



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 4, 2024 – 10:24 pm GMT

PDB ID : 8PQN
Title : NQO1 bound to RBS-10
Authors : Pous, J.; Jose-Duran, F.; Mayor-Ruiz, C.; Riera, A.
Deposited on : 2023-07-11
Resolution : 3.80 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 : 4.02b-467
Mogul : 1.8.4, 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

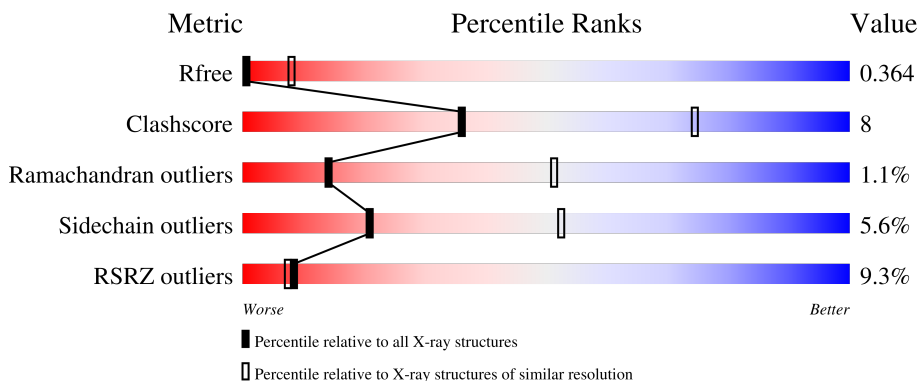
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

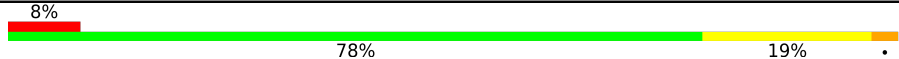

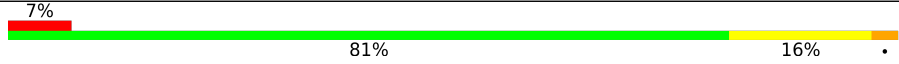
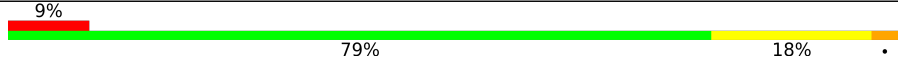
The reported resolution of this entry is 3.80 Å.

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	1212 (4.00-3.60)
Clashscore	141614	1288 (4.00-3.60)
Ramachandran outliers	138981	1243 (4.00-3.60)
Sidechain outliers	138945	1237 (4.00-3.60)
RSRZ outliers	127900	1121 (4.00-3.60)

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	273	
1	B	273	
1	C	273	
1	D	273	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8856 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 NAD(P)H dehydrogenase [quinone] 1.

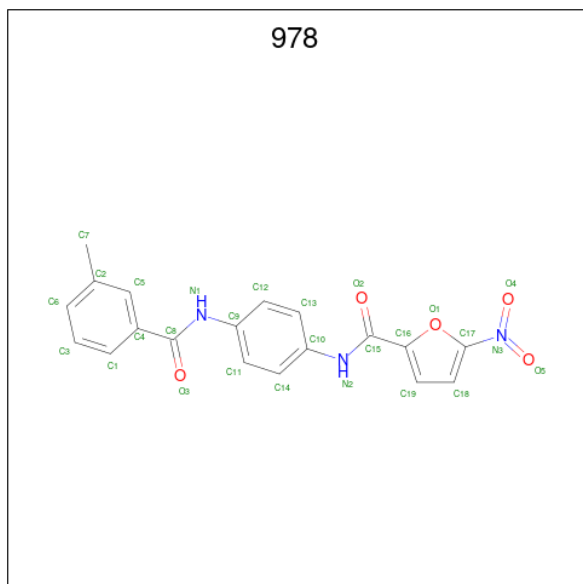
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	273	Total	C	N	O	S	0	0	0
			2175	1414	365	389	7			
1	B	273	Total	C	N	O	S	0	0	0
			2175	1414	365	389	7			
1	C	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			
1	D	273	Total	C	N	O	S	0	0	0
			2173	1412	365	389	7			

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

- Molecule 3 is {N}-[4-[(3-methylphenyl)carbonylamino]phenyl]-5-nitro-furan-2-carboxamide (three-letter code: 978) (formula: C₁₉H₁₅N₃O₅) (labeled as "Ligand of Interest" by depositor).

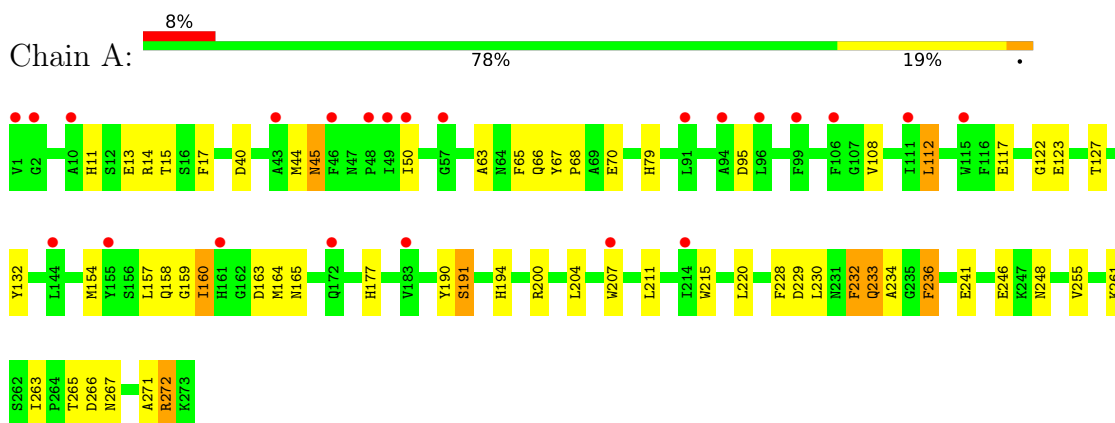


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	B	1	27	19	3	5	0	0
3	C	1	27	19	3	5	0	0

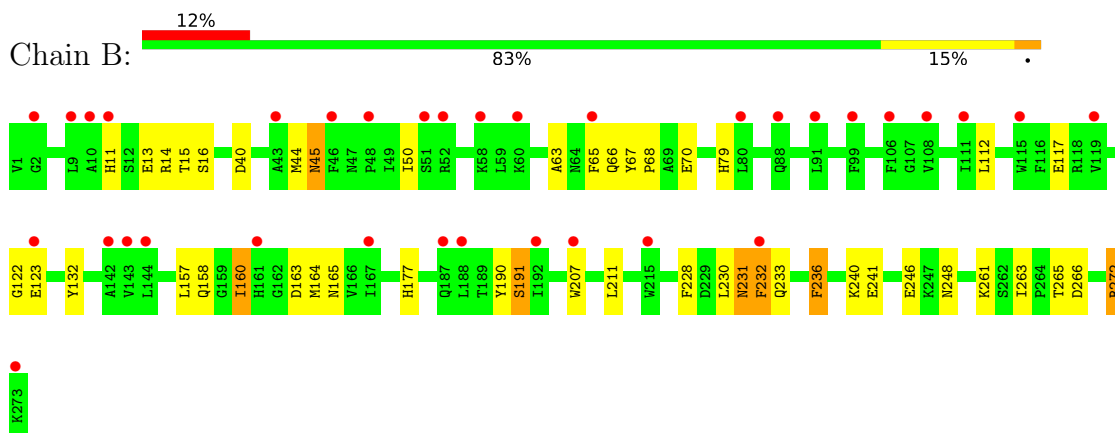
3 Residue-property plots [i](#)

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.

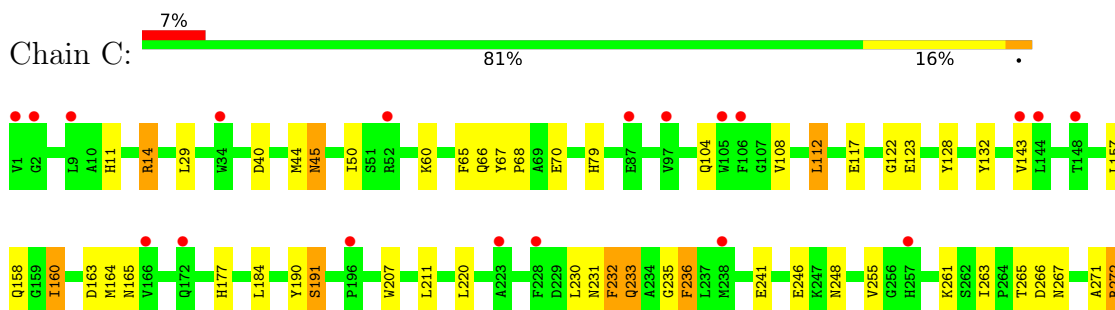
- Molecule 1: NAD(P)H dehydrogenase [quinone] 1



- Molecule 1: NAD(P)H dehydrogenase [quinone] 1



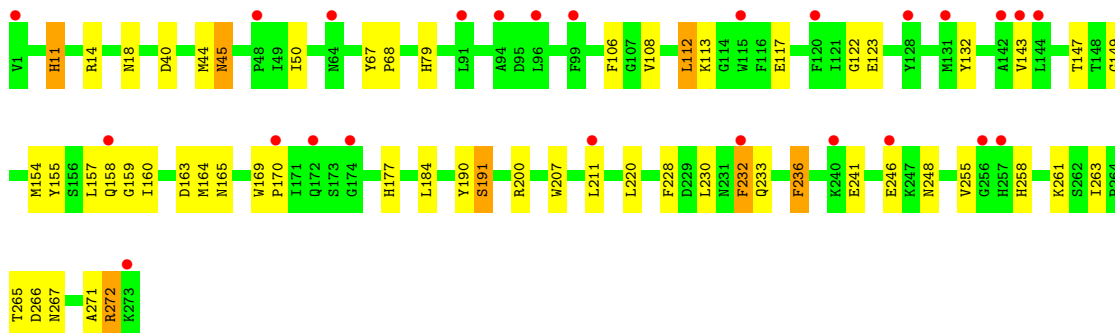
- Molecule 1: NAD(P)H dehydrogenase [quinone] 1



K273

● Molecule 1: NAD(P)H dehydrogenase [quinone] 1

Chain D: 9% 79% 18%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	70.75Å 178.37Å 210.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.13 – 3.80 68.04 – 3.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (68.13-3.80) 99.9 (68.04-3.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.11 (at 3.77Å)	Xtrriage
Refinement program	REFMAC 5.8.0405	Depositor
R, R_{free}	0.261 , 0.362 0.266 , 0.364	Depositor DCC
R_{free} test set	687 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	130.0	Xtrriage
Anisotropy	0.973	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 184.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8856	wwPDB-VP
Average B, all atoms (Å ²)	215.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 978, 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	0.27	0/2233	0.52	0/3015
1	B	0.26	0/2233	0.51	0/3015
1	C	0.26	0/2231	0.51	0/3012
1	D	0.26	0/2231	0.50	0/3012
All	All	0.26	0/8928	0.51	0/12054

There are no bond length outliers.

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	2175	0	2179	44	1
1	B	2175	0	2179	37	0
1	C	2173	0	2172	42	0
1	D	2173	0	2172	39	0
2	A	53	0	31	2	0
2	D	53	0	31	10	0
3	B	27	0	0	5	0
3	C	27	0	0	0	0
All	All	8856	0	8764	141	1

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 8.

All (141) 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)
3:B:301:978:O4	3:B:301:978:N3	1.57	1.34
1:D:155:TYR:OH	2:D:301:FAD:HM82	1.90	0.71
1:D:50:ILE:HD11	1:D:117:GLU:HB3	1.75	0.69
1:C:50:ILE:HD11	1:C:117:GLU:HB3	1.74	0.67
1:D:200:ARG:NH1	2:D:301:FAD:H1B	2.13	0.64
1:A:50:ILE:HD11	1:A:117:GLU:HB3	1.78	0.63
1:A:200:ARG:HD2	2:A:301:FAD:H8A	1.81	0.63
1:B:50:ILE:HD11	1:B:117:GLU:HB3	1.82	0.62
1:B:232:PHE:CD2	1:C:232:PHE:HB2	2.36	0.60
1:B:232:PHE:HB2	1:C:232:PHE:CD2	2.37	0.60
2:A:301:FAD:PA	1:B:66:GLN:HE22	2.25	0.60
1:D:232:PHE:CZ	1:D:236:PHE:HE2	2.22	0.58
1:C:232:PHE:CE2	1:C:236:PHE:HE2	2.21	0.58
1:A:228:PHE:CG	1:B:160:ILE:HG23	2.38	0.58
1:A:232:PHE:CE2	1:A:236:PHE:HE2	2.21	0.58
1:B:165:ASN:HD21	1:B:266:ASP:H	1.50	0.57
1:A:194:HIS:CE1	1:C:232:PHE:C	2.77	0.57
1:A:207:TRP:O	1:A:211:LEU:HG	2.05	0.57
1:C:263:ILE:HD11	1:D:263:ILE:HD11	1.86	0.57
1:A:65:PHE:HB3	1:B:13:GLU:CD	2.25	0.56
1:A:194:HIS:ND1	1:C:235:GLY:HA2	2.21	0.56
1:D:232:PHE:CE2	1:D:236:PHE:HE2	2.22	0.56
1:A:165:ASN:HD21	1:A:266:ASP:H	1.54	0.56
1:D:200:ARG:HH11	2:D:301:FAD:H1B	1.69	0.55
1:A:232:PHE:CZ	1:A:236:PHE:HE2	2.25	0.55
1:A:66:GLN:HE22	3:B:301:978:C1	2.20	0.55
1:B:232:PHE:HB2	1:C:232:PHE:CE2	2.41	0.55
1:B:232:PHE:CE2	1:B:236:PHE:HE2	2.24	0.55
1:D:232:PHE:CZ	1:D:236:PHE:CE2	2.95	0.55
1:C:232:PHE:CZ	1:C:236:PHE:HE2	2.24	0.54
1:D:165:ASN:HD21	1:D:266:ASP:H	1.54	0.54
1:A:44:MET:O	1:A:45:ASN:C	2.45	0.54
1:B:207:TRP:O	1:B:211:LEU:HG	2.08	0.54
1:A:13:GLU:CD	1:B:65:PHE:HB3	2.28	0.54
1:C:246:GLU:O	1:C:261:LYS:NZ	2.39	0.54
1:A:228:PHE:CD2	1:B:160:ILE:HG23	2.44	0.53
1:A:160:ILE:HG23	1:B:228:PHE:CG	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:ASN:HD21	1:C:266:ASP:H	1.56	0.53
1:C:232:PHE:CZ	1:C:236:PHE:CE2	2.96	0.53
1:B:44:MET:O	1:B:45:ASN:C	2.46	0.53
1:B:232:PHE:CD1	1:C:232:PHE:CD1	2.97	0.52
1:B:232:PHE:CZ	1:B:236:PHE:HE2	2.27	0.52
1:D:106:PHE:CE1	2:D:301:FAD:HM72	2.44	0.52
1:B:16:SER:HA	3:B:301:978:C3	2.40	0.52
1:D:246:GLU:O	1:D:261:LYS:NZ	2.43	0.52
1:A:232:PHE:CZ	1:A:236:PHE:CE2	2.98	0.51
1:C:44:MET:O	1:C:45:ASN:C	2.48	0.51
1:A:63:ALA:O	1:B:15:THR:HG21	2.10	0.51
3:B:301:978:O4	3:B:301:978:C17	2.49	0.51
1:D:44:MET:O	1:D:45:ASN:C	2.49	0.51
1:B:240:LYS:CD	1:C:60:LYS:HG2	2.41	0.50
1:C:272:ARG:N	1:C:272:ARG:HD3	2.26	0.50
1:B:230:LEU:HA	1:B:236:PHE:CE1	2.47	0.50
1:D:255:VAL:HG23	1:D:267:ASN:HD22	1.76	0.50
1:A:67:TYR:N	1:A:68:PRO:CD	2.75	0.50
1:D:207:TRP:O	1:D:211:LEU:HG	2.12	0.50
1:B:67:TYR:N	1:B:68:PRO:CD	2.74	0.49
1:B:231:ASN:HD22	1:C:128:TYR:HB2	1.78	0.49
1:A:160:ILE:HG23	1:B:228:PHE:CD2	2.48	0.49
1:B:65:PHE:CE1	1:B:70:GLU:HG3	2.47	0.49
1:B:232:PHE:CZ	1:B:236:PHE:CE2	3.00	0.49
1:A:15:THR:HG21	1:B:63:ALA:O	2.13	0.49
1:A:163:ASP:OD1	1:A:164:MET:N	2.46	0.49
1:C:67:TYR:N	1:C:68:PRO:CD	2.76	0.49
1:D:272:ARG:N	1:D:272:ARG:HD3	2.28	0.49
1:B:163:ASP:OD1	1:B:164:MET:N	2.46	0.48
1:D:230:LEU:HA	1:D:236:PHE:CE1	2.48	0.48
1:D:157:LEU:HD11	1:D:265:THR:HG21	1.94	0.48
1:C:255:VAL:HG23	1:C:267:ASN:HD22	1.78	0.48
1:A:246:GLU:O	1:A:261:LYS:NZ	2.44	0.48
1:D:147:THR:O	2:D:301:FAD:O4'	2.31	0.48
1:B:272:ARG:N	1:B:272:ARG:HD3	2.29	0.48
1:D:163:ASP:OD1	1:D:164:MET:N	2.45	0.47
1:A:132:TYR:HA	1:A:177:HIS:O	2.15	0.47
1:A:230:LEU:HA	1:A:236:PHE:CE1	2.50	0.47
1:A:272:ARG:N	1:A:272:ARG:HD3	2.30	0.47
1:B:246:GLU:O	1:B:261:LYS:NZ	2.44	0.47
1:C:66:GLN:NE2	2:D:301:FAD:O1A	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:67:TYR:N	1:C:68:PRO:HD2	2.30	0.47
1:A:66:GLN:HE22	3:B:301:978:C4	2.28	0.47
1:A:95:ASP:HB3	1:A:215:TRP:CZ3	2.50	0.47
1:C:65:PHE:CE1	1:C:70:GLU:HG3	2.50	0.47
1:D:67:TYR:N	1:D:68:PRO:CD	2.78	0.46
1:A:157:LEU:HD11	1:A:265:THR:HG21	1.98	0.46
1:B:67:TYR:N	1:B:68:PRO:HD2	2.31	0.46
1:B:165:ASN:ND2	1:B:266:ASP:H	2.12	0.46
1:A:255:VAL:HG23	1:A:267:ASN:HD22	1.81	0.46
1:C:163:ASP:OD1	1:C:164:MET:N	2.49	0.46
1:A:233:GLN:HG3	1:A:234:ALA:N	2.32	0.45
1:B:122:GLY:O	1:B:123:GLU:HB3	2.16	0.45
1:C:190:TYR:O	1:C:191:SER:C	2.55	0.45
1:D:67:TYR:N	1:D:68:PRO:HD2	2.32	0.45
1:C:157:LEU:HD11	1:C:265:THR:HG21	1.97	0.45
1:D:149:GLY:HA3	2:D:301:FAD:H9	1.98	0.45
1:A:67:TYR:N	1:A:68:PRO:HD2	2.31	0.45
1:C:207:TRP:O	1:C:211:LEU:HG	2.17	0.45
1:C:230:LEU:HA	1:C:236:PHE:CE1	2.51	0.45
1:B:157:LEU:HD11	1:B:265:THR:HG21	1.98	0.44
1:C:122:GLY:O	1:C:123:GLU:HB3	2.17	0.44
1:A:65:PHE:CE1	1:A:70:GLU:HG3	2.52	0.44
1:A:190:TYR:O	1:A:191:SER:C	2.55	0.44
1:C:160:ILE:HG23	1:D:228:PHE:CG	2.52	0.44
1:C:220:LEU:HD23	1:C:271:ALA:HB2	1.98	0.44
1:A:108:VAL:HG11	1:A:112:LEU:HD13	2.00	0.44
1:B:240:LYS:HD2	1:C:60:LYS:HG2	1.98	0.44
1:A:122:GLY:O	1:A:123:GLU:HB3	2.18	0.44
1:D:190:TYR:O	1:D:191:SER:C	2.56	0.44
1:C:14:ARG:H	1:C:14:ARG:HD2	1.81	0.43
1:D:18:ASN:ND2	2:D:301:FAD:O2P	2.51	0.43
1:D:122:GLY:O	1:D:123:GLU:HB3	2.17	0.43
1:B:132:TYR:HA	1:B:177:HIS:O	2.18	0.43
1:D:132:TYR:HA	1:D:177:HIS:O	2.18	0.43
1:A:263:ILE:HD11	1:B:263:ILE:HD11	2.00	0.43
1:D:154:MET:O	1:D:159:GLY:HA3	2.18	0.43
1:A:154:MET:O	1:A:159:GLY:HA3	2.19	0.43
1:D:232:PHE:N	1:D:232:PHE:CD1	2.87	0.43
1:D:165:ASN:ND2	1:D:266:ASP:H	2.15	0.43
1:A:17:PHE:CE1	1:A:204:LEU:HD21	2.54	0.43
1:A:95:ASP:HB3	1:A:215:TRP:CH2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LEU:HD23	1:A:271:ALA:HB2	2.01	0.42
1:C:132:TYR:HA	1:C:177:HIS:O	2.19	0.42
1:C:143:VAL:HG22	1:C:184:LEU:HB2	2.02	0.42
1:D:11:HIS:CD2	2:D:301:FAD:O2P	2.73	0.42
1:C:108:VAL:HG11	1:C:112:LEU:HD13	2.02	0.42
1:D:232:PHE:CE2	1:D:236:PHE:CE2	3.07	0.41
1:C:104:GLN:O	1:D:113:LYS:NZ	2.51	0.41
1:A:165:ASN:ND2	1:A:266:ASP:H	2.17	0.41
1:B:190:TYR:O	1:B:191:SER:C	2.58	0.41
1:C:29:LEU:CD1	1:C:211:LEU:HD13	2.51	0.41
1:D:220:LEU:HD23	1:D:271:ALA:HB2	2.01	0.41
1:D:143:VAL:HG22	1:D:184:LEU:HB2	2.03	0.41
1:D:200:ARG:HH11	2:D:301:FAD:C1B	2.34	0.41
1:A:229:ASP:O	1:A:236:PHE:HA	2.21	0.41
1:C:232:PHE:CE2	1:C:236:PHE:CE2	3.05	0.41
1:C:232:PHE:CD1	1:C:232:PHE:N	2.88	0.41
1:D:169:TRP:HB3	1:D:170:PRO:HD3	2.03	0.41
1:A:194:HIS:HE1	1:C:233:GLN:N	2.19	0.41
1:C:108:VAL:CG1	1:C:112:LEU:HB3	2.51	0.40
1:A:232:PHE:CD1	1:A:232:PHE:N	2.89	0.40
1:C:263:ILE:HD12	1:D:258:HIS:CG	2.56	0.40
1:D:108:VAL:CG1	1:D:112:LEU:HB3	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:ASN:ND2	1:A:45:ASN:ND2[4_555]	2.04	0.16

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	271/273 (99%)	238 (88%)	30 (11%)	3 (1%)	14	51
1	B	271/273 (99%)	236 (87%)	32 (12%)	3 (1%)	14	51
1	C	271/273 (99%)	237 (88%)	31 (11%)	3 (1%)	14	51
1	D	271/273 (99%)	237 (88%)	31 (11%)	3 (1%)	14	51
All	All	1084/1092 (99%)	948 (88%)	124 (11%)	12 (1%)	14	51

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	B	45	ASN
1	C	45	ASN
1	D	45	ASN
1	A	191	SER
1	B	191	SER
1	C	191	SER
1	D	191	SER
1	D	40	ASP
1	A	40	ASP
1	B	40	ASP
1	C	40	ASP

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	227/227 (100%)	214 (94%)	13 (6%)	20	52
1	B	227/227 (100%)	214 (94%)	13 (6%)	20	52
1	C	226/227 (100%)	213 (94%)	13 (6%)	20	51
1	D	226/227 (100%)	214 (95%)	12 (5%)	22	54
All	All	906/908 (100%)	855 (94%)	51 (6%)	21	52

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	HIS
1	A	14	ARG
1	A	79	HIS
1	A	112	LEU
1	A	127	THR
1	A	158	GLN
1	A	160	ILE
1	A	232	PHE
1	A	233	GLN
1	A	236	PHE
1	A	241	GLU
1	A	248	ASN
1	A	272	ARG
1	B	11	HIS
1	B	14	ARG
1	B	79	HIS
1	B	112	LEU
1	B	158	GLN
1	B	160	ILE
1	B	231	ASN
1	B	232	PHE
1	B	233	GLN
1	B	236	PHE
1	B	241	GLU
1	B	248	ASN
1	B	272	ARG
1	C	11	HIS
1	C	14	ARG
1	C	79	HIS
1	C	112	LEU
1	C	158	GLN
1	C	160	ILE
1	C	231	ASN
1	C	232	PHE
1	C	233	GLN
1	C	236	PHE
1	C	241	GLU
1	C	248	ASN
1	C	272	ARG
1	D	11	HIS
1	D	14	ARG
1	D	79	HIS
1	D	112	LEU

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Mol	Chain	Res	Type
1	D	158	GLN
1	D	160	ILE
1	D	232	PHE
1	D	233	GLN
1	D	236	PHE
1	D	241	GLU
1	D	248	ASN
1	D	272	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	104	GLN
1	A	213	ASN
1	A	267	ASN
1	A	268	GLN
1	B	64	ASN
1	B	66	GLN
1	B	104	GLN
1	B	213	ASN
1	B	231	ASN
1	B	267	ASN
1	B	268	GLN
1	C	104	GLN
1	C	213	ASN
1	C	231	ASN
1	C	267	ASN
1	C	268	GLN
1	D	11	HIS
1	D	66	GLN
1	D	104	GLN
1	D	267	ASN
1	D	268	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	D	301	-	53,58,58	0.61	0	68,89,89	0.72	2 (2%)
3	978	B	301	-	25,29,29	4.84	8 (32%)	32,40,40	1.70	5 (15%)
2	FAD	A	301	-	53,58,58	0.58	0	68,89,89	0.75	2 (2%)
3	978	C	301	-	25,29,29	4.44	8 (32%)	32,40,40	1.84	6 (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	FAD	D	301	-	-	6/30/50/50	0/6/6/6
3	978	B	301	-	-	11/15/20/20	0/3/3/3
2	FAD	A	301	-	-	19/30/50/50	0/6/6/6
3	978	C	301	-	-	5/15/20/20	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	301	978	O4-N3	14.74	1.57	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	301	978	O5-N3	12.29	1.52	1.22
3	B	301	978	O5-N3	11.83	1.50	1.22
3	C	301	978	C16-C15	-11.16	1.32	1.49
3	B	301	978	C16-C15	-10.30	1.34	1.49
3	C	301	978	O4-N3	9.69	1.45	1.22
3	B	301	978	C17-N3	-8.04	1.32	1.46
3	C	301	978	C17-N3	-7.60	1.33	1.46
3	C	301	978	C4-C8	-4.90	1.40	1.50
3	B	301	978	C4-C8	-4.53	1.40	1.50
3	C	301	978	C10-N2	-3.85	1.33	1.41
3	B	301	978	C9-N1	-3.36	1.34	1.41
3	C	301	978	C9-N1	-3.33	1.34	1.41
3	B	301	978	C10-N2	-3.20	1.35	1.41
3	C	301	978	C7-C2	-3.16	1.39	1.51
3	B	301	978	C7-C2	-2.84	1.41	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	301	978	C10-N2-C15	-5.80	111.52	126.58
3	C	301	978	C10-N2-C15	-5.75	111.64	126.58
3	C	301	978	C9-N1-C8	-5.73	111.69	126.58
3	B	301	978	C16-C15-N2	4.82	121.29	113.89
3	C	301	978	C16-C15-N2	4.21	120.36	113.89
3	B	301	978	C18-C17-N3	-3.67	122.55	127.43
2	D	301	FAD	C5A-C6A-N6A	2.31	123.86	120.35
3	B	301	978	C4-C8-N1	-2.25	110.97	115.92
2	A	301	FAD	C5A-C6A-N6A	2.24	123.76	120.35
3	C	301	978	O2-C15-C16	-2.22	116.25	121.08
3	C	301	978	O3-C8-C4	-2.19	117.03	120.94
3	C	301	978	C4-C8-N1	2.16	120.67	115.92
2	D	301	FAD	C4-N3-C2	-2.15	121.67	125.64
3	B	301	978	O2-C15-C16	-2.11	116.48	121.08
2	A	301	FAD	C4-N3-C2	-2.10	121.76	125.64

There are no chirality outliers.

All (41) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	FAD	C5B-O5B-PA-O2A
2	A	301	FAD	O4B-C4B-C5B-O5B
2	A	301	FAD	C1'-C2'-C3'-O3'

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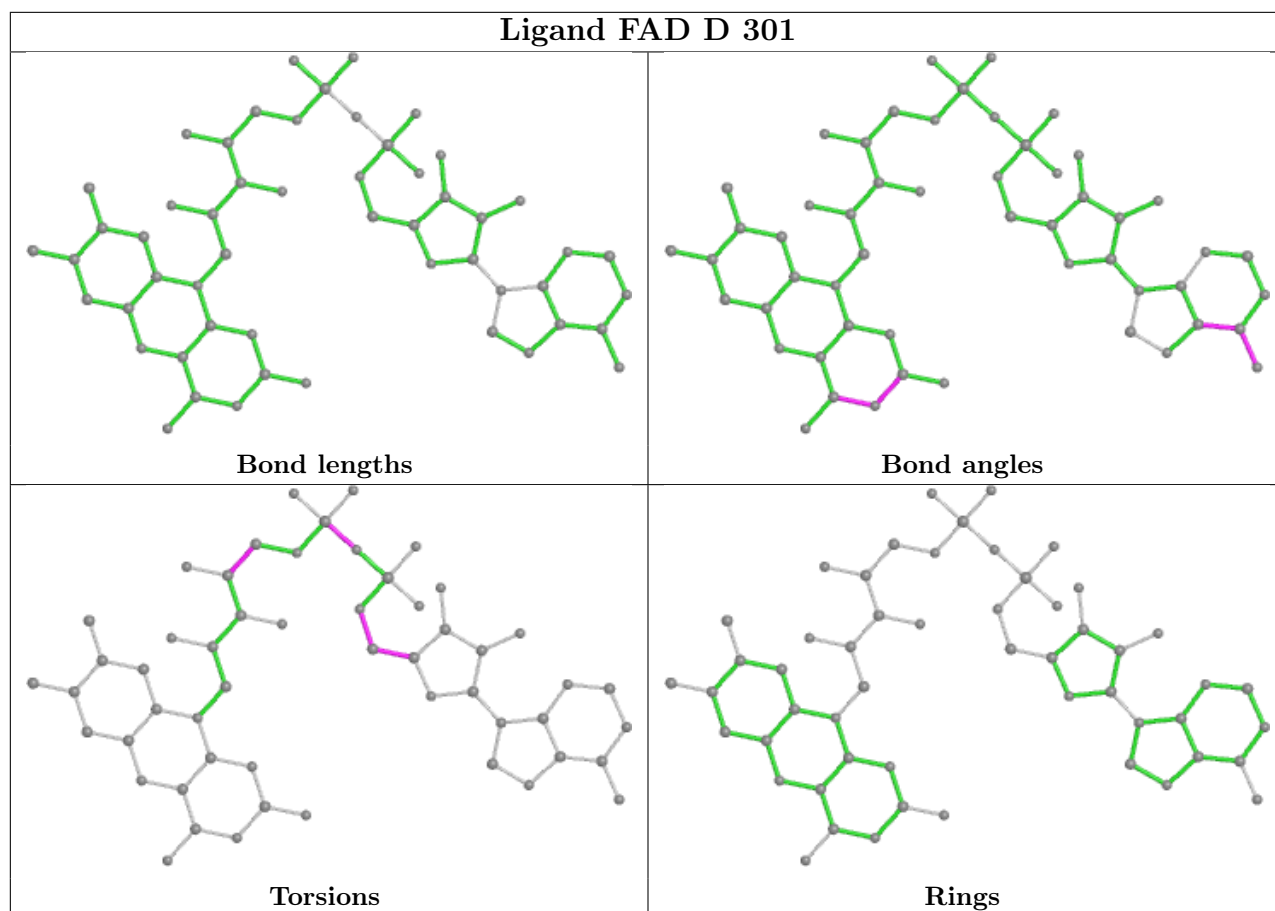
Mol	Chain	Res	Type	Atoms
2	A	301	FAD	C1'-C2'-C3'-C4'
2	A	301	FAD	C3'-C4'-C5'-O5'
2	A	301	FAD	O4'-C4'-C5'-O5'
2	A	301	FAD	C5'-O5'-P-O2P
2	D	301	FAD	C3B-C4B-C5B-O5B
2	D	301	FAD	C3'-C4'-C5'-O5'
2	D	301	FAD	O4'-C4'-C5'-O5'
3	B	301	978	O2-C15-C16-C19
3	C	301	978	O2-C15-C16-C19
3	B	301	978	C1-C4-C8-O3
3	B	301	978	C1-C4-C8-N1
3	B	301	978	C5-C4-C8-N1
3	B	301	978	C5-C4-C8-O3
2	A	301	FAD	O2'-C2'-C3'-O3'
2	A	301	FAD	O2'-C2'-C3'-C4'
3	C	301	978	C14-C10-N2-C15
3	C	301	978	C12-C9-N1-C8
3	C	301	978	C13-C10-N2-C15
3	C	301	978	C11-C9-N1-C8
3	B	301	978	C14-C10-N2-C15
3	B	301	978	C13-C10-N2-C15
3	B	301	978	C11-C9-N1-C8
2	D	301	FAD	O4B-C4B-C5B-O5B
3	B	301	978	C12-C9-N1-C8
2	A	301	FAD	C2'-C3'-C4'-O4'
2	A	301	FAD	C4'-C5'-O5'-P
2	D	301	FAD	C4B-C5B-O5B-PA
2	A	301	FAD	C5B-O5B-PA-O3P
2	A	301	FAD	C5'-O5'-P-O3P
3	B	301	978	C18-C17-N3-O4
3	B	301	978	C18-C17-N3-O5
2	A	301	FAD	C5B-O5B-PA-O1A
2	A	301	FAD	O3'-C3'-C4'-O4'
2	A	301	FAD	C2'-C3'-C4'-C5'
2	A	301	FAD	O3'-C3'-C4'-C5'
2	A	301	FAD	PA-O3P-P-O1P
2	D	301	FAD	PA-O3P-P-O2P
2	A	301	FAD	C5'-O5'-P-O1P

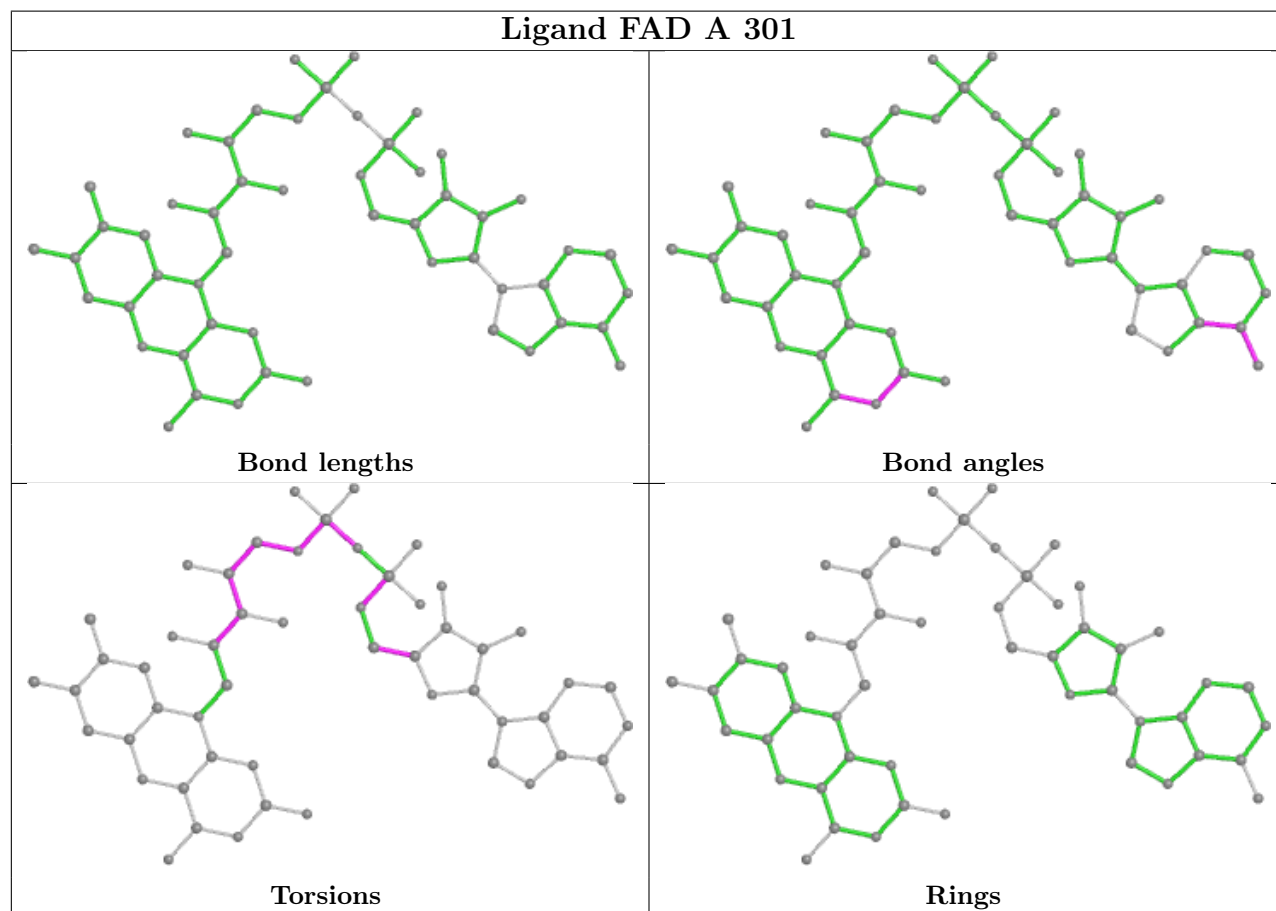
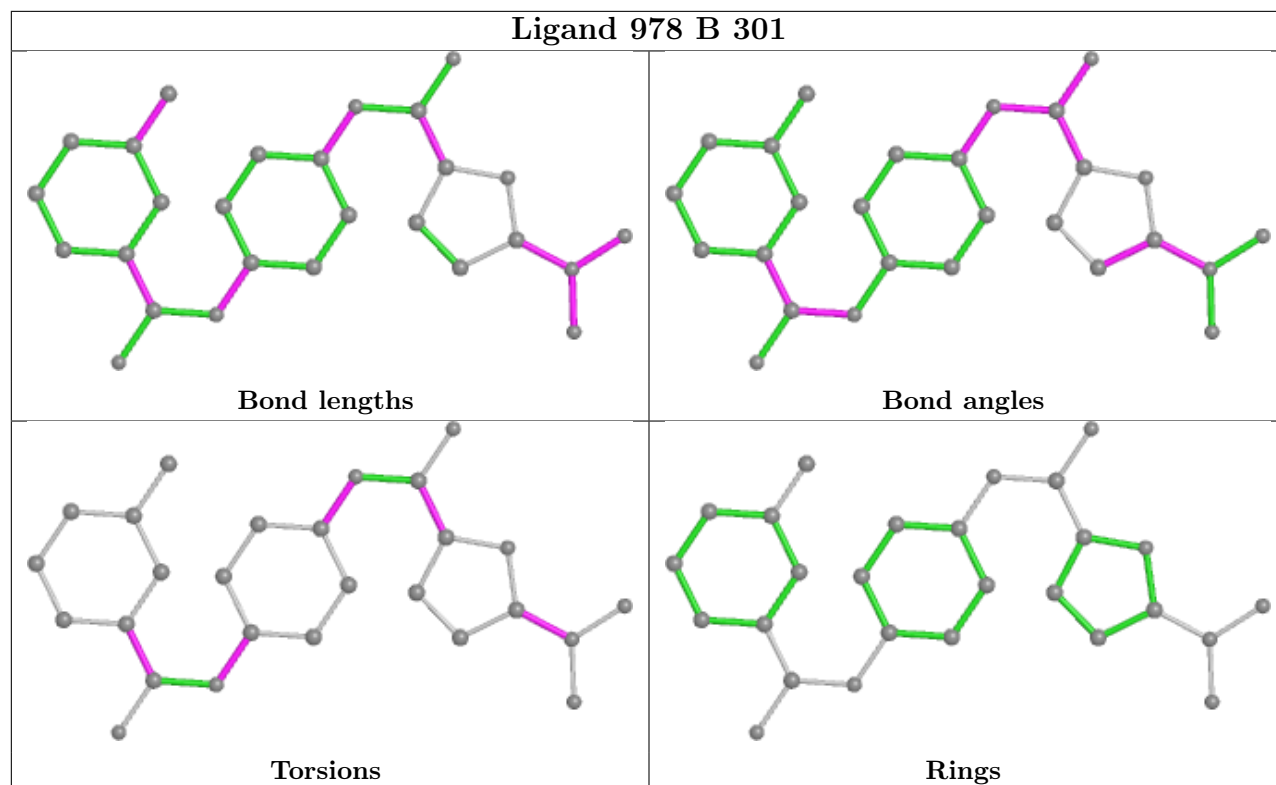
There are no ring outliers.

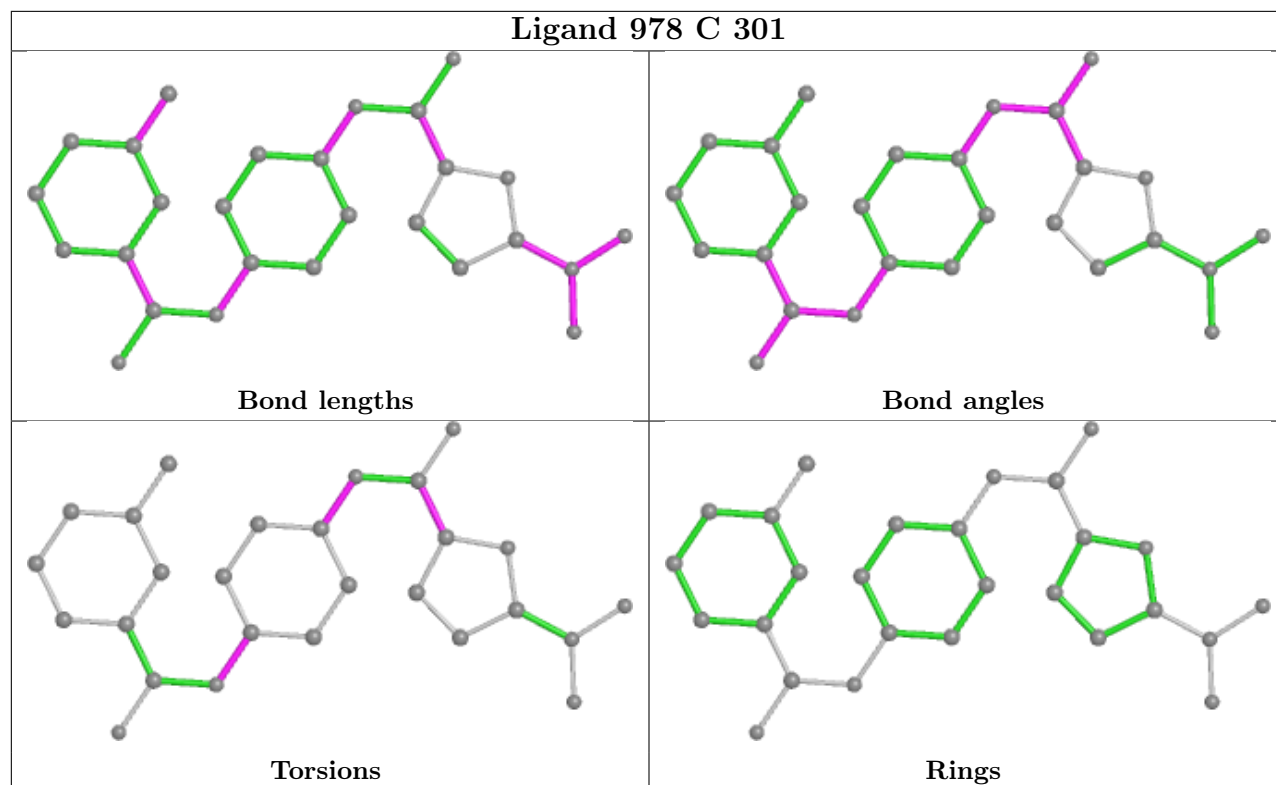
3 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	FAD	10	0
3	B	301	978	5	0
2	A	301	FAD	2	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	273/273 (100%)	0.60	23 (8%) 11 9	128, 202, 260, 314	0
1	B	273/273 (100%)	0.65	34 (12%) 3 4	130, 209, 265, 331	0
1	C	273/273 (100%)	0.52	20 (7%) 15 12	129, 214, 284, 314	0
1	D	273/273 (100%)	0.58	25 (9%) 9 7	131, 231, 293, 325	0
All	All	1092/1092 (100%)	0.59	102 (9%) 8 7	128, 214, 279, 331	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	142	ALA	5.0
1	A	1	VAL	4.7
1	A	43	ALA	4.7
1	B	10	ALA	4.5
1	C	1	VAL	4.4
1	A	94	ALA	4.2
1	D	1	VAL	4.2
1	A	48	PRO	4.2
1	A	50	ILE	4.1
1	A	49	ILE	3.9
1	C	238	MET	3.8
1	A	111	ILE	3.6
1	B	91	LEU	3.6
1	A	144	LEU	3.5
1	B	48	PRO	3.5
1	B	207	TRP	3.5
1	B	9	LEU	3.4
1	C	87	GLU	3.4
1	D	158	GLN	3.4
1	C	2	GLY	3.3
1	C	9	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	207	TRP	3.3
1	D	144	LEU	3.2
1	D	96	LEU	3.2
1	D	170	PRO	3.2
1	B	273	LYS	3.2
1	B	11	HIS	3.2
1	C	223	ALA	3.1
1	D	172	GLN	3.1
1	A	2	GLY	3.1
1	D	143	VAL	3.0
1	B	2	GLY	3.0
1	D	91	LEU	3.0
1	A	106	PHE	3.0
1	C	257	HIS	3.0
1	B	115	TRP	2.9
1	A	46	PHE	2.9
1	B	187	GLN	2.9
1	B	99	PHE	2.9
1	B	232	PHE	2.9
1	D	48	PRO	2.9
1	D	64	ASN	2.9
1	D	240	LYS	2.8
1	B	188	LEU	2.8
1	B	88	GLN	2.8
1	A	214	ILE	2.8
1	A	161	HIS	2.8
1	A	155	TYR	2.7
1	A	172	GLN	2.7
1	C	196	PRO	2.7
1	B	60	LYS	2.7
1	B	144	LEU	2.6
1	B	215	TRP	2.6
1	C	52	ARG	2.6
1	D	128	TYR	2.6
1	D	94	ALA	2.6
1	B	51	SER	2.6
1	B	123	GLU	2.6
1	B	43	ALA	2.6
1	D	273	LYS	2.6
1	A	91	LEU	2.6
1	C	273	LYS	2.6
1	B	52	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	166	VAL	2.5
1	A	10	ALA	2.5
1	A	99	PHE	2.5
1	A	183	VAL	2.5
1	C	148	THR	2.4
1	D	115	TRP	2.4
1	D	246	GLU	2.4
1	C	105	TRP	2.4
1	B	108	VAL	2.3
1	B	65	PHE	2.3
1	D	120	PHE	2.3
1	B	46	PHE	2.3
1	B	106	PHE	2.3
1	B	167	ILE	2.2
1	C	106	PHE	2.2
1	B	161	HIS	2.2
1	A	57	GLY	2.2
1	B	111	ILE	2.2
1	A	115	TRP	2.2
1	C	144	LEU	2.2
1	B	143	VAL	2.2
1	C	97	VAL	2.2
1	D	174	GLY	2.2
1	B	192	ILE	2.2
1	D	257	HIS	2.2
1	A	96	LEU	2.2
1	D	232	PHE	2.1
1	C	34	TRP	2.1
1	C	228	PHE	2.1
1	C	172	GLN	2.1
1	D	131	MET	2.1
1	D	99	PHE	2.1
1	D	142	ALA	2.1
1	D	211	LEU	2.1
1	C	143	VAL	2.0
1	D	256	GLY	2.0
1	B	80	LEU	2.0
1	B	58	LYS	2.0
1	B	119	VAL	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](#)

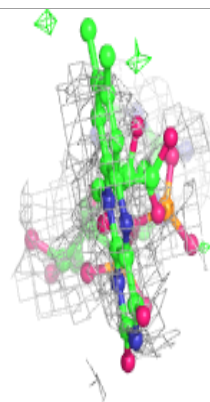
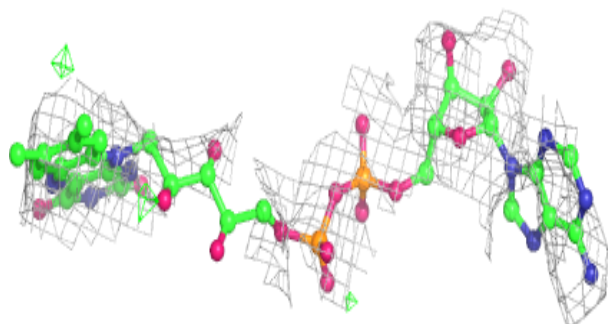
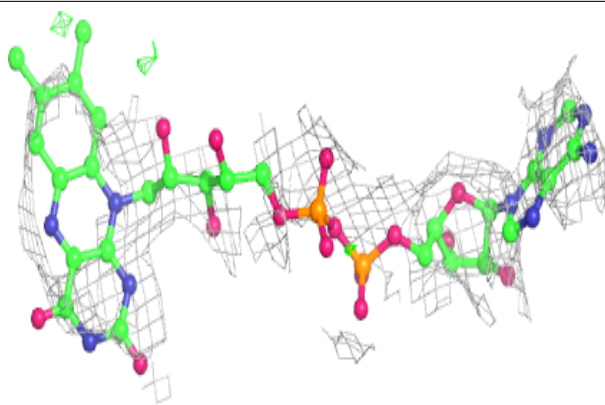
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FAD	D	301	53/53	0.69	0.33	169,251,290,293	0
3	978	B	301	27/27	0.74	0.38	151,198,234,248	0
2	FAD	A	301	53/53	0.78	0.39	168,235,289,305	0
3	978	C	301	27/27	0.79	0.23	160,219,242,243	0

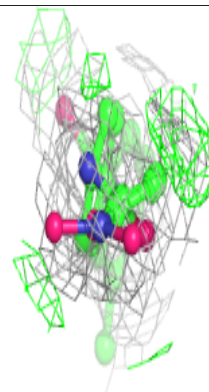
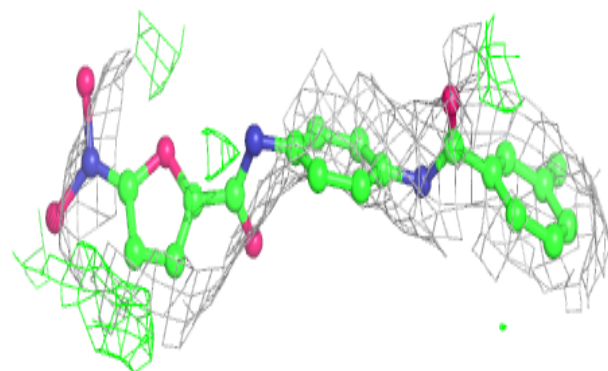
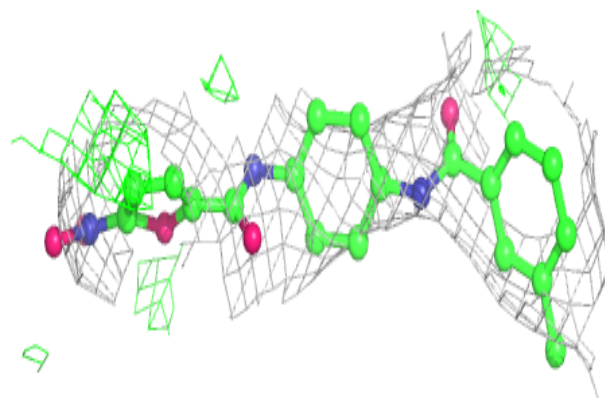
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD D 301:

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

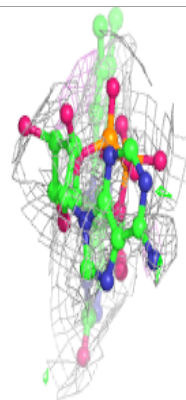
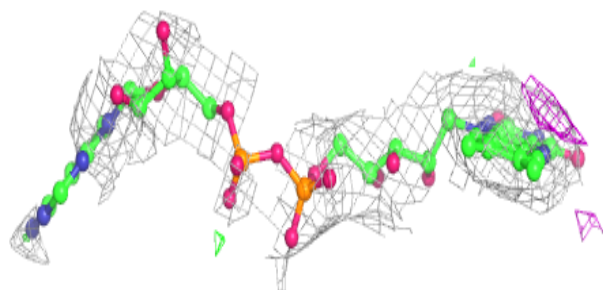
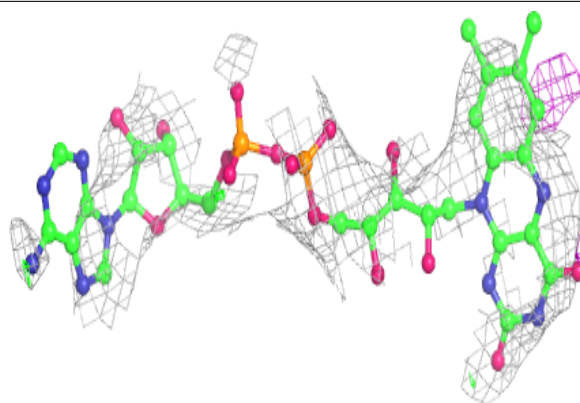
**Electron density around 978 B 301:**

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

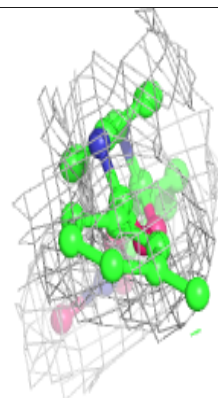
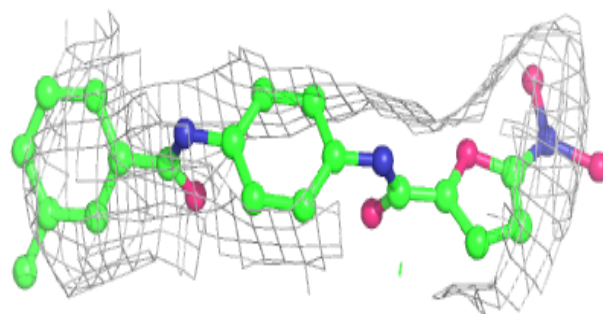
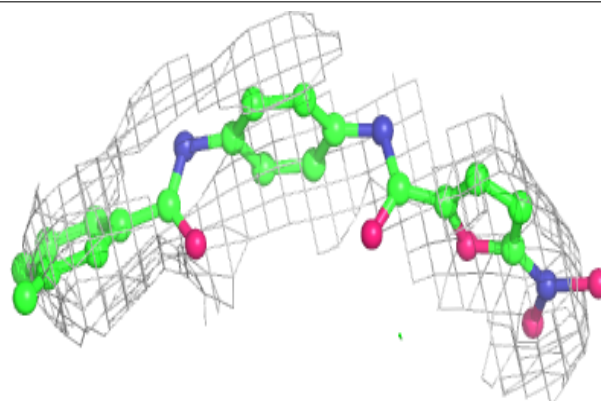


Electron density around FAD A 301:

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

**Electron density around 978 C 301:**

$2mF_o-DF_c$ (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.