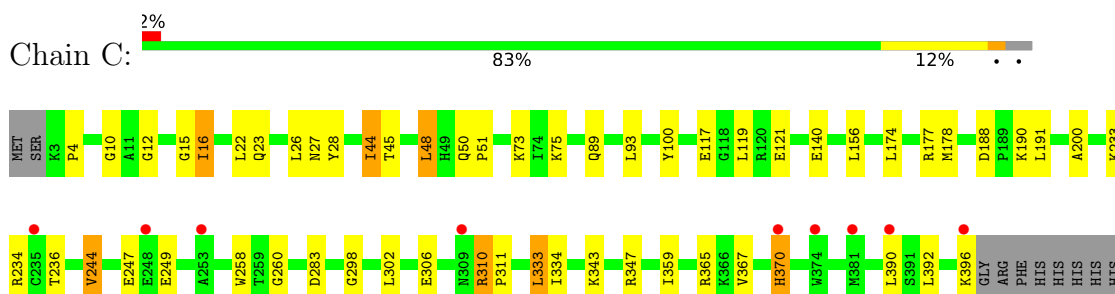
- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



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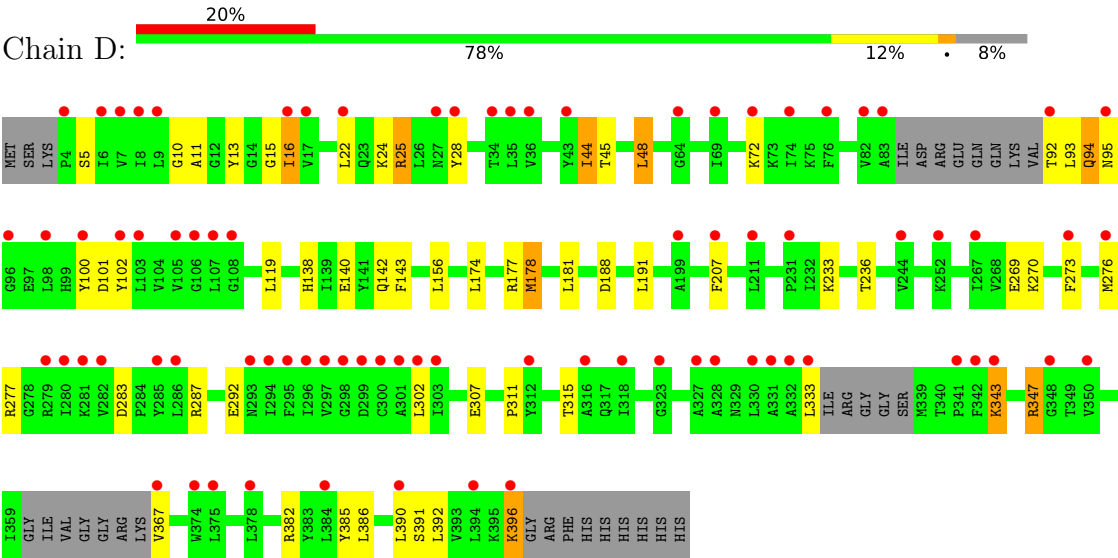


- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



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● Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.79Å 114.48Å 130.59Å 90.00° 92.04° 90.00°	Depositor
Resolution (Å)	64.61 – 2.50 64.61 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (64.61-2.50) 99.5 (64.61-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.217 , 0.269 0.219 , 0.269	Depositor DCC
R_{free} test set	3717 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	69.2	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 62.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12454	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.83	3/3111 (0.1%)	0.99	10/4214 (0.2%)
1	B	0.81	4/3120 (0.1%)	0.88	7/4226 (0.2%)
1	C	0.77	5/3098 (0.2%)	0.85	4/4197 (0.1%)
1	D	0.73	7/2945 (0.2%)	0.95	13/3991 (0.3%)
All	All	0.79	19/12274 (0.2%)	0.92	34/16628 (0.2%)

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

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	28	TYR	CA-CB	-13.06	1.32	1.53
1	D	269	GLU	CA-CB	-11.26	1.35	1.53
1	A	308	ASN	N-CA	11.19	1.57	1.46
1	D	102	TYR	CA-CB	9.22	1.74	1.53
1	C	370	HIS	CB-CG	-8.82	1.37	1.50
1	B	28	TYR	CB-CG	7.87	1.69	1.51
1	B	378	LEU	CB-CG	-7.17	1.39	1.53
1	A	347	ARG	CB-CG	-6.98	1.31	1.52
1	C	310	ARG	CA-CB	-6.90	1.44	1.53
1	B	190	LYS	CG-CD	6.86	1.73	1.52
1	D	25	ARG	CA-CB	-6.81	1.43	1.53
1	C	306	GLU	CA-CB	6.29	1.64	1.53
1	D	101	ASP	CA-CB	6.28	1.63	1.53
1	B	343	LYS	CA-CB	6.18	1.61	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	95	ASN	CA-CB	6.02	1.61	1.53
1	D	347	ARG	CB-CG	-5.90	1.34	1.52
1	D	276	MET	CA-CB	-5.81	1.45	1.53
1	C	75	LYS	CB-CG	-5.34	1.36	1.52
1	A	252	LYS	CG-CD	5.25	1.68	1.52

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	307	GLU	O-C-N	-21.51	88.00	122.41
1	D	28	TYR	CB-CA-C	16.21	135.81	109.70
1	A	307	GLU	CB-CA-C	-14.34	82.46	110.27
1	D	95	ASN	CA-CB-CG	-12.71	99.89	112.60
1	D	28	TYR	CA-CB-CG	10.52	132.83	113.90
1	A	307	GLU	CA-C-N	10.51	138.41	121.18
1	A	307	GLU	C-N-CA	10.51	138.41	121.18
1	D	28	TYR	N-CA-CB	-10.45	93.07	109.69
1	A	310	ARG	CB-CG-CD	-10.03	88.24	111.30
1	A	247	GLU	CB-CA-C	9.70	125.64	109.84
1	D	343	LYS	CA-CB-CG	8.60	131.30	114.10
1	C	247	GLU	CB-CA-C	7.76	123.15	109.72
1	C	28	TYR	CB-CA-C	7.64	121.31	109.84
1	D	307	GLU	CA-CB-CG	-7.29	99.51	114.10
1	B	190	LYS	CB-CG-CD	-6.93	95.36	111.30
1	B	347	ARG	CG-CD-NE	-6.88	96.85	112.00
1	C	347	ARG	CB-CG-CD	6.80	126.95	111.30
1	D	72	LYS	CB-CA-C	6.73	121.53	110.90
1	B	28	TYR	CB-CG-CD2	6.65	130.78	120.80
1	B	396	LYS	CB-CG-CD	6.63	126.55	111.30
1	D	95	ASN	CB-CA-C	-6.49	100.36	111.46
1	B	28	TYR	CB-CG-CD1	-6.33	111.31	120.80
1	D	207	PHE	CA-CB-CG	-6.24	107.56	113.80
1	D	270	LYS	CB-CA-C	-6.15	97.45	110.31
1	D	307	GLU	CB-CG-CD	-5.96	102.47	112.60
1	C	306	GLU	N-CA-CB	-5.92	100.75	110.40
1	A	238	GLU	CB-CG-CD	-5.88	102.60	112.60
1	D	273	PHE	CA-CB-CG	5.79	119.58	113.80
1	D	277	ARG	CA-CB-CG	-5.73	102.65	114.10
1	A	252	LYS	CB-CG-CD	-5.64	98.32	111.30
1	A	307	GLU	CA-CB-CG	-5.47	103.16	114.10
1	B	395	LYS	CA-CB-CG	5.11	124.31	114.10
1	B	138	HIS	N-CA-C	-5.09	105.81	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	308	ASN	N-CA-CB	-5.09	101.87	111.43

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	307	GLU	Mainchain

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	3049	0	3084	38	0
1	B	3055	0	3090	35	0
1	C	3039	0	3075	24	0
1	D	2886	0	2904	26	0
2	A	53	0	31	2	0
2	B	53	0	31	1	0
2	C	53	0	31	1	0
2	D	53	0	31	2	0
3	A	65	0	0	0	0
3	B	77	0	0	3	0
3	C	53	0	0	0	0
3	D	18	0	0	0	0
All	All	12454	0	12277	121	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 5.

All (121) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:GLU:HB3	1:C:234:ARG:HG2	1.57	0.86
1:B:178:MET:HA	1:B:178:MET:HE3	1.70	0.74
1:D:178:MET:HE3	1:D:178:MET:HA	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:MET:HA	1:A:178:MET:HE3	1.70	0.73
1:A:390:LEU:HA	1:A:393:VAL:HG22	1.71	0.72
1:C:26:LEU:O	1:C:73:LYS:NZ	2.22	0.72
1:A:302:LEU:HD11	1:A:311:PRO:HB3	1.72	0.70
1:D:386:LEU:HD23	1:D:386:LEU:O	1.93	0.67
1:B:386:LEU:O	1:B:386:LEU:HD23	1.95	0.66
1:A:4:PRO:HD2	1:A:31:ALA:HA	1.77	0.66
1:A:71:GLU:H	1:A:71:GLU:CD	2.06	0.62
1:D:5:SER:OG	1:D:100:TYR:HA	2.01	0.61
1:A:5:SER:OG	1:A:100:TYR:HA	2.01	0.60
1:D:315:THR:HB	2:D:601:FAD:O2	2.01	0.60
1:D:13:TYR:CD2	2:D:601:FAD:O4'	2.54	0.60
1:D:94:GLN:NE2	1:D:94:GLN:HA	2.17	0.58
1:C:396:LYS:HD2	1:C:396:LYS:C	2.29	0.58
1:B:178:MET:HE1	1:B:181:LEU:HD23	1.85	0.58
1:B:5:SER:OG	1:B:100:TYR:HA	2.05	0.57
1:C:45:THR:HA	1:C:48:LEU:HD22	1.86	0.57
1:B:338:SER:HB3	1:C:200:ALA:HB2	1.88	0.56
1:B:45:THR:HA	1:B:48:LEU:HD22	1.88	0.56
1:C:283:ASP:OD1	1:C:283:ASP:C	2.49	0.56
1:D:178:MET:HE1	1:D:181:LEU:HD23	1.87	0.56
1:A:188:ASP:HB3	1:A:191:LEU:HD12	1.88	0.56
1:D:386:LEU:HD23	1:D:386:LEU:C	2.29	0.55
1:A:71:GLU:OE2	1:A:71:GLU:N	2.22	0.55
1:D:45:THR:HA	1:D:48:LEU:HD22	1.87	0.55
1:D:283:ASP:OD1	1:D:283:ASP:C	2.50	0.55
1:B:283:ASP:C	1:B:283:ASP:OD1	2.50	0.54
1:A:178:MET:HE1	1:A:181:LEU:HD23	1.90	0.54
1:A:225[A]:GLU:OE2	1:A:227:LYS:NZ	2.30	0.53
1:D:302:LEU:HD11	1:D:311:PRO:HB3	1.90	0.53
1:B:188:ASP:HB3	1:B:191:LEU:HD12	1.90	0.53
1:B:386:LEU:HD23	1:B:386:LEU:C	2.34	0.53
1:C:188:ASP:HB3	1:C:191:LEU:HD12	1.90	0.53
1:C:302:LEU:HD11	1:C:311:PRO:HB3	1.90	0.53
1:A:44:ILE:CD1	1:A:44:ILE:N	2.72	0.52
1:D:188:ASP:HB3	1:D:191:LEU:HD12	1.90	0.52
1:A:112:GLU:OE2	1:A:263:ARG:HD2	2.10	0.51
1:A:283:ASP:OD1	1:A:283:ASP:C	2.53	0.51
1:B:89:GLN:HA	1:B:100:TYR:CZ	2.46	0.51
1:B:306:GLU:HG3	3:B:762:HOH:O	2.10	0.49
1:B:178:MET:HA	1:B:178:MET:CE	2.39	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:44:ILE:CD1	1:D:44:ILE:N	2.75	0.49
1:A:6:ILE:HD11	1:A:334:ILE:HD11	1.93	0.49
1:C:117:GLU:OE1	1:C:234:ARG:HD3	2.12	0.49
1:C:244:VAL:HG21	1:C:249:GLU:OE1	2.13	0.49
1:B:315:THR:HB	2:B:601:FAD:O2	2.12	0.49
1:B:326:VAL:HA	3:B:730:HOH:O	2.12	0.49
1:B:302:LEU:HD11	1:B:311:PRO:HB3	1.95	0.48
1:A:112:GLU:HG3	1:A:114:PHE:CD1	2.49	0.47
1:C:258:TRP:CZ2	1:C:260:GLY:HA3	2.49	0.47
1:C:298:GLY:HA3	2:C:601:FAD:O2P	2.14	0.47
1:D:287:ARG:NH1	1:D:292:GLU:HG3	2.30	0.47
1:A:112:GLU:HG3	1:A:114:PHE:CE1	2.49	0.47
1:D:385:TYR:C	1:D:385:TYR:CD1	2.93	0.46
1:B:143:PHE:CE2	1:B:178:MET:HE2	2.51	0.46
1:B:338:SER:CB	1:C:200:ALA:HB2	2.45	0.46
1:A:385:TYR:C	1:A:385:TYR:CD1	2.94	0.46
1:A:44:ILE:N	1:A:44:ILE:HD13	2.31	0.45
1:C:89:GLN:HA	1:C:100:TYR:CZ	2.51	0.45
1:A:143:PHE:CE2	1:A:178:MET:HE2	2.51	0.45
1:A:178:MET:HE3	1:A:181:LEU:HB3	1.99	0.45
1:A:178:MET:HA	1:A:178:MET:CE	2.42	0.45
1:D:386:LEU:C	1:D:386:LEU:CD2	2.90	0.45
1:C:4:PRO:HG2	1:C:334:ILE:HD12	1.98	0.44
1:C:44:ILE:N	1:C:44:ILE:CD1	2.80	0.44
1:A:114:PHE:CD1	1:A:114:PHE:N	2.85	0.44
1:A:71:GLU:CD	1:A:71:GLU:N	2.72	0.44
1:A:143:PHE:CE2	1:A:178:MET:CE	3.00	0.44
1:D:396:LYS:O	1:D:396:LYS:HD3	2.17	0.44
1:B:140:GLU:OE2	1:B:177:ARG:NH2	2.51	0.44
1:C:12:GLY:O	1:C:16:ILE:HB	2.17	0.44
1:A:316:ALA:HB2	2:A:601:FAD:H2'	2.00	0.43
1:D:143:PHE:CE2	1:D:178:MET:HE2	2.53	0.43
1:B:44:ILE:N	1:B:44:ILE:CD1	2.81	0.43
1:D:138:HIS:O	1:D:142:GLN:HG2	2.18	0.43
1:C:50:GLN:HB2	1:C:51:PRO:CD	2.48	0.43
1:B:382:ARG:HG2	1:B:382:ARG:HH11	1.83	0.43
1:C:140:GLU:OE2	1:C:177:ARG:NH2	2.52	0.43
1:A:12:GLY:O	1:A:16:ILE:HB	2.18	0.43
1:A:298:GLY:HA3	2:A:601:FAD:O2P	2.18	0.43
1:B:6:ILE:HD11	1:B:334:ILE:HD11	2.00	0.43
1:B:385:TYR:CD1	1:B:385:TYR:C	2.96	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:178:MET:HA	1:D:178:MET:CE	2.40	0.43
1:A:138:HIS:O	1:A:142:GLN:HG2	2.20	0.42
1:B:138:HIS:O	1:B:142:GLN:HG2	2.20	0.42
1:A:140:GLU:OE2	1:A:177:ARG:NH2	2.52	0.42
1:A:143:PHE:CZ	1:A:178:MET:HE1	2.55	0.42
1:B:178:MET:HE3	1:B:181:LEU:HB3	2.02	0.42
1:A:3:LYS:HD2	1:A:3:LYS:HA	1.89	0.42
1:A:143:PHE:CZ	1:A:178:MET:CE	3.03	0.42
1:B:143:PHE:CE2	1:B:178:MET:CE	3.03	0.42
1:C:10:GLY:O	1:C:15:GLY:HA3	2.19	0.42
1:B:79:ASP:OD2	1:B:95:ASN:HB2	2.20	0.42
1:B:89:GLN:HA	1:B:100:TYR:CE1	2.55	0.42
1:C:119:LEU:HD12	1:C:119:LEU:HA	1.93	0.42
1:B:375:LEU:O	1:B:379:ILE:HG13	2.20	0.41
1:A:5:SER:HA	1:A:32:ASP:HB2	2.02	0.41
1:A:375:LEU:O	1:A:379:ILE:HG13	2.21	0.41
1:D:140:GLU:OE2	1:D:177:ARG:NH2	2.53	0.41
1:B:136:ARG:HD2	3:B:712:HOH:O	2.20	0.41
1:B:211:LEU:O	1:B:212:VAL:C	2.64	0.41
1:D:11:ALA:HB1	1:D:16:ILE:HD13	2.02	0.41
1:A:50:GLN:HB2	1:A:51:PRO:CD	2.50	0.41
1:A:211:LEU:O	1:A:212:VAL:C	2.64	0.41
1:D:10:GLY:O	1:D:15:GLY:HA3	2.21	0.41
1:B:16:ILE:CG2	1:B:17:VAL:N	2.84	0.41
1:D:44:ILE:N	1:D:44:ILE:HD13	2.35	0.41
1:D:178:MET:HE3	1:D:181:LEU:HB3	2.03	0.41
1:C:333:LEU:HD12	1:C:333:LEU:HA	1.97	0.41
1:A:154:ASP:OD1	1:A:154:ASP:N	2.50	0.40
1:B:10:GLY:O	1:B:15:GLY:HA3	2.21	0.40
1:C:26:LEU:C	1:C:73:LYS:NZ	2.79	0.40
1:D:119:LEU:HD12	1:D:119:LEU:HA	1.94	0.40
1:B:50:GLN:HB2	1:B:51:PRO:CD	2.52	0.40
1:C:23:GLN:HA	1:C:73:LYS:HE2	2.03	0.40
1:A:305:ASN:O	1:A:309:ASN:N	2.53	0.40
1:B:11:ALA:HB1	1:B:16:ILE:HD13	2.03	0.40
1:B:16:ILE:HG22	1:B:17:VAL:N	2.35	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	394/405 (97%)	380 (96%)	14 (4%)	0	100	100
1	B	395/405 (98%)	384 (97%)	11 (3%)	0	100	100
1	C	392/405 (97%)	381 (97%)	11 (3%)	0	100	100
1	D	366/405 (90%)	353 (96%)	13 (4%)	0	100	100
All	All	1547/1620 (96%)	1498 (97%)	49 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	320/329 (97%)	303 (95%)	17 (5%)	20	42
1	B	321/329 (98%)	298 (93%)	23 (7%)	13	28
1	C	319/329 (97%)	296 (93%)	23 (7%)	13	28
1	D	304/329 (92%)	281 (92%)	23 (8%)	12	26
All	All	1264/1316 (96%)	1178 (93%)	86 (7%)	14	31

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	22	LEU

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Mol	Chain	Res	Type
1	A	27	ASN
1	A	44	ILE
1	A	73	LYS
1	A	90	LYS
1	A	93	LEU
1	A	99	HIS
1	A	121	GLU
1	A	156	LEU
1	A	174	LEU
1	A	178	MET
1	A	236	THR
1	A	333	LEU
1	A	359	ILE
1	A	367	VAL
1	A	390	LEU
1	B	16	ILE
1	B	22	LEU
1	B	28	TYR
1	B	44	ILE
1	B	48	LEU
1	B	92	THR
1	B	93	LEU
1	B	156	LEU
1	B	174	LEU
1	B	178	MET
1	B	236	THR
1	B	274	GLU
1	B	292	GLU
1	B	333	LEU
1	B	343	LYS
1	B	347	ARG
1	B	359	ILE
1	B	367	VAL
1	B	378	LEU
1	B	390	LEU
1	B	391	SER
1	B	392	LEU
1	B	396	LYS
1	C	16	ILE
1	C	22	LEU
1	C	27	ASN
1	C	44	ILE

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Mol	Chain	Res	Type
1	C	48	LEU
1	C	93	LEU
1	C	121	GLU
1	C	156	LEU
1	C	174	LEU
1	C	178	MET
1	C	190	LYS
1	C	233	LYS
1	C	236	THR
1	C	244	VAL
1	C	310	ARG
1	C	333	LEU
1	C	343	LYS
1	C	359	ILE
1	C	365	ARG
1	C	367	VAL
1	C	370	HIS
1	C	390	LEU
1	C	392	LEU
1	D	16	ILE
1	D	22	LEU
1	D	24	LYS
1	D	25	ARG
1	D	44	ILE
1	D	48	LEU
1	D	92	THR
1	D	93	LEU
1	D	94	GLN
1	D	156	LEU
1	D	174	LEU
1	D	178	MET
1	D	233	LYS
1	D	236	THR
1	D	333	LEU
1	D	343	LYS
1	D	347	ARG
1	D	367	VAL
1	D	382	ARG
1	D	390	LEU
1	D	391	SER
1	D	392	LEU
1	D	396	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	291	HIS
1	A	370	HIS
1	C	99	HIS
1	C	122	HIS
1	C	356	ASN
1	D	94	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
2	FAD	C	601	-	58,58,58	1.46	9 (15%)	85,89,89	1.77	19 (22%)
2	FAD	D	601	-	58,58,58	1.71	12 (20%)	85,89,89	1.47	18 (21%)
2	FAD	B	601	-	58,58,58	1.64	9 (15%)	85,89,89	1.58	16 (18%)
2	FAD	A	601	-	58,58,58	1.39	8 (13%)	85,89,89	1.67	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	C	601	-	-	5/34/50/50	0/6/6/6
2	FAD	D	601	-	-	6/34/50/50	0/6/6/6
2	FAD	B	601	-	-	1/34/50/50	0/6/6/6
2	FAD	A	601	-	-	1/34/50/50	0/6/6/6

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	FAD	C9A-C5X	5.46	1.49	1.41
2	D	601	FAD	C5A-C4A	5.17	1.48	1.39
2	A	601	FAD	C9A-C5X	4.93	1.49	1.41
2	B	601	FAD	C9A-C5X	4.78	1.48	1.41
2	C	601	FAD	C9A-C5X	4.18	1.47	1.41
2	B	601	FAD	C4'-C3'	-4.12	1.46	1.53
2	B	601	FAD	C5A-C4A	3.91	1.46	1.39
2	D	601	FAD	C8-C7	3.80	1.50	1.40
2	D	601	FAD	C8A-N7A	3.70	1.38	1.31
2	C	601	FAD	C5A-C4A	3.59	1.45	1.39
2	B	601	FAD	C5X-N5	-3.56	1.32	1.39
2	C	601	FAD	C5A-N7A	-3.28	1.33	1.39
2	B	601	FAD	C8A-N7A	3.28	1.38	1.31
2	D	601	FAD	PA-O3P	3.09	1.62	1.59
2	A	601	FAD	C8A-N7A	3.01	1.37	1.31
2	B	601	FAD	C8-C7	3.00	1.48	1.40
2	C	601	FAD	C8-C7	2.94	1.48	1.40
2	C	601	FAD	C5X-N5	-2.94	1.34	1.39
2	C	601	FAD	C4A-N9A	-2.92	1.31	1.37
2	D	601	FAD	C4A-N9A	-2.92	1.31	1.37
2	D	601	FAD	P-O3P	2.91	1.62	1.59
2	C	601	FAD	P-O3P	2.91	1.62	1.59
2	B	601	FAD	C5A-N7A	-2.82	1.33	1.39
2	A	601	FAD	C4A-N9A	-2.64	1.32	1.37
2	A	601	FAD	C4'-C3'	-2.63	1.48	1.53
2	A	601	FAD	C5X-N5	-2.56	1.34	1.39
2	B	601	FAD	PA-O3P	2.51	1.62	1.59
2	A	601	FAD	C8-C7	2.47	1.46	1.40
2	C	601	FAD	C4-N3	-2.42	1.34	1.38
2	D	601	FAD	C5A-N7A	-2.38	1.34	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	601	FAD	C5A-C6A	2.32	1.47	1.41
2	A	601	FAD	C5A-C4A	2.31	1.43	1.39
2	D	601	FAD	C4X-N5	2.31	1.35	1.30
2	D	601	FAD	C4-N3	-2.31	1.34	1.38
2	A	601	FAD	O4B-C1B	2.21	1.47	1.42
2	C	601	FAD	C5A-C6A	2.19	1.47	1.41
2	D	601	FAD	C1'-C2'	2.08	1.55	1.52
2	B	601	FAD	C2-N3	-2.07	1.34	1.39

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	FAD	C5A-C4A-N3A	-5.38	119.31	126.72
2	A	601	FAD	N3A-C2A-N1A	-4.84	121.25	128.58
2	B	601	FAD	C5A-C4A-N3A	-4.75	120.18	126.72
2	A	601	FAD	C5A-C4A-N3A	-4.49	120.53	126.72
2	C	601	FAD	N3A-C4A-N9A	4.32	134.51	127.17
2	D	601	FAD	C5A-C4A-N3A	-4.28	120.82	126.72
2	C	601	FAD	C4A-N9A-C8A	4.02	109.96	105.74
2	C	601	FAD	C4A-C5A-N7A	-3.96	106.06	110.58
2	B	601	FAD	N3A-C4A-N9A	3.82	133.67	127.17
2	C	601	FAD	N9A-C8A-N7A	-3.77	108.59	113.94
2	C	601	FAD	C2A-N3A-C4A	3.72	120.92	111.83
2	A	601	FAD	C2A-N3A-C4A	3.67	120.80	111.83
2	C	601	FAD	C5A-N7A-C8A	3.66	109.20	103.45
2	A	601	FAD	O2A-PA-O3P	-3.63	97.46	107.27
2	C	601	FAD	N3A-C2A-N1A	-3.61	123.11	128.58
2	D	601	FAD	N3A-C2A-N1A	-3.58	123.16	128.58
2	B	601	FAD	O4-C4-C4X	-3.48	117.35	126.53
2	A	601	FAD	O4-C4-C4X	-3.39	117.57	126.53
2	C	601	FAD	O4-C4-C4X	-3.34	117.72	126.53
2	A	601	FAD	N3A-C4A-N9A	3.33	132.84	127.17
2	A	601	FAD	O2P-P-O1P	3.22	127.44	112.44
2	A	601	FAD	C4A-N9A-C8A	3.03	108.92	105.74
2	B	601	FAD	O2A-PA-O3P	-3.02	99.11	107.27
2	D	601	FAD	N3A-C4A-N9A	3.00	132.26	127.17
2	D	601	FAD	C2A-N1A-C6A	2.99	123.64	118.73
2	B	601	FAD	C4A-C5A-N7A	-2.98	107.17	110.58
2	B	601	FAD	C4A-N9A-C8A	2.97	108.85	105.74
2	B	601	FAD	O2P-P-O3P	-2.92	99.39	107.27
2	A	601	FAD	O2-C2-N1	-2.91	116.97	121.80
2	D	601	FAD	C2A-N3A-C4A	2.90	118.92	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	N9A-C8A-N7A	-2.89	109.83	113.94
2	A	601	FAD	N9A-C8A-N7A	-2.88	109.86	113.94
2	C	601	FAD	C4X-C10-N1	-2.81	117.69	124.59
2	B	601	FAD	N3A-C2A-N1A	-2.79	124.36	128.58
2	D	601	FAD	C4A-C5A-N7A	-2.79	107.40	110.58
2	B	601	FAD	C4'-C3'-C2'	-2.69	109.09	113.57
2	B	601	FAD	C2A-N3A-C4A	2.64	118.27	111.83
2	B	601	FAD	C5A-N7A-C8A	2.61	107.55	103.45
2	C	601	FAD	C10-N1-C2	2.59	122.45	116.85
2	D	601	FAD	C4A-N9A-C8A	2.58	108.45	105.74
2	C	601	FAD	C4'-C3'-C2'	-2.58	109.28	113.57
2	A	601	FAD	C2A-N1A-C6A	2.56	122.94	118.73
2	B	601	FAD	O4-C4-N3	2.56	124.93	120.11
2	C	601	FAD	O3B-C3B-C2B	-2.55	103.63	111.82
2	B	601	FAD	O2P-P-O1P	2.48	123.97	112.44
2	C	601	FAD	C2B-C3B-C4B	2.39	107.23	102.61
2	D	601	FAD	C4X-C10-N10	2.36	119.86	116.48
2	A	601	FAD	C4A-N9A-C1B	-2.36	121.12	126.63
2	D	601	FAD	O2-C2-N1	-2.35	117.90	121.80
2	A	601	FAD	C4'-C3'-C2'	-2.33	109.69	113.57
2	D	601	FAD	C4A-N9A-C1B	-2.32	121.22	126.63
2	D	601	FAD	C4X-C10-N1	-2.31	118.93	124.59
2	A	601	FAD	C2B-C3B-C4B	2.27	106.99	102.61
2	C	601	FAD	O4B-C1B-N9A	2.26	112.42	108.09
2	A	601	FAD	C4-N3-C2	-2.24	121.66	125.64
2	A	601	FAD	C4X-C4-N3	2.21	118.88	113.25
2	D	601	FAD	C4-C4X-N5	2.21	121.26	118.21
2	D	601	FAD	C5X-C9A-N10	2.17	119.93	117.97
2	C	601	FAD	C9A-N10-C10	-2.15	117.47	120.75
2	B	601	FAD	C4X-C10-N1	-2.14	119.33	124.59
2	D	601	FAD	O2P-P-O3P	-2.14	101.49	107.27
2	A	601	FAD	C4X-C10-N10	2.12	119.52	116.48
2	D	601	FAD	O4B-C1B-N9A	2.12	112.16	108.09
2	C	601	FAD	C4-N3-C2	-2.09	121.93	125.64
2	C	601	FAD	C6A-C5A-N7A	2.08	136.10	132.09
2	A	601	FAD	C9A-N10-C10	-2.06	117.61	120.75
2	C	601	FAD	C4X-C4-N3	2.06	118.48	113.25
2	D	601	FAD	C4-N3-C2	-2.02	122.05	125.64
2	D	601	FAD	C9A-C5X-N5	-2.02	120.31	122.45
2	D	601	FAD	O4-C4-C4X	-2.02	121.21	126.53
2	B	601	FAD	C4X-C10-N10	2.01	119.36	116.48

There are no chirality outliers.

All (13) torsion outliers are listed below:

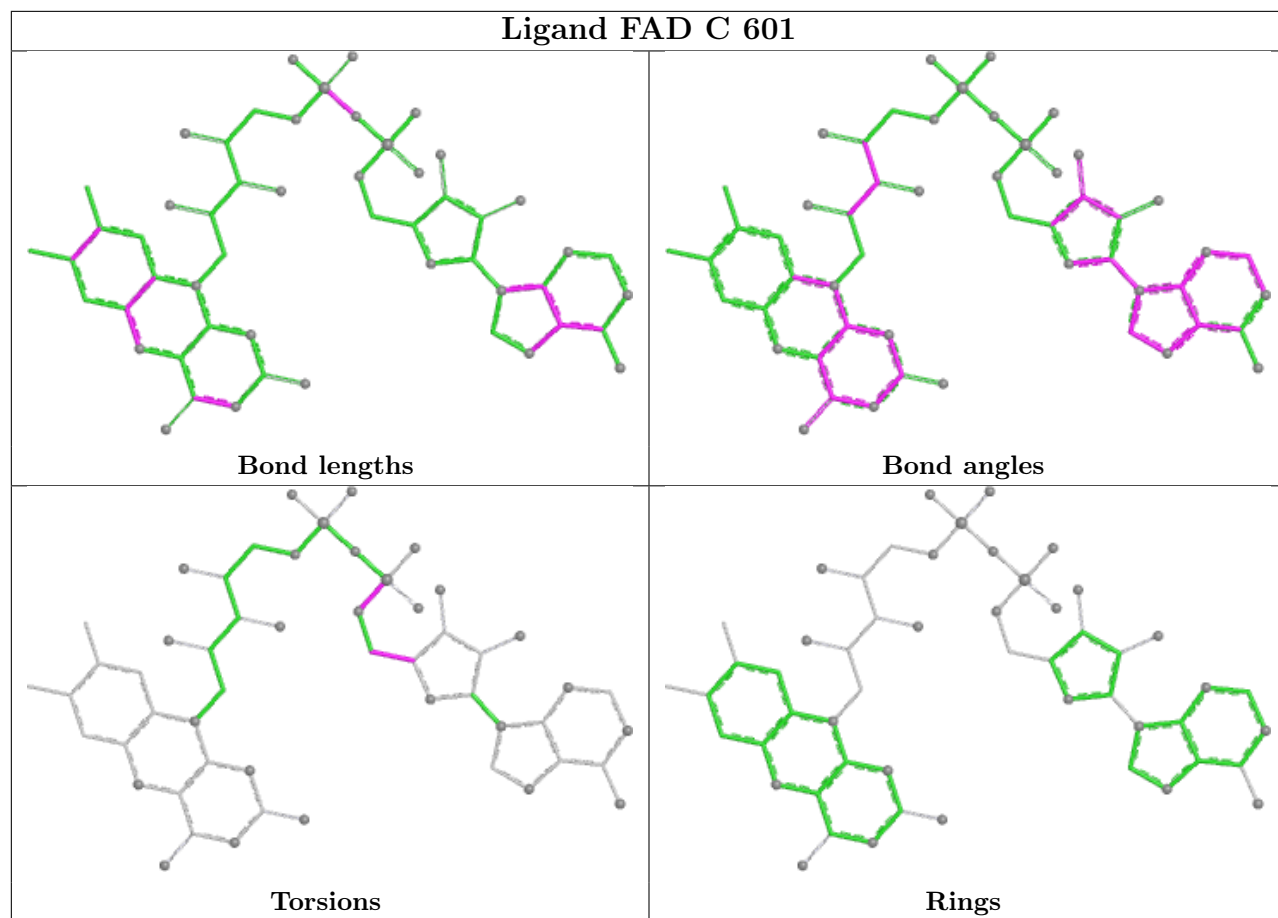
Mol	Chain	Res	Type	Atoms
2	C	601	FAD	C5B-O5B-PA-O1A
2	C	601	FAD	C5B-O5B-PA-O2A
2	D	601	FAD	C1'-C2'-C3'-C4'
2	D	601	FAD	C5'-O5'-P-O3P
2	C	601	FAD	O4B-C4B-C5B-O5B
2	B	601	FAD	PA-O3P-P-O5'
2	D	601	FAD	C1'-C2'-C3'-O3'
2	C	601	FAD	C5B-O5B-PA-O3P
2	C	601	FAD	C3B-C4B-C5B-O5B
2	D	601	FAD	O2'-C2'-C3'-O3'
2	D	601	FAD	O2'-C2'-C3'-C4'
2	A	601	FAD	O4B-C4B-C5B-O5B
2	D	601	FAD	N10-C1'-C2'-O2'

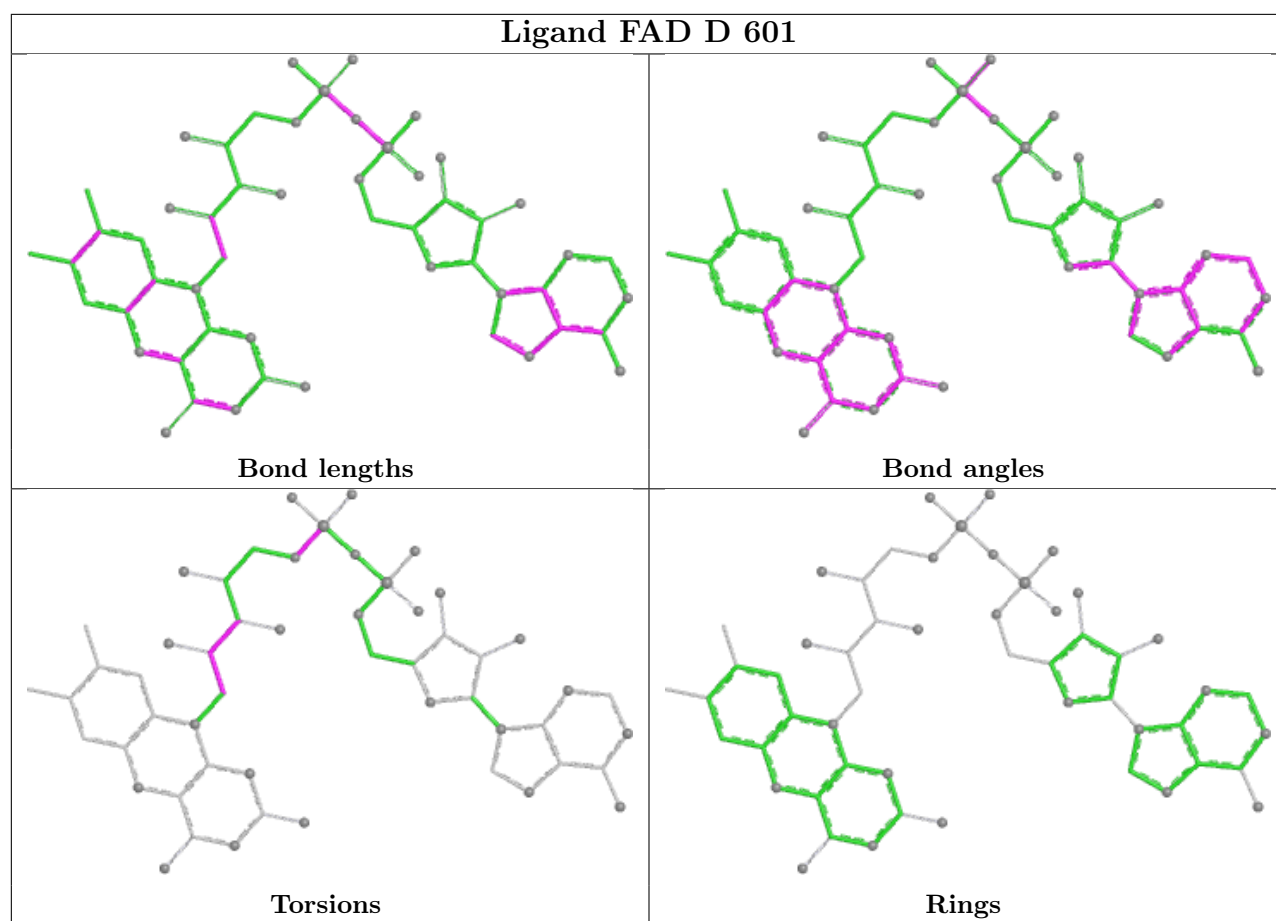
There are no ring outliers.

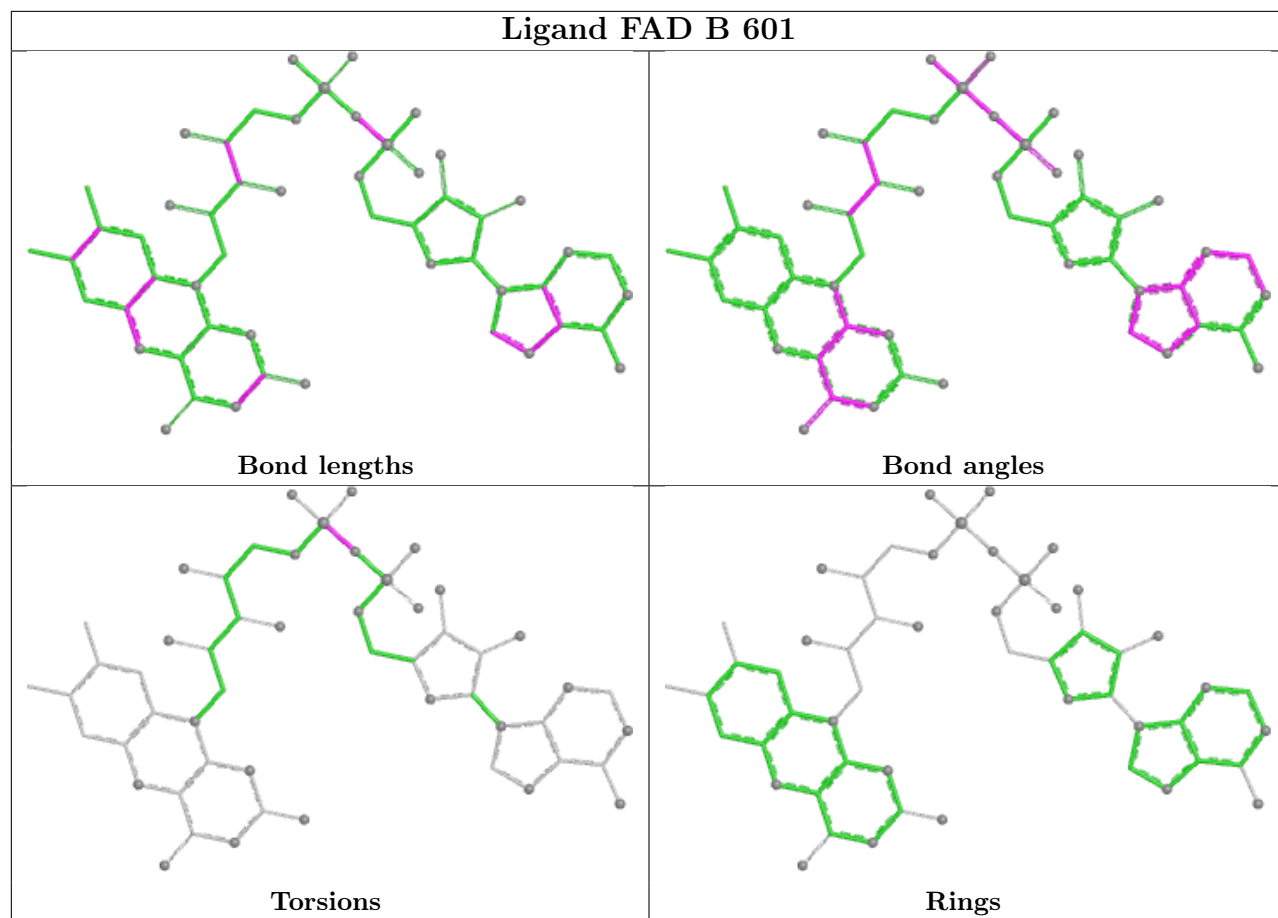
4 monomers are involved in 6 short contacts:

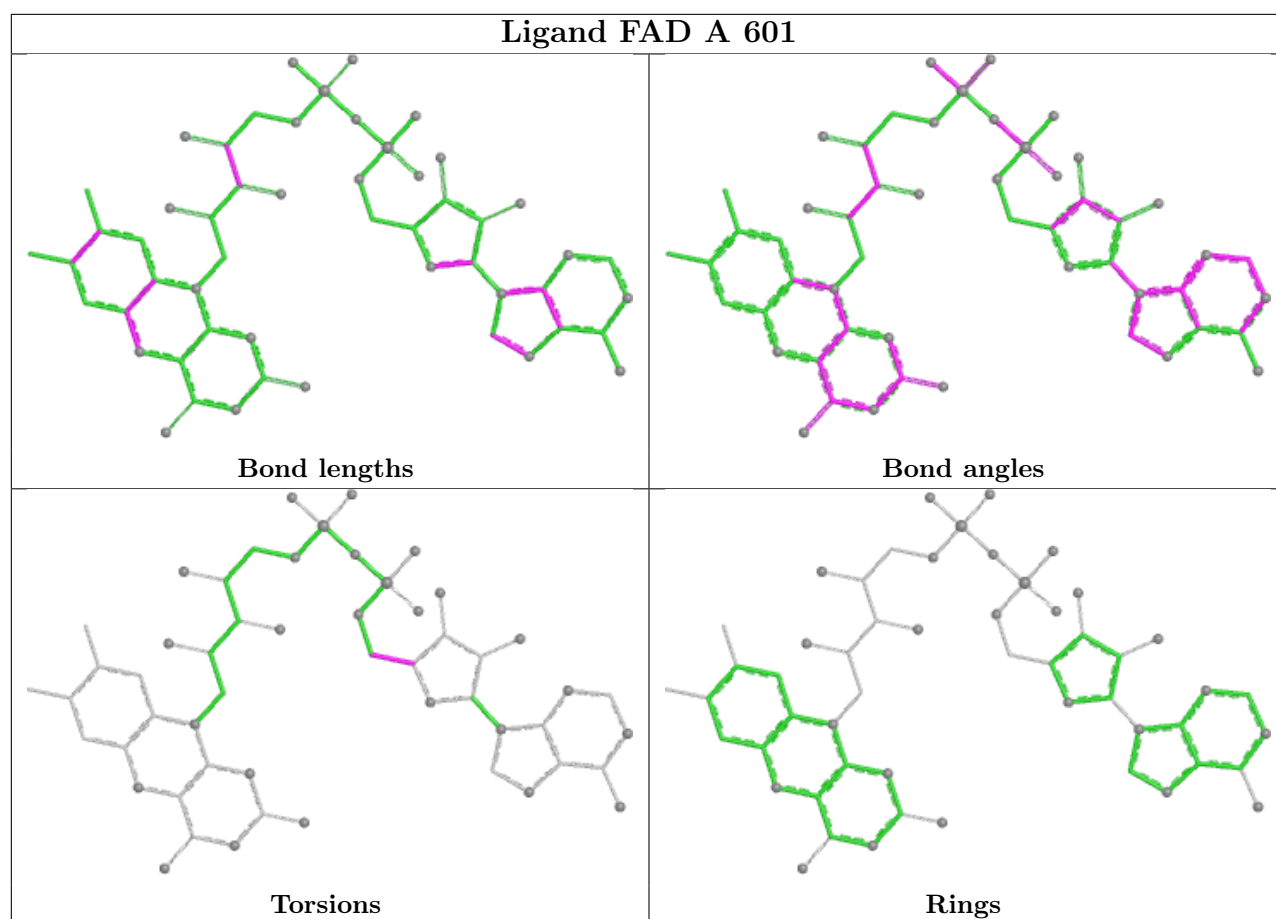
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	FAD	1	0
2	D	601	FAD	2	0
2	B	601	FAD	1	0
2	A	601	FAD	2	0

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.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	395/405 (97%)	0.23	6 (1%) 72 68	31, 65, 102, 130	20 (5%)
1	B	395/405 (97%)	0.41	20 (5%) 33 29	33, 69, 102, 126	11 (2%)
1	C	394/405 (97%)	0.34	9 (2%) 61 57	34, 74, 110, 135	17 (4%)
1	D	373/405 (92%)	1.20	81 (21%) 2 2	39, 98, 138, 164	21 (5%)
All	All	1557/1620 (96%)	0.54	116 (7%) 20 18	31, 74, 120, 164	69 (4%)

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	83	ALA	6.1
1	D	302	LEU	5.2
1	D	295	PHE	5.0
1	D	106	GLY	4.8
1	D	331	ALA	4.6
1	D	4	PRO	4.4
1	B	341	PRO	4.3
1	D	280	ILE	4.3
1	D	296	ILE	4.3
1	D	102	TYR	4.2
1	D	367	VAL	4.1
1	D	301	ALA	4.1
1	D	7	VAL	4.0
1	D	34	THR	4.0
1	D	211	LEU	3.9
1	D	282	VAL	3.9
1	D	76	PHE	3.8
1	C	370	HIS	3.7
1	D	332	ALA	3.7
1	D	267	ILE	3.7
1	D	318	ILE	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	92	THR	3.5
1	D	327	ALA	3.4
1	B	28	TYR	3.4
1	B	317	GLN	3.3
1	D	35	LEU	3.3
1	D	36	VAL	3.3
1	D	64	GLY	3.2
1	D	273	PHE	3.2
1	D	312	TYR	3.2
1	D	333	LEU	3.1
1	D	303	ILE	3.1
1	C	390	LEU	3.1
1	D	286	LEU	3.1
1	D	297	VAL	3.1
1	D	390	LEU	3.0
1	D	6	ILE	3.0
1	C	248	GLU	3.0
1	D	374	TRP	3.0
1	D	341	PRO	2.9
1	B	294	ILE	2.9
1	C	235	CYS	2.8
1	B	333	LEU	2.8
1	D	107	LEU	2.8
1	C	381	MET	2.8
1	D	342	PHE	2.8
1	D	348	GLY	2.8
1	D	96	GLY	2.8
1	D	350	VAL	2.8
1	A	374	TRP	2.7
1	D	330	LEU	2.7
1	D	199	ALA	2.7
1	D	299	ASP	2.7
1	B	295	PHE	2.7
1	D	16	ILE	2.7
1	D	82	VAL	2.7
1	D	231	PRO	2.7
1	C	374	TRP	2.7
1	D	298	GLY	2.6
1	A	84	ILE	2.6
1	D	396	LYS	2.6
1	D	69	ILE	2.6
1	D	375	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	98	LEU	2.6
1	D	316	ALA	2.6
1	B	34	THR	2.5
1	D	285	TYR	2.5
1	B	33	ILE	2.5
1	D	276	MET	2.5
1	A	31	ALA	2.5
1	D	279	ARG	2.4
1	D	281	LYS	2.4
1	D	343	LYS	2.4
1	D	244	VAL	2.4
1	D	100	TYR	2.4
1	D	207	PHE	2.4
1	D	252	LYS	2.4
1	D	300	CYS	2.3
1	B	207	PHE	2.3
1	D	8	ILE	2.3
1	B	99	HIS	2.3
1	C	309	ASN	2.3
1	B	103	LEU	2.3
1	D	103	LEU	2.3
1	D	105	VAL	2.3
1	B	289	PRO	2.3
1	B	330	LEU	2.2
1	C	396	LYS	2.2
1	D	293	ASN	2.2
1	B	368	TYR	2.2
1	A	3	LYS	2.2
1	D	9	LEU	2.2
1	D	384	LEU	2.2
1	D	394	LEU	2.2
1	B	324	GLU	2.2
1	B	297	VAL	2.2
1	D	22	LEU	2.2
1	B	27	ASN	2.2
1	C	253	ALA	2.2
1	D	328	ALA	2.2
1	D	28	TYR	2.1
1	D	74	ILE	2.1
1	B	274	GLU	2.1
1	B	374	TRP	2.1
1	D	27	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	108	GLY	2.1
1	D	95	ASN	2.1
1	D	323	GLY	2.1
1	D	294	ILE	2.1
1	A	207	PHE	2.0
1	A	102	TYR	2.0
1	D	43	TYR	2.0
1	B	273	PHE	2.0
1	D	17	VAL	2.0
1	D	378	LEU	2.0
1	D	72	LYS	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

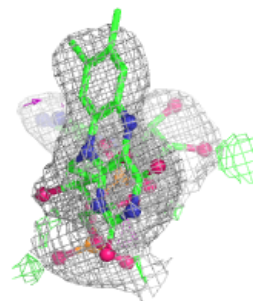
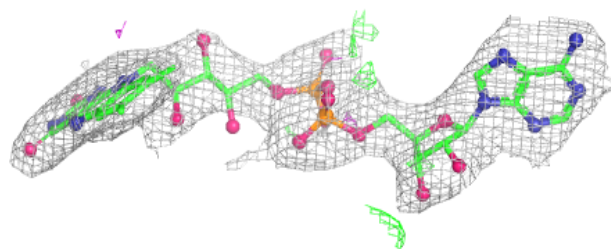
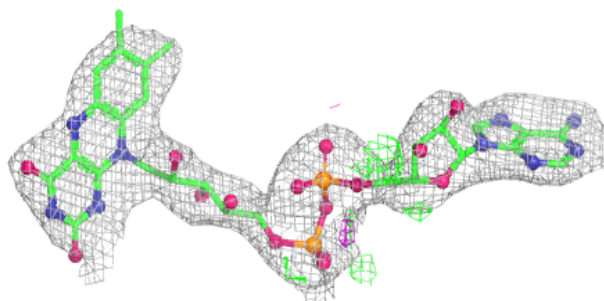
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FAD	D	601	53/53	0.92	0.12	63,80,106,121	0
2	FAD	B	601	53/53	0.97	0.07	28,46,58,63	0
2	FAD	C	601	53/53	0.97	0.07	44,54,75,88	0
2	FAD	A	601	53/53	0.97	0.07	39,49,59,64	0

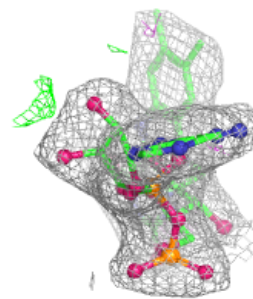
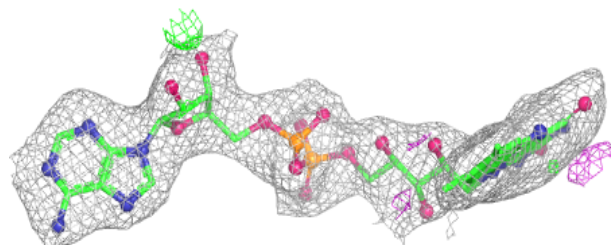
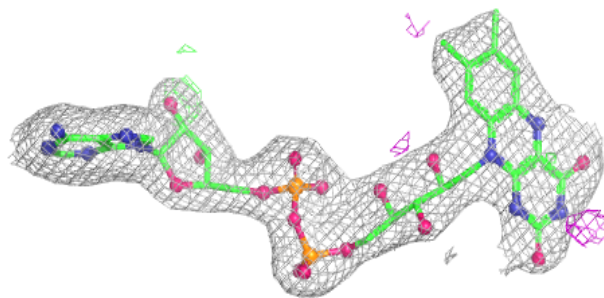
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD D 601:

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

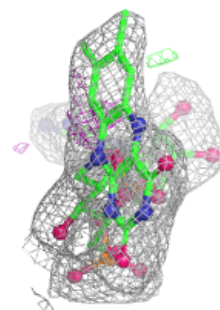
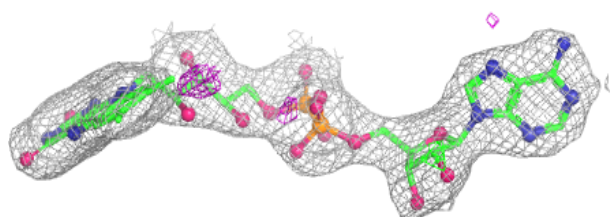
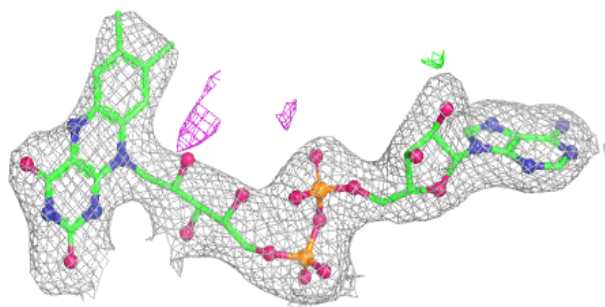
**Electron density around FAD B 601:**

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

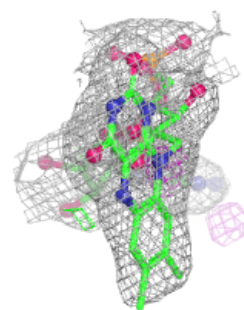
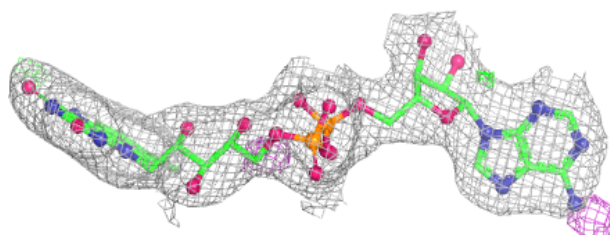
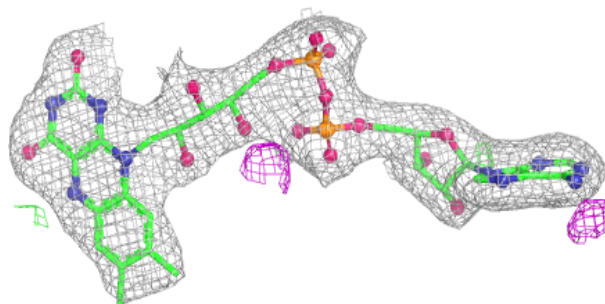


Electron density around FAD C 601:

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

**Electron density around FAD A 601:**

$2mF_o-DF_c$ (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.