Full wwPDB X-ray Structure Validation Report (i)

Dec 16, 2023 – 01:02 PM EST

PDB ID	:	4P8K
Title	:	Crystal structure of M. tuberculosis DprE1 in complex with the non-covalent inhibitor Ty38c
Authors	:	Neres, J.; Pojer, F.; Cole, S.T.
Deposited on	:	2014-03-31
Resolution	:	2.49 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

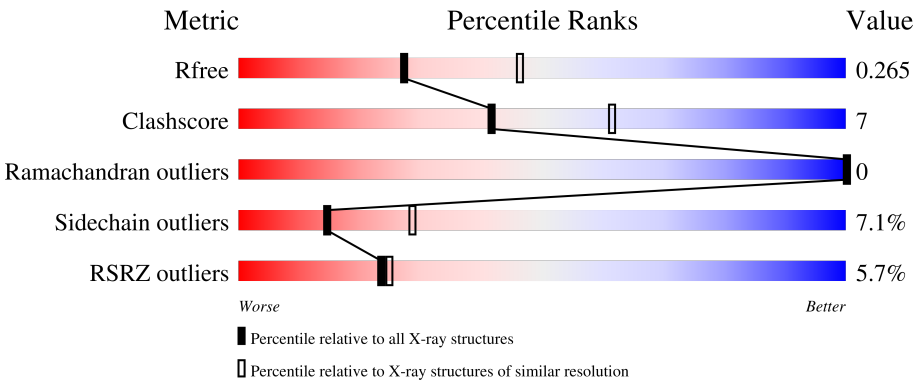
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.49 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	480	5% 76% 16% • 7%
1	B	480	5% 69% 16% • 13%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6907 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 Probable decaprenylphosphoryl-beta-D-ribose oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	448	3442	2192	605	635	10	0	1	0
1	B	419	3218	2040	571	597	10	0	0	0

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	GLY	-	expression tag	UNP P72056
A	-17	SER	-	expression tag	UNP P72056
A	-16	SER	-	expression tag	UNP P72056
A	-15	HIS	-	expression tag	UNP P72056
A	-14	HIS	-	expression tag	UNP P72056
A	-13	HIS	-	expression tag	UNP P72056
A	-12	HIS	-	expression tag	UNP P72056
A	-11	HIS	-	expression tag	UNP P72056
A	-10	HIS	-	expression tag	UNP P72056
A	-9	SER	-	expression tag	UNP P72056
A	-8	SER	-	expression tag	UNP P72056
A	-7	GLY	-	expression tag	UNP P72056
A	-6	LEU	-	expression tag	UNP P72056
A	-5	VAL	-	expression tag	UNP P72056
A	-4	PRO	-	expression tag	UNP P72056
A	-3	ARG	-	expression tag	UNP P72056
A	-2	GLY	-	expression tag	UNP P72056
A	-1	SER	-	expression tag	UNP P72056
A	0	HIS	-	expression tag	UNP P72056
B	-18	GLY	-	expression tag	UNP P72056
B	-17	SER	-	expression tag	UNP P72056
B	-16	SER	-	expression tag	UNP P72056
B	-15	HIS	-	expression tag	UNP P72056
B	-14	HIS	-	expression tag	UNP P72056
B	-13	HIS	-	expression tag	UNP P72056

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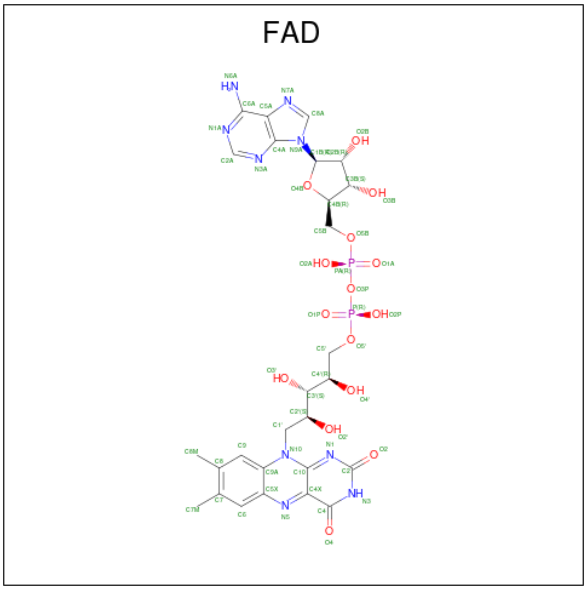
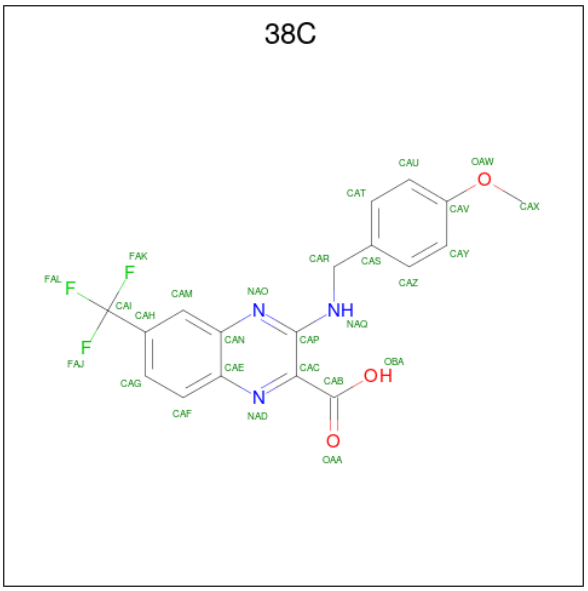
Chain	Residue	Modelled	Actual	Comment	Reference
B	-12	HIS	-	expression tag	UNP P72056
B	-11	HIS	-	expression tag	UNP P72056
B	-10	HIS	-	expression tag	UNP P72056
B	-9	SER	-	expression tag	UNP P72056
B	-8	SER	-	expression tag	UNP P72056
B	-7	GLY	-	expression tag	UNP P72056
B	-6	LEU	-	expression tag	UNP P72056
B	-5	VAL	-	expression tag	UNP P72056
B	-4	PRO	-	expression tag	UNP P72056
B	-3	ARG	-	expression tag	UNP P72056
B	-2	GLY	-	expression tag	UNP P72056
B	-1	SER	-	expression tag	UNP P72056
B	0	HIS	-	expression tag	UNP P72056

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



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

- Molecule 3 is 3-[(4-methoxybenzyl)amino]-6-(trifluoromethyl)quinoxaline-2-carboxylic acid (three-letter code: 38C) (formula: $C_{18}H_{14}F_3N_3O_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			27	18	3	3	3		
3	B	1	Total	C	F	N	O	0	0
			27	18	3	3	3		

- Molecule 4 is water.

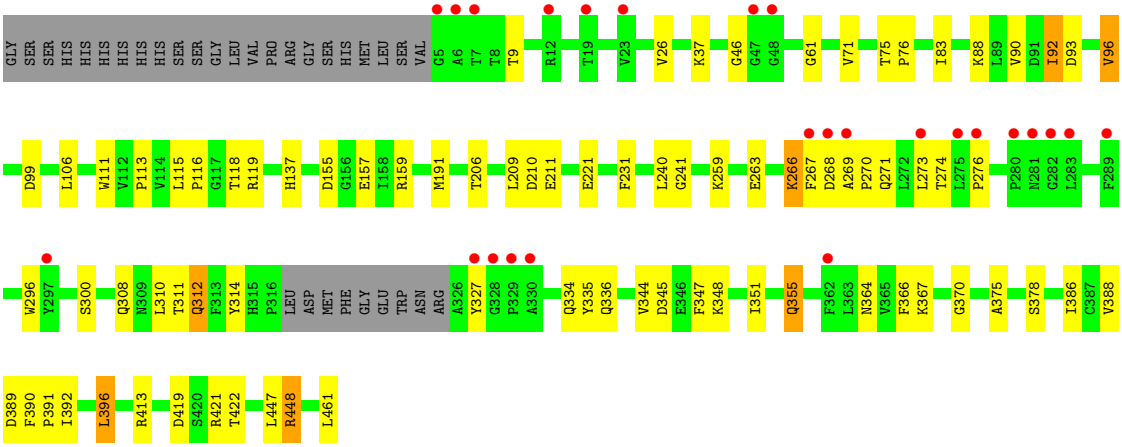
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total	O	0	0
			51	51		
4	B	36	Total	O	0	0
			36	36		

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.

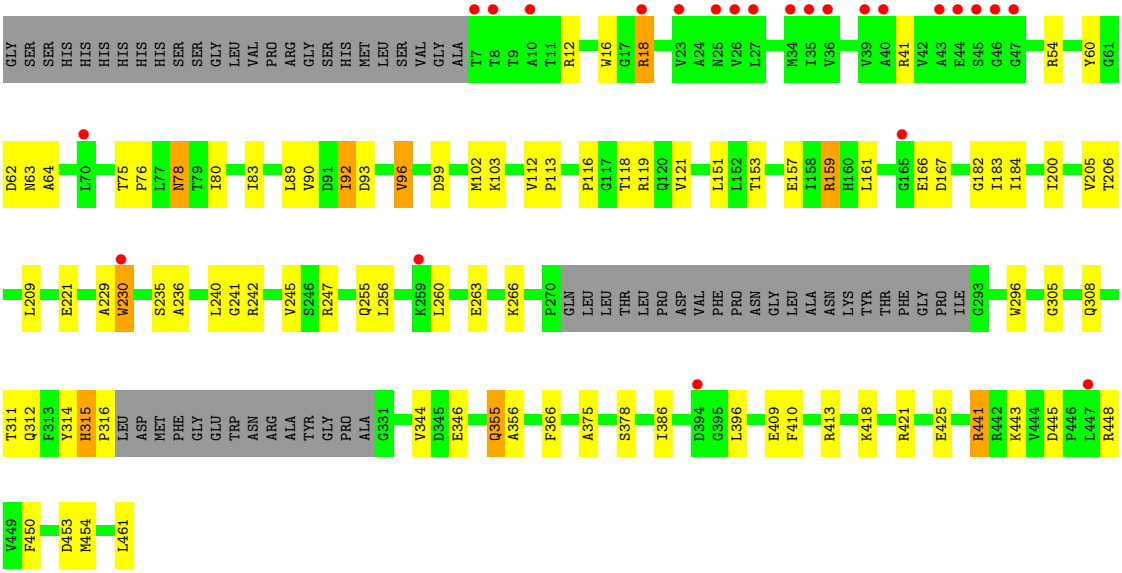
• Molecule 1: Probable decaprenylphosphoryl-beta-D-ribose oxidase

Chain A:



• Molecule 1: Probable decaprenylphosphoryl-beta-D-ribose oxidase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	76.78Å 82.88Å 80.54Å 90.00° 103.00° 90.00°	Depositor
Resolution (Å)	42.10 – 2.49 42.14 – 2.49	Depositor EDS
% Data completeness (in resolution range)	98.5 (42.10-2.49) 98.5 (42.14-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.59 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.221 , 0.265 0.221 , 0.265	Depositor DCC
R_{free} test set	1728 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	40.3	Xtriage
Anisotropy	0.064	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 33.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6907	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 38C, 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	0.46	1/3529 (0.0%)	0.52	0/4797
1	B	0.45	3/3291 (0.1%)	0.52	0/4466
All	All	0.46	4/6820 (0.1%)	0.52	0/9263

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	230	TRP	CD2-CE2	5.15	1.47	1.41
1	B	296	TRP	CD2-CE2	5.14	1.47	1.41
1	A	296	TRP	CD2-CE2	5.12	1.47	1.41
1	B	16	TRP	CD2-CE2	5.03	1.47	1.41

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3442	0	3415	44	0
1	B	3218	0	3193	51	0
2	A	53	0	31	0	0
2	B	53	0	31	1	0
3	A	27	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	27	0	13	1	0
4	A	51	0	0	2	0
4	B	36	0	0	2	0
All	All	6907	0	6696	95	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 7.

All (95) 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:B:159:ARG:HG2	1:B:159:ARG:HH11	1.16	1.08
1:B:18:ARG:HG3	1:B:18:ARG:HH11	1.34	0.91
1:A:269:ALA:N	1:A:270:PRO:HD3	1.93	0.83
1:A:274:THR:HG22	1:A:276:PRO:HD2	1.66	0.77
1:B:236:ALA:HA	1:B:355:GLN:HE22	1.51	0.73
1:B:93:ASP:O	1:B:96:VAL:HG13	1.89	0.72
1:A:263:GLU:HB3	1:A:266:LYS:HB3	1.73	0.71
1:A:273:LEU:O	1:A:273:LEU:HG	1.90	0.70
1:B:159:ARG:HG2	1:B:159:ARG:NH1	1.93	0.69
1:B:236:ALA:HA	1:B:355:GLN:NE2	2.08	0.67
1:B:78:ASN:HD22	1:B:78:ASN:H	1.41	0.66
1:A:119:ARG:HE	1:A:310:LEU:HD21	1.61	0.66
1:A:221:GLU:HG2	4:A:637:HOH:O	1.99	0.61
1:A:99:ASP:HB2	1:A:119:ARG:HB3	1.80	0.61
1:B:240:LEU:HD22	1:B:355:GLN:HG2	1.81	0.61
1:A:83:ILE:HG12	1:A:90:VAL:HG12	1.83	0.60
1:A:269:ALA:HB1	1:A:311:THR:CG2	2.31	0.60
1:A:93:ASP:O	1:A:96:VAL:HG13	2.02	0.59
1:A:269:ALA:N	1:A:270:PRO:CD	2.65	0.58
1:B:12:ARG:HH12	1:B:18:ARG:HH12	1.52	0.57
1:B:83:ILE:HG12	1:B:90:VAL:HG12	1.86	0.57
1:B:18:ARG:HG3	1:B:18:ARG:NH1	2.08	0.57
1:A:266:LYS:HE3	1:A:268:ASP:HB3	1.88	0.56
1:A:116:PRO:HG2	1:A:118:THR:O	2.06	0.56
1:A:270:PRO:O	1:A:271:GLN:HG3	2.06	0.55
1:B:102:MET:HG3	4:B:623:HOH:O	2.07	0.55
1:A:375:ALA:HB3	1:A:378:SER:HB2	1.87	0.55
1:B:221:GLU:HG2	4:B:614:HOH:O	2.07	0.54
1:B:99:ASP:HB2	1:B:119:ARG:HB3	1.90	0.54
1:B:184:ILE:HD12	2:B:501:FAD:C2A	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:THR:O	1:B:242:ARG:HB2	2.07	0.54
1:B:256:LEU:HD23	1:B:260:LEU:HB3	1.89	0.53
1:A:206:THR:HB	1:A:211:GLU:HB3	1.91	0.53
1:B:62:ASP:HB2	1:B:453:ASP:HB2	1.90	0.53
1:A:308:GLN:HB3	1:A:312:GLN:HG3	1.89	0.53
1:A:345:ASP:HA	1:A:348:LYS:HD3	1.89	0.53
1:B:75:THR:N	1:B:76:PRO:HD2	2.25	0.52
1:B:308:GLN:HB3	1:B:312:GLN:HG3	1.91	0.52
1:A:268:ASP:C	1:A:270:PRO:HD3	2.30	0.52
1:B:441:ARG:HH12	1:B:445:ASP:HB3	1.75	0.51
1:A:61:GLY:HA3	1:A:419:ASP:O	2.10	0.51
1:B:355:GLN:HG3	1:B:356:ALA:N	2.26	0.50
1:A:99:ASP:OD2	1:A:119:ARG:NH2	2.44	0.49
1:B:119:ARG:HE	1:B:315:HIS:HE1	1.60	0.49
1:B:60:TYR:HB2	1:B:418:LYS:HA	1.94	0.49
1:A:75:THR:N	1:A:76:PRO:HD2	2.28	0.49
1:B:441:ARG:NH1	1:B:445:ASP:HB3	2.27	0.49
1:A:448:ARG:HG2	1:A:461:LEU:HG	1.95	0.49
1:B:166:GLU:O	1:B:167:ASP:HB2	2.12	0.48
1:A:137:HIS:HA	1:A:370:GLY:HA3	1.97	0.47
1:B:92:ILE:HB	1:B:96:VAL:HG11	1.96	0.47
1:B:230:TRP:HB2	3:B:502:38C:H8	1.97	0.46
1:B:229:ALA:HB3	1:B:366:PHE:HB3	1.97	0.46
1:A:210:ASP:OD1	1:A:348:LYS:NZ	2.34	0.46
1:A:347:PHE:O	1:A:351:ILE:HG12	2.16	0.46
1:B:80:ILE:HG23	1:B:92:ILE:HG22	1.98	0.45
1:B:441:ARG:HH21	1:B:461:LEU:HD23	1.82	0.45
1:B:200:ILE:HG23	1:B:305:GLY:HA2	1.98	0.45
1:A:355:GLN:HG3	4:A:634:HOH:O	2.15	0.45
1:B:315:HIS:HA	1:B:316:PRO:HA	1.75	0.45
1:A:115:LEU:O	1:A:314:TYR:OH	2.26	0.45
1:A:106:LEU:HD21	1:A:111:TRP:CZ3	2.52	0.45
1:B:54:ARG:HG3	1:B:64:ALA:HB1	1.98	0.44
1:B:235:SER:O	1:B:355:GLN:NE2	2.51	0.44
1:A:26:VAL:HA	1:A:71:VAL:O	2.17	0.44
1:B:209:LEU:HB2	1:B:241:GLY:HA3	1.99	0.44
1:B:245:VAL:HG12	1:B:247:ARG:HG3	1.99	0.44
1:B:375:ALA:HB3	1:B:378:SER:HB2	2.00	0.44
1:A:267:PHE:CZ	1:A:269:ALA:HB3	2.53	0.44
1:B:153:THR:HG21	1:B:159:ARG:HD3	2.00	0.44
1:B:63:ASN:HB3	1:B:454:MET:SD	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:ILE:CG2	1:B:305:GLY:HA2	2.48	0.43
1:A:106:LEU:HD21	1:A:111:TRP:HZ3	1.83	0.43
1:B:200:ILE:HD11	1:B:255:GLN:OE1	2.18	0.43
1:B:78:ASN:HD22	1:B:78:ASN:N	2.06	0.43
1:B:112:VAL:HA	1:B:113:PRO:HD3	1.86	0.43
1:A:231:PHE:CE2	1:A:364:ASN:HB3	2.53	0.42
1:A:46:GLY:HA2	1:A:447:LEU:HD13	2.02	0.42
1:B:182:GLY:HA2	1:B:450:PHE:CZ	2.54	0.42
1:A:390:PHE:HA	1:A:391:PRO:HD2	1.93	0.41
1:A:419:ASP:OD2	1:A:422:THR:HG22	2.21	0.41
1:A:392:ILE:HA	1:A:396:LEU:HD12	2.01	0.41
1:B:346:GLU:HG2	1:B:410:PHE:CD1	2.56	0.41
1:A:92:ILE:HB	1:A:96:VAL:HG11	2.03	0.41
1:A:269:ALA:HB1	1:A:311:THR:HG23	2.01	0.41
1:A:336:GLN:HA	1:A:386:ILE:O	2.21	0.41
1:A:209:LEU:HB2	1:A:241:GLY:HA3	2.03	0.40
1:A:335:TYR:HB3	1:A:388:VAL:HG23	2.02	0.40
1:A:113:PRO:HG3	1:A:191:MET:HA	2.03	0.40
1:B:159:ARG:HH11	1:B:159:ARG:CG	2.05	0.40
1:B:263:GLU:HB3	1:B:266:LYS:HB2	2.02	0.40
1:A:366:PHE:C	1:A:366:PHE:CD1	2.95	0.40
1:B:18:ARG:NH1	1:B:18:ARG:CG	2.77	0.40
1:B:116:PRO:HG2	1:B:118:THR:O	2.22	0.40
1:B:151:LEU:HD22	1:B:161:LEU:HD11	2.03	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	445/480 (93%)	436 (98%)	9 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	413/480 (86%)	407 (98%)	6 (2%)	0	100	100
All	All	858/960 (89%)	843 (98%)	15 (2%)	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	357/384 (93%)	334 (94%)	23 (6%)	17	33
1	B	335/384 (87%)	309 (92%)	26 (8%)	12	24
All	All	692/768 (90%)	643 (93%)	49 (7%)	14	28

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	37	LYS
1	A	88	LYS
1	A	92	ILE
1	A	96	VAL
1	A	155	ASP
1	A	157	GLU
1	A	159	ARG
1	A	240	LEU
1	A	259	LYS
1	A	266	LYS
1	A	300	SER
1	A	312	GLN
1	A	327	TYR
1	A	334	GLN
1	A	344	VAL
1	A	355	GLN
1	A	367	LYS

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Mol	Chain	Res	Type
1	A	389	ASP
1	A	396	LEU
1	A	413	ARG
1	A	421	ARG
1	A	448	ARG
1	B	18	ARG
1	B	41	ARG
1	B	78	ASN
1	B	89	LEU
1	B	92	ILE
1	B	96	VAL
1	B	103	LYS
1	B	121	VAL
1	B	157	GLU
1	B	159	ARG
1	B	183	ILE
1	B	205	VAL
1	B	311	THR
1	B	314	TYR
1	B	315	HIS
1	B	344	VAL
1	B	355	GLN
1	B	386	ILE
1	B	396	LEU
1	B	409	GLU
1	B	413	ARG
1	B	421	ARG
1	B	425	GLU
1	B	441	ARG
1	B	443	LYS
1	B	448	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	355	GLN
1	B	78	ASN
1	B	315	HIS
1	B	355	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	FAD	A	501	-	53,58,58	1.34	5 (9%)	68,89,89	1.30	11 (16%)
3	38C	B	502	-	29,29,29	1.33	3 (10%)	41,42,42	1.32	5 (12%)
2	FAD	B	501	-	53,58,58	1.35	6 (11%)	68,89,89	1.26	9 (13%)
3	38C	A	502	-	29,29,29	1.31	2 (6%)	41,42,42	1.12	1 (2%)

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	A	501	-	-	2/30/50/50	0/6/6/6
3	38C	B	502	-	-	2/17/17/17	0/3/3/3
2	FAD	B	501	-	-	2/30/50/50	0/6/6/6
3	38C	A	502	-	-	6/17/17/17	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	FAD	C9A-C5X	5.55	1.50	1.41
2	B	501	FAD	C9A-C5X	5.51	1.50	1.41
3	B	502	38C	CAP-NAO	3.99	1.36	1.31
3	A	502	38C	CAP-NAO	3.70	1.36	1.31
2	B	501	FAD	C8-C7	3.60	1.49	1.40
2	A	501	FAD	C8-C7	3.58	1.49	1.40
3	B	502	38C	CAR-NAQ	-3.49	1.33	1.45
3	A	502	38C	CAR-NAQ	-3.44	1.33	1.45
2	A	501	FAD	C4X-N5	2.61	1.35	1.30
2	B	501	FAD	C5A-C4A	2.51	1.47	1.40
2	A	501	FAD	C5A-C4A	2.51	1.47	1.40
2	B	501	FAD	C4X-N5	2.50	1.35	1.30
2	A	501	FAD	C4-N3	-2.35	1.34	1.38
2	B	501	FAD	C4-N3	-2.13	1.34	1.38
3	B	502	38C	CAC-NAD	2.10	1.35	1.32
2	B	501	FAD	C10-N10	2.08	1.41	1.37

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	502	38C	CAS-CAR-NAQ	3.68	123.34	113.77
2	A	501	FAD	N3A-C2A-N1A	-3.64	122.99	128.68
2	B	501	FAD	N3A-C2A-N1A	-3.52	123.17	128.68
2	A	501	FAD	C4-C4X-N5	3.13	122.69	118.23
2	A	501	FAD	P-O3P-PA	-2.97	122.64	132.83
2	B	501	FAD	C4-C4X-N5	2.90	122.36	118.23
2	B	501	FAD	C4A-C5A-N7A	-2.79	106.49	109.40
2	B	501	FAD	P-O3P-PA	-2.63	123.80	132.83
2	B	501	FAD	C4X-C10-N1	-2.62	118.66	124.73
2	A	501	FAD	C4A-C5A-N7A	-2.61	106.68	109.40
2	A	501	FAD	C4X-C10-N1	-2.57	118.77	124.73
2	B	501	FAD	O4-C4-C4X	-2.53	119.89	126.60
3	B	502	38C	CAR-NAQ-CAP	2.51	126.44	123.14
2	A	501	FAD	C10-N1-C2	2.46	121.83	116.90
2	B	501	FAD	C10-N1-C2	2.43	121.77	116.90
3	A	502	38C	CAC-CAP-NAO	-2.38	119.31	121.63
2	A	501	FAD	O4-C4-C4X	-2.29	120.52	126.60
2	A	501	FAD	C4X-C4-N3	2.26	118.93	113.19
3	B	502	38C	CAC-CAP-NAQ	-2.25	117.90	121.69
2	A	501	FAD	C4X-C10-N10	2.20	119.69	116.48
2	B	501	FAD	O2-C2-N1	-2.19	118.20	121.83
3	B	502	38C	CAC-CAP-NAO	-2.18	119.51	121.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	FAD	C4X-C4-N3	2.17	118.70	113.19
2	A	501	FAD	O2-C2-N1	-2.04	118.45	121.83
2	A	501	FAD	C2A-N1A-C6A	2.01	122.19	118.75
3	B	502	38C	CAP-NAO-CAN	2.01	120.43	116.56

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	38C	CAC-CAP-NAQ-CAR
3	A	502	38C	CAY-CAV-OAW-CAX
3	A	502	38C	CAU-CAV-OAW-CAX
3	B	502	38C	CAY-CAV-OAW-CAX
3	B	502	38C	CAU-CAV-OAW-CAX
3	A	502	38C	NAO-CAP-NAQ-CAR
2	B	501	FAD	PA-O3P-P-O2P
2	A	501	FAD	PA-O3P-P-O2P
3	A	502	38C	OBA-CAB-CAC-NAD
3	A	502	38C	OAA-CAB-CAC-NAD
2	B	501	FAD	PA-O3P-P-O1P
2	A	501	FAD	PA-O3P-P-O1P

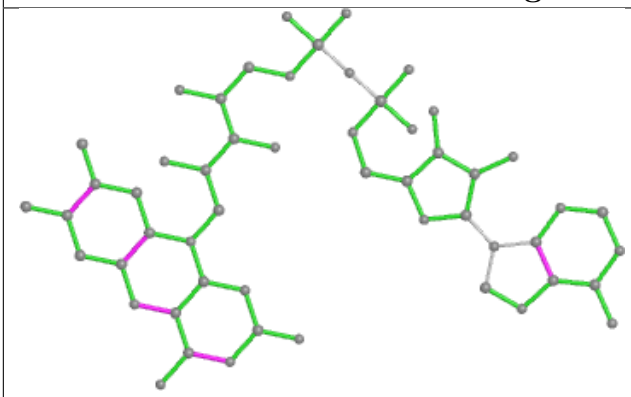
There are no ring outliers.

2 monomers are involved in 2 short contacts:

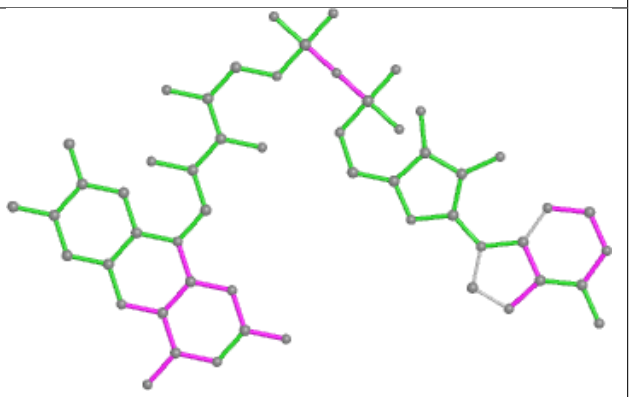
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	38C	1	0
2	B	501	FAD	1	0

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 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.

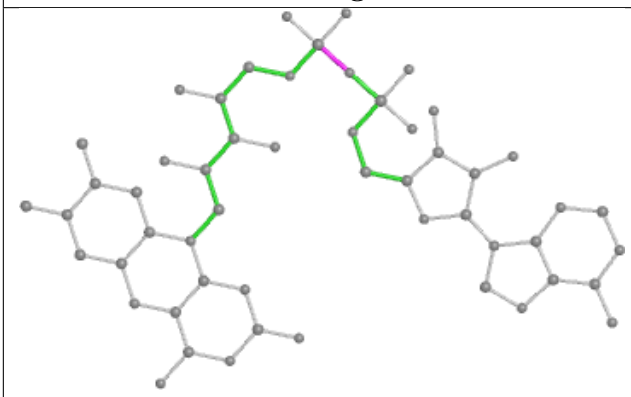
Ligand FAD A 501



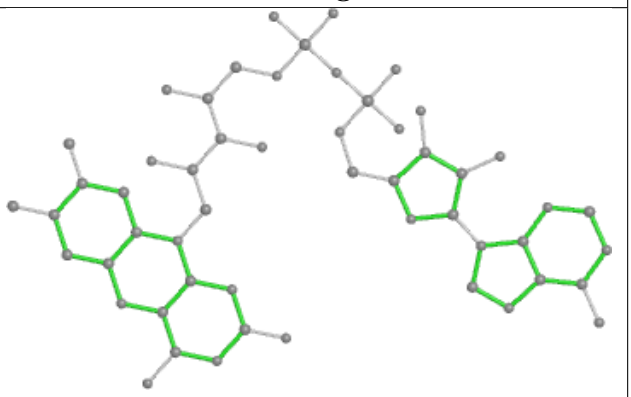
Bond lengths



Bond angles

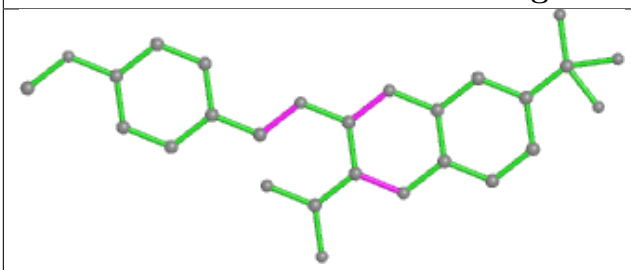


Torsions

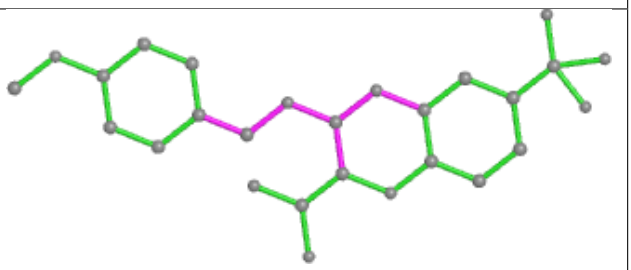


Rings

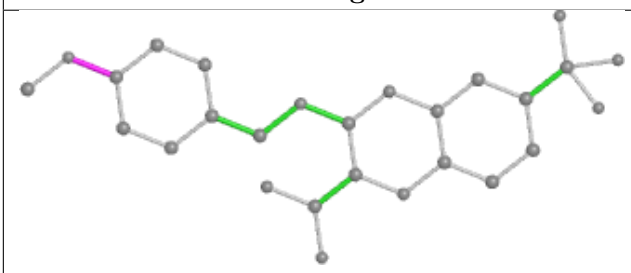
Ligand 38C B 502



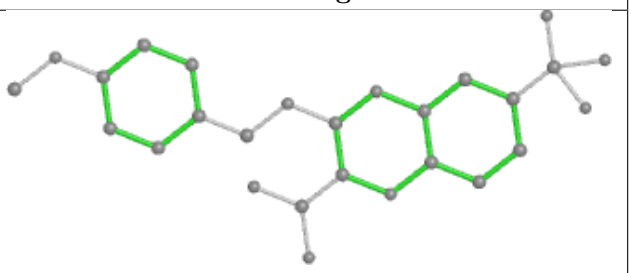
Bond lengths



Bond angles

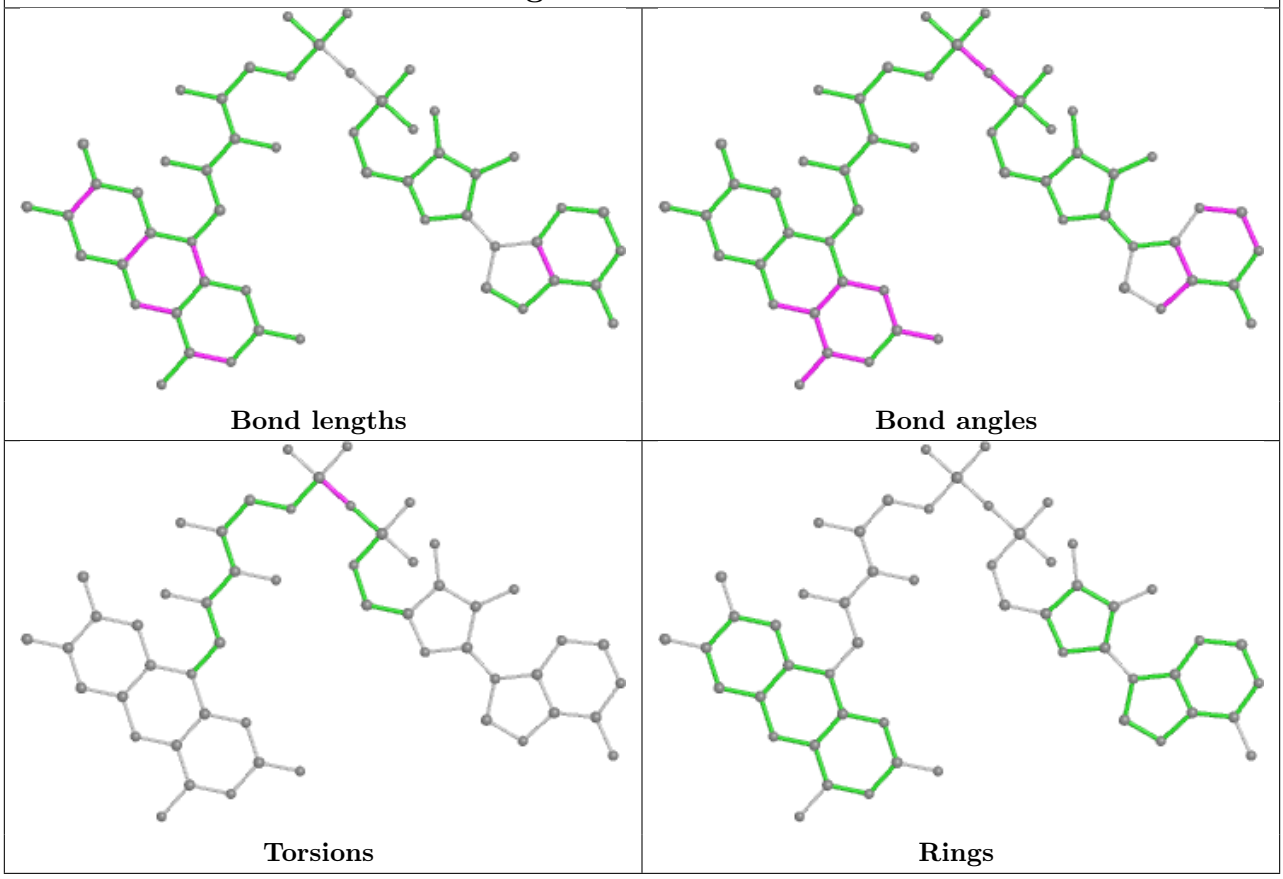


Torsions

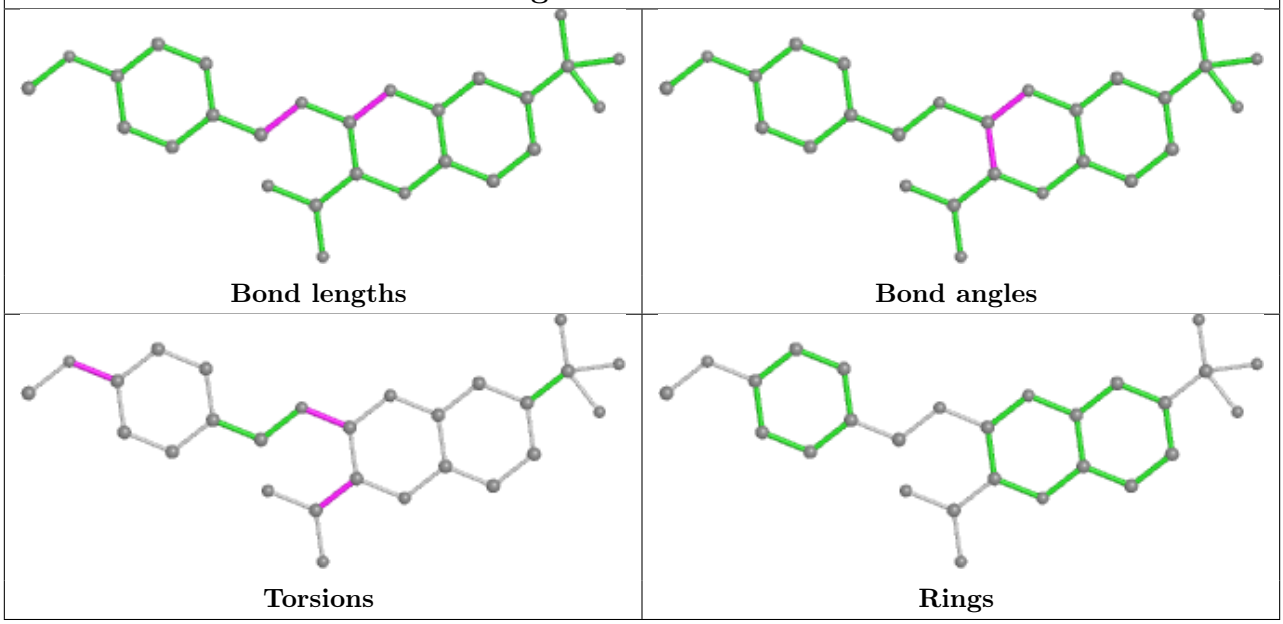


Rings

Ligand FAD B 501



Ligand 38C A 502



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	448/480 (93%)	0.12	25 (5%)	24	25	22, 36, 73, 94	0
1	B	419/480 (87%)	0.29	24 (5%)	23	25	28, 45, 71, 89	0
All	All	867/960 (90%)	0.20	49 (5%)	23	25	22, 41, 71, 94	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	330	ALA	5.4
1	A	269	ALA	5.4
1	B	7	THR	5.4
1	A	362	PHE	5.3
1	A	327	TYR	4.8
1	A	328	GLY	4.6
1	B	46	GLY	4.4
1	A	329	PRO	4.4
1	B	47	GLY	4.2
1	A	283	LEU	4.1
1	A	282	GLY	3.8
1	A	273	LEU	3.8
1	A	268	ASP	3.6
1	A	48	GLY	3.4
1	B	230	TRP	3.4
1	B	165	GLY	3.2
1	A	289	PHE	2.9
1	A	47	GLY	2.9
1	B	8	THR	2.8
1	A	12	ARG	2.8
1	A	275	LEU	2.8
1	B	34	MET	2.8
1	B	18	ARG	2.8
1	B	44	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	36	VAL	2.7
1	A	6	ALA	2.6
1	B	27	LEU	2.6
1	B	45	SER	2.4
1	B	259	LYS	2.4
1	B	35	ILE	2.4
1	B	10	ALA	2.3
1	A	281	ASN	2.3
1	A	5	GLY	2.3
1	A	280	PRO	2.3
1	B	70	LEU	2.3
1	B	447	LEU	2.2
1	B	43	ALA	2.2
1	A	297	TYR	2.2
1	B	39	VAL	2.2
1	A	7	THR	2.2
1	B	25	ASN	2.2
1	B	394	ASP	2.2
1	B	26	VAL	2.2
1	B	40	ALA	2.1
1	A	19	THR	2.1
1	A	23	VAL	2.1
1	B	23	VAL	2.1
1	A	267	PHE	2.0
1	A	276	PRO	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)

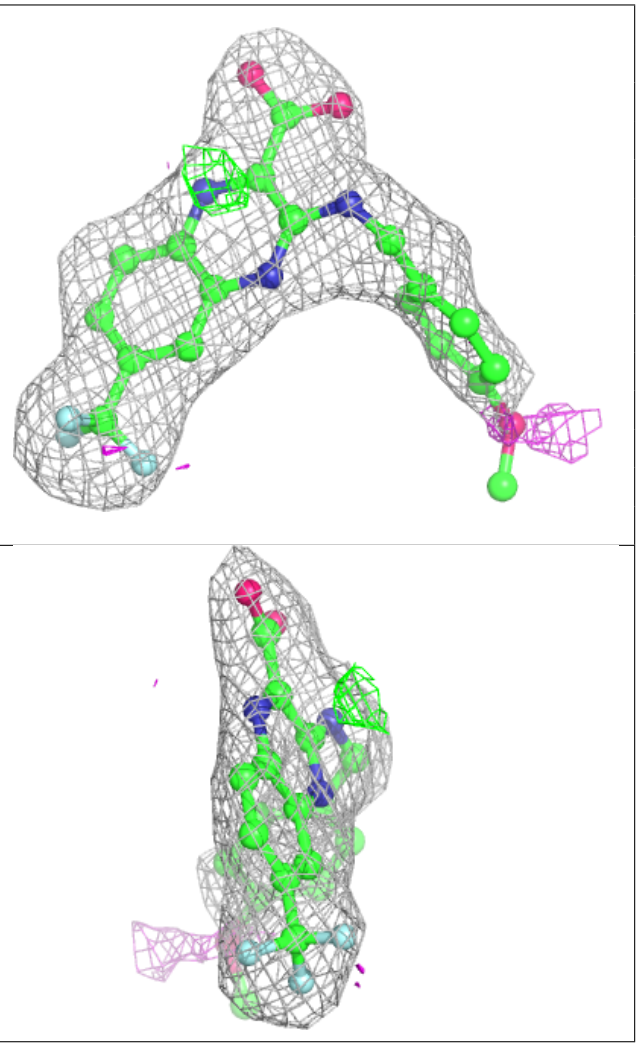
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	38C	B	502	27/27	0.89	0.24	50,52,64,65	0
3	38C	A	502	27/27	0.92	0.19	41,43,56,56	0
2	FAD	A	501	53/53	0.96	0.16	27,30,32,33	0
2	FAD	B	501	53/53	0.96	0.16	34,35,36,37	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.

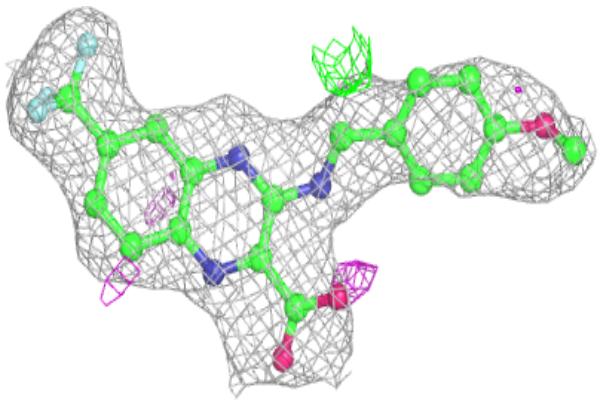
Electron density around 38C B 502:

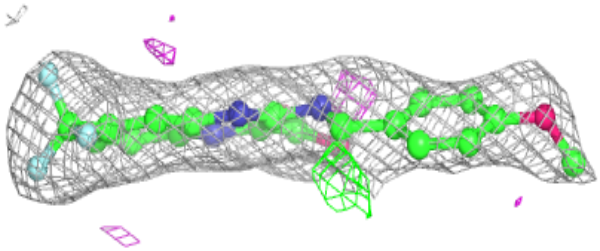
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

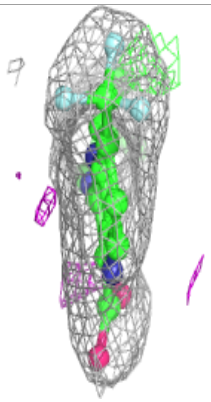


Electron density around 38C A 502:

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

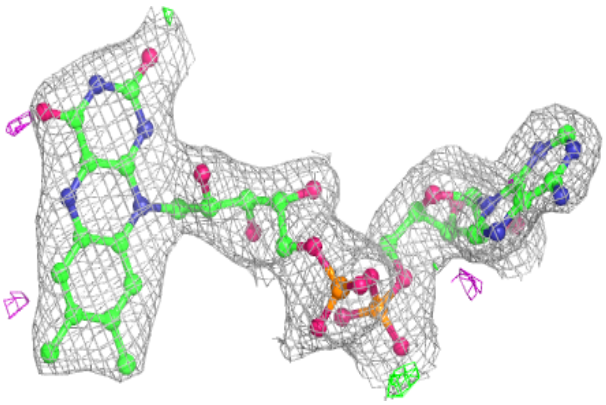


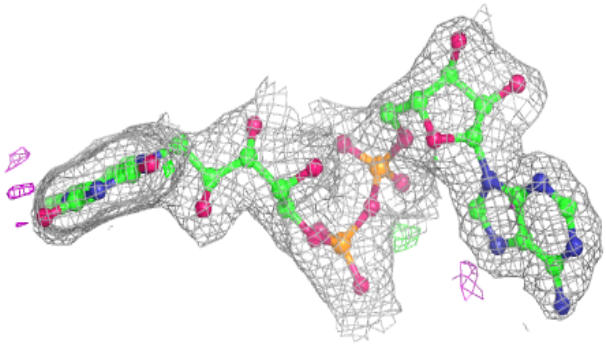


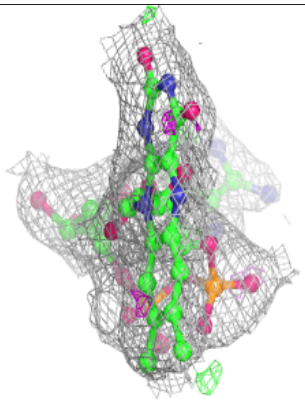


Electron density around FAD A 501:

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

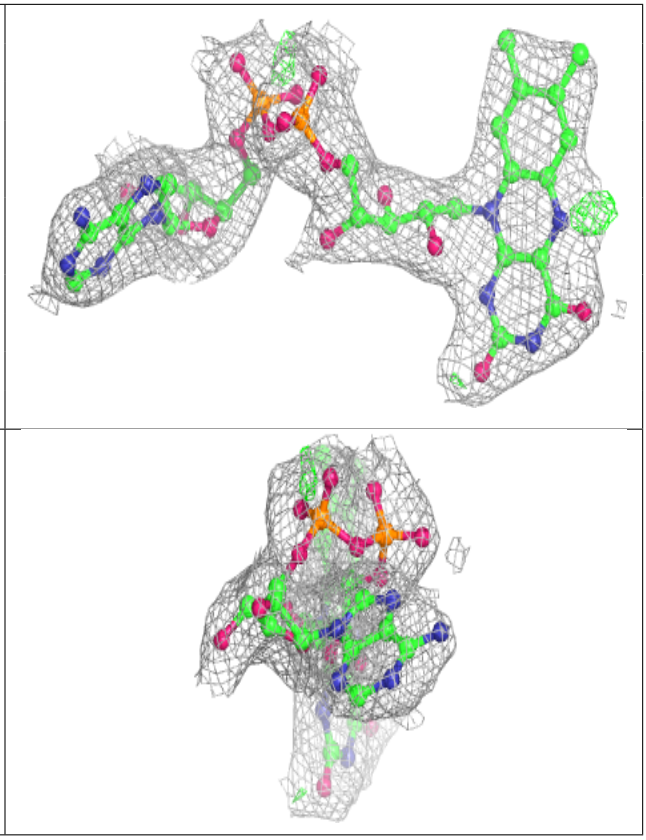






Electron density around FAD B 501:

$2mF_o$-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.