Full wwPDB X-ray Structure Validation Report (i)

Aug 21, 2020 – 08:59 PM BST

PDB ID	:	6K8I
Title	:	Crystal structure of Arabidopsis thaliana CRY2
Authors	:	Ma, L.; Wang, X.; Guan, Z.; Yin, P.
Deposited on	:	2019-06-12
Resolution	:	2.70 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 (i) symbol.

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

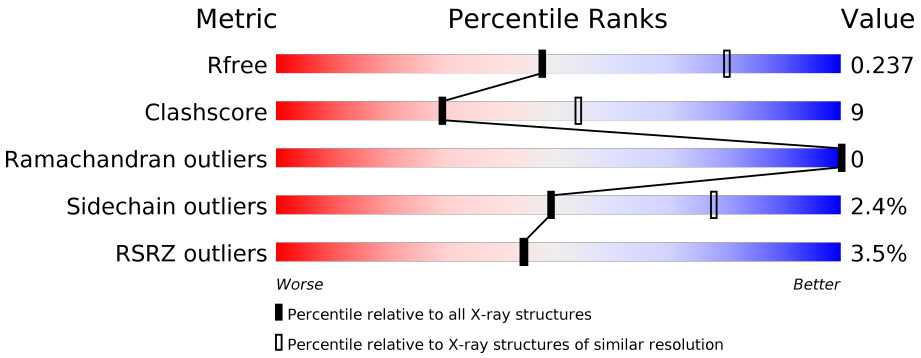
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	612	3% 58% 16% • 25%
1	B	612	2% 57% 18% 25%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7608 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.

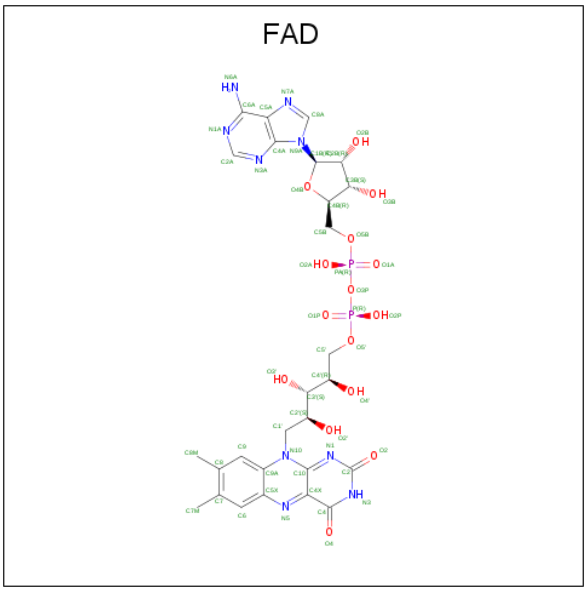
- Molecule 1 is a protein called Cryptochrome-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	457	Total	C	N	O	S	0	1	0
			3728	2412	632	668	16			
1	B	461	Total	C	N	O	S	0	0	0
			3752	2427	638	671	16			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	GLN	LYS	See Sequence details	UNP Q96524
A	95	PRO	ALA	See Sequence details	UNP Q96524
A	366	GLY	ALA	See Sequence details	UNP Q96524
A	534	VAL	ALA	See Sequence details	UNP Q96524
B	78	GLN	LYS	See Sequence details	UNP Q96524
B	95	PRO	ALA	See Sequence details	UNP Q96524
B	366	GLY	ALA	See Sequence details	UNP Q96524
B	534	VAL	ALA	See Sequence details	UNP Q96524

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by author).



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		

- Molecule 3 is water.

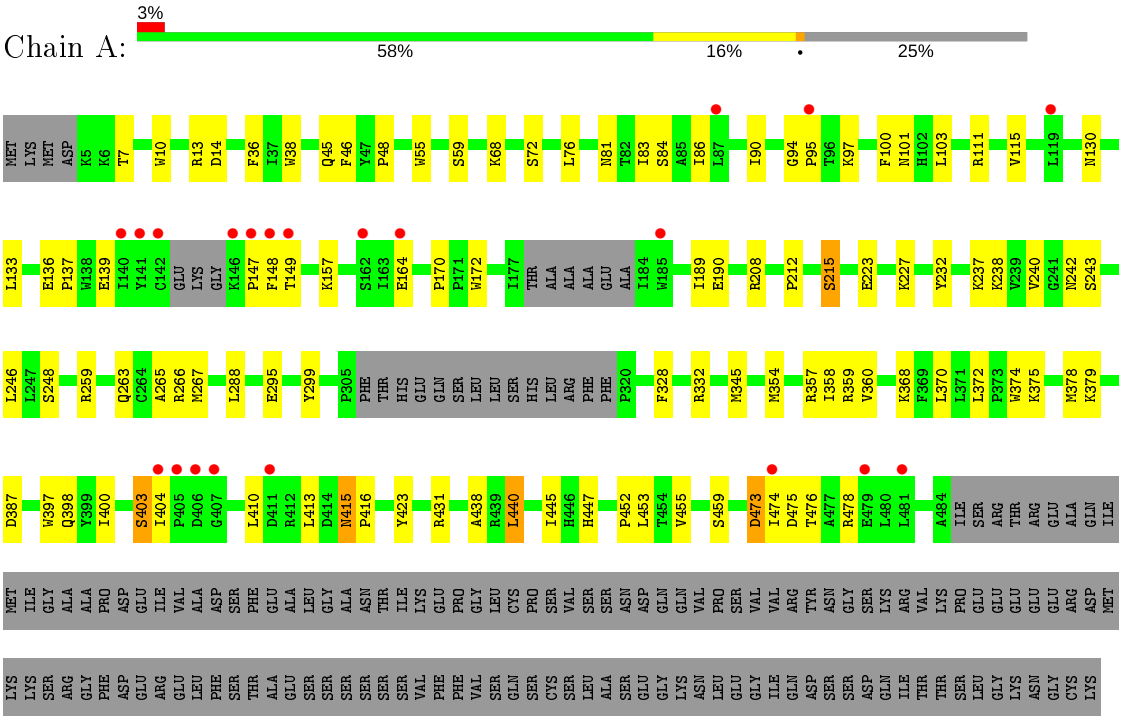
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	9	Total	O	0	0
			9	9		
3	B	13	Total	O	0	0
			13	13		

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.

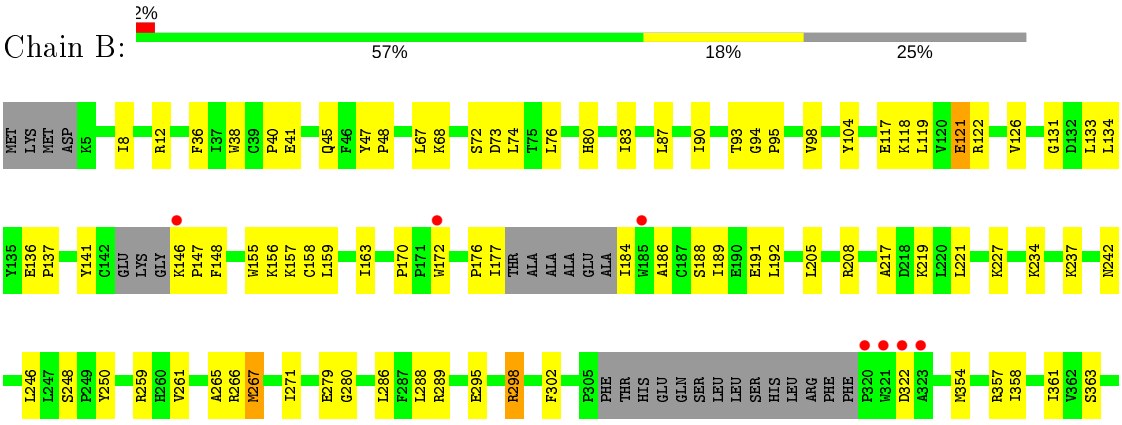
- Molecule 1: Cryptochrome-2

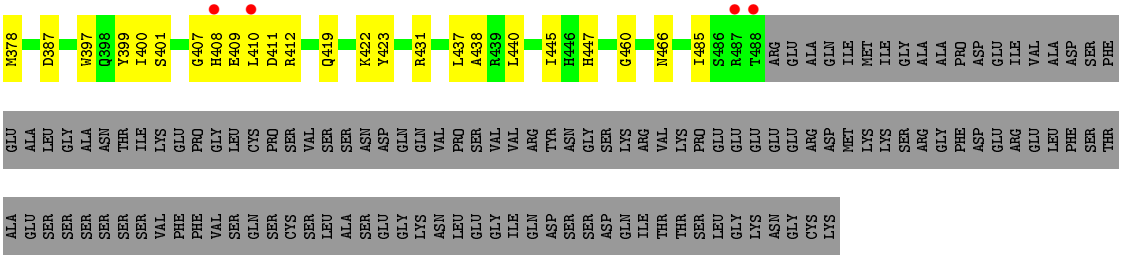
Chain A:



- Molecule 1: Cryptochrome-2

Chain B:




M378 D387 W397 Q398 Y399 I400 S401 G407 H408 E409 L410 D411 R412 Q419 K422 Y423 R431 L437 A438 R439 L440 I445 H446 H447 G460 N466 I485 S486 R487 T488
GLU ALA LEU GLY ALA ASN THR ILE LYS GLU PRO GLY LEU CYS PRO SER VAL SER SER ASN ASP GLN GLN VAL PRO SER VAL VAL ARG TYR ASN GLY SER LYS ARG VAL LYS PRO GLU GLU GLU GLU GLU ARG ASP MET LYS LYS SER ARG GLY PHE ASP GLU ARG GLU LEU PHE SER THR
ALA GLU SER SER SER SER SER VAL PHE PHE VAL SER GLN SER CYS SER LEU ALA SER GLU GLY LYS ASN LEU GLU GLY ILE GLN ASP SER SER ASP GLN ILE THR THR SER LEU GLY LYS ASN GLY CYS LYS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	123.12Å 160.37Å 139.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.03 – 2.70 49.04 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.03-2.70) 99.8 (49.04-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.26 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???)	Depositor
R, R_{free}	0.185 , 0.237 0.185 , 0.237	Depositor DCC
R_{free} test set	1948 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å^2)	61.6	Xtriage
Anisotropy	0.681	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 54.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7608	wwPDB-VP
Average B, all atoms (Å^2)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.44	0/3839	0.60	1/5219 (0.0%)
1	B	0.47	0/3863	0.63	3/5251 (0.1%)
All	All	0.45	0/7702	0.62	4/10470 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	422	LYS	CD-CE-NZ	-5.52	99.01	111.70
1	A	459	SER	C-N-CA	-5.43	110.89	122.30
1	B	121	GLU	CA-CB-CG	-5.37	101.58	113.40
1	B	422	LYS	CB-CG-CD	-5.12	98.27	111.60

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	3728	0	3657	71	0
1	B	3752	0	3690	77	0
2	A	53	0	30	5	0
2	B	53	0	30	3	0
3	A	9	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	13	0	0	0	0
All	All	7608	0	7407	142	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 9.

All (142) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:701:FAD:O4B	2:A:701:FAD:C1B	1.63	1.14
1:A:266:ARG:NH1	1:B:172:TRP:HE3	1.46	1.13
1:A:266:ARG:NH1	1:B:172:TRP:CE3	2.21	1.09
1:A:266:ARG:CZ	1:B:172:TRP:HE3	1.75	0.99
1:A:189:ILE:HD12	1:A:190:GLU:H	1.32	0.91
1:A:266:ARG:CZ	1:B:172:TRP:CE3	2.58	0.86
1:B:460:GLY:O	1:B:466:ASN:ND2	2.11	0.83
1:B:234:LYS:HD3	1:B:237:LYS:HE3	1.62	0.80
1:A:189:ILE:HD12	1:A:190:GLU:N	2.01	0.76
1:A:431:ARG:HG2	1:A:438:ALA:HA	1.69	0.75
1:B:431:ARG:HG2	1:B:438:ALA:HA	1.70	0.74
1:B:409:GLU:HB3	1:B:411:ASP:OD1	1.89	0.73
1:B:36:PHE:HB3	1:B:76:LEU:HD23	1.72	0.70
1:B:41:GLU:OE2	1:B:80:HIS:ND1	2.16	0.69
1:A:398:GLN:HB3	1:A:404:ILE:HG12	1.74	0.68
1:B:83:ILE:HD12	1:B:83:ILE:H	1.58	0.67
1:A:242:ASN:HB3	1:A:246:LEU:HD21	1.78	0.65
1:B:131:GLY:HA2	1:B:259:ARG:HH22	1.59	0.65
1:A:475:ASP:O	1:A:478:ARG:HB2	1.98	0.64
1:A:90:ILE:HG23	1:A:95:PRO:HD2	1.79	0.63
1:B:205:LEU:O	1:B:208:ARG:HB2	2.00	0.62
1:A:238:LYS:HD2	1:A:240:VAL:O	2.00	0.61
1:A:232:TYR:OH	2:A:701:FAD:O1A	2.16	0.61
1:A:398:GLN:HG2	1:A:403:SER:OG	2.01	0.61
1:A:370:LEU:HB3	1:A:372:LEU:HD13	1.81	0.61
1:B:411:ASP:HA	1:B:485:ILE:HD11	1.81	0.61
1:B:98:VAL:HG11	1:B:119:LEU:HD13	1.81	0.60
1:B:248:SER:HB3	2:B:701:FAD:H5'2	1.82	0.60
1:B:188:SER:O	1:B:192:LEU:HD13	2.02	0.60
1:A:13:ARG:CZ	1:A:103:LEU:HD13	2.32	0.59
1:A:172:TRP:CZ3	1:B:266:ARG:HD3	2.38	0.59
1:B:184:ILE:HG13	1:B:186:ALA:H	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:THR:HG23	1:A:97:LYS:HB3	1.85	0.58
1:A:440:LEU:HD13	1:A:445:ILE:HA	1.85	0.58
1:A:45:GLN:H	1:A:45:GLN:CD	2.07	0.58
1:A:111:ARG:O	1:A:115:VAL:HG13	2.05	0.56
1:A:215:SER:HB2	1:B:271:ILE:HD12	1.88	0.56
1:A:267:MET:SD	1:B:170:PRO:HG3	2.44	0.56
1:A:266:ARG:NH2	1:B:172:TRP:HA	2.22	0.55
1:B:159:LEU:HD23	1:B:163:ILE:HD12	1.88	0.55
1:A:147:PRO:HB2	1:A:149:THR:HG23	1.88	0.55
1:B:357:ARG:HD3	1:B:423:TYR:CE2	2.42	0.55
1:B:148:PHE:CZ	1:B:157:LYS:HG2	2.42	0.54
1:B:45:GLN:CD	1:B:45:GLN:H	2.10	0.54
1:B:156:LYS:NZ	1:B:289:ARG:HH12	2.06	0.54
1:A:368:LYS:HE2	1:A:410:LEU:HA	1.91	0.53
1:B:354:MET:HE3	1:B:358:ILE:HG22	1.90	0.53
1:B:408:HIS:NE2	1:B:412:ARG:HB3	2.23	0.53
1:A:170:PRO:HG3	1:B:267:MET:SD	2.49	0.53
1:A:14:ASP:OD2	1:A:101:ASN:ND2	2.33	0.53
1:A:38:TRP:HH2	1:A:48:PRO:HG2	1.74	0.52
1:A:133:LEU:HB3	1:A:295:GLU:OE2	2.10	0.52
1:A:354:MET:HG3	1:A:359:ARG:HG2	1.90	0.52
1:B:155:TRP:O	1:B:158:CYS:HB2	2.10	0.52
1:B:242:ASN:HB3	1:B:246:LEU:HD11	1.91	0.51
1:B:93:THR:OG1	1:B:95:PRO:HD3	2.10	0.51
1:A:415:ASN:HD22	1:A:416:PRO:HD2	1.74	0.51
1:B:134:LEU:N	1:B:295:GLU:OE2	2.44	0.51
1:A:10:TRP:HB3	1:A:100:PHE:HB3	1.92	0.50
1:B:38:TRP:HH2	1:B:48:PRO:HG2	1.76	0.50
1:B:265:ALA:HB3	1:B:288:LEU:HD21	1.93	0.50
1:A:86:ILE:O	1:A:90:ILE:HD12	2.11	0.50
1:B:408:HIS:HD2	1:B:409:GLU:O	1.94	0.50
1:B:146:LYS:HD2	1:B:146:LYS:N	2.27	0.50
1:A:83:ILE:CG1	1:A:115:VAL:HG12	2.42	0.50
1:B:217:ALA:HB2	1:B:250:TYR:CD1	2.47	0.49
1:B:94:GLY:N	1:B:95:PRO:HD3	2.26	0.49
1:A:259:ARG:O	1:A:263:GLN:HG2	2.13	0.49
1:A:374:TRP:CZ3	1:A:375:LYS:HG2	2.47	0.49
1:B:90:ILE:HA	1:B:95:PRO:CD	2.42	0.49
1:B:8:ILE:HD11	1:B:95:PRO:HG3	1.94	0.49
1:B:399:TYR:CD2	1:B:407:GLY:HA2	2.48	0.48
1:A:81:ASN:HB3	1:A:84:SER:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLU:OE1	1:A:227:LYS:HG3	2.14	0.48
1:B:401:SER:O	1:B:410:LEU:HD21	2.14	0.47
1:A:357:ARG:HD2	1:A:423:TYR:CE1	2.49	0.47
1:A:130:ASN:HB2	1:A:136:GLU:OE2	2.14	0.47
1:B:440:LEU:HD13	1:B:445:ILE:HA	1.97	0.47
2:A:701:FAD:O5B	2:A:701:FAD:H8A	2.14	0.47
1:B:361:ILE:HG12	1:B:419:GLN:NE2	2.29	0.47
1:B:363:SER:OG	1:B:400:ILE:HG13	2.15	0.46
1:A:370:LEU:CB	1:A:372:LEU:HD13	2.45	0.46
1:B:98:VAL:HG13	1:B:126:VAL:HG13	1.96	0.46
1:A:328:PHE:HA	1:A:372:LEU:HD21	1.97	0.46
1:A:452:PRO:HG2	1:A:455:VAL:CG2	2.45	0.46
1:A:413:LEU:HB3	1:A:478:ARG:HD2	1.98	0.46
2:B:701:FAD:H8A	2:B:701:FAD:O5B	2.16	0.46
1:B:148:PHE:HZ	1:B:157:LYS:HG2	1.80	0.46
1:B:437:LEU:HD22	1:B:440:LEU:HD12	1.98	0.46
1:B:40:PRO:HB3	1:B:47:TYR:CE1	2.51	0.45
1:B:146:LYS:HB3	1:B:147:PRO:HD2	1.98	0.45
1:A:360:VAL:HG13	1:A:400:ILE:HG22	1.98	0.45
1:B:279:GLU:HG2	1:B:280:GLY:N	2.31	0.45
1:B:387:ASP:O	2:B:701:FAD:N3	2.41	0.45
1:A:453:LEU:H	1:A:453:LEU:HD12	1.81	0.45
1:B:189:ILE:HA	1:B:192:LEU:HD13	1.98	0.45
1:B:133:LEU:HD21	1:B:298:ARG:HD3	1.99	0.45
1:B:408:HIS:CD2	1:B:409:GLU:O	2.69	0.45
1:A:45:GLN:HG2	1:A:46:PHE:CD2	2.52	0.45
1:B:242:ASN:HB3	1:B:246:LEU:HD21	1.99	0.45
1:B:104:TYR:OH	1:B:136:GLU:OE1	2.22	0.45
1:B:72:SER:OG	1:B:73:ASP:N	2.48	0.45
1:A:378:MET:HE2	1:A:397:TRP:CG	2.52	0.45
1:A:332:ARG:HA	1:A:345:MET:HE2	1.99	0.44
1:B:90:ILE:HA	1:B:95:PRO:HD2	1.99	0.44
1:A:136:GLU:HB2	1:A:139:GLU:HG3	2.00	0.44
1:A:357:ARG:CD	1:A:423:TYR:CZ	3.01	0.44
1:B:87:LEU:HD22	1:B:122:ARG:HH21	1.81	0.44
1:B:137:PRO:HB2	1:B:302:PHE:CD2	2.53	0.44
1:B:156:LYS:HZ1	1:B:289:ARG:HH12	1.64	0.44
1:A:387:ASP:O	2:A:701:FAD:N3	2.42	0.43
1:A:431:ARG:NH1	1:A:445:ILE:HD12	2.34	0.43
1:A:248:SER:HB3	2:A:701:FAD:H5'2	2.00	0.43
1:A:94:GLY:N	1:A:95:PRO:HD3	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:LYS:HA	1:A:72:SER:O	2.18	0.43
1:B:363:SER:HB2	1:B:397:TRP:CZ3	2.53	0.43
1:B:117:GLU:O	1:B:121:GLU:OE1	2.37	0.43
1:A:354:MET:HE3	1:A:358:ILE:HG22	1.99	0.43
1:B:67:LEU:HD12	1:B:74:LEU:HB2	2.00	0.43
1:A:36:PHE:HB3	1:A:76:LEU:HD23	2.01	0.42
1:B:68:LYS:HA	1:B:72:SER:O	2.19	0.42
1:B:188:SER:O	1:B:192:LEU:CD1	2.67	0.42
1:A:473[A]:ASP:OD1	1:A:476:THR:HG23	2.19	0.42
1:B:87:LEU:CD2	1:B:122:ARG:HE	2.32	0.42
1:B:205:LEU:O	1:B:208:ARG:CB	2.66	0.42
1:B:188:SER:OG	1:B:191:GLU:HG3	2.19	0.41
1:A:415:ASN:HD22	1:A:416:PRO:CD	2.33	0.41
1:B:141:TYR:CE2	1:B:146:LYS:N	2.88	0.41
1:A:55:TRP:CE3	1:A:212:PRO:HG3	2.54	0.41
1:A:137:PRO:HB3	1:A:299:TYR:HA	2.01	0.41
1:B:221:LEU:HD13	1:B:261:VAL:HG22	2.03	0.41
1:B:83:ILE:H	1:B:83:ILE:CD1	2.29	0.41
1:A:238:LYS:HE3	1:A:243:SER:HB3	2.03	0.41
1:A:164:GLU:OE1	1:A:164:GLU:HA	2.21	0.41
1:A:237:LYS:HA	1:A:237:LYS:HD3	1.76	0.41
1:A:189:ILE:CD1	1:A:190:GLU:N	2.80	0.41
1:A:265:ALA:HB3	1:A:288:LEU:HD21	2.03	0.40
1:A:148:PHE:CE2	1:A:157:LYS:HG2	2.56	0.40
1:A:90:ILE:HA	1:A:95:PRO:CD	2.52	0.40
1:B:176:PRO:O	1:B:177:ILE:HG23	2.21	0.40
1:A:474:ILE:O	1:A:478:ARG:HG2	2.22	0.40
1:B:122:ARG:HG2	1:B:122:ARG:HH11	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	450/612 (74%)	432 (96%)	18 (4%)	0	100	100
1	B	453/612 (74%)	427 (94%)	26 (6%)	0	100	100
All	All	903/1224 (74%)	859 (95%)	44 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	399/531 (75%)	389 (98%)	10 (2%)	47	76
1	B	402/531 (76%)	392 (98%)	10 (2%)	47	76
All	All	801/1062 (75%)	781 (98%)	20 (2%)	49	76

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	SER
1	A	208	ARG
1	A	215	SER
1	A	379	LYS
1	A	403	SER
1	A	415	ASN
1	A	440	LEU
1	A	447	HIS
1	A	473[A]	ASP
1	A	473[B]	ASP
1	B	12	ARG
1	B	118	LYS
1	B	219	LYS
1	B	227	LYS
1	B	267	MET
1	B	286	LEU
1	B	298	ARG
1	B	322	ASP

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Mol	Chain	Res	Type
1	B	378	MET
1	B	447	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	415	ASN
1	B	408	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no monosaccharides 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	701	-	51,58,58	4.62	21 (41%)	60,89,89	2.46	14 (23%)
2	FAD	A	701	-	51,58,58	4.93	21 (41%)	60,89,89	2.43	15 (25%)

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	701	-	-	4/30/50/50	0/6/6/6
2	FAD	A	701	-	-	6/30/50/50	0/6/6/6

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	FAD	O4B-C1B	16.20	1.63	1.41
2	B	701	FAD	O4B-C1B	14.87	1.61	1.41
2	B	701	FAD	C2B-C1B	-14.43	1.31	1.53
2	A	701	FAD	C2B-C1B	-14.07	1.32	1.53
2	A	701	FAD	C10-N1	11.13	1.47	1.33
2	B	701	FAD	C10-N1	10.93	1.47	1.33
2	A	701	FAD	C9A-N10	10.65	1.52	1.38
2	A	701	FAD	C5X-N5	9.36	1.50	1.35
2	A	701	FAD	C4X-N5	8.96	1.46	1.33
2	B	701	FAD	C9A-N10	8.39	1.49	1.38
2	B	701	FAD	C5X-N5	8.33	1.49	1.35
2	B	701	FAD	C4X-N5	8.12	1.44	1.33
2	A	701	FAD	C4-N3	7.96	1.46	1.33
2	B	701	FAD	C4-N3	7.88	1.46	1.33
2	A	701	FAD	C4X-C10	7.80	1.46	1.38
2	A	701	FAD	C4-C4X	6.58	1.52	1.41
2	A	701	FAD	C2-N3	6.26	1.50	1.38
2	B	701	FAD	C4X-C10	6.25	1.45	1.38
2	B	701	FAD	C4-C4X	6.06	1.51	1.41
2	A	701	FAD	O4B-C4B	-5.99	1.31	1.45
2	B	701	FAD	C2-N3	5.86	1.49	1.38
2	B	701	FAD	O4B-C4B	-5.84	1.31	1.45
2	A	701	FAD	C2-N1	5.08	1.48	1.38
2	B	701	FAD	C2-N1	4.63	1.47	1.38
2	A	701	FAD	O2B-C2B	3.87	1.52	1.43
2	A	701	FAD	C1'-N10	3.76	1.52	1.48
2	A	701	FAD	C2A-N3A	3.55	1.37	1.32
2	B	701	FAD	C1'-N10	3.49	1.51	1.48
2	B	701	FAD	C2A-N3A	3.43	1.37	1.32
2	B	701	FAD	O2B-C2B	3.23	1.50	1.43
2	A	701	FAD	O3B-C3B	-3.18	1.35	1.43
2	B	701	FAD	O3B-C3B	-3.07	1.35	1.43
2	B	701	FAD	C6A-N6A	2.81	1.44	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	FAD	O4-C4	-2.61	1.18	1.24
2	B	701	FAD	O2'-C2'	-2.61	1.37	1.43
2	A	701	FAD	C6A-N6A	2.56	1.43	1.34
2	A	701	FAD	O2'-C2'	-2.40	1.38	1.43
2	B	701	FAD	O4'-C4'	-2.38	1.38	1.43
2	B	701	FAD	C5A-C4A	-2.35	1.34	1.40
2	A	701	FAD	C5A-C4A	-2.19	1.35	1.40
2	A	701	FAD	O4-C4	-2.19	1.19	1.24
2	A	701	FAD	O4'-C4'	-2.08	1.39	1.43

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	701	FAD	C5A-C6A-N6A	8.80	133.72	120.35
2	B	701	FAD	C5A-C6A-N6A	8.31	132.98	120.35
2	A	701	FAD	C4-N3-C2	6.08	120.27	115.14
2	B	701	FAD	C7M-C7-C8	6.00	133.04	120.74
2	B	701	FAD	C4-N3-C2	5.80	120.04	115.14
2	A	701	FAD	N6A-C6A-N1A	-5.74	106.65	118.57
2	B	701	FAD	C7M-C7-C6	-5.59	106.96	120.34
2	A	701	FAD	C1'-N10-C9A	5.59	122.69	118.29
2	B	701	FAD	N3A-C2A-N1A	-5.53	120.04	128.68
2	B	701	FAD	C1'-N10-C10	5.46	123.30	118.41
2	B	701	FAD	N6A-C6A-N1A	-5.33	107.51	118.57
2	A	701	FAD	N3A-C2A-N1A	-5.23	120.50	128.68
2	A	701	FAD	C7M-C7-C8	4.58	130.13	120.74
2	B	701	FAD	C5X-C9A-N10	4.38	120.89	117.72
2	A	701	FAD	C7M-C7-C6	-4.23	110.23	120.34
2	A	701	FAD	C5X-C9A-N10	3.89	120.53	117.72
2	A	701	FAD	C5'-C4'-C3'	-3.66	105.14	112.20
2	A	701	FAD	C8M-C8-C7	-3.62	113.32	120.74
2	B	701	FAD	C4X-N5-C5X	3.60	120.37	116.77
2	A	701	FAD	C4X-C4-N3	-3.59	118.52	123.43
2	B	701	FAD	P-O3P-PA	-2.98	122.59	132.83
2	B	701	FAD	C4X-C4-N3	-2.98	119.35	123.43
2	A	701	FAD	C4X-N5-C5X	2.93	119.70	116.77
2	A	701	FAD	C9A-N10-C10	-2.81	118.23	121.91
2	B	701	FAD	C5'-C4'-C3'	-2.75	106.89	112.20
2	B	701	FAD	C9A-N10-C10	-2.72	118.35	121.91
2	A	701	FAD	P-O3P-PA	-2.42	124.52	132.83
2	B	701	FAD	C3B-C2B-C1B	2.21	104.31	100.98
2	A	701	FAD	C8M-C8-C9	2.13	125.44	120.34

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	701	FAD	C5B-O5B-PA-O1A
2	B	701	FAD	C5B-O5B-PA-O2A
2	B	701	FAD	C4'-C5'-O5'-P
2	A	701	FAD	C4'-C5'-O5'-P
2	B	701	FAD	C5B-O5B-PA-O3P
2	A	701	FAD	P-O3P-PA-O2A
2	A	701	FAD	C5B-O5B-PA-O3P
2	A	701	FAD	P-O3P-PA-O1A
2	A	701	FAD	C5B-O5B-PA-O1A
2	A	701	FAD	C5'-O5'-P-O1P

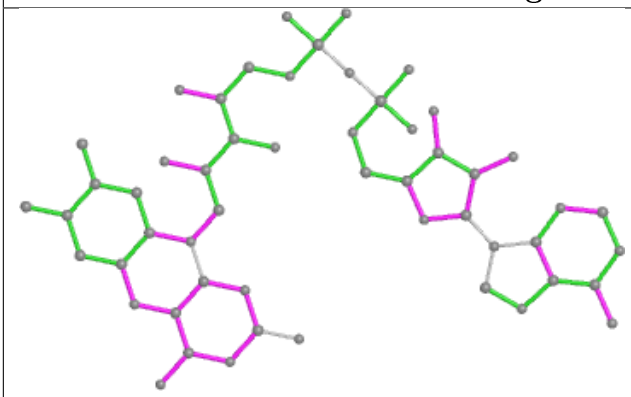
There are no ring outliers.

2 monomers are involved in 8 short contacts:

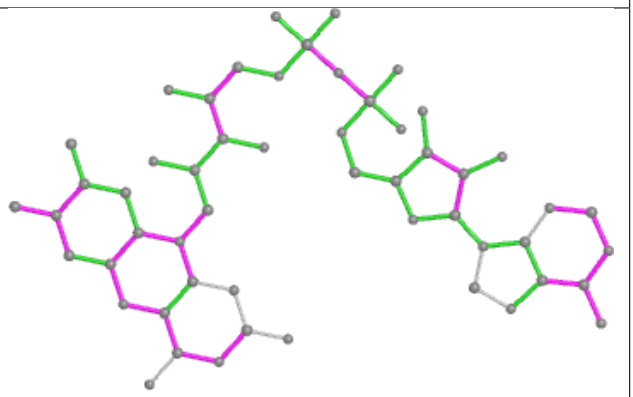
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	FAD	3	0
2	A	701	FAD	5	0

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

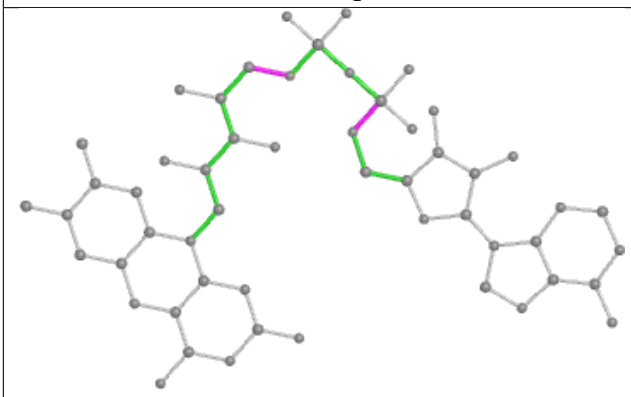
Ligand FAD B 701



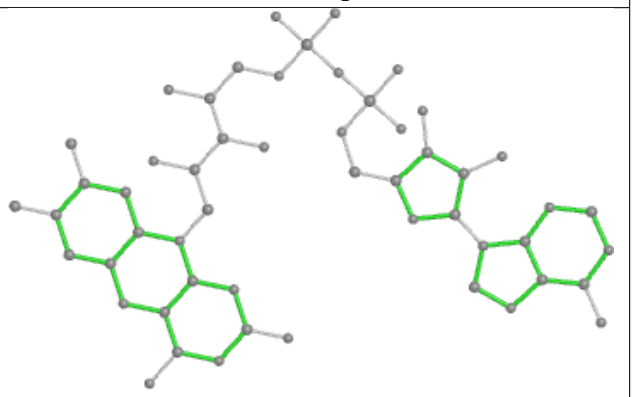
Bond lengths



Bond angles

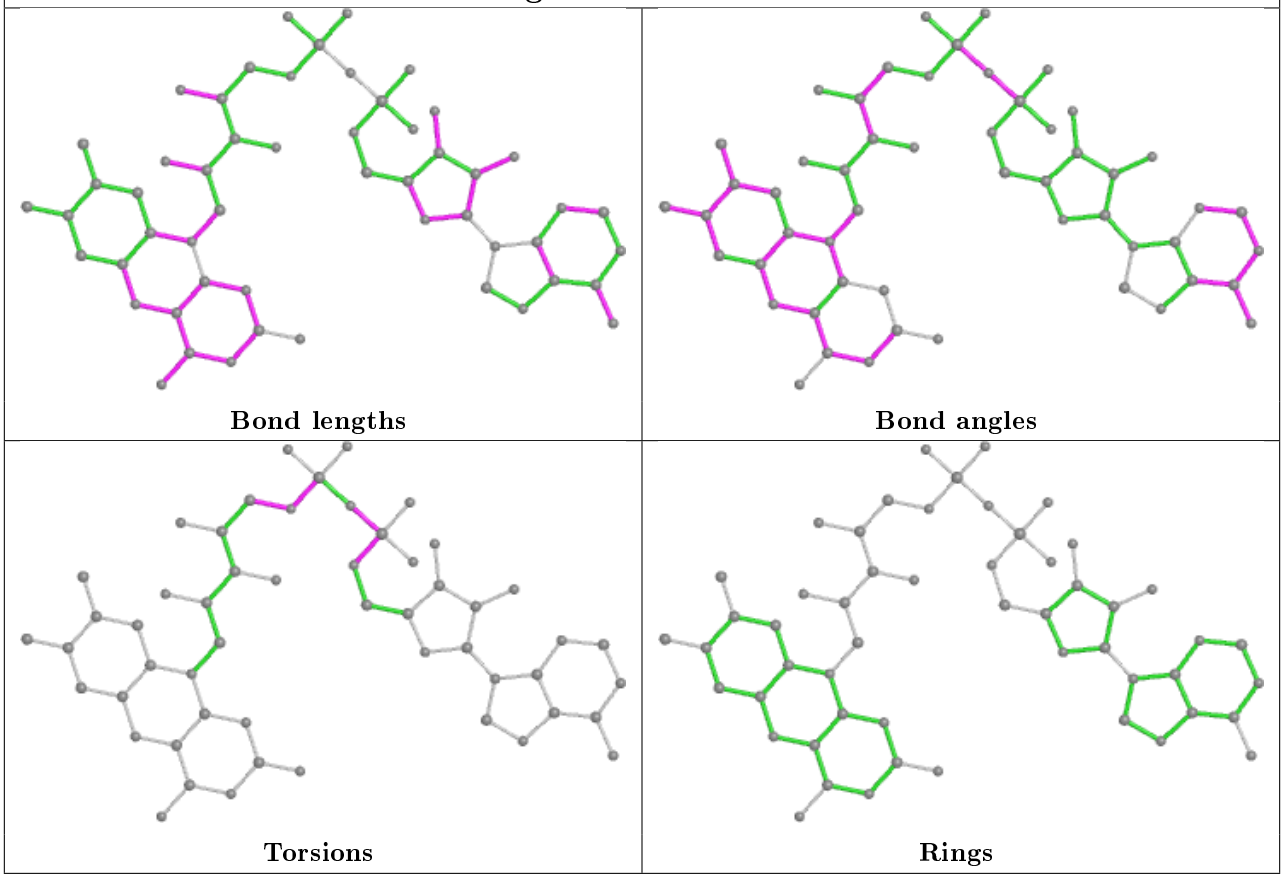


Torsions



Rings

Ligand FAD A 701



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	457/612 (74%)	0.17	21 (4%)	32	31	41, 69, 121, 165	0
1	B	461/612 (75%)	0.02	11 (2%)	59	60	42, 62, 115, 147	0
All	All	918/1224 (75%)	0.09	32 (3%)	44	44	41, 65, 120, 165	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	147	PRO	8.0
1	A	141	TYR	6.0
1	B	320	PRO	5.6
1	B	410	LEU	5.1
1	A	405	PRO	4.9
1	A	406	ASP	4.2
1	B	146	LYS	4.2
1	A	142	CYS	4.1
1	B	488	THR	4.0
1	B	487	ARG	3.6
1	A	140	ILE	3.6
1	A	481	LEU	3.6
1	B	185	TRP	3.5
1	B	321	TRP	3.5
1	A	162	SER	3.5
1	A	407	GLY	3.5
1	A	87	LEU	3.4
1	A	95	PRO	3.3
1	A	404	ILE	3.2
1	A	148	PHE	3.1
1	B	322	ASP	3.0
1	A	146	LYS	2.9
1	A	185	TRP	2.7
1	B	323	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	411	ASP	2.6
1	B	408	HIS	2.5
1	B	172	TRP	2.5
1	A	164	GLU	2.2
1	A	479	GLU	2.2
1	A	474	ILE	2.2
1	A	149	THR	2.0
1	A	119	LEU	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 monosaccharides in this entry.

6.4 Ligands (i)

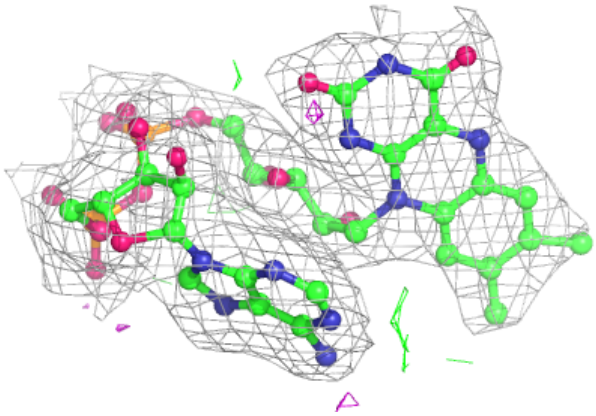
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

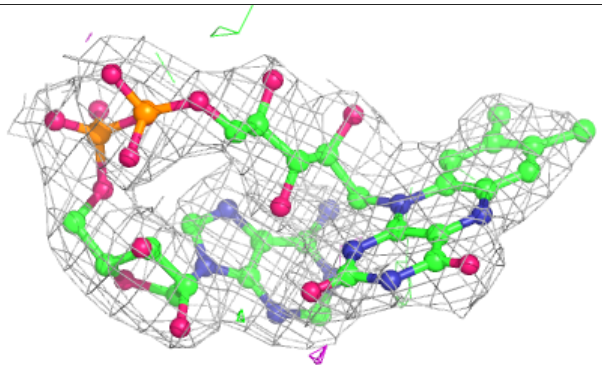
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FAD	B	701	53/53	0.98	0.20	36,48,57,63	0
2	FAD	A	701	53/53	0.98	0.17	44,52,63,70	0

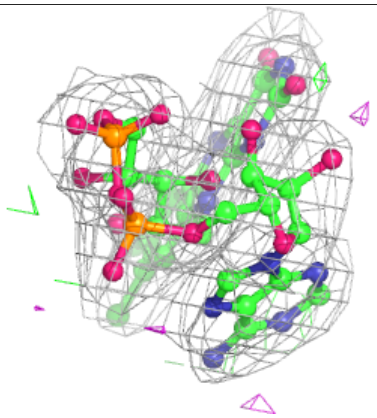
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD B 701:

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

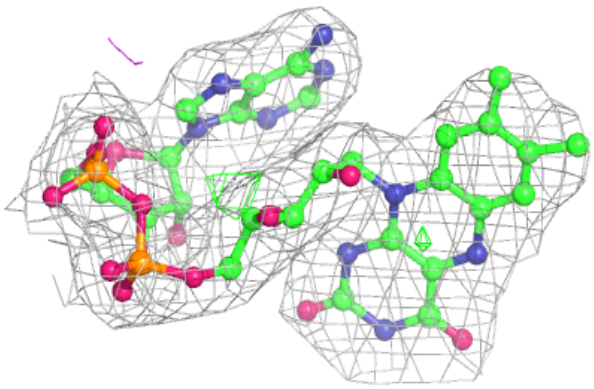


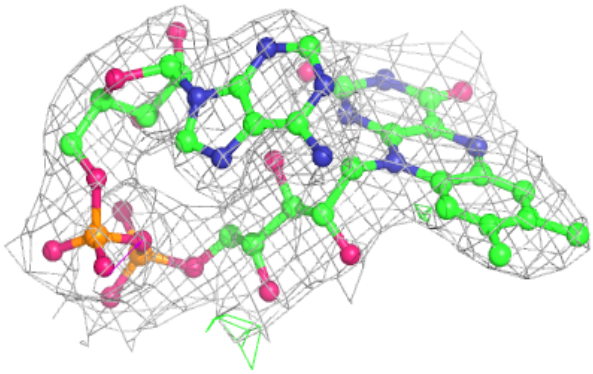


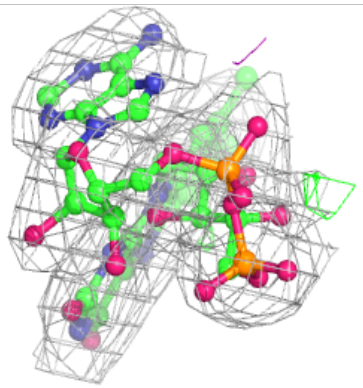


Electron density around FAD A 701:

$2mF_o$-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.