



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:46 PM UTC

PDB ID : 3EY9 / pdb_00003ey9
Title : Structural basis for membrane binding and catalytic activation of the peripheral membrane enzyme pyruvate oxidase from Escherichia coli
Authors : Neumann, P.; Weidner, A.; Pech, A.; Stubbs, M.T.; Tittmann, K.
Deposited on : 2008-10-20
Resolution : 2.90 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

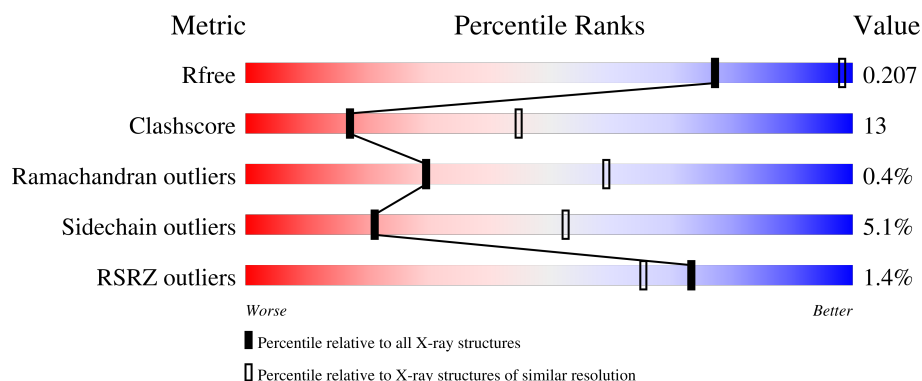
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	572	% 73% 25% .
1	B	572	% 74% 23% .

2 Entry composition [i](#)

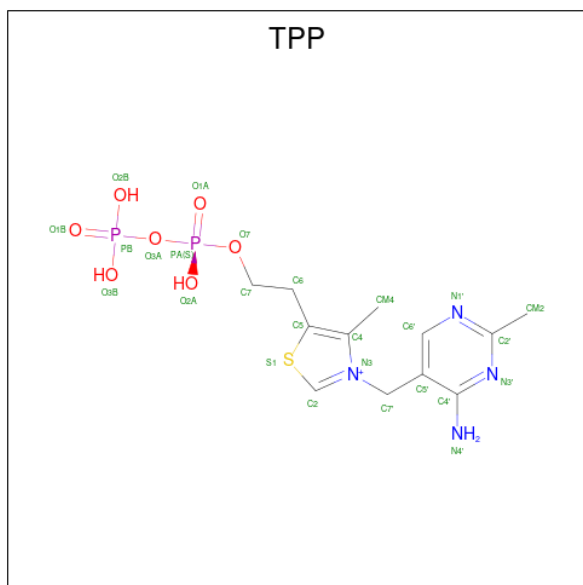
There are 5 unique types of molecules in this entry. The entry contains 8854 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pyruvate dehydrogenase [cytochrome].

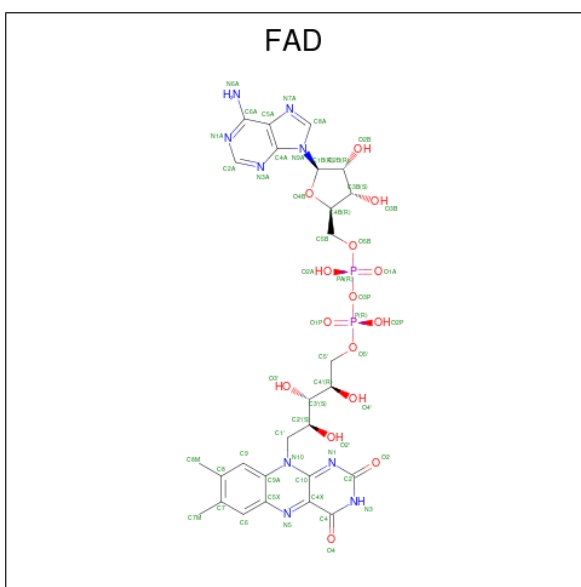
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	571	Total	C	N	O	S	0	0	0
			4342	2750	753	810	29			
1	B	571	Total	C	N	O	S	0	0	0
			4342	2750	753	810	29			

- Molecule 2 is THIAMINE DIPHOSPHATE (CCD ID: TPP) (formula: $C_{12}H_{19}N_4O_7P_2S$).



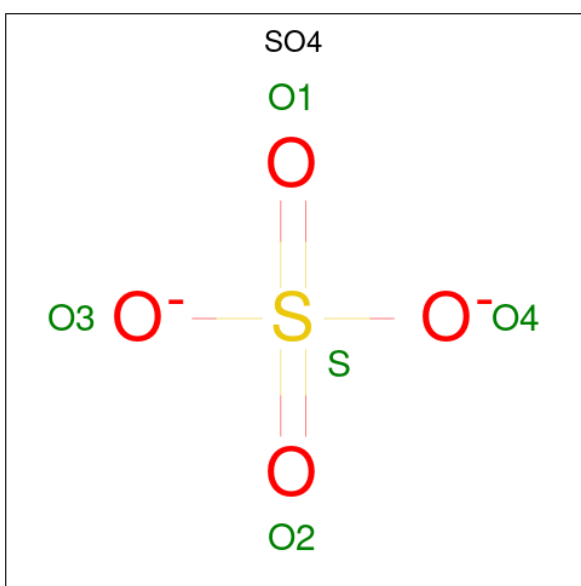
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0
			26	12	4	7	2	1	
2	B	1	Total	C	N	O	P	S	0
			26	12	4	7	2	1	

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



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

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



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

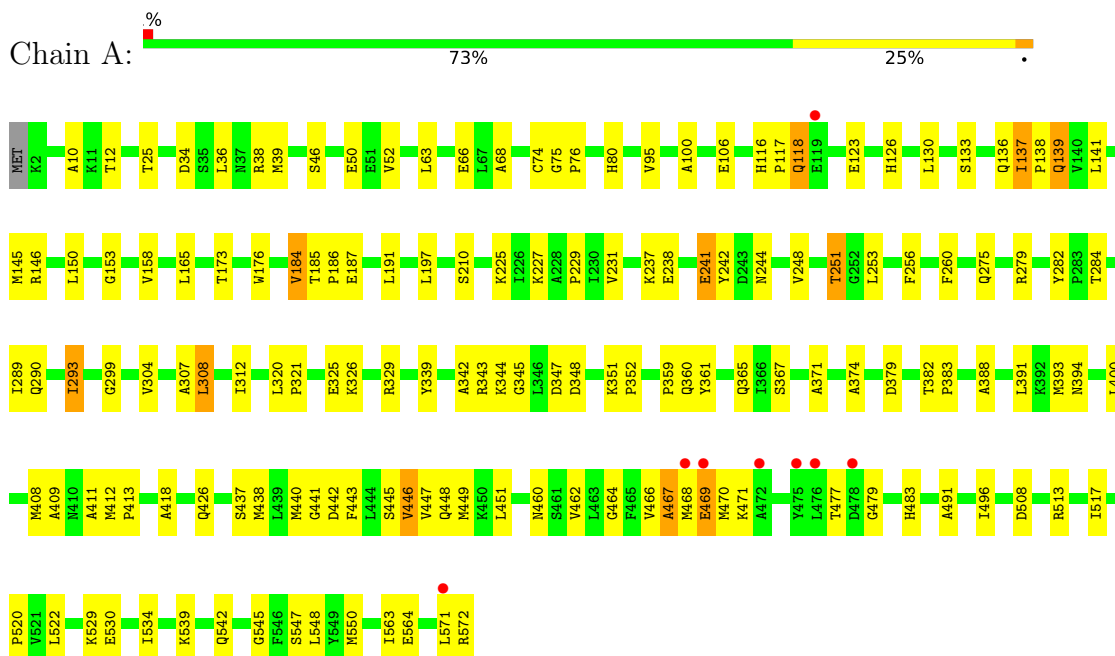
- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Mg 1	0	0
5	B	1	Total 1	Mg 1	0	0

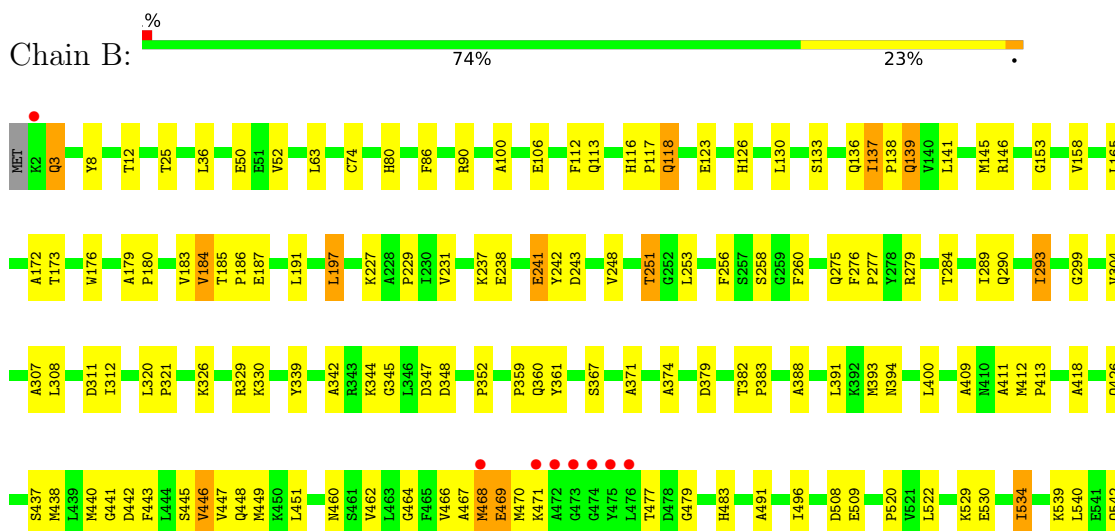
3 Residue-property plots

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: Pyruvate dehydrogenase [cytochrome]



• Molecule 1: Pyruvate dehydrogenase [cytochrome]



G545	F546	S547	L548	Y549	M550	I563	R572
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4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	151.37Å 151.37Å 153.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.97 – 2.90 29.97 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.97-2.90) 98.7 (29.97-2.90)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	0.01	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.75 (at 2.90Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.182 , 0.216 0.177 , 0.207	Depositor DCC
R_{free} test set	1983 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	73.5	Xtriage
Anisotropy	0.427	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h,l,k 0.000 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8854	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.46	0/4430	0.85	8/6001 (0.1%)
1	B	0.48	1/4430 (0.0%)	0.86	5/6001 (0.1%)
All	All	0.47	1/8860 (0.0%)	0.86	13/12002 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	137	ILE	CA-CB	5.67	1.57	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	137	ILE	N-CA-CB	6.78	114.77	110.50
1	B	137	ILE	N-CA-CB	6.33	114.49	110.50
1	B	534	ILE	N-CA-C	6.12	114.90	107.61
1	A	275	GLN	N-CA-C	-5.70	103.81	112.99
1	B	275	GLN	N-CA-C	-5.38	104.42	112.54
1	B	451	LEU	CA-C-N	5.38	126.56	119.84
1	B	451	LEU	C-N-CA	5.38	126.56	119.84
1	A	244	ASN	CA-C-N	5.29	124.96	119.56
1	A	244	ASN	C-N-CA	5.29	124.96	119.56
1	A	451	LEU	CA-C-N	5.09	126.21	119.84
1	A	451	LEU	C-N-CA	5.09	126.21	119.84
1	A	282	TYR	CA-C-N	5.06	125.06	119.90
1	A	282	TYR	C-N-CA	5.06	125.06	119.90

There are no chirality outliers.

There are no planarity outliers.

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	4342	0	4348	113	0
1	B	4342	0	4348	118	0
2	A	26	0	16	5	0
2	B	26	0	16	6	0
3	A	53	0	31	2	0
3	B	53	0	31	3	0
4	A	5	0	0	1	0
4	B	5	0	0	1	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	8854	0	8790	223	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 13.

All (223) 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:A:466:VAL:HG21	1:A:529:LYS:HB3	1.45	0.98
1:A:547:SER:H	1:A:550:MET:HE3	1.28	0.97
1:B:547:SER:H	1:B:550:MET:HE3	1.30	0.96
1:B:466:VAL:HG21	1:B:529:LYS:HB3	1.45	0.94
1:B:445:SER:HB3	1:B:449:MET:HE3	1.49	0.94
1:A:445:SER:HB3	1:A:449:MET:HE3	1.51	0.92
1:A:371:ALA:H	1:A:426:GLN:NE2	1.69	0.89
1:B:141:LEU:HG	1:B:145:MET:HE2	1.55	0.88
1:A:371:ALA:H	1:A:426:GLN:HE22	1.19	0.87
1:A:123:GLU:OE2	1:B:116:HIS:HD2	1.57	0.86
1:B:371:ALA:H	1:B:426:GLN:NE2	1.73	0.86
1:A:141:LEU:HG	1:A:145:MET:HE2	1.57	0.85
1:B:371:ALA:H	1:B:426:GLN:HE22	1.27	0.82
1:B:330:LYS:HZ3	1:B:572:ARG:HD3	1.45	0.81
1:A:116:HIS:HD2	1:B:123:GLU:OE2	1.63	0.80
1:A:539:LYS:HA	1:A:542:GLN:HE21	1.48	0.79
1:A:227:LYS:HG2	1:A:329:ARG:HB3	1.64	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:GLN:HG2	1:B:130:LEU:HB2	1.66	0.77
1:B:227:LYS:HG2	1:B:329:ARG:HB3	1.66	0.76
1:B:330:LYS:HZ3	1:B:572:ARG:HB3	1.50	0.76
1:A:118:GLN:HG2	1:A:130:LEU:HB2	1.66	0.75
1:A:483:HIS:H	1:B:448:GLN:NE2	1.85	0.74
1:B:539:LYS:HA	1:B:542:GLN:HE21	1.51	0.73
1:B:136:GLN:HA	1:B:139:GLN:HE21	1.53	0.73
1:A:136:GLN:HA	1:A:139:GLN:HE21	1.54	0.72
1:B:445:SER:HB3	1:B:449:MET:CE	2.20	0.71
1:B:547:SER:N	1:B:550:MET:HE3	2.06	0.71
1:A:448:GLN:NE2	1:B:483:HIS:H	1.89	0.71
1:A:547:SER:N	1:A:550:MET:HE3	2.05	0.70
1:A:445:SER:HB3	1:A:449:MET:CE	2.21	0.69
1:A:225:LYS:HG3	1:A:325:GLU:HG2	1.73	0.69
1:A:123:GLU:OE2	1:B:116:HIS:CD2	2.44	0.69
1:B:382:THR:HG23	2:B:611:TPP:O2B	1.94	0.68
1:B:330:LYS:NZ	1:B:572:ARG:HD3	2.09	0.67
1:A:443:PHE:HE2	1:A:491:ALA:HA	1.62	0.65
1:B:237:LYS:HE3	1:B:388:ALA:HA	1.78	0.65
1:A:237:LYS:HE3	1:A:388:ALA:HA	1.79	0.64
1:A:133:SER:O	1:A:136:GLN:HG2	1.97	0.64
1:A:304:VAL:HG11	1:A:307:ALA:HB2	1.79	0.64
1:A:371:ALA:N	1:A:426:GLN:HE22	1.95	0.64
1:A:184:VAL:HG11	1:A:293:ILE:HD11	1.80	0.63
1:B:443:PHE:HE2	1:B:491:ALA:HA	1.62	0.63
1:B:52:VAL:HG13	1:B:413:PRO:HB3	1.80	0.62
1:B:467:ALA:C	1:B:469:GLU:H	2.08	0.62
1:B:330:LYS:NZ	1:B:572:ARG:HB3	2.15	0.62
1:B:133:SER:O	1:B:136:GLN:HG2	1.99	0.61
1:B:184:VAL:HG11	1:B:293:ILE:HD11	1.81	0.61
1:A:137:ILE:HG22	1:A:138:PRO:HD3	1.82	0.61
1:A:52:VAL:HG13	1:A:413:PRO:HB3	1.83	0.61
1:A:312:ILE:HG12	3:A:612:FAD:C2A	2.31	0.60
1:A:116:HIS:CD2	1:B:123:GLU:OE2	2.51	0.60
1:A:383:PRO:HD3	2:A:611:TPP:O3B	2.01	0.60
1:A:352:PRO:HD3	1:A:361:TYR:CE1	2.37	0.59
1:B:63:LEU:HD12	1:B:418:ALA:HA	1.84	0.59
1:B:304:VAL:HG11	1:B:307:ALA:HB2	1.84	0.59
1:A:447:VAL:HG23	1:A:496:ILE:HD11	1.85	0.58
1:B:371:ALA:N	1:B:426:GLN:HE22	2.00	0.58
1:B:3:GLN:HE22	1:B:8:TYR:HA	1.68	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:ILE:HG22	1:B:138:PRO:HD3	1.84	0.58
1:B:290:GLN:HE22	1:B:299:GLY:H	1.52	0.58
1:B:352:PRO:HD3	1:B:361:TYR:CE1	2.39	0.58
1:A:469:GLU:HG3	1:A:539:LYS:CE	2.34	0.58
1:B:447:VAL:HG23	1:B:496:ILE:HD11	1.85	0.57
1:A:137:ILE:CG2	1:A:138:PRO:HD3	2.34	0.57
1:B:367:SER:HB2	1:B:391:LEU:HD23	1.86	0.57
1:A:367:SER:HB2	1:A:391:LEU:HD23	1.87	0.57
2:B:611:TPP:HM41	2:B:611:TPP:H6'	1.87	0.57
1:B:118:GLN:CG	1:B:130:LEU:HB2	2.34	0.56
1:A:483:HIS:H	1:B:448:GLN:HE21	1.52	0.56
1:B:534:ILE:O	1:B:539:LYS:HE3	2.05	0.56
1:B:137:ILE:CG2	1:B:138:PRO:HD3	2.35	0.56
1:B:469:GLU:HG3	1:B:539:LYS:CE	2.35	0.56
1:B:136:GLN:HA	1:B:139:GLN:NE2	2.21	0.55
1:A:290:GLN:HE22	1:A:299:GLY:H	1.54	0.55
1:A:63:LEU:HD12	1:A:418:ALA:HA	1.89	0.55
1:B:248:VAL:HG12	1:B:248:VAL:O	2.06	0.55
1:B:251:THR:HG21	1:B:260:PHE:HB2	1.89	0.55
1:A:279:ARG:NH1	1:B:106:GLU:OE2	2.38	0.55
1:A:409:ALA:HB2	1:A:438:MET:HB3	1.90	0.54
2:B:611:TPP:HM41	2:B:611:TPP:C6'	2.38	0.54
1:A:136:GLN:HA	1:A:139:GLN:NE2	2.21	0.54
1:A:460:ASN:CG	1:A:462:VAL:HG12	2.34	0.53
1:A:251:THR:HG21	1:A:260:PHE:HB2	1.89	0.53
1:A:468:MET:O	1:A:534:ILE:HD12	2.08	0.53
1:B:496:ILE:HG23	1:B:520:PRO:HB2	1.90	0.53
1:A:441:GLY:HA3	1:B:437:SER:O	2.09	0.53
1:B:307:ALA:C	1:B:308:LEU:HD12	2.34	0.53
1:A:443:PHE:CE2	1:A:491:ALA:HA	2.44	0.53
1:A:345:GLY:HA2	1:A:348:ASP:OD1	2.09	0.53
1:A:545:GLY:HA2	1:B:165:LEU:HD12	1.90	0.53
1:A:308:LEU:N	1:A:308:LEU:HD12	2.25	0.52
1:A:229:PRO:HB2	1:A:248:VAL:HG21	1.91	0.52
1:B:460:ASN:CG	1:B:462:VAL:HG12	2.34	0.52
1:A:106:GLU:OE2	1:B:279:ARG:NH1	2.42	0.52
1:B:382:THR:N	1:B:383:PRO:CD	2.72	0.52
1:A:534:ILE:O	1:A:539:LYS:HE3	2.09	0.52
1:B:522:LEU:C	1:B:522:LEU:HD23	2.35	0.52
1:B:312:ILE:HG12	3:B:612:FAD:C2A	2.40	0.52
1:A:382:THR:N	1:A:383:PRO:CD	2.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:SER:O	1:B:441:GLY:HA3	2.10	0.51
1:B:443:PHE:CE2	1:B:491:ALA:HA	2.45	0.51
1:A:12:THR:CG2	1:A:145:MET:HE1	2.40	0.51
1:A:496:ILE:HG23	1:A:520:PRO:HB2	1.92	0.51
1:B:229:PRO:HB2	1:B:248:VAL:HG21	1.93	0.51
1:B:383:PRO:HD3	2:B:611:TPP:O1B	2.11	0.51
1:A:467:ALA:HB1	1:A:470:MET:C	2.36	0.51
1:B:467:ALA:HB1	1:B:470:MET:C	2.36	0.51
1:A:248:VAL:HG12	1:A:248:VAL:O	2.10	0.50
1:A:446:VAL:HG22	1:A:496:ILE:HD11	1.92	0.50
1:A:522:LEU:C	1:A:522:LEU:HD23	2.35	0.50
1:B:409:ALA:HB2	1:B:438:MET:HB3	1.92	0.50
1:A:118:GLN:CG	1:A:130:LEU:HB2	2.38	0.49
1:A:371:ALA:HB3	1:A:374:ALA:HB2	1.94	0.49
2:A:611:TPP:H7'2	4:B:621:SO4:O3	2.12	0.49
1:B:469:GLU:HG3	1:B:539:LYS:NZ	2.26	0.49
1:B:345:GLY:HA2	1:B:348:ASP:OD1	2.12	0.49
1:A:186:PRO:HG2	1:A:191:LEU:HD21	1.94	0.49
1:B:412:MET:HB3	1:B:413:PRO:HD3	1.95	0.49
1:B:446:VAL:HG22	1:B:496:ILE:HD11	1.94	0.49
1:A:469:GLU:HG3	1:A:539:LYS:NZ	2.27	0.49
1:B:241:GLU:O	1:B:241:GLU:HG3	2.12	0.48
1:A:379:ASP:N	1:A:411:ALA:HB2	2.29	0.48
1:A:448:GLN:HE21	1:B:483:HIS:H	1.60	0.48
1:A:412:MET:HB3	1:A:413:PRO:HD3	1.96	0.48
1:B:12:THR:CG2	1:B:145:MET:HE1	2.44	0.48
1:A:279:ARG:HD3	1:B:106:GLU:OE1	2.14	0.48
1:A:227:LYS:HD3	1:A:326:LYS:O	2.13	0.48
1:B:197:LEU:HD22	1:B:289:ILE:HD11	1.96	0.48
3:A:612:FAD:H9	3:A:612:FAD:H1'1	1.59	0.47
1:A:293:ILE:HD13	1:A:293:ILE:O	2.15	0.47
1:A:238:GLU:HB2	1:A:393:MET:HE2	1.97	0.47
1:B:227:LYS:N	1:B:227:LYS:HD2	2.30	0.47
1:A:467:ALA:HB2	1:A:471:LYS:HA	1.97	0.46
1:A:197:LEU:HD22	1:A:289:ILE:HD11	1.97	0.46
1:B:227:LYS:HD3	1:B:326:LYS:O	2.16	0.46
1:B:8:TYR:CE1	1:B:172:ALA:HB1	2.51	0.46
1:B:50:GLU:HB2	1:B:80:HIS:HB3	1.98	0.46
1:A:241:GLU:HG3	1:A:241:GLU:O	2.15	0.45
1:B:437:SER:HA	1:B:440:MET:HB2	1.97	0.45
1:B:468:MET:O	1:B:534:ILE:HD12	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:ASP:O	1:A:38:ARG:HG3	2.17	0.45
1:A:50:GLU:HB2	1:A:80:HIS:HB3	1.99	0.45
1:A:165:LEU:HD12	1:B:545:GLY:HA2	1.99	0.45
1:A:347:ASP:O	1:A:348:ASP:C	2.59	0.45
1:B:12:THR:HG21	1:B:145:MET:HE1	1.98	0.45
1:B:74:CYS:HA	1:B:100:ALA:O	2.17	0.45
1:B:186:PRO:HG2	1:B:191:LEU:HD21	1.98	0.45
1:B:464:GLY:HA2	2:B:611:TPP:O2B	2.17	0.45
1:A:12:THR:HG21	1:A:145:MET:HE1	1.99	0.45
2:A:611:TPP:HN42	2:A:611:TPP:H2	1.81	0.45
1:B:238:GLU:HB2	1:B:393:MET:HE2	1.99	0.45
1:B:293:ILE:HD13	1:B:293:ILE:O	2.17	0.45
1:A:367:SER:HB2	1:A:391:LEU:HA	1.98	0.45
1:A:571:LEU:O	1:A:572:ARG:C	2.60	0.45
1:A:227:LYS:HD2	1:A:227:LYS:N	2.32	0.44
1:B:371:ALA:HB3	1:B:374:ALA:HB2	1.99	0.44
1:A:442:ASP:O	1:A:445:SER:HB2	2.18	0.44
1:A:361:TYR:O	1:A:365:GLN:HG2	2.18	0.44
1:B:379:ASP:N	1:B:411:ALA:HB2	2.32	0.44
1:A:75:GLY:HA3	1:A:76:PRO:HD3	1.83	0.44
1:A:359:PRO:HG2	1:A:530:GLU:HG3	2.00	0.44
1:B:320:LEU:HB2	1:B:321:PRO:HD3	1.99	0.44
1:A:477:THR:C	1:A:479:GLY:H	2.24	0.44
1:A:477:THR:C	1:A:479:GLY:N	2.76	0.44
1:B:308:LEU:HD12	1:B:308:LEU:N	2.32	0.44
1:B:477:THR:C	1:B:479:GLY:H	2.25	0.43
1:A:145:MET:HE3	1:A:176:TRP:CZ3	2.54	0.43
1:B:477:THR:C	1:B:479:GLY:N	2.75	0.43
1:A:117:PRO:HB2	1:A:158:VAL:CG2	2.49	0.43
1:B:467:ALA:HB2	1:B:471:LYS:HA	1.99	0.43
1:A:74:CYS:HA	1:A:100:ALA:O	2.18	0.43
1:A:307:ALA:C	1:A:308:LEU:HD12	2.43	0.43
1:B:367:SER:HB2	1:B:391:LEU:HA	2.01	0.43
4:A:621:SO4:O3	2:B:611:TPP:H7'2	2.18	0.43
1:A:320:LEU:HB2	1:A:321:PRO:HD3	2.01	0.43
1:A:339:TYR:O	1:A:343:ARG:HG3	2.19	0.43
1:A:467:ALA:C	1:A:469:GLU:H	2.26	0.43
1:B:347:ASP:O	1:B:348:ASP:C	2.62	0.43
1:A:68:ALA:O	1:A:95:VAL:HA	2.19	0.42
1:A:348:ASP:O	1:A:351:LYS:HE3	2.18	0.42
1:A:437:SER:HA	1:A:440:MET:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:611:TPP:H6'	2:A:611:TPP:HM41	2.02	0.42
1:A:464:GLY:HA2	2:A:611:TPP:O2B	2.20	0.42
1:B:145:MET:HE3	1:B:176:TRP:CZ3	2.54	0.42
1:B:179:ALA:HA	1:B:180:PRO:HD2	1.94	0.42
1:B:258:SER:HB3	1:B:339:TYR:HA	2.01	0.42
1:B:359:PRO:HG2	1:B:530:GLU:HG3	2.02	0.42
1:A:10:ALA:CB	1:A:39:MET:HE1	2.50	0.42
1:A:513:ARG:HD2	1:A:517:ILE:HD11	2.01	0.42
1:B:253:LEU:HD22	1:B:548:LEU:HB2	2.01	0.42
3:B:612:FAD:H1'1	3:B:612:FAD:H9	1.70	0.42
1:A:238:GLU:CB	1:A:393:MET:HE2	2.50	0.42
1:B:242:TYR:HB2	1:B:394:ASN:HA	2.01	0.42
1:A:126:HIS:CE1	1:A:153:GLY:HA3	2.55	0.41
1:A:253:LEU:HD22	1:A:548:LEU:HB2	2.01	0.41
1:A:469:GLU:OE1	1:A:539:LYS:HD3	2.20	0.41
1:A:513:ARG:O	1:A:517:ILE:HG13	2.20	0.41
1:B:290:GLN:NE2	1:B:299:GLY:H	2.16	0.41
1:B:137:ILE:N	1:B:138:PRO:CD	2.83	0.41
1:A:150:LEU:HD23	1:A:150:LEU:HA	1.83	0.41
1:A:342:ALA:C	1:A:344:LYS:H	2.27	0.41
1:B:352:PRO:HD3	1:B:361:TYR:CZ	2.56	0.41
1:A:251:THR:HG23	1:A:256:PHE:O	2.20	0.41
1:B:50:GLU:HB2	1:B:80:HIS:CB	2.51	0.41
1:B:540:LEU:HD23	1:B:540:LEU:HA	1.75	0.41
1:A:106:GLU:OE1	1:B:279:ARG:HD3	2.21	0.41
1:B:126:HIS:CE1	1:B:153:GLY:HA3	2.56	0.41
1:B:251:THR:HG23	1:B:256:PHE:O	2.19	0.41
1:B:112:PHE:O	1:B:113:GLN:HB2	2.21	0.41
1:A:137:ILE:N	1:A:138:PRO:CD	2.84	0.41
1:B:311:ASP:HA	3:B:612:FAD:N1A	2.36	0.41
1:B:242:TYR:O	1:B:243:ASP:C	2.64	0.41
1:B:86:PHE:O	1:B:90:ARG:HG2	2.20	0.40
1:B:117:PRO:HB2	1:B:158:VAL:HG21	2.03	0.40
1:B:469:GLU:OE1	1:B:539:LYS:HD3	2.21	0.40
1:B:342:ALA:C	1:B:344:LYS:H	2.28	0.40
1:B:442:ASP:O	1:B:445:SER:HB2	2.20	0.40
1:B:467:ALA:C	1:B:469:GLU:N	2.73	0.40
1:A:408:MET:HE2	1:A:408:MET:HB3	1.94	0.40
1:B:276:PHE:HA	1:B:277:PRO:HD3	1.91	0.40
1:A:242:TYR:HB2	1:A:394:ASN:HA	2.04	0.40
1:A:352:PRO:HD3	1:A:361:TYR:CZ	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:GLN:CD	1:B:139:GLN:H	2.28	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	569/572 (100%)	537 (94%)	30 (5%)	2 (0%)	30	59
1	B	569/572 (100%)	537 (94%)	30 (5%)	2 (0%)	30	59
All	All	1138/1144 (100%)	1074 (94%)	60 (5%)	4 (0%)	30	59

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	468	MET
1	A	469	GLU
1	B	469	GLU
1	A	467	ALA

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	457/458 (100%)	433 (95%)	24 (5%)	20	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	457/458 (100%)	434 (95%)	23 (5%)	22	53
All	All	914/916 (100%)	867 (95%)	47 (5%)	21	53

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	THR
1	A	36	LEU
1	A	46	SER
1	A	66	GLU
1	A	118	GLN
1	A	139	GLN
1	A	146	ARG
1	A	173	THR
1	A	184	VAL
1	A	185	THR
1	A	187	GLU
1	A	210	SER
1	A	231	VAL
1	A	241	GLU
1	A	251	THR
1	A	284	THR
1	A	293	ILE
1	A	308	LEU
1	A	360	GLN
1	A	400	LEU
1	A	446	VAL
1	A	508	ASP
1	A	563	ILE
1	A	564	GLU
1	B	3	GLN
1	B	25	THR
1	B	36	LEU
1	B	118	GLN
1	B	139	GLN
1	B	146	ARG
1	B	173	THR
1	B	183	VAL
1	B	184	VAL
1	B	185	THR
1	B	187	GLU
1	B	197	LEU

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Mol	Chain	Res	Type
1	B	231	VAL
1	B	241	GLU
1	B	251	THR
1	B	284	THR
1	B	293	ILE
1	B	360	GLN
1	B	400	LEU
1	B	446	VAL
1	B	508	ASP
1	B	509	GLU
1	B	563	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	92	HIS
1	A	116	HIS
1	A	126	HIS
1	A	139	GLN
1	A	151	ASN
1	A	181	GLN
1	A	261	HIS
1	A	275	GLN
1	A	290	GLN
1	A	301	HIS
1	A	364	GLN
1	A	365	GLN
1	A	426	GLN
1	A	448	GLN
1	A	542	GLN
1	B	3	GLN
1	B	92	HIS
1	B	116	HIS
1	B	126	HIS
1	B	139	GLN
1	B	151	ASN
1	B	181	GLN
1	B	261	HIS
1	B	275	GLN
1	B	290	GLN
1	B	301	HIS

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Mol	Chain	Res	Type
1	B	364	GLN
1	B	365	GLN
1	B	426	GLN
1	B	448	GLN
1	B	542	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](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 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	TPP	B	611	5	26,27,27	1.62	4 (15%)	38,40,40	1.87	9 (23%)
4	SO4	B	621	-	4,4,4	0.25	0	6,6,6	0.35	0
4	SO4	A	621	-	4,4,4	0.21	0	6,6,6	0.33	0
3	FAD	B	612	-	58,58,58	1.15	5 (8%)	85,89,89	1.63	19 (22%)
2	TPP	A	611	5	26,27,27	1.73	5 (19%)	38,40,40	1.90	10 (26%)
3	FAD	A	612	-	58,58,58	1.16	6 (10%)	85,89,89	1.59	16 (18%)

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	TPP	B	611	5	-	2/17/17/17	0/2/2/2
3	FAD	B	612	-	-	10/34/50/50	0/6/6/6
2	TPP	A	611	5	-	2/17/17/17	0/2/2/2
3	FAD	A	612	-	-	7/34/50/50	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	611	TPP	C5-C4	6.41	1.48	1.35
2	B	611	TPP	C5-C4	6.09	1.48	1.35
3	A	612	FAD	C9A-N10	-4.16	1.33	1.41
3	B	612	FAD	C9A-N10	-3.24	1.35	1.41
2	A	611	TPP	PA-O3A	3.08	1.62	1.59
3	B	612	FAD	P-O3P	2.98	1.62	1.59
2	A	611	TPP	C4-N3	-2.58	1.34	1.39
2	B	611	TPP	PA-O3A	2.50	1.62	1.59
3	A	612	FAD	C4X-N5	2.46	1.36	1.30
2	B	611	TPP	C4-N3	-2.36	1.35	1.39
3	A	612	FAD	C5A-N7A	-2.31	1.34	1.39
3	A	612	FAD	P-O3P	2.25	1.61	1.59
3	B	612	FAD	C8A-N7A	2.25	1.36	1.31
3	B	612	FAD	C4X-N5	2.19	1.35	1.30
3	A	612	FAD	C8A-N7A	2.15	1.35	1.31
2	A	611	TPP	C5'-C4'	-2.09	1.39	1.42
3	B	612	FAD	C5X-N5	-2.08	1.35	1.39
3	A	612	FAD	C8A-N9A	-2.04	1.34	1.37
2	B	611	TPP	C2-S1	2.03	1.75	1.69
2	A	611	TPP	C2-S1	2.01	1.75	1.69

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	611	TPP	C2-S1-C5	-6.33	87.03	91.22
2	B	611	TPP	C2-S1-C5	-5.79	87.38	91.22
3	B	612	FAD	C5A-C4A-N3A	-5.21	119.54	126.72
3	B	612	FAD	N3A-C2A-N1A	-5.00	121.01	128.58
3	A	612	FAD	C5A-C4A-N3A	-4.85	120.03	126.72
3	A	612	FAD	N3A-C2A-N1A	-4.69	121.48	128.58
3	A	612	FAD	N9A-C8A-N7A	-3.84	108.49	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	612	FAD	N9A-C8A-N7A	-3.79	108.56	113.94
3	B	612	FAD	C2A-N3A-C4A	3.72	120.93	111.83
2	A	611	TPP	S1-C2-N3	3.36	116.55	112.30
3	A	612	FAD	C2A-N3A-C4A	3.29	119.88	111.83
2	A	611	TPP	C6'-C5'-C4'	3.25	119.54	115.55
3	A	612	FAD	N3A-C4A-N9A	3.21	132.62	127.17
2	B	611	TPP	C5'-C7'-N3	-3.16	106.15	112.97
3	B	612	FAD	C4A-C5A-N7A	-3.12	107.01	110.58
3	A	612	FAD	C4A-N9A-C8A	3.08	108.97	105.74
3	B	612	FAD	N3A-C4A-N9A	3.01	132.29	127.17
2	A	611	TPP	C6'-N1'-C2'	2.97	120.94	116.07
3	B	612	FAD	C9A-C5X-N5	-2.94	119.34	122.45
2	B	611	TPP	C6'-N1'-C2'	2.90	120.84	116.07
2	B	611	TPP	C6'-C5'-C4'	2.87	119.08	115.55
2	A	611	TPP	C5'-C7'-N3	-2.86	106.80	112.97
2	B	611	TPP	S1-C2-N3	2.85	115.91	112.30
3	A	612	FAD	C9A-C5X-N5	-2.82	119.46	122.45
2	A	611	TPP	C5'-C6'-N1'	-2.80	119.28	123.83
3	B	612	FAD	C5A-N7A-C8A	2.78	107.83	103.45
3	A	612	FAD	C5A-N7A-C8A	2.70	107.70	103.45
3	A	612	FAD	C4A-C5A-N7A	-2.70	107.50	110.58
2	B	611	TPP	C6-C5-S1	-2.69	114.60	120.90
2	B	611	TPP	C5'-C6'-N1'	-2.62	119.56	123.83
2	B	611	TPP	CM4-C4-N3	2.62	126.61	120.57
3	A	612	FAD	C5'-C4'-C3'	-2.61	107.29	112.22
2	B	611	TPP	N1'-C2'-N3'	-2.59	121.22	125.53
3	B	612	FAD	C4A-N9A-C8A	2.57	108.43	105.74
3	A	612	FAD	C10-C4X-N5	-2.54	119.62	124.81
2	A	611	TPP	N1'-C2'-N3'	-2.53	121.33	125.53
3	B	612	FAD	C4-N3-C2	-2.49	121.21	125.64
3	A	612	FAD	C4-C4X-N5	2.48	121.63	118.21
3	B	612	FAD	C10-C4X-N5	-2.46	119.78	124.81
3	B	612	FAD	C4X-C10-N1	-2.39	118.72	124.59
3	A	612	FAD	C4X-C10-N10	2.32	119.81	116.48
3	B	612	FAD	C4X-C4-N3	2.32	119.16	113.25
3	B	612	FAD	O4B-C1B-N9A	2.31	112.53	108.09
2	A	611	TPP	N4'-C4'-N3'	2.30	120.13	117.03
3	B	612	FAD	C4X-C10-N10	2.29	119.76	116.48
3	A	612	FAD	C4X-C4-N3	2.28	119.06	113.25
3	B	612	FAD	C4'-C3'-C2'	-2.21	109.89	113.57
2	A	611	TPP	C6-C5-S1	-2.21	115.73	120.90
3	A	612	FAD	C4-N3-C2	-2.20	121.74	125.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	612	FAD	O4-C4-C4X	-2.20	120.73	126.53
2	A	611	TPP	CM2-C2'-N3'	2.19	120.41	117.13
3	B	612	FAD	C5A-C4A-N9A	2.16	108.16	105.81
3	A	612	FAD	O4-C4-C4X	-2.06	121.10	126.53
3	B	612	FAD	C10-N1-C2	2.04	121.27	116.85

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	611	TPP	PA-O3A-PB-O3B
3	A	612	FAD	O4B-C4B-C5B-O5B
3	A	612	FAD	O4'-C4'-C5'-O5'
3	B	612	FAD	C3'-C4'-C5'-O5'
3	B	612	FAD	O4'-C4'-C5'-O5'
3	A	612	FAD	C3B-C4B-C5B-O5B
3	B	612	FAD	O4B-C4B-C5B-O5B
3	B	612	FAD	C3B-C4B-C5B-O5B
3	B	612	FAD	O3'-C3'-C4'-C5'
3	B	612	FAD	C2'-C3'-C4'-C5'
3	B	612	FAD	C2'-C3'-C4'-O4'
3	A	612	FAD	C3'-C4'-C5'-O5'
3	B	612	FAD	O3'-C3'-C4'-O4'
3	A	612	FAD	P-O3P-PA-O5B
3	B	612	FAD	P-O3P-PA-O5B
2	A	611	TPP	PA-O3A-PB-O3B
2	B	611	TPP	PA-O3A-PB-O2B
2	A	611	TPP	C5-C6-C7-O7
3	A	612	FAD	N10-C1'-C2'-O2'
3	A	612	FAD	P-O3P-PA-O2A
3	B	612	FAD	P-O3P-PA-O2A

There are no ring outliers.

6 monomers are involved in 16 short contacts:

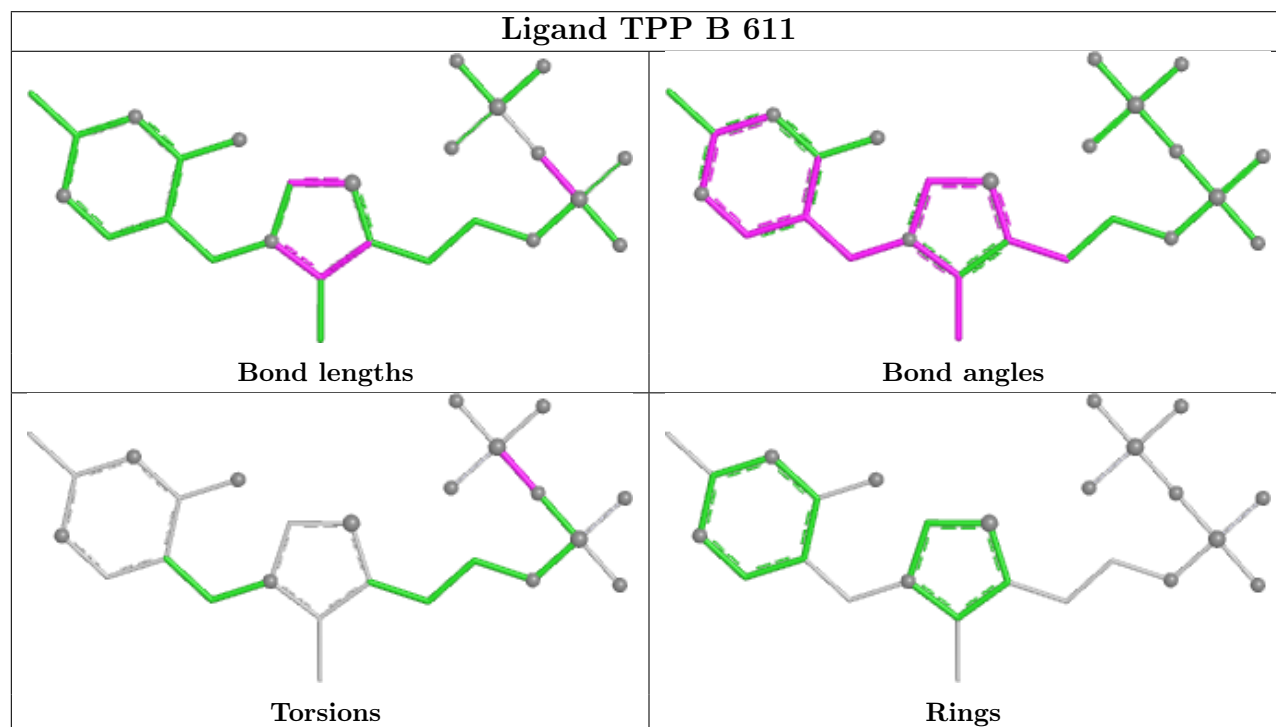
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	611	TPP	6	0
4	B	621	SO4	1	0
4	A	621	SO4	1	0
3	B	612	FAD	3	0
2	A	611	TPP	5	0

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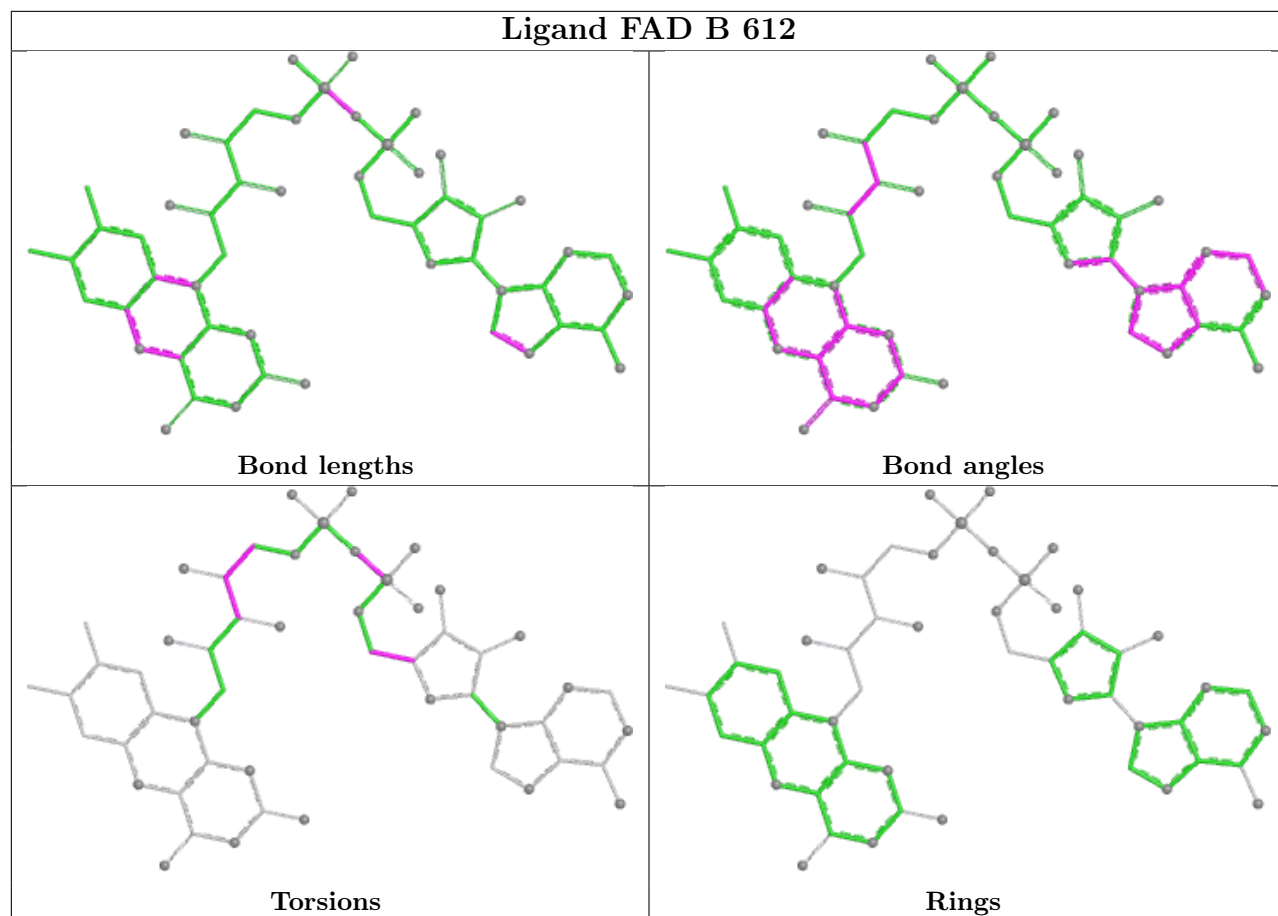
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	612	FAD	2	0

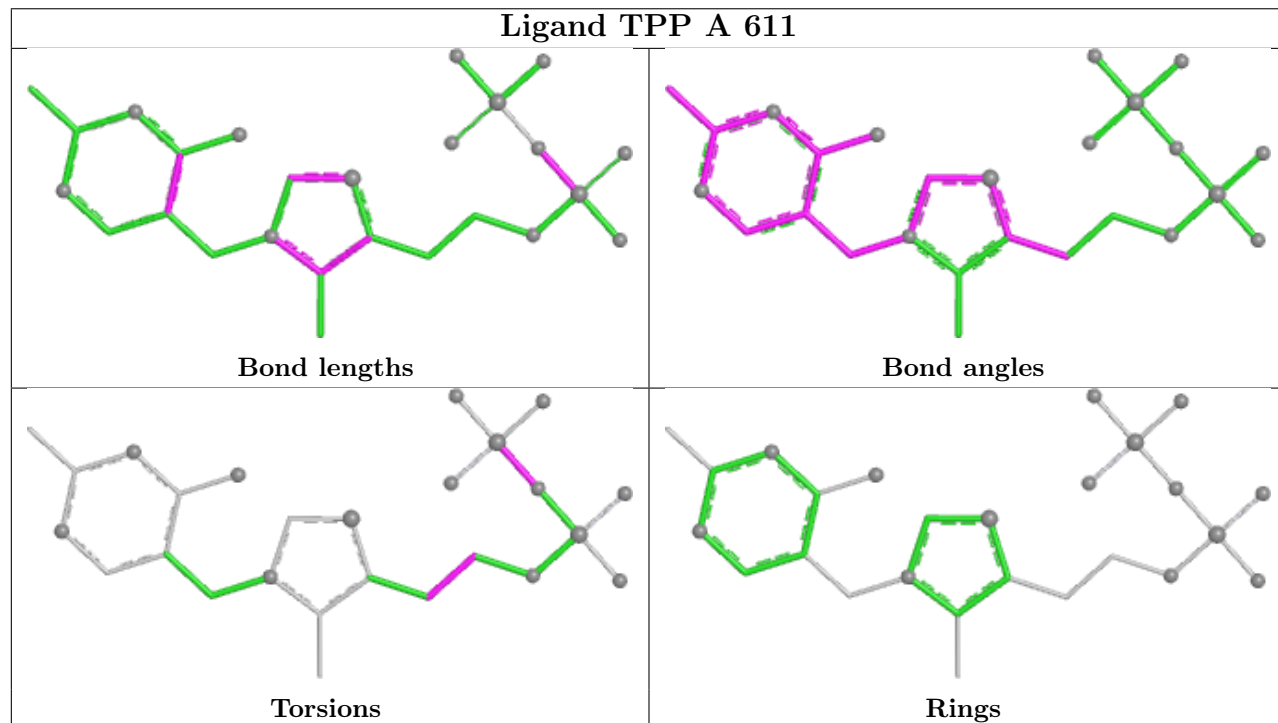
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

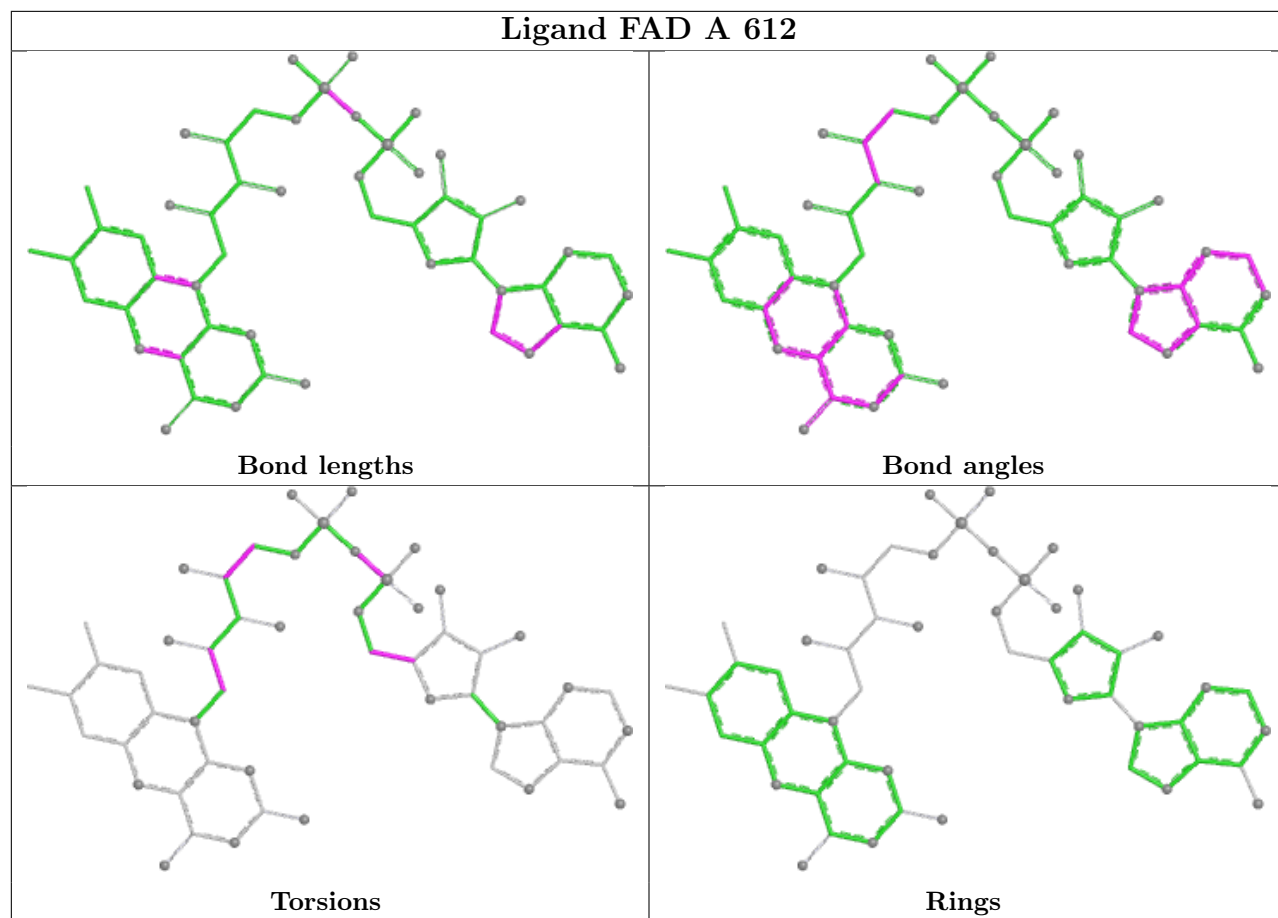


Ligand FAD B 612



Ligand TPP A 611





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	571/572 (99%)	-0.44	8 (1%) 73 65	50, 77, 113, 227	0
1	B	571/572 (99%)	-0.39	8 (1%) 73 65	49, 75, 114, 221	0
All	All	1142/1144 (99%)	-0.41	16 (1%) 73 65	49, 76, 114, 227	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	468	MET	5.5
1	A	468	MET	5.4
1	B	473	GLY	4.3
1	A	476	LEU	4.0
1	B	475	TYR	3.1
1	A	571	LEU	2.9
1	B	476	LEU	2.8
1	B	474	GLY	2.7
1	A	119	GLU	2.7
1	A	475	TYR	2.4
1	B	2	LYS	2.4
1	A	472	ALA	2.2
1	A	478	ASP	2.2
1	A	469	GLU	2.1
1	B	472	ALA	2.1
1	B	471	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](#)

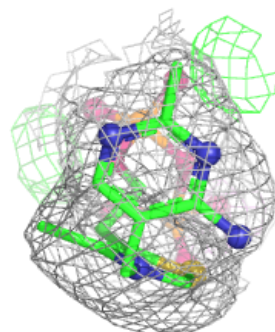
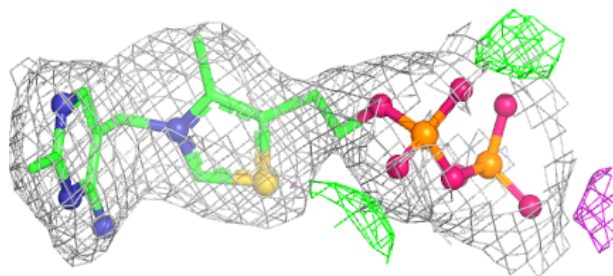
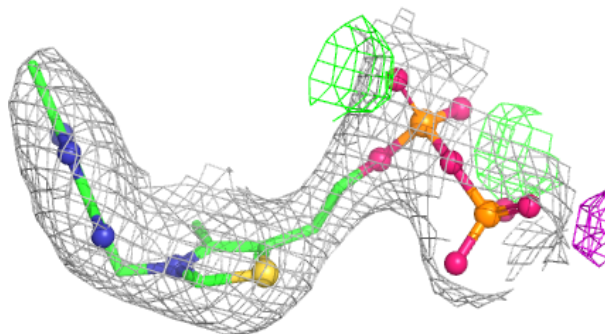
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	A	621	5/5	0.90	0.08	74,82,95,108	0
2	TPP	A	611	26/26	0.96	0.07	46,60,79,82	0
3	FAD	A	612	53/53	0.97	0.06	46,65,79,93	0
2	TPP	B	611	26/26	0.97	0.07	34,64,74,122	0
4	SO4	B	621	5/5	0.97	0.05	68,71,82,87	0
5	MG	A	616	1/1	0.97	0.05	57,57,57,57	0
5	MG	B	617	1/1	0.97	0.12	59,59,59,59	0
3	FAD	B	612	53/53	0.98	0.05	49,66,84,91	0

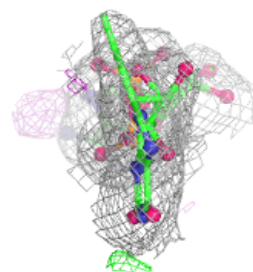
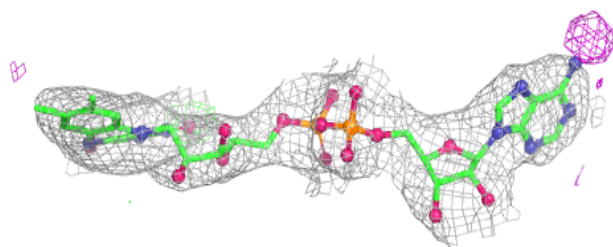
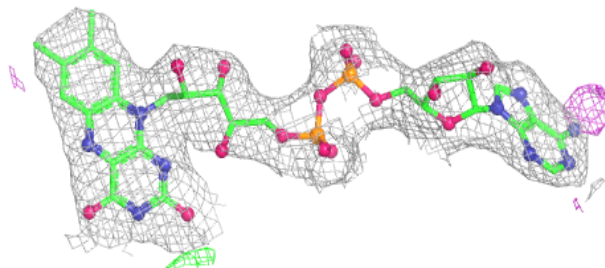
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TPP A 611:

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

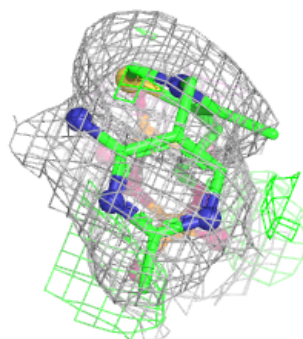
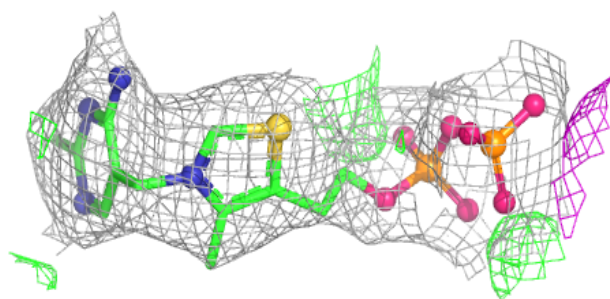
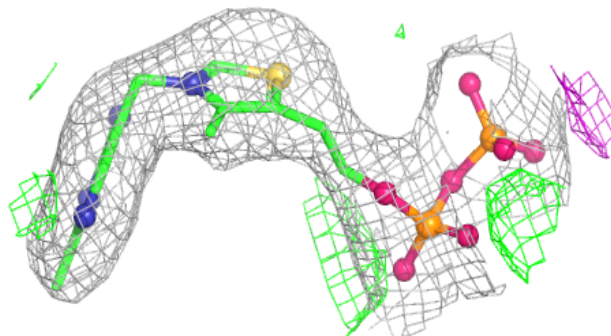
**Electron density around FAD A 612:**

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

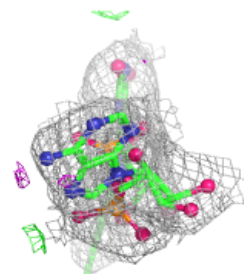
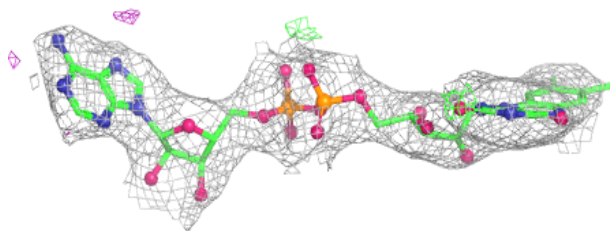
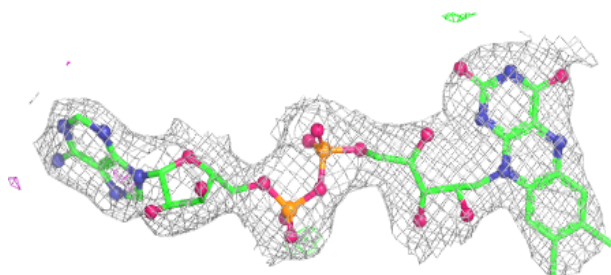


Electron density around TPP B 611:

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

**Electron density around FAD B 612:**

$2mF_o-DF_c$ (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.