



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2023 – 10:48 PM EDT

PDB ID : 3O73  
Title : Crystal structure of quinone reductase 2 in complex with the indolequinone MAC627  
Authors : Dufour, M.; Yan, C.; Colucci, M.A.; Siegel, D.; Li, Y.; De Matteis, C.I.; Ross, D.; Moody, C.J.  
Deposited on : 2010-07-30  
Resolution : 2.00 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](mailto: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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35  
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) : 2.35

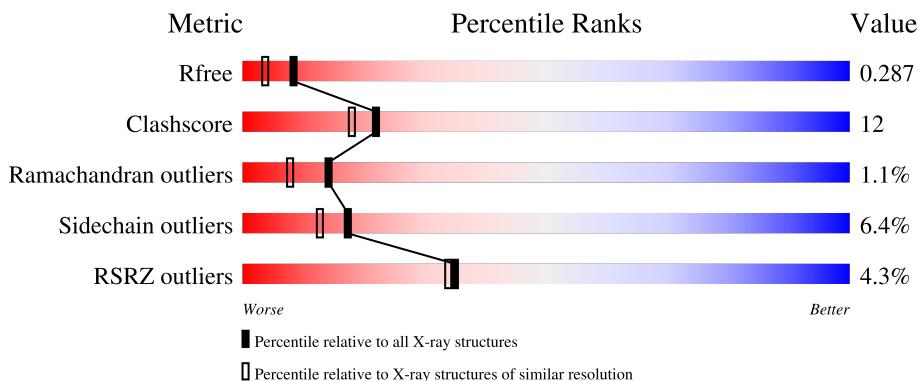
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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  $\geq 3$ , 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	237	
1	B	237	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	O73	A	239	-	-	X	-
4	O73	A	240	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3876 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 Ribosyldihyronicotinamide dehydrogenase [quinone].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	230	1823	1174	304	337	8	0	0	0
1	B	230	1823	1174	304	337	8	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	231	HIS	-	expression tag	UNP P16083
A	232	HIS	-	expression tag	UNP P16083
A	233	HIS	-	expression tag	UNP P16083
A	234	HIS	-	expression tag	UNP P16083
A	235	HIS	-	expression tag	UNP P16083
A	236	HIS	-	expression tag	UNP P16083
B	231	HIS	-	expression tag	UNP P16083
B	232	HIS	-	expression tag	UNP P16083
B	233	HIS	-	expression tag	UNP P16083
B	234	HIS	-	expression tag	UNP P16083
B	235	HIS	-	expression tag	UNP P16083
B	236	HIS	-	expression tag	UNP P16083

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

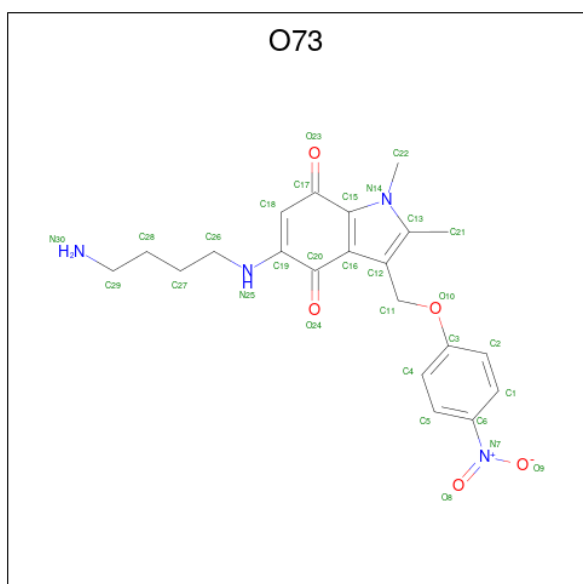
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C<sub>27</sub>H<sub>33</sub>N<sub>9</sub>O<sub>15</sub>P<sub>2</sub>).



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

- Molecule 4 is 5-[(4-aminobutyl)amino]-1,2-dimethyl-3-[(4-nitrophenoxy)methyl]-1H-indole-4,7-dione (three-letter code: O73) (formula: C<sub>21</sub>H<sub>24</sub>N<sub>4</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	30	21	4	5	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	30	21	4	5	0	0

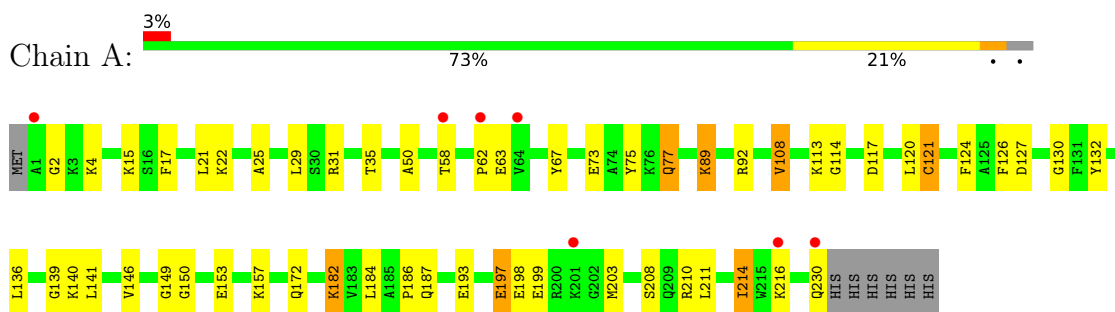
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	31	Total	O	0	0
			31	31		
5	B	31	Total	O	0	0
			31	31		

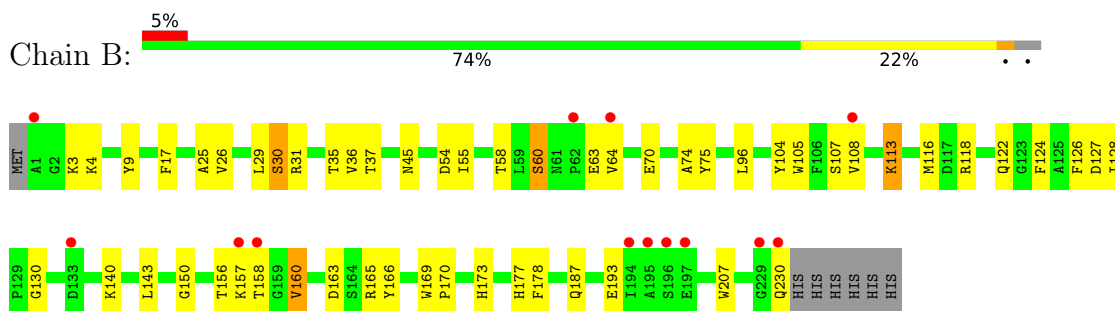
### 3 Residue-property plots [i](#)

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: Ribosyldihyronicotinamide dehydrogenase [quinone]



- Molecule 1: Ribosyldihyronicotinamide dehydrogenase [quinone]



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.08Å 82.05Å 106.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.45 – 2.00 27.56 – 2.00	Depositor EDS
% Data completeness (in resolution range)	81.1 (35.45-2.00) 81.1 (27.56-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.90 (at 1.99Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.220 , 0.288 0.227 , 0.287	Depositor DCC
$R_{free}$ test set	1406 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.1	Xtrriage
Anisotropy	0.238	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 44.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3876	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, FAD, O73

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	1.01	4/1873 (0.2%)	0.86	0/2542
1	B	1.03	3/1873 (0.2%)	0.94	1/2542 (0.0%)
All	All	1.02	7/3746 (0.2%)	0.90	1/5084 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	9	TYR	CZ-OH	5.97	1.48	1.37
1	A	121	CYS	CB-SG	-5.88	1.72	1.81
1	B	9	TYR	CD2-CE2	5.73	1.48	1.39
1	B	9	TYR	CD1-CE1	5.68	1.47	1.39
1	A	197	GLU	CG-CD	5.58	1.60	1.51
1	A	108	VAL	CB-CG2	-5.55	1.41	1.52
1	A	114	GLY	N-CA	5.30	1.54	1.46

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	143	LEU	CB-CG-CD1	-5.14	102.27	111.00

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	1823	0	1779	45	0
1	B	1823	0	1779	39	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	53	0	31	3	0
3	B	53	0	31	5	0
4	A	60	0	48	28	0
5	A	31	0	0	2	0
5	B	31	0	0	0	0
All	All	3876	0	3668	88	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 12.

All (88) 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:113:LYS:HA	1:B:116:MET:HE3	1.42	1.01
4:A:240:O73:H28A	3:B:238:FAD:HM72	1.40	0.99
1:A:31:ARG:HG2	1:A:31:ARG:HH11	1.29	0.95
1:B:108:VAL:HG21	1:B:116:MET:CE	2.01	0.90
1:A:17:PHE:HB2	3:A:238:FAD:H51A	1.53	0.90
1:B:108:VAL:HG21	1:B:116:MET:HE1	1.57	0.84
1:A:139:GLY:H	1:A:182:LYS:NZ	1.75	0.82
1:B:113:LYS:HA	1:B:116:MET:CE	2.09	0.81
4:A:240:O73:H21B	1:B:150:GLY:CA	2.12	0.80
1:A:121:CYS:HA	4:A:240:O73:H29	1.63	0.79
4:A:239:O73:H22	5:A:259:HOH:O	1.82	0.78
1:A:29:LEU:HD21	1:A:214:ILE:HD11	1.68	0.76
4:A:240:O73:H21B	1:B:150:GLY:N	2.01	0.74
1:A:182:LYS:N	1:A:182:LYS:HD3	2.02	0.73
1:B:113:LYS:CA	1:B:116:MET:HE3	2.17	0.73
1:A:139:GLY:H	1:A:182:LYS:HZ3	1.37	0.72
1:A:31:ARG:HG2	1:A:31:ARG:NH1	1.98	0.70
4:A:240:O73:H28A	3:B:238:FAD:C7M	2.18	0.69
1:A:150:GLY:CA	4:A:239:O73:H21B	2.24	0.68
1:B:17:PHE:HB2	3:B:238:FAD:H51A	1.76	0.68
1:A:193:GLU:HB3	4:A:239:O73:O9	1.94	0.67
1:A:92:ARG:O	1:A:140:LYS:NZ	2.28	0.66
4:A:240:O73:H5	1:B:193:GLU:HB2	1.76	0.66
1:B:108:VAL:HG21	1:B:116:MET:HE2	1.76	0.64
4:A:240:O73:H21B	1:B:150:GLY:HA2	1.80	0.64

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:187:GLN:HE21	1:A:210:ARG:HH11	1.47	0.63
1:A:187:GLN:NE2	1:A:210:ARG:HH11	1.98	0.61
1:B:26:VAL:O	1:B:30:SER:HB2	2.00	0.61
1:A:193:GLU:CB	4:A:239:O73:O9	2.49	0.61
1:A:126:PHE:CZ	4:A:240:O73:H26	2.35	0.60
1:B:169:TRP:HB3	1:B:170:PRO:HD3	1.83	0.60
1:B:157:LYS:HE2	1:B:166:TYR:OH	2.02	0.59
4:A:240:O73:C28	3:B:238:FAD:HM72	2.24	0.59
1:B:54:ASP:OD2	1:B:118:ARG:HD2	2.03	0.59
1:B:25:ALA:O	1:B:29:LEU:HD12	2.04	0.58
1:A:73:GLU:O	1:A:77:GLN:HG2	2.02	0.57
4:A:240:O73:O10	4:A:240:O73:O24	2.22	0.57
1:A:149:GLY:C	4:A:239:O73:H11A	2.26	0.56
4:A:239:O73:H26	1:B:126:PHE:HZ	1.70	0.56
4:A:239:O73:H28	1:B:122:GLN:H	1.72	0.55
1:A:139:GLY:H	1:A:182:LYS:HZ1	1.54	0.55
1:A:15:LYS:HD3	1:B:64:VAL:HG22	1.90	0.53
4:A:239:O73:H22A	1:B:178:PHE:CE1	2.46	0.51
1:B:54:ASP:OD2	1:B:118:ARG:CD	2.59	0.51
1:B:36:VAL:HG12	1:B:37:THR:N	2.26	0.51
1:A:120:LEU:O	4:A:240:O73:H29	2.11	0.51
1:B:122:GLN:NE2	1:B:128:ILE:HG12	2.27	0.50
1:A:50:ALA:HB1	1:A:67:TYR:CZ	2.48	0.49
1:A:150:GLY:HA2	4:A:239:O73:H21B	1.95	0.48
1:B:54:ASP:C	1:B:55:ILE:HG13	2.34	0.48
1:B:55:ILE:HG23	1:B:74:ALA:HB2	1.95	0.47
1:A:172:GLN:HE22	1:A:186:PRO:HD3	1.79	0.47
1:A:4:LYS:HG2	1:A:35:THR:HB	1.95	0.47
1:B:187:GLN:HG2	1:B:207:TRP:CE3	2.50	0.47
1:B:127:ASP:C	1:B:128:ILE:HD13	2.35	0.47
1:B:60:SER:HB2	1:B:70:GLU:OE1	2.15	0.47
4:A:240:O73:H28	1:B:105:TRP:CZ2	2.51	0.46
1:A:17:PHE:HB2	3:A:238:FAD:C5B	2.36	0.46
4:A:240:O73:O23	4:A:240:O73:H22	2.15	0.46
1:A:29:LEU:CD2	1:A:214:ILE:HD11	2.43	0.46
4:A:239:O73:H26	1:B:126:PHE:CZ	2.51	0.45
1:A:230:GLN:HA	1:A:230:GLN:OE1	2.17	0.45
1:A:75:TYR:CE1	1:A:124:PHE:HB2	2.53	0.44
1:A:132:TYR:HE2	1:B:160:VAL:HG13	1.83	0.43
1:A:117:ASP:O	4:A:240:O73:H29A	2.18	0.43
1:B:127:ASP:O	1:B:128:ILE:HD13	2.18	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:TYR:HA	3:B:238:FAD:C5X	2.48	0.43
1:B:163:ASP:OD2	1:B:165:ARG:HG3	2.18	0.43
1:A:25:ALA:HA	1:A:211:LEU:HD13	2.01	0.43
1:B:3:LYS:HD3	1:B:3:LYS:HA	1.84	0.43
1:A:121:CYS:HA	4:A:240:O73:C29	2.42	0.43
1:A:199:GLU:O	1:A:203:MET:HG3	2.19	0.43
1:A:126:PHE:HA	1:A:136:LEU:HG	2.01	0.43
1:A:17:PHE:CE2	1:A:21:LEU:HD11	2.54	0.42
3:A:238:FAD:HM82	4:A:239:O73:H29	2.01	0.42
1:B:4:LYS:HD3	1:B:35:THR:HG21	2.01	0.42
1:A:126:PHE:CZ	4:A:240:O73:C26	3.01	0.42
1:A:141:LEU:HD22	1:A:184:LEU:HD21	2.02	0.42
1:A:31:ARG:HH11	1:A:31:ARG:CG	2.13	0.42
1:B:173:HIS:HA	1:B:177:HIS:HB3	2.01	0.42
1:A:126:PHE:HZ	4:A:240:O73:C26	2.33	0.42
1:A:113:LYS:CE	1:B:107:SER:HB3	2.50	0.42
1:A:197:GLU:HG2	1:A:198:GLU:OE2	2.20	0.42
1:A:25:ALA:HA	1:A:211:LEU:CD1	2.50	0.41
1:B:75:TYR:CE1	1:B:124:PHE:HB2	2.55	0.41
1:A:89:LYS:HD3	5:A:272:HOH:O	2.21	0.41
1:A:139:GLY:N	1:A:182:LYS:HZ3	2.12	0.41
1:A:108:VAL:H	1:A:108:VAL:HG22	1.45	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	228/237 (96%)	214 (94%)	10 (4%)	4 (2%)	<b>8</b> <b>3</b>
1	B	228/237 (96%)	219 (96%)	8 (4%)	1 (0%)	34 30
All	All	456/474 (96%)	433 (95%)	18 (4%)	5 (1%)	<b>14</b> <b>8</b>

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	GLY
1	A	63	GLU
1	A	62	PRO
1	B	130	GLY
1	A	130	GLY

### 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	194/201 (96%)	182 (94%)	12 (6%)	18	13
1	B	194/201 (96%)	181 (93%)	13 (7%)	16	11
All	All	388/402 (96%)	363 (94%)	25 (6%)	17	13

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

Mol	Chain	Res	Type
1	A	22	LYS
1	A	58	THR
1	A	77	GLN
1	A	89	LYS
1	A	127	ASP
1	A	146	VAL
1	A	153	GLU
1	A	157	LYS
1	A	182	LYS
1	A	208	SER
1	A	214	ILE
1	A	216	LYS
1	B	30	SER
1	B	31	ARG
1	B	45	ASN
1	B	58	THR
1	B	60	SER
1	B	63	GLU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	96	LEU
1	B	113	LYS
1	B	140	LYS
1	B	156	THR
1	B	158	THR
1	B	160	VAL
1	B	230	GLN

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	172	GLN
1	A	187	GLN
1	B	88	GLN
1	B	225	HIS

### 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](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	O73	A	240	-	30,32,32	2.18	6 (20%)	33,45,45	2.30	11 (33%)
4	O73	A	239	-	30,32,32	2.28	5 (16%)	33,45,45	1.71	9 (27%)
3	FAD	B	238	-	53,58,58	1.37	6 (11%)	68,89,89	1.93	14 (20%)
3	FAD	A	238	-	53,58,58	1.41	7 (13%)	68,89,89	1.51	12 (17%)

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
4	O73	A	240	-	-	3/11/31/31	0/3/3/3
4	O73	A	239	-	-	3/11/31/31	0/3/3/3
3	FAD	B	238	-	-	1/30/50/50	0/6/6/6
3	FAD	A	238	-	-	2/30/50/50	0/6/6/6

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	239	O73	C16-C15	7.50	1.47	1.40
4	A	240	O73	C16-C15	6.88	1.46	1.40
4	A	240	O73	C13-C12	5.72	1.51	1.39
4	A	239	O73	C13-C12	5.43	1.50	1.39
4	A	239	O73	C16-C12	5.06	1.48	1.39
3	B	238	FAD	C4X-N5	4.67	1.39	1.30
4	A	240	O73	C13-N14	-3.89	1.36	1.39
3	B	238	FAD	C2A-N3A	3.87	1.38	1.32
3	A	238	FAD	C4X-N5	3.81	1.38	1.30
3	A	238	FAD	C2A-N3A	3.76	1.38	1.32
4	A	240	O73	C16-C12	3.60	1.46	1.39
4	A	239	O73	C13-N14	-3.51	1.36	1.39
3	A	238	FAD	C10-N1	3.42	1.40	1.33
3	B	238	FAD	C1'-C2'	3.40	1.57	1.52
3	A	238	FAD	C2A-N1A	2.84	1.39	1.33
3	B	238	FAD	C6-C7	2.45	1.43	1.39
4	A	239	O73	C18-C17	-2.38	1.39	1.44
4	A	240	O73	C18-C19	2.32	1.41	1.36
3	A	238	FAD	C4X-C4	-2.30	1.36	1.44
3	B	238	FAD	C2A-N1A	2.25	1.38	1.33
4	A	240	O73	C18-C17	-2.16	1.40	1.44
3	A	238	FAD	C1'-C2'	2.09	1.55	1.52

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	238	FAD	C9A-N10	-2.09	1.37	1.41
3	B	238	FAD	C4X-C10	-2.00	1.38	1.44

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	238	FAD	N3A-C2A-N1A	-6.40	118.68	128.68
4	A	240	O73	C20-C19-N25	5.86	119.08	114.17
3	A	238	FAD	N3A-C2A-N1A	-5.54	120.02	128.68
4	A	240	O73	C16-C15-N14	5.00	113.16	108.24
4	A	240	O73	C11-C12-C13	4.53	132.58	127.17
3	B	238	FAD	C9A-C5X-N5	-4.33	117.72	122.43
4	A	240	O73	C26-N25-C19	-4.11	114.20	124.34
4	A	240	O73	C12-C13-N14	4.08	110.56	106.86
3	B	238	FAD	O3'-C3'-C2'	-4.02	99.11	108.81
3	A	238	FAD	C1B-N9A-C4A	-3.95	119.70	126.64
3	B	238	FAD	P-O3P-PA	-3.91	119.42	132.83
4	A	239	O73	C20-C19-N25	3.85	117.40	114.17
4	A	240	O73	C27-C26-N25	3.74	122.89	112.21
4	A	239	O73	C22-N14-C15	3.62	128.46	124.49
4	A	239	O73	C1-C6-N7	3.61	122.09	119.38
3	B	238	FAD	C4X-C10-N10	3.45	121.53	116.48
3	A	238	FAD	O4'-C4'-C5'	-3.37	102.35	109.92
3	B	238	FAD	C4-C4X-N5	3.36	123.01	118.23
4	A	240	O73	C11-O10-C3	3.32	125.86	117.65
3	B	238	FAD	C10-N1-C2	3.30	123.50	116.90
3	A	238	FAD	O4-C4-C4X	-3.14	118.28	126.60
3	A	238	FAD	C5X-C9A-N10	3.12	121.18	117.95
4	A	239	O73	C12-C13-N14	3.12	109.69	106.86
3	B	238	FAD	C10-C4X-N5	-3.08	118.31	124.86
3	B	238	FAD	C5X-N5-C4X	2.88	122.86	118.07
4	A	239	O73	C16-C15-N14	2.86	111.06	108.24
4	A	239	O73	C22-N14-C13	-2.85	121.30	124.09
3	B	238	FAD	C5X-C9A-N10	2.73	120.78	117.95
4	A	240	O73	C22-N14-C15	2.63	127.37	124.49
3	A	238	FAD	O5B-PA-O1A	-2.58	98.99	109.07
4	A	240	O73	C18-C19-N25	-2.57	120.71	124.88
3	B	238	FAD	O2-C2-N3	2.56	123.62	118.65
4	A	239	O73	C11-O10-C3	2.55	123.94	117.65
3	A	238	FAD	C4X-C4-N3	2.45	119.42	113.19
3	A	238	FAD	C9-C9A-N10	-2.40	118.59	121.84
4	A	240	O73	C5-C6-N7	2.35	121.14	119.38

*Continued on next page...*



Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	240	O73	O8-N7-C6	2.26	122.00	118.80
3	B	238	FAD	C4X-C4-N3	2.21	118.80	113.19
3	A	238	FAD	C1'-N10-C9A	2.19	124.16	120.51
3	A	238	FAD	C4X-C10-N10	2.18	119.67	116.48
3	A	238	FAD	C5'-C4'-C3'	2.11	116.29	112.20
3	B	238	FAD	C4-N3-C2	-2.10	121.76	125.64
3	A	238	FAD	C4-N3-C2	-2.07	121.82	125.64
4	A	239	O73	C18-C19-N25	-2.03	121.58	124.88
4	A	239	O73	O23-C17-C18	-2.03	118.40	121.79
3	B	238	FAD	C9-C9A-N10	-2.01	119.12	121.84

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	238	FAD	O4B-C4B-C5B-O5B
4	A	239	O73	C2-C3-O10-C11
4	A	240	O73	C2-C3-O10-C11
3	B	238	FAD	O4B-C4B-C5B-O5B
4	A	240	O73	C4-C3-O10-C11
4	A	239	O73	C4-C3-O10-C11
4	A	239	O73	C26-C27-C28-C29
4	A	240	O73	C27-C28-C29-N30
3	A	238	FAD	C4'-C5'-O5'-P

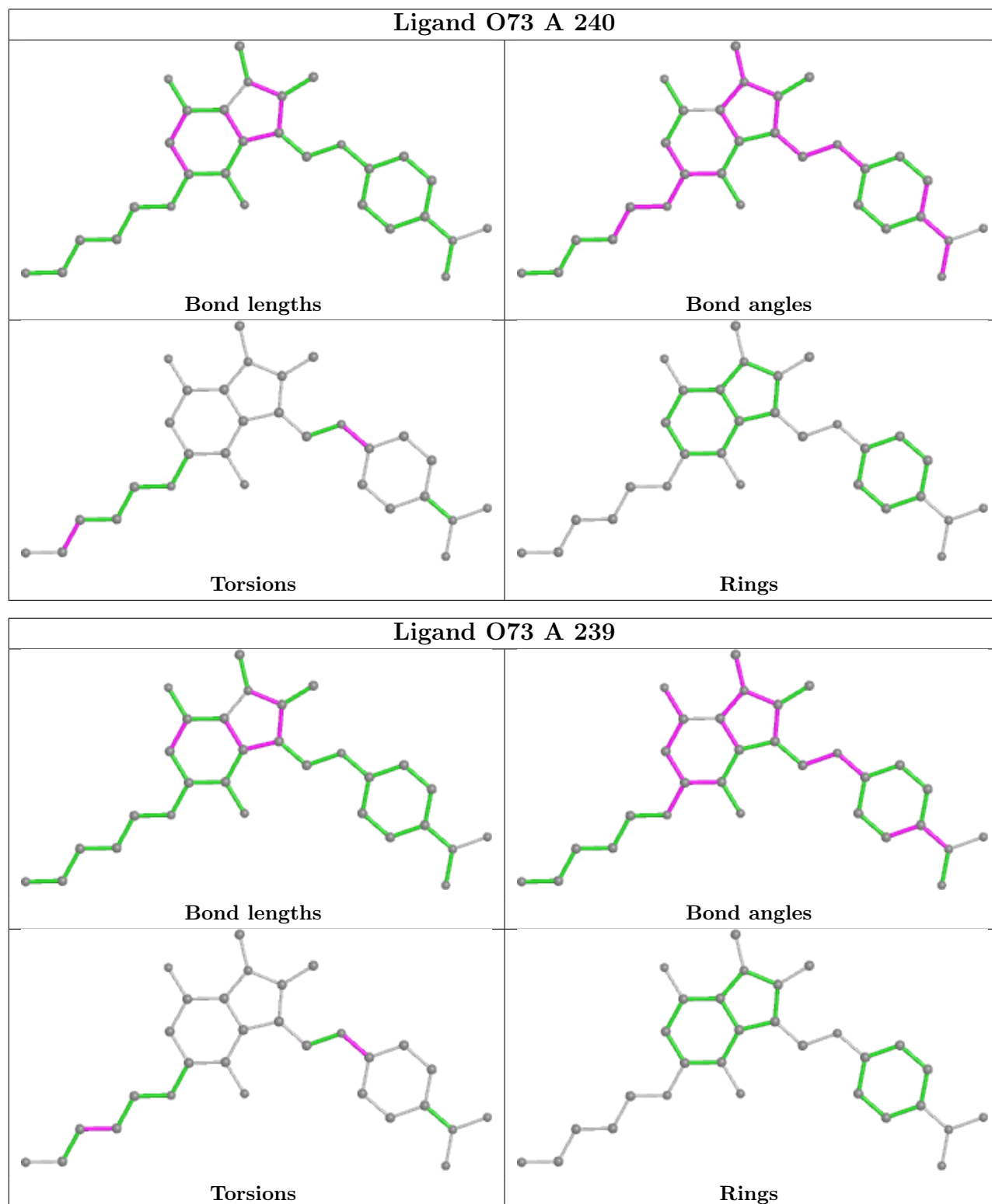
There are no ring outliers.

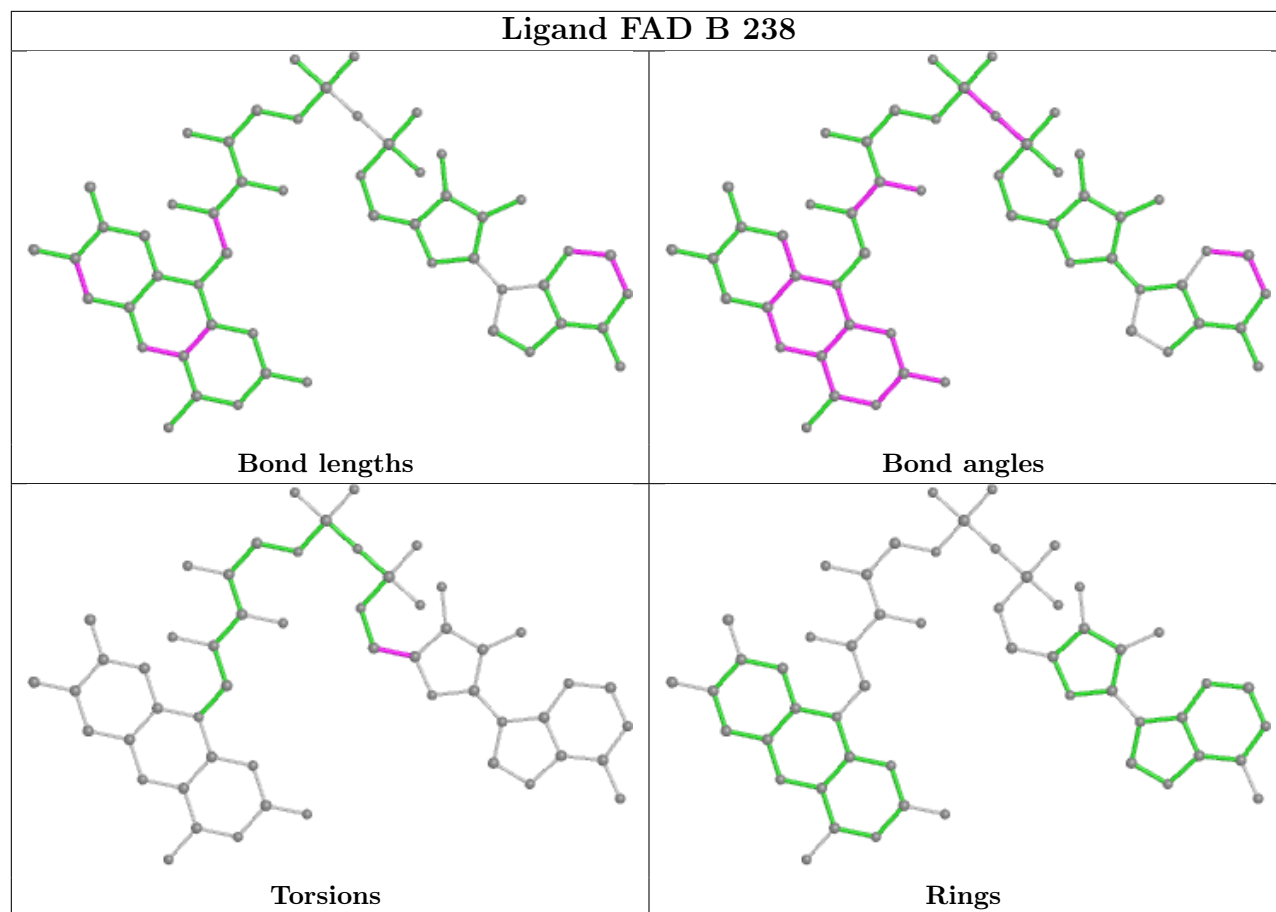
4 monomers are involved in 32 short contacts:

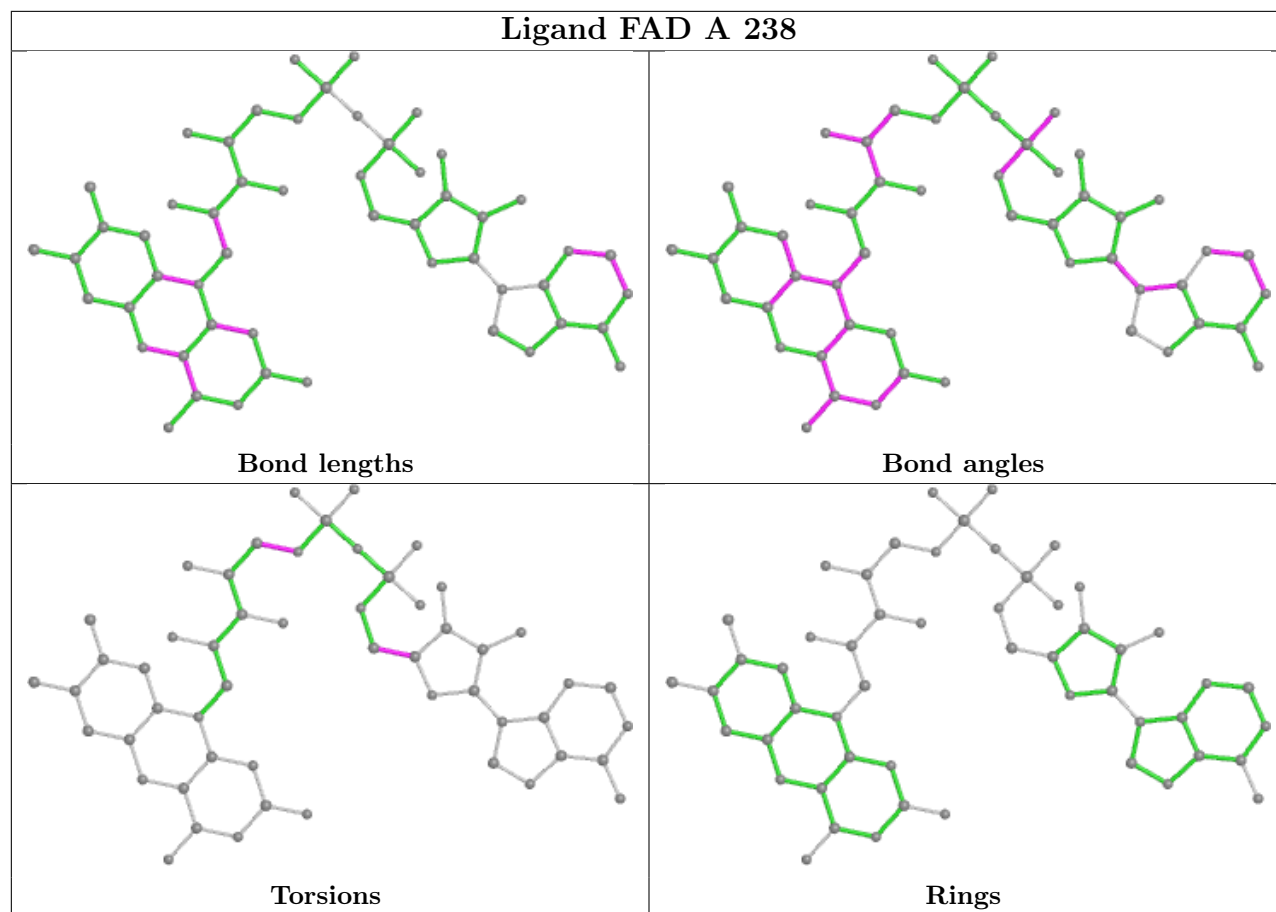
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	240	O73	17	0
4	A	239	O73	11	0
3	B	238	FAD	5	0
3	A	238	FAD	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	230/237 (97%)	0.19	7 (3%) 50 49	16, 35, 54, 65	0
1	B	230/237 (97%)	0.29	13 (5%) 23 23	21, 35, 55, 64	0
All	All	460/474 (97%)	0.24	20 (4%) 35 34	16, 35, 55, 65	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	ALA	8.5
1	B	230	GLN	4.7
1	B	196	SER	3.8
1	B	62	PRO	3.7
1	A	1	ALA	3.0
1	B	194	ILE	3.0
1	B	197	GLU	2.9
1	B	195	ALA	2.7
1	B	158	THR	2.7
1	A	201	LYS	2.6
1	A	62	PRO	2.5
1	B	64	VAL	2.4
1	B	157	LYS	2.4
1	B	133	ASP	2.3
1	A	230	GLN	2.3
1	A	216	LYS	2.1
1	A	58	THR	2.1
1	B	229	GLY	2.1
1	A	64	VAL	2.0
1	B	108	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](#)

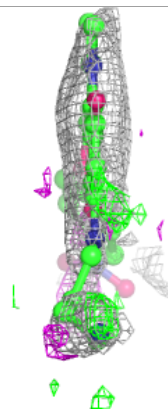
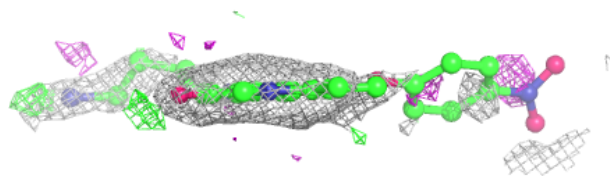
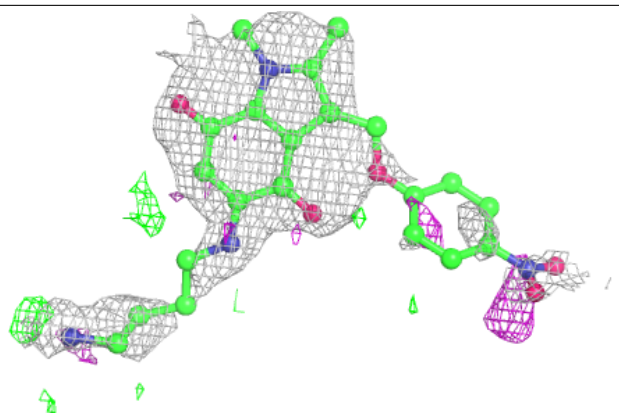
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	O73	A	239	30/30	0.76	0.34	51,58,92,93	0
4	O73	A	240	30/30	0.82	0.29	52,57,85,87	0
3	FAD	A	238	53/53	0.94	0.12	19,31,48,50	0
3	FAD	B	238	53/53	0.94	0.14	21,31,54,54	0
2	ZN	A	237	1/1	0.99	0.07	36,36,36,36	0
2	ZN	B	237	1/1	0.99	0.05	38,38,38,38	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.

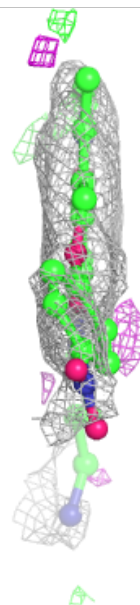
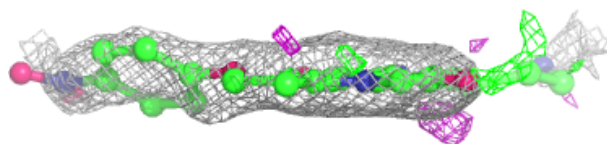
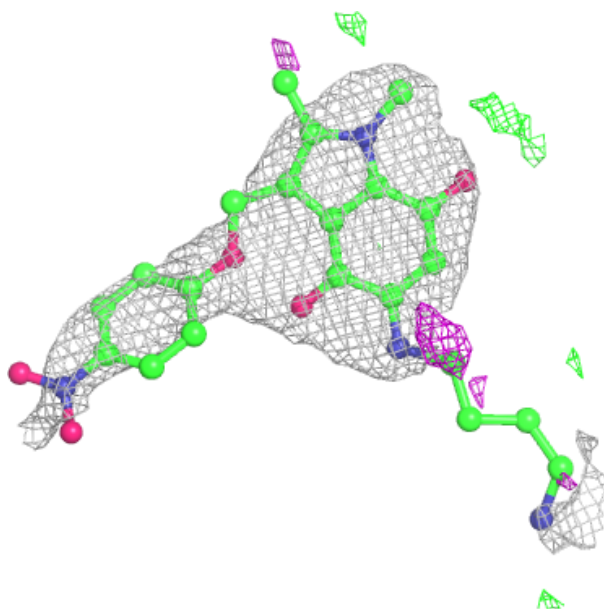
**Electron density around O73 A 239:**

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



**Electron density around O73 A 240:**

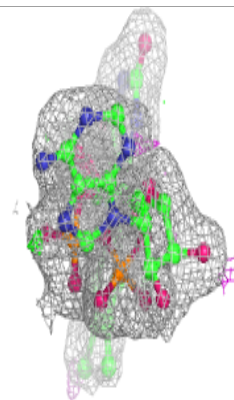
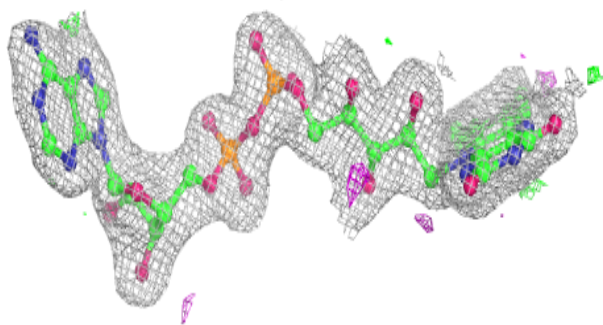
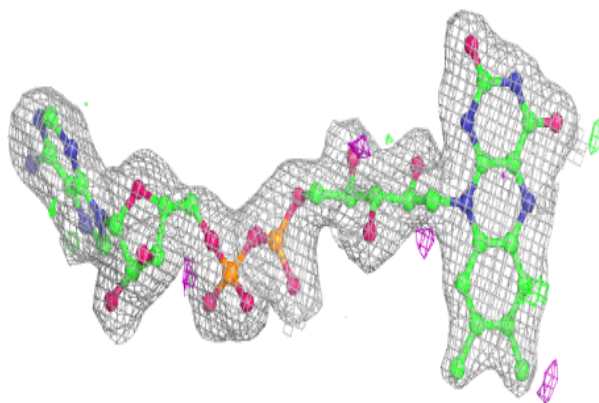
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



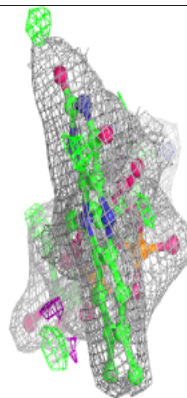
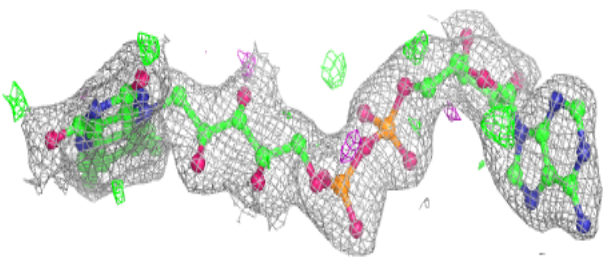
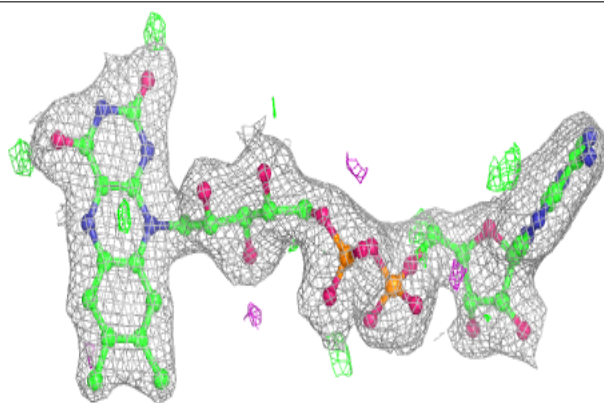


**Electron density around FAD A 238:**

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

**Electron density around FAD B 238:**

$2mF_o-DF_c$  (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.