



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:47 PM UTC

PDB ID : 3LSI / pdb_00003lsi
Title : Pyranose 2-oxidase T169A, tetragonal
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Deposited on : 2010-02-12
Resolution : 1.90 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

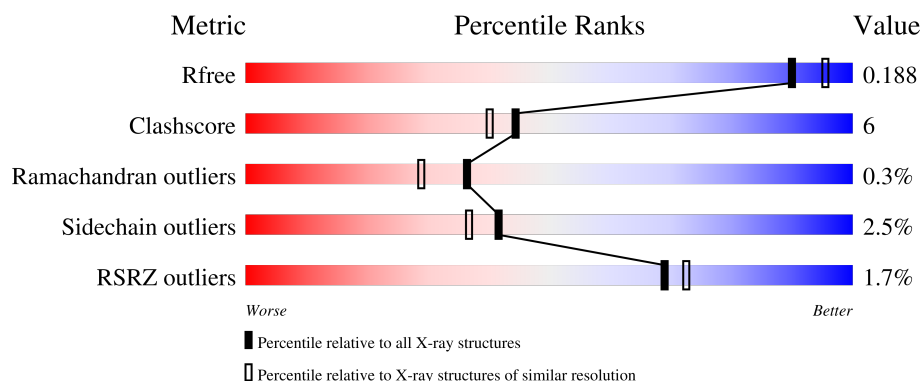
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

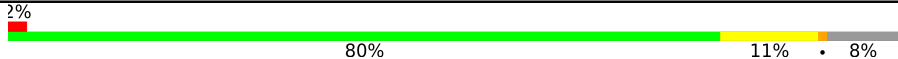
The reported resolution of this entry is 1.90 Å.

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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	623	
1	B	623	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10182 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 Pyranose 2-oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	575	Total	C	N	O	S	0	0	0
			4532	2862	776	870	24			
1	B	576	Total	C	N	O	S	0	0	0
			4540	2867	777	871	25			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	ALA	THR	engineered mutation	UNP Q7ZA32
B	169	ALA	THR	engineered mutation	UNP Q7ZA32

- 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		

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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		

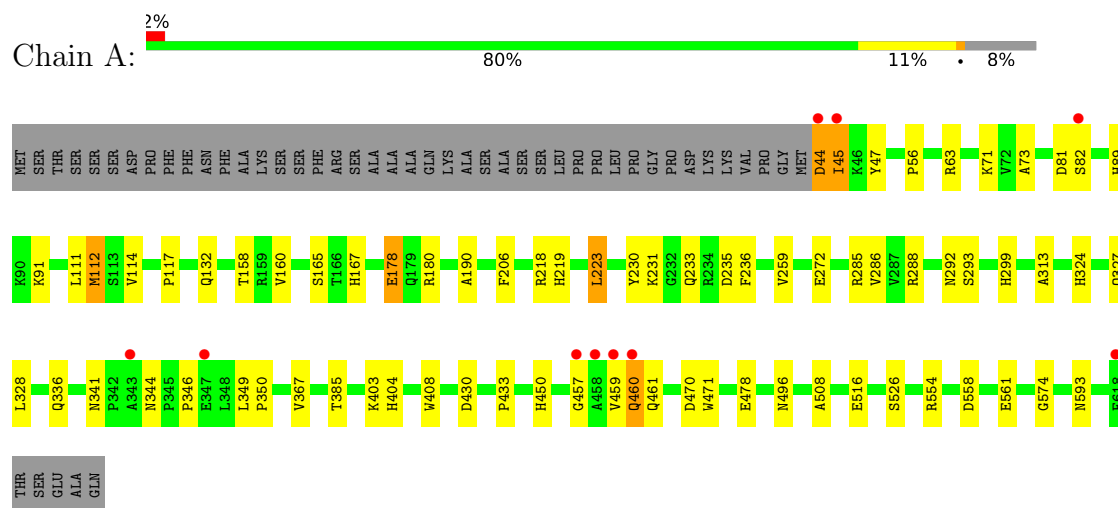
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	507	Total	O	0	0
			507	507		
3	B	497	Total	O	0	0
			497	497		

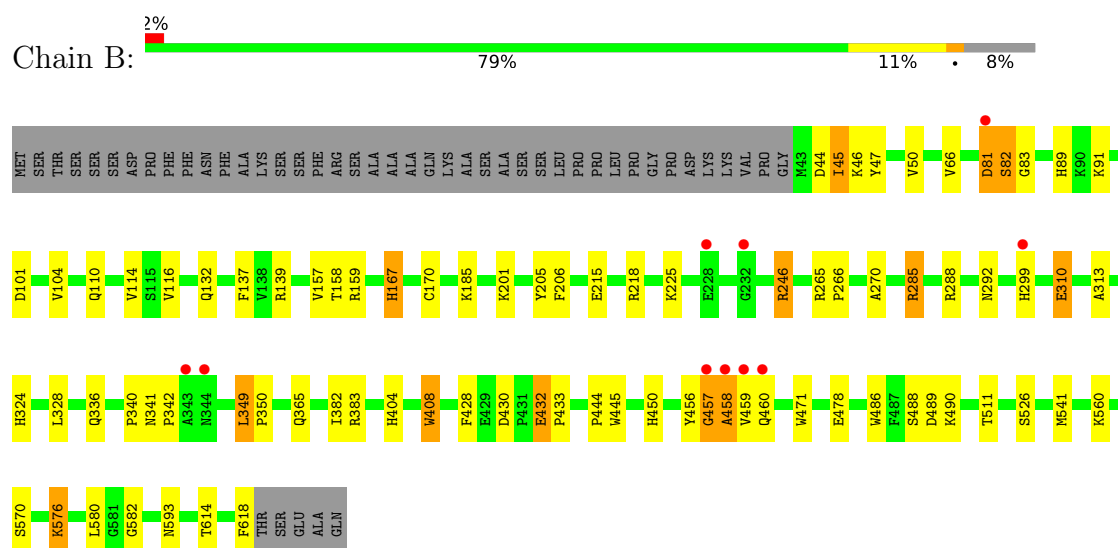
3 Residue-property plots [i](#)

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

• Molecule 1: Pyranose 2-oxidase



• Molecule 1: Pyranose 2-oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	101.30Å 101.30Å 249.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.10 – 1.90 29.10 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.10-1.90) 99.5 (29.10-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.164 , 0.213 (Not available) , 0.188	Depositor DCC
R_{free} test set	3090 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10182	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.81 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.3323e-03.*

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

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	1.25	5/4648 (0.1%)	1.12	10/6320 (0.2%)
1	B	1.23	5/4656 (0.1%)	1.16	17/6330 (0.3%)
All	All	1.24	10/9304 (0.1%)	1.14	27/12650 (0.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	582	GLY	C-O	6.07	1.29	1.23
1	B	488	SER	C-O	-5.93	1.16	1.23
1	B	81	ASP	CA-C	5.71	1.59	1.52
1	B	114	VAL	CA-C	5.69	1.59	1.52
1	A	367	VAL	CA-CB	-5.51	1.47	1.54
1	A	117	PRO	N-CA	-5.29	1.41	1.47
1	A	508	ALA	C-O	5.29	1.31	1.23
1	A	82	SER	N-CA	5.28	1.52	1.46
1	B	349	LEU	CA-C	-5.13	1.49	1.53
1	A	190	ALA	CA-CB	5.12	1.61	1.53

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	116	VAL	CA-C-N	-8.59	111.96	120.21
1	B	116	VAL	C-N-CA	-8.59	111.96	120.21
1	A	82	SER	CB-CA-C	-6.88	98.12	110.37
1	B	432	GLU	CA-C-N	6.82	126.58	119.76
1	B	432	GLU	C-N-CA	6.82	126.58	119.76
1	A	82	SER	N-CA-C	6.56	120.89	113.02
1	B	457	GLY	N-CA-C	6.56	121.89	114.67
1	B	428	PHE	CA-C-N	-6.37	113.02	122.83
1	B	428	PHE	C-N-CA	-6.37	113.02	122.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	614	THR	CA-C-N	-6.22	113.37	119.78
1	B	614	THR	C-N-CA	-6.22	113.37	119.78
1	A	114	VAL	N-CA-C	6.07	116.23	110.53
1	A	457	GLY	N-CA-C	5.78	123.23	114.61
1	B	44	ASP	N-CA-C	-5.74	102.91	110.43
1	A	112	MET	CG-SD-CE	-5.71	88.34	100.90
1	B	170	CYS	N-CA-C	5.57	119.66	112.86
1	A	180	ARG	CA-C-O	5.56	124.22	119.66
1	A	554	ARG	N-CA-C	5.47	117.75	110.53
1	B	66	VAL	N-CA-C	-5.47	105.19	110.72
1	B	167	HIS	N-CA-C	5.41	121.73	113.19
1	A	496	ASN	N-CA-C	5.40	118.89	111.54
1	B	383	ARG	NE-CZ-NH2	5.35	124.01	119.20
1	A	341	ASN	CA-C-N	5.24	125.54	119.47
1	A	341	ASN	C-N-CA	5.24	125.54	119.47
1	B	82	SER	N-CA-C	5.18	121.83	110.80
1	B	270	ALA	CA-C-N	-5.08	114.51	119.64
1	B	270	ALA	C-N-CA	-5.08	114.51	119.64

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	4532	0	4377	55	0
1	B	4540	0	4386	54	0
2	A	53	0	30	1	0
2	B	53	0	30	3	0
3	A	507	0	0	13	0
3	B	497	0	0	14	0
All	All	10182	0	8823	108	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 6.

All (108) 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:460:GLN:HB3	3:A:1821:HOH:O	1.57	1.03
1:A:459:VAL:HA	3:A:1442:HOH:O	1.61	1.00
1:B:101:ASP:O	1:B:104:VAL:HG12	1.65	0.96
1:A:459:VAL:HG12	1:A:460:GLN:O	1.66	0.94
1:A:132:GLN:HG2	3:A:1850:HOH:O	1.65	0.94
1:B:365:GLN:HE22	1:B:459:VAL:HG22	1.38	0.85
1:B:132:GLN:HG2	3:B:1822:HOH:O	1.80	0.82
1:A:478:GLU:HB3	3:A:1931:HOH:O	1.78	0.82
1:B:132:GLN:CG	3:B:1822:HOH:O	2.29	0.80
1:B:457:GLY:HA3	3:B:1373:HOH:O	1.82	0.80
1:A:45:ILE:O	1:A:45:ILE:CG2	2.34	0.75
1:B:382:ILE:HD13	3:B:1450:HOH:O	1.86	0.74
1:A:132:GLN:CG	3:A:1850:HOH:O	2.29	0.70
1:A:45:ILE:O	1:A:45:ILE:HG22	1.91	0.69
1:A:167:HIS:CD2	2:A:801:FAD:HM82	2.27	0.69
1:B:225:LYS:HE2	3:B:1650:HOH:O	1.94	0.68
1:A:285:ARG:NH1	1:A:299:HIS:HD2	1.92	0.68
1:B:576:LYS:HE2	3:B:1942:HOH:O	1.93	0.68
1:A:459:VAL:HG12	1:A:460:GLN:N	2.10	0.66
1:B:167:HIS:CD2	2:B:801:FAD:HM82	2.31	0.66
1:A:167:HIS:ND1	3:A:1300:HOH:O	2.28	0.66
1:A:44:ASP:HB2	1:A:71:LYS:HZ3	1.64	0.63
1:A:404:HIS:HE1	3:A:1385:HOH:O	1.81	0.63
1:B:218:ARG:HD2	3:B:1026:HOH:O	1.99	0.61
1:B:404:HIS:HE1	3:B:1539:HOH:O	1.83	0.61
1:B:246:ARG:O	1:B:246:ARG:HG3	2.01	0.60
1:B:299:HIS:HB2	1:B:310:GLU:OE2	2.02	0.60
1:A:63:ARG:HD3	1:A:259:VAL:O	2.02	0.59
1:A:112:MET:N	1:A:112:MET:HE2	2.18	0.59
1:A:285:ARG:NH1	1:A:299:HIS:CD2	2.70	0.58
1:A:111:LEU:C	1:A:112:MET:HE2	2.29	0.58
1:A:230:TYR:O	1:A:233:GLN:HG3	2.04	0.57
1:A:403:LYS:HE2	3:A:1973:HOH:O	2.04	0.56
1:B:341:ASN:HD22	1:B:342:PRO:HD2	1.72	0.55
1:A:81:ASP:O	1:B:81:ASP:O	2.25	0.54
1:A:450:HIS:O	1:A:470:ASP:HB2	2.07	0.54
1:B:201:LYS:HE2	1:B:205:TYR:OH	2.07	0.53
1:A:285:ARG:HH12	1:A:299:HIS:CD2	2.26	0.53
1:A:89:HIS:CE1	1:A:91:LYS:HB2	2.44	0.53
1:A:44:ASP:HB2	1:A:71:LYS:NZ	2.23	0.52
1:B:285:ARG:NH2	3:B:1974:HOH:O	2.31	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:VAL:HG21	1:B:324:HIS:HE1	1.75	0.52
1:A:459:VAL:CG1	1:A:460:GLN:N	2.72	0.52
1:B:132:GLN:NE2	3:B:1822:HOH:O	2.38	0.52
1:A:558:ASP:OD2	1:A:561:GLU:HG3	2.10	0.51
1:A:158:THR:HG22	1:A:160:VAL:HG22	1.93	0.51
1:B:471:TRP:CH2	1:B:526:SER:HA	2.45	0.51
1:A:91:LYS:NZ	3:A:1361:HOH:O	2.43	0.51
1:B:81:ASP:OD1	1:B:81:ASP:N	2.42	0.50
1:B:576:LYS:CE	3:B:1942:HOH:O	2.56	0.50
1:B:456:TYR:CD2	1:B:541:MET:HE2	2.47	0.49
1:A:47:TYR:O	1:A:313:ALA:HA	2.11	0.49
1:B:47:TYR:O	1:B:313:ALA:HA	2.12	0.49
1:B:167:HIS:HD2	2:B:801:FAD:HM82	1.77	0.49
1:A:230:TYR:O	1:A:231:LYS:C	2.57	0.48
1:B:285:ARG:NE	3:B:1974:HOH:O	2.25	0.48
1:B:576:LYS:N	1:B:576:LYS:HD2	2.28	0.48
1:A:218:ARG:HD2	3:A:1032:HOH:O	2.14	0.47
1:B:432:GLU:HB2	1:B:433:PRO:HD2	1.97	0.47
1:B:408:TRP:C	1:B:408:TRP:CD1	2.93	0.47
1:A:89:HIS:ND1	1:A:91:LYS:HB2	2.30	0.47
1:A:461:GLN:HG3	3:B:1832:HOH:O	2.14	0.46
1:B:89:HIS:CE1	1:B:91:LYS:HB2	2.50	0.46
1:B:328:LEU:O	1:B:328:LEU:HD23	2.15	0.46
1:B:444:PRO:HD2	1:B:445:TRP:CZ3	2.51	0.46
1:A:56:PRO:HD3	1:A:165:SER:HB3	1.97	0.46
1:A:223:LEU:HD13	1:A:236:PHE:HB3	1.97	0.46
1:A:223:LEU:CD1	1:A:236:PHE:HB3	2.46	0.46
1:B:50:VAL:HG13	1:B:313:ALA:HB2	1.97	0.45
1:B:137:PHE:CE2	1:B:139:ARG:HG3	2.51	0.45
1:A:403:LYS:CE	3:A:1973:HOH:O	2.61	0.45
1:B:618:PHE:CD1	1:B:618:PHE:C	2.93	0.45
1:B:299:HIS:CB	1:B:310:GLU:OE2	2.64	0.45
1:A:288:ARG:HD2	1:A:292:ASN:OD1	2.17	0.45
1:A:336:GLN:NE2	1:A:344:ASN:O	2.50	0.45
1:B:218:ARG:HG3	1:B:430:ASP:OD2	2.17	0.45
1:B:489:ASP:O	1:B:490:LYS:HD3	2.17	0.45
1:B:288:ARG:HD2	1:B:292:ASN:OD1	2.16	0.44
1:A:346:PRO:HG2	1:A:350:PRO:HA	1.99	0.44
1:B:46:LYS:HD3	1:B:47:TYR:N	2.33	0.44
1:A:218:ARG:HG3	1:A:430:ASP:OD2	2.18	0.44
1:A:459:VAL:O	3:A:1649:HOH:O	2.21	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:461:GLN:CB	3:B:1832:HOH:O	2.66	0.44
1:B:570:SER:HB3	1:B:580:LEU:O	2.18	0.44
1:A:408:TRP:C	1:A:408:TRP:CD1	2.96	0.43
1:B:265:ARG:HA	1:B:266:PRO:C	2.43	0.43
1:A:293:SER:HA	1:A:574:GLY:O	2.18	0.43
1:A:324:HIS:HD2	1:A:327:GLN:OE1	2.02	0.43
1:B:489:ASP:OD1	1:B:490:LYS:HG2	2.18	0.43
1:B:349:LEU:N	1:B:350:PRO:CD	2.81	0.43
1:A:45:ILE:O	1:A:45:ILE:HG23	2.16	0.43
1:B:101:ASP:O	1:B:104:VAL:CG1	2.51	0.43
1:A:178:GLU:O	1:A:178:GLU:HG3	2.18	0.43
1:A:44:ASP:CB	1:A:71:LYS:NZ	2.82	0.43
1:A:219:HIS:HB2	1:A:433:PRO:HA	2.01	0.43
1:B:457:GLY:O	1:B:458:ALA:C	2.62	0.42
1:B:45:ILE:HD13	1:B:45:ILE:N	2.35	0.42
1:B:341:ASN:HD22	1:B:342:PRO:CD	2.31	0.42
1:B:433:PRO:O	1:B:450:HIS:HA	2.19	0.42
1:A:404:HIS:CE1	3:A:1385:HOH:O	2.62	0.42
1:B:159:ARG:HA	2:B:801:FAD:O2B	2.19	0.42
1:A:47:TYR:CD2	1:A:73:ALA:HB2	2.55	0.42
1:B:478:GLU:HG3	1:B:511:THR:OG1	2.21	0.41
1:B:81:ASP:O	1:B:83:GLY:N	2.45	0.40
1:B:110:GLN:HG3	1:B:158:THR:HG23	2.03	0.40
1:A:471:TRP:CH2	1:A:526:SER:HA	2.56	0.40
1:B:340:PRO:HD3	1:B:486:TRP:CG	2.57	0.40
1:A:328:LEU:HD23	1:A:328:LEU:C	2.45	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	573/623 (92%)	552 (96%)	20 (4%)	1 (0%)	43	36
1	B	574/623 (92%)	557 (97%)	15 (3%)	2 (0%)	36	29
All	All	1147/1246 (92%)	1109 (97%)	35 (3%)	3 (0%)	36	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	82	SER
1	B	458	ALA
1	A	45	ILE

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	502/541 (93%)	490 (98%)	12 (2%)	43	38
1	B	503/541 (93%)	490 (97%)	13 (3%)	40	35
All	All	1005/1082 (93%)	980 (98%)	25 (2%)	42	37

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ASP
1	A	178	GLU
1	A	206	PHE
1	A	223	LEU
1	A	235	ASP
1	A	272	GLU
1	A	286	VAL
1	A	349	LEU
1	A	385	THR
1	A	460	GLN
1	A	516	GLU
1	A	593	ASN
1	B	45	ILE

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Mol	Chain	Res	Type
1	B	185	LYS
1	B	206	PHE
1	B	215	GLU
1	B	246	ARG
1	B	285	ARG
1	B	310	GLU
1	B	336	GLN
1	B	408	TRP
1	B	460	GLN
1	B	560	LYS
1	B	576	LYS
1	B	593	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	ASN
1	A	257	ASN
1	A	263	GLN
1	A	299	HIS
1	A	324	HIS
1	A	341	ASN
1	A	404	HIS
1	A	419	HIS
1	B	233	GLN
1	B	263	GLN
1	B	324	HIS
1	B	341	ASN
1	B	344	ASN
1	B	365	GLN
1	B	404	HIS
1	B	419	HIS
1	B	461	GLN
1	B	496	ASN
1	B	611	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	B	801	-	58,58,58	1.24	6 (10%)	85,89,89	2.31	27 (31%)
2	FAD	A	801	-	58,58,58	1.20	6 (10%)	85,89,89	2.40	30 (35%)

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	801	-	-	2/34/50/50	0/6/6/6
2	FAD	A	801	-	-	0/34/50/50	0/6/6/6

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	FAD	P-O3P	3.88	1.63	1.59
2	A	801	FAD	C8M-C8	2.97	1.56	1.51
2	A	801	FAD	C5'-C4'	2.85	1.55	1.51
2	B	801	FAD	C1'-C2'	2.55	1.56	1.52
2	B	801	FAD	C10-N1	2.51	1.38	1.33
2	A	801	FAD	C2A-N1A	2.43	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	FAD	C5A-N7A	-2.32	1.34	1.39
2	B	801	FAD	C2A-N3A	2.28	1.38	1.33
2	A	801	FAD	C4X-N5	2.14	1.35	1.30
2	B	801	FAD	C4-N3	-2.12	1.34	1.38
2	A	801	FAD	O2B-C2B	-2.09	1.37	1.43
2	A	801	FAD	O4'-C4'	2.02	1.47	1.43

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	C4A-N9A-C8A	8.24	114.39	105.74
2	B	801	FAD	C4A-N9A-C8A	7.74	113.87	105.74
2	A	801	FAD	N9A-C8A-N7A	-6.58	104.60	113.94
2	B	801	FAD	N9A-C8A-N7A	-6.48	104.75	113.94
2	B	801	FAD	N3A-C2A-N1A	-5.97	119.54	128.58
2	B	801	FAD	O4-C4-C4X	-5.77	111.30	126.53
2	A	801	FAD	C2B-C1B-N9A	5.60	127.21	113.30
2	A	801	FAD	O4-C4-C4X	-5.04	113.22	126.53
2	A	801	FAD	N3A-C4A-N9A	5.02	135.71	127.17
2	A	801	FAD	N3A-C2A-N1A	-4.92	121.14	128.58
2	A	801	FAD	C1B-N9A-C8A	-4.58	116.92	127.09
2	B	801	FAD	C4-C4X-N5	4.53	124.47	118.21
2	B	801	FAD	C2B-C1B-N9A	4.48	124.44	113.30
2	B	801	FAD	N3A-C4A-N9A	4.39	134.63	127.17
2	A	801	FAD	O2A-PA-O3P	4.29	118.86	107.27
2	A	801	FAD	C5A-C4A-N3A	-4.06	121.13	126.72
2	B	801	FAD	O4B-C1B-N9A	3.94	115.66	108.09
2	A	801	FAD	C5A-N7A-C8A	3.76	109.36	103.45
2	B	801	FAD	C1B-N9A-C8A	-3.63	119.03	127.09
2	A	801	FAD	C5B-C4B-C3B	3.56	128.04	115.21
2	A	801	FAD	C4-C4X-N5	3.54	123.10	118.21
2	A	801	FAD	O4B-C1B-N9A	3.42	114.66	108.09
2	A	801	FAD	O3B-C3B-C4B	3.40	120.86	111.08
2	B	801	FAD	O4B-C4B-C3B	3.33	111.76	105.15
2	B	801	FAD	C5A-C4A-N9A	-3.19	102.33	105.81
2	B	801	FAD	O4B-C4B-C5B	3.18	119.53	109.33
2	B	801	FAD	O4-C4-N3	3.14	126.02	120.11
2	B	801	FAD	O2B-C2B-C3B	3.11	121.79	111.82
2	A	801	FAD	O2B-C2B-C1B	3.09	120.75	110.10
2	B	801	FAD	C5A-N7A-C8A	3.01	108.18	103.45
2	A	801	FAD	O3'-C3'-C2'	2.97	115.68	108.93
2	A	801	FAD	C7M-C7-C6	-2.84	114.58	119.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	FAD	O4B-C4B-C3B	2.80	110.71	105.15
2	A	801	FAD	C2A-N3A-C4A	2.76	118.56	111.83
2	B	801	FAD	C2A-N3A-C4A	2.70	118.42	111.83
2	B	801	FAD	O3P-PA-O1A	-2.69	102.61	110.70
2	B	801	FAD	C5A-C4A-N3A	-2.67	123.04	126.72
2	B	801	FAD	O2B-C2B-C1B	2.60	119.05	110.10
2	A	801	FAD	C10-C4X-N5	-2.46	119.79	124.81
2	B	801	FAD	C10-N1-C2	2.45	122.15	116.85
2	A	801	FAD	C5A-C4A-N9A	-2.44	103.15	105.81
2	A	801	FAD	O4-C4-N3	2.35	124.53	120.11
2	B	801	FAD	C4'-C3'-C2'	-2.31	109.72	113.57
2	B	801	FAD	O2A-PA-O3P	2.29	113.47	107.27
2	A	801	FAD	O4B-C4B-C5B	2.20	116.39	109.33
2	B	801	FAD	O3B-C3B-C4B	2.18	117.34	111.08
2	A	801	FAD	C10-N1-C2	2.13	121.47	116.85
2	B	801	FAD	C5B-C4B-C3B	2.10	122.79	115.21
2	A	801	FAD	O2P-P-O3P	-2.09	101.64	107.27
2	A	801	FAD	C9A-C5X-N5	-2.08	120.25	122.45
2	A	801	FAD	O4'-C4'-C5'	2.07	114.55	109.99
2	A	801	FAD	C3B-C2B-C1B	2.07	105.38	101.46
2	A	801	FAD	C4X-C10-N1	-2.06	119.53	124.59
2	B	801	FAD	O4B-C1B-C2B	2.06	111.02	106.62
2	B	801	FAD	C4X-C4-N3	2.04	118.46	113.25
2	B	801	FAD	C2A-N1A-C6A	2.02	122.05	118.73
2	A	801	FAD	O2'-C2'-C1'	-2.00	101.96	110.20

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	801	FAD	PA-O3P-P-O5'
2	B	801	FAD	O4B-C4B-C5B-O5B

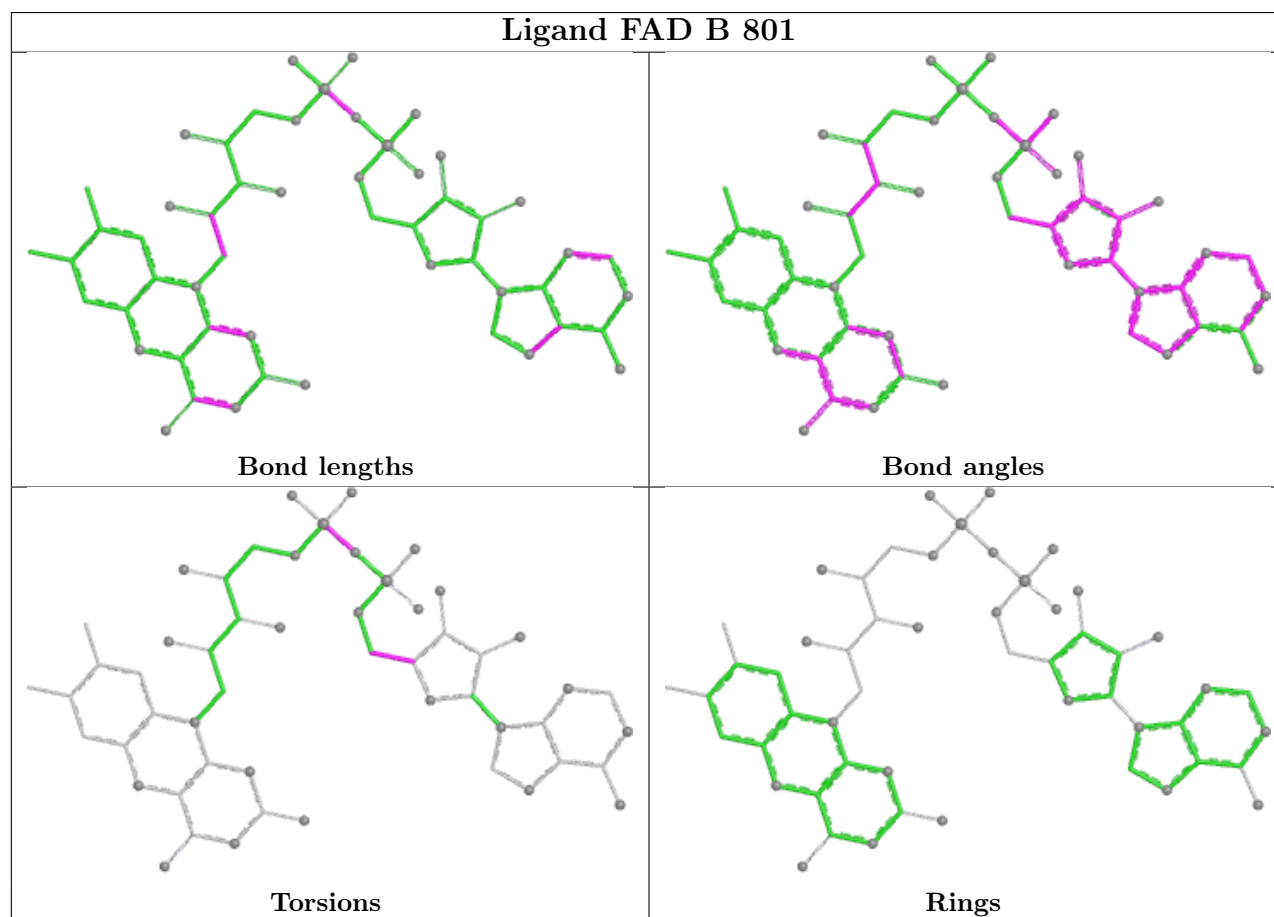
There are no ring outliers.

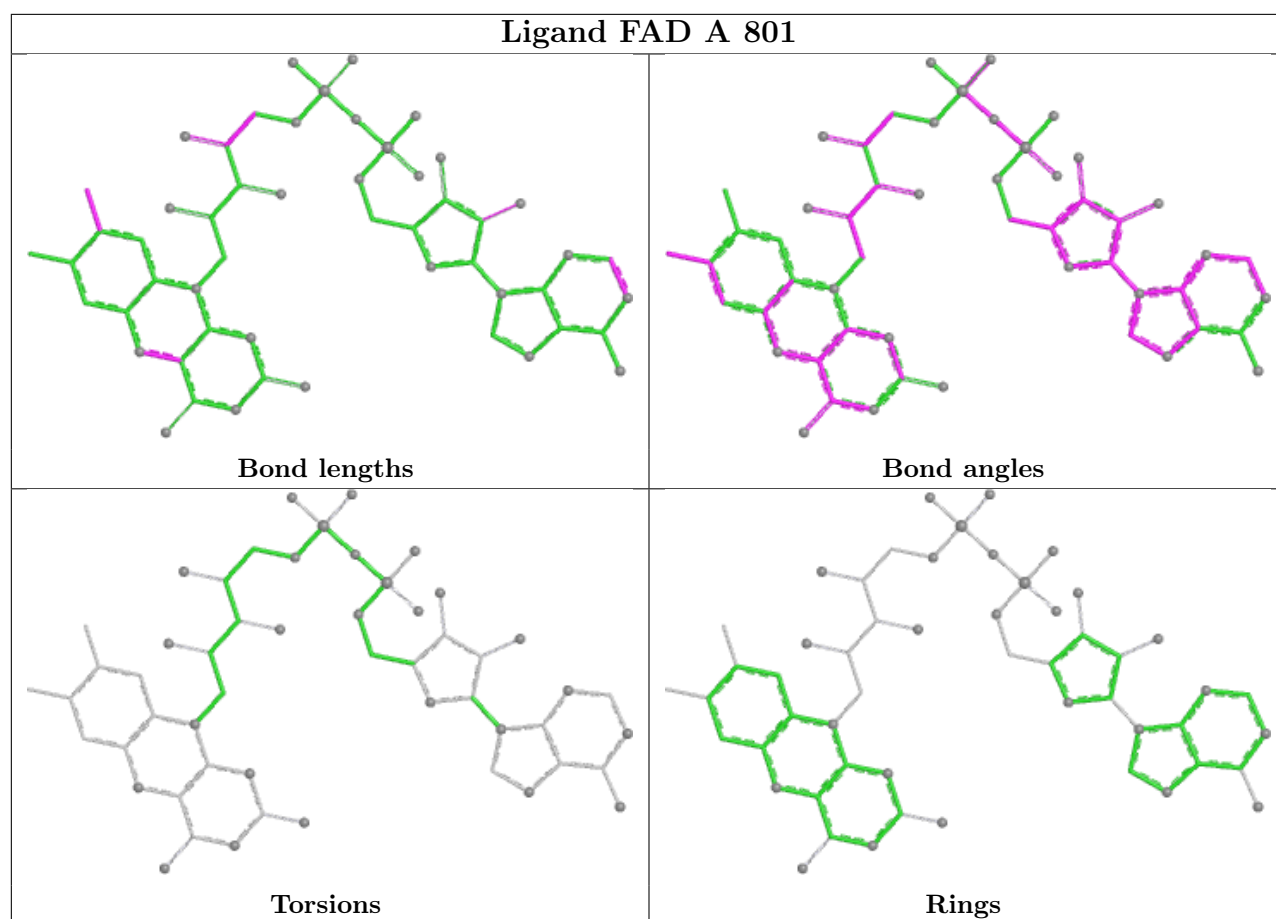
2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	FAD	3	0
2	A	801	FAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	575/623 (92%)	-0.32	10 (1%) 69 72	12, 17, 33, 55	0
1	B	576/623 (92%)	-0.27	10 (1%) 69 72	13, 19, 35, 58	0
All	All	1151/1246 (92%)	-0.30	20 (1%) 69 72	12, 18, 34, 58	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	459	VAL	5.9
1	A	45	ILE	5.7
1	B	458	ALA	5.0
1	B	459	VAL	4.8
1	A	458	ALA	4.6
1	A	457	GLY	3.8
1	B	232	GLY	3.8
1	B	457	GLY	3.5
1	A	44	ASP	3.4
1	B	343	ALA	3.3
1	B	344	ASN	3.1
1	B	460	GLN	2.9
1	A	343	ALA	2.8
1	A	347	GLU	2.5
1	A	460	GLN	2.4
1	A	618	PHE	2.4
1	B	81	ASP	2.3
1	A	82	SER	2.3
1	B	228	GLU	2.1
1	B	299	HIS	2.1

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](#)

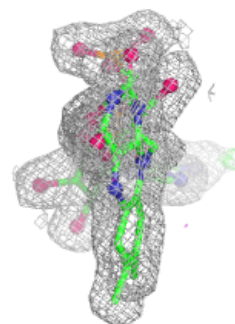
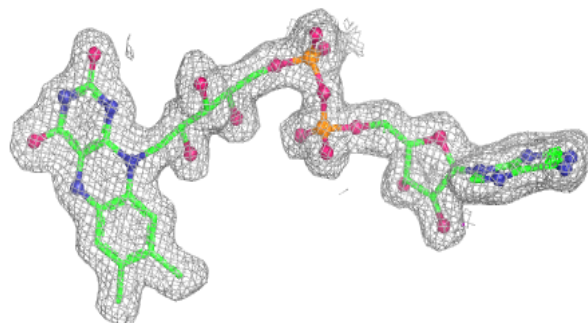
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	FAD	A	801	53/53	0.98	0.04	9,13,17,18	0
2	FAD	B	801	53/53	0.98	0.05	10,14,18,19	0

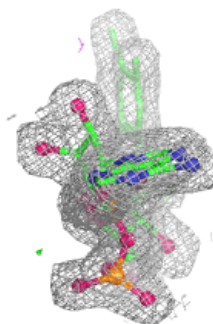
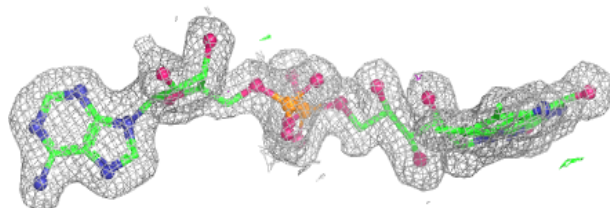
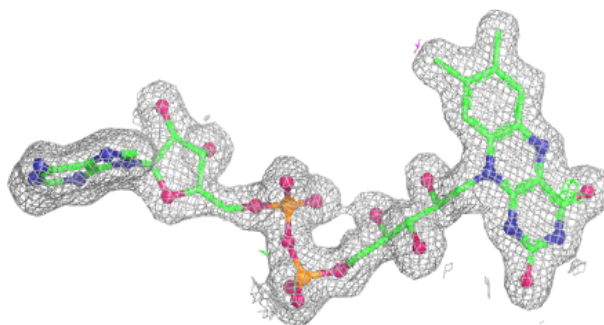
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD A 801:

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

**Electron density around FAD B 801:**

$2mF_o-DF_c$ (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.