



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 03:38 pm BST

PDB ID : 1ZX1
Title : Human quinone oxidoreductase 2 (NQO2) in complex with the cytostatic pro-drug CB1954
Authors : Jansson, A.; Wu, X.; Kavanagh, K.; Kerr, D.; Knox, R.; Walton, R.; Gunther, U.; Ludwig, C.; Edwards, A.; Arrowsmith, C.; Sundstrom, M.; von Delft, F.; Oppermann, U.; Structural Genomics Consortium (SGC)
Deposited on : 2005-06-06
Resolution : 2.16 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 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 : 2.11
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) : 2.11

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3870 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.

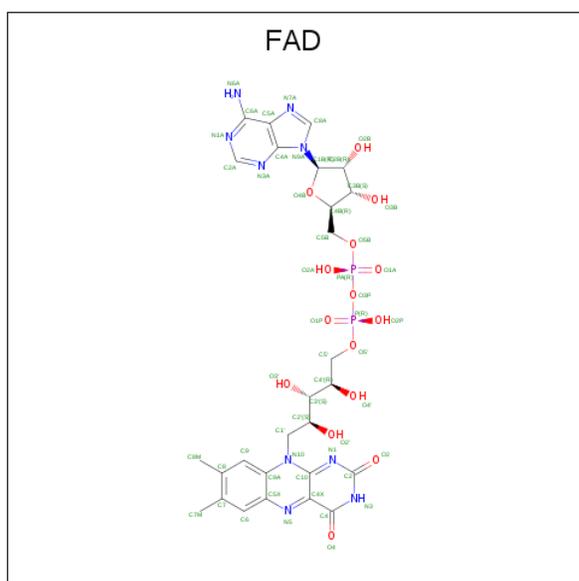
- Molecule 1 is a protein called NRH dehydrogenase [quinone] 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	227	Total 1786	C 1155	N 296	O 327	S 8	0	1	0
1	B	229	Total 1784	C 1153	N 297	O 326	S 8	0	0	0

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

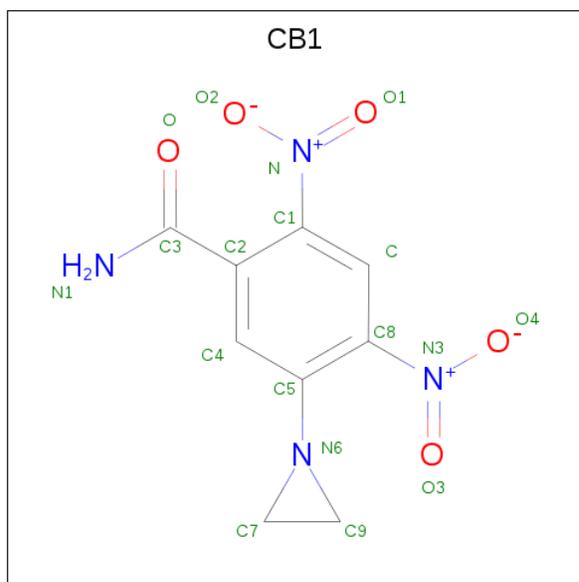
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total 1	Zn 1	0	0
2	A	1	Total 1	Zn 1	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 4 is 5-(AZIRIDIN-1-YL)-2,4-DINITROBENZAMIDE (three-letter code: CB1) (formula: C₉H₈N₄O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			18	9	4	5		
4	B	1	Total	C	N	O	0	0
			18	9	4	5		

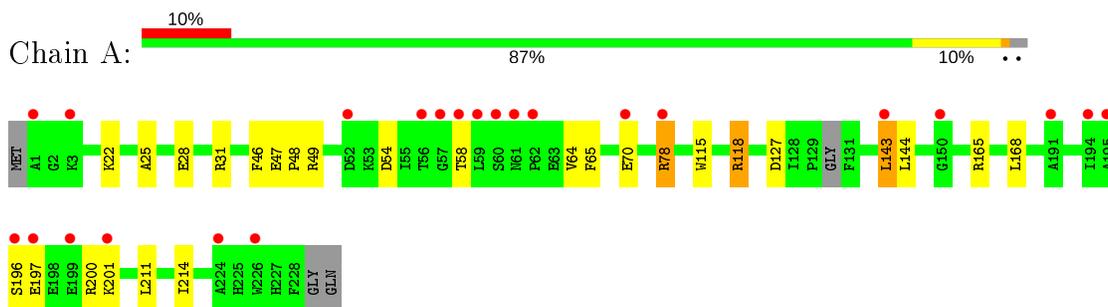
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	64	Total	O	0	1
			65	65		
5	B	90	Total	O	0	1
			91	91		

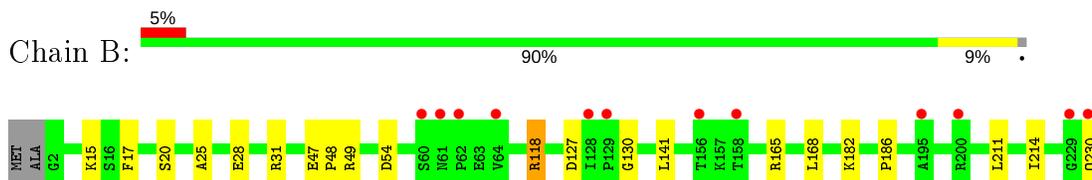
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: NRH dehydrogenase [quinone] 2



- Molecule 1: NRH dehydrogenase [quinone] 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.33Å 84.34Å 106.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.16 45.05 – 2.16	Depositor EDS
% Data completeness (in resolution range)	96.8 (50.00-2.16) 96.8 (45.05-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.53 (at 2.16Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.174 , 0.207 0.176 , 0.208	Depositor DCC
R_{free} test set	1365 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	33.1	Xtrriage
Anisotropy	0.034	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3870	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.40% 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: ZN, CB1, 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	0.76	1/1838 (0.1%)	0.78	4/2499 (0.2%)
1	B	0.75	0/1834	0.78	2/2494 (0.1%)
All	All	0.75	1/3672 (0.0%)	0.78	6/4993 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	196	SER	CB-OG	5.99	1.50	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	ARG	NE-CZ-NH2	-12.13	114.24	120.30
1	B	118	ARG	NE-CZ-NH2	-11.39	114.60	120.30
1	B	118	ARG	NE-CZ-NH1	10.93	125.76	120.30
1	A	118	ARG	NE-CZ-NH1	9.22	124.91	120.30
1	A	78	ARG	NE-CZ-NH2	-6.72	116.94	120.30
1	A	78	ARG	NE-CZ-NH1	6.61	123.60	120.30

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	1786	0	1722	15	0
1	B	1784	0	1708	11	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	53	0	31	1	0
3	B	53	0	31	3	0
4	A	18	0	8	1	0
4	B	18	0	8	1	0
5	A	65	0	0	1	0
5	B	91	0	0	1	0
All	All	3870	0	3508	29	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 4.

All (29) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:LEU:HD23	1:A:143:LEU:C	1.85	0.96
1:A:143:LEU:CD2	1:A:143:LEU:C	2.42	0.88
1:A:143:LEU:HD23	1:A:144:LEU:N	1.95	0.81
4:A:233:CB1:O2	4:A:233:CB1:O	2.06	0.73
1:A:65:PHE:CD1	1:A:70:GLU:OE1	2.53	0.61
1:A:54:ASP:OD2	1:A:118:ARG:HD2	2.01	0.60
1:B:54:ASP:OD2	1:B:118:ARG:HD2	2.02	0.59
5:A:293:HOH:O	3:B:234:FAD:HM72	2.03	0.58
1:A:143:LEU:HD22	1:A:143:LEU:O	2.06	0.55
4:B:235:CB1:O	4:B:235:CB1:O2	2.27	0.52
1:A:143:LEU:CD2	1:A:143:LEU:O	2.58	0.52
1:B:28:GLU:OE2	1:B:31:ARG:NH1	2.43	0.51
1:B:49:ARG:O	1:B:118:ARG:HD3	2.11	0.50
1:A:49:ARG:O	1:A:118:ARG:HD3	2.12	0.49
1:B:17:PHE:HB2	3:B:234:FAD:H51A	1.94	0.48
1:B:141:LEU:HD23	1:B:182:LYS:HB2	1.96	0.47
1:A:28:GLU:OE2	1:A:31:ARG:NH1	2.48	0.47
1:A:64:VAL:HG22	1:B:15:LYS:HD2	1.98	0.46
3:A:232:FAD:HM72	5:B:310:HOH:O	2.16	0.46
1:A:46:PHE:CE2	1:A:115:TRP:HA	2.52	0.45
1:A:197:GLU:HG2	1:A:200:ARG:HH11	1.81	0.45
1:B:20:SER:OG	3:B:234:FAD:N7A	2.45	0.43
1:B:47:GLU:HA	1:B:48:PRO:HD3	1.90	0.42
1:B:165:ARG:HA	1:B:168:LEU:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:ARG:HA	1:A:168:LEU:HD12	2.02	0.41
1:A:47:GLU:HA	1:A:48:PRO:HD3	1.90	0.41
1:B:25:ALA:HA	1:B:211:LEU:CD1	2.51	0.41
1:A:25:ALA:HA	1:A:211:LEU:HD13	2.03	0.40
1:B:168:LEU:HD13	1:B:186:PRO:HG3	2.02	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	224/231 (97%)	218 (97%)	6 (3%)	0	100	100
1	B	227/231 (98%)	221 (97%)	5 (2%)	1 (0%)	34	29
All	All	451/462 (98%)	439 (97%)	11 (2%)	1 (0%)	47	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	186/195 (95%)	179 (96%)	7 (4%)	33	31
1	B	183/195 (94%)	180 (98%)	3 (2%)	62	67
All	All	369/390 (95%)	359 (97%)	10 (3%)	44	46

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	58	THR
1	A	78	ARG
1	A	127	ASP
1	A	143	LEU
1	A	201	LYS
1	A	214	ILE
1	B	127	ASP
1	B	214	ILE
1	B	230	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	45	ASN

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 carbohydrates 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
3	FAD	A	232	-	51,58,58	1.51	7 (13%)	60,89,89	2.06	11 (18%)
4	CB1	B	235	-	17,19,19	4.77	8 (47%)	18,28,28	1.34	3 (16%)
4	CB1	A	233	-	17,19,19	4.78	9 (52%)	18,28,28	1.72	4 (22%)
3	FAD	B	234	-	51,58,58	1.43	7 (13%)	60,89,89	1.75	9 (15%)

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
3	FAD	A	232	-	-	1/30/50/50	0/6/6/6
4	CB1	B	235	-	-	0/12/18/18	0/2/2/2
4	CB1	A	233	-	-	2/12/18/18	0/2/2/2
3	FAD	B	234	-	-	2/30/50/50	0/6/6/6

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	233	CB1	O1-N	13.82	1.46	1.22
4	B	235	CB1	O1-N	12.64	1.44	1.22
4	B	235	CB1	O3-N3	10.80	1.41	1.22
4	A	233	CB1	O3-N3	9.34	1.38	1.22
4	B	235	CB1	C8-N3	-6.39	1.34	1.45
4	A	233	CB1	C8-N3	-6.09	1.34	1.45
4	A	233	CB1	C1-N	-5.79	1.35	1.45
4	B	235	CB1	C1-N	-5.28	1.36	1.45
3	A	232	FAD	C1'-N10	4.81	1.53	1.48
3	B	234	FAD	C10-N1	4.67	1.39	1.33
3	A	232	FAD	C2A-N3A	4.25	1.39	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	234	FAD	C2A-N3A	4.21	1.38	1.32
3	B	234	FAD	C4X-N5	4.19	1.39	1.33
3	A	232	FAD	C4X-N5	4.11	1.39	1.33
3	A	232	FAD	C10-N1	3.59	1.37	1.33
4	B	235	CB1	C9-N6	3.41	1.49	1.46
4	B	235	CB1	C7-N6	3.26	1.48	1.46
4	A	233	CB1	C5-N6	-3.03	1.34	1.41
4	A	233	CB1	C7-N6	2.95	1.48	1.46
3	B	234	FAD	C5X-N5	2.94	1.40	1.35
4	B	235	CB1	C5-N6	-2.78	1.35	1.41
3	A	232	FAD	C2A-N1A	2.77	1.39	1.33
3	B	234	FAD	C2A-N1A	2.63	1.38	1.33
4	B	235	CB1	C2-C3	2.58	1.53	1.50
4	A	233	CB1	C2-C3	2.57	1.53	1.50
4	A	233	CB1	C9-N6	2.47	1.48	1.46
4	A	233	CB1	C9-C7	2.45	1.53	1.47
3	A	232	FAD	C4X-C10	2.39	1.41	1.38
3	A	232	FAD	C5X-N5	2.26	1.39	1.35
3	B	234	FAD	C1'-N10	2.16	1.50	1.48
3	B	234	FAD	C4-N3	2.06	1.36	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	232	FAD	C1'-N10-C9A	8.94	125.33	118.29
3	A	232	FAD	C4-N3-C2	6.90	120.97	115.14
3	B	234	FAD	C4-N3-C2	6.30	120.46	115.14
3	B	234	FAD	N3A-C2A-N1A	-5.61	119.91	128.68
3	A	232	FAD	N3A-C2A-N1A	-5.56	119.99	128.68
3	B	234	FAD	C1'-N10-C9A	5.50	122.62	118.29
4	A	233	CB1	C9-N6-C7	4.60	62.12	60.53
3	B	234	FAD	C4X-N5-C5X	3.66	120.43	116.77
3	A	232	FAD	C4X-C4-N3	-3.38	118.81	123.43
3	A	232	FAD	C5X-C9A-N10	3.22	120.05	117.72
4	B	235	CB1	C4-C2-C3	-3.04	114.03	121.53
3	B	234	FAD	C5X-C9A-N10	3.04	119.92	117.72
3	A	232	FAD	C9A-N10-C10	-2.98	118.00	121.91
4	A	233	CB1	C9-C7-N6	-2.98	58.84	59.73
3	B	234	FAD	C4-C4X-N5	2.53	121.48	118.60
3	A	232	FAD	C1B-N9A-C4A	-2.44	122.35	126.64
3	B	234	FAD	C10-C4X-N5	-2.39	119.60	121.26
4	B	235	CB1	C-C8-C5	2.36	121.33	118.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	233	CB1	C4-C2-C3	-2.30	115.87	121.53
4	A	233	CB1	C7-C9-N6	-2.29	59.04	59.73
3	B	234	FAD	P-O3P-PA	-2.27	125.04	132.83
3	A	232	FAD	O2'-C2'-C3'	2.26	114.60	109.10
3	A	232	FAD	C4X-N5-C5X	2.21	118.98	116.77
4	B	235	CB1	C4-C5-N6	-2.11	119.40	122.52
3	B	234	FAD	C4A-C5A-N7A	-2.09	107.22	109.40
3	A	232	FAD	C1'-N10-C10	-2.05	116.57	118.41
3	A	232	FAD	C4A-C5A-N7A	-2.00	107.31	109.40

There are no chirality outliers.

All (5) torsion outliers are listed below:

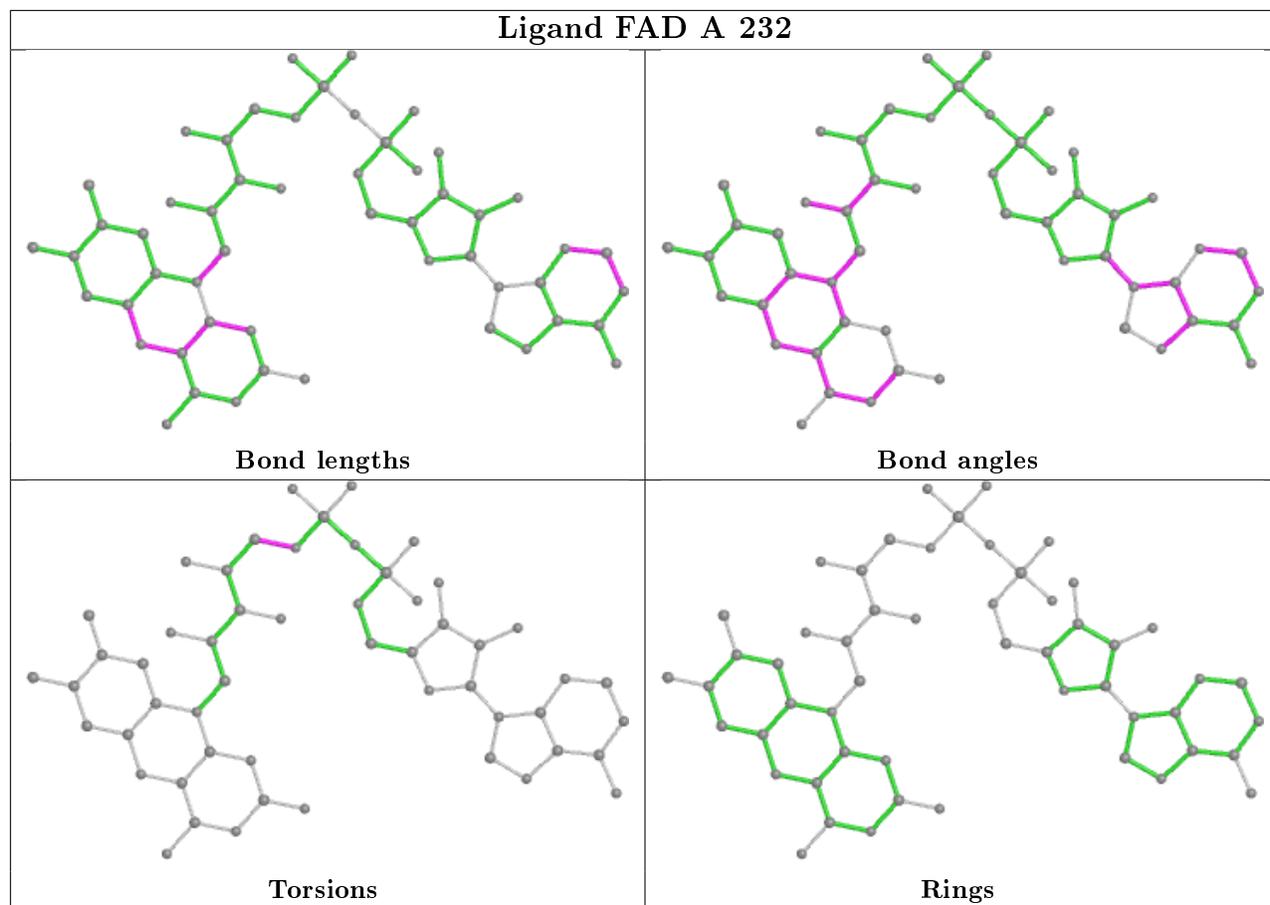
Mol	Chain	Res	Type	Atoms
4	A	233	CB1	C4-C5-N6-C7
4	A	233	CB1	C8-C5-N6-C7
3	B	234	FAD	O4B-C4B-C5B-O5B
3	B	234	FAD	C3B-C4B-C5B-O5B
3	A	232	FAD	C4'-C5'-O5'-P

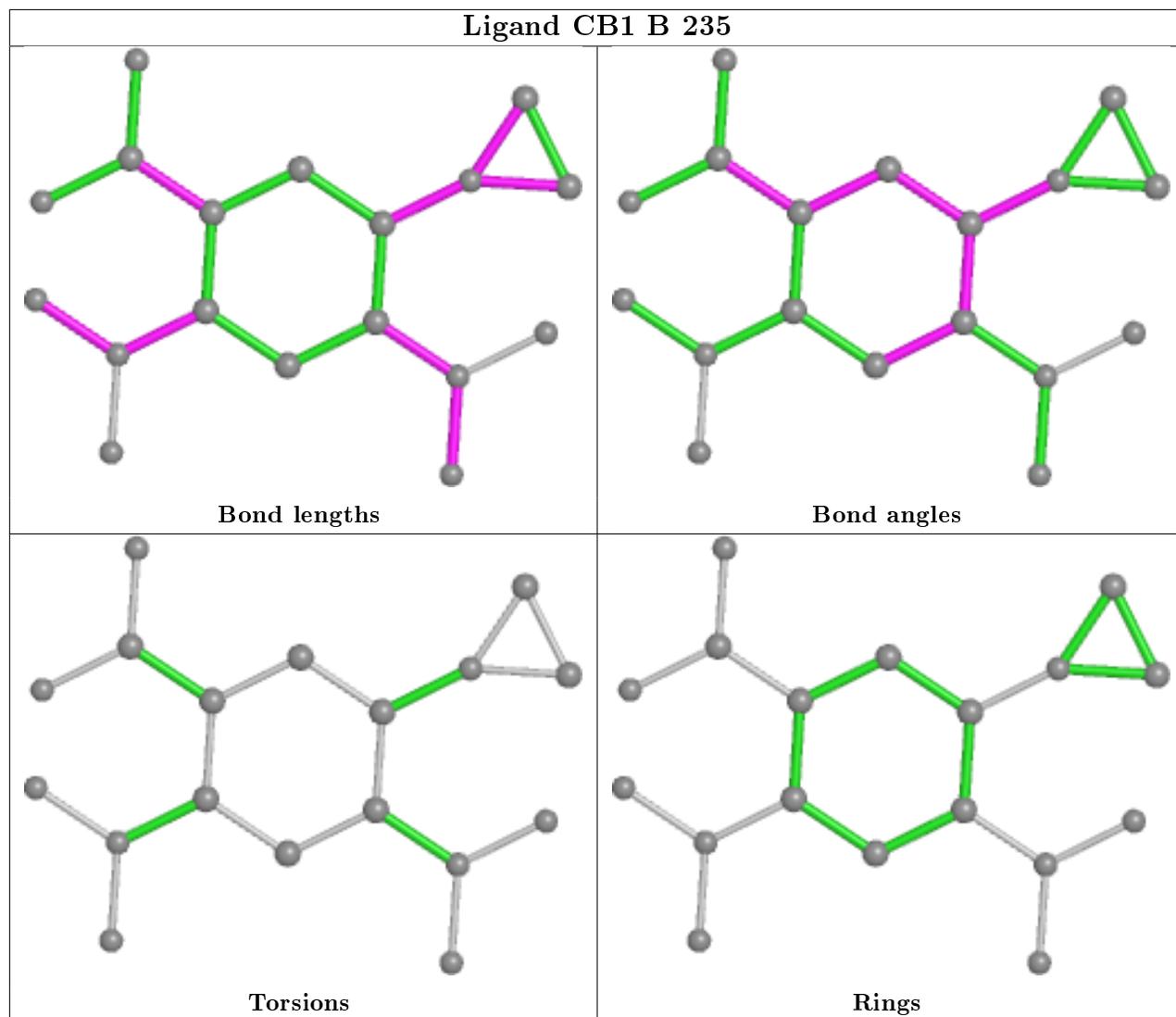
There are no ring outliers.

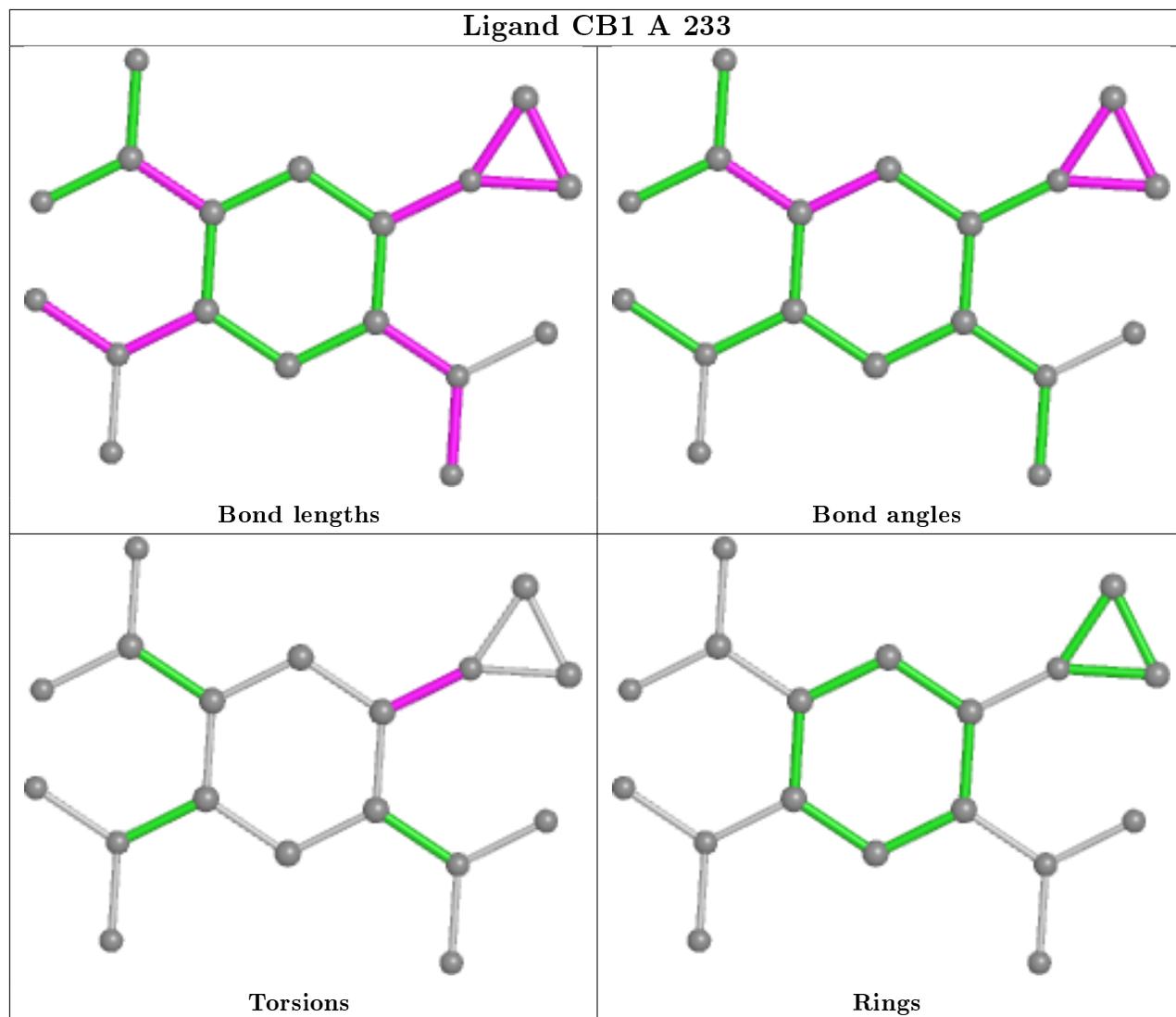
4 monomers are involved in 6 short contacts:

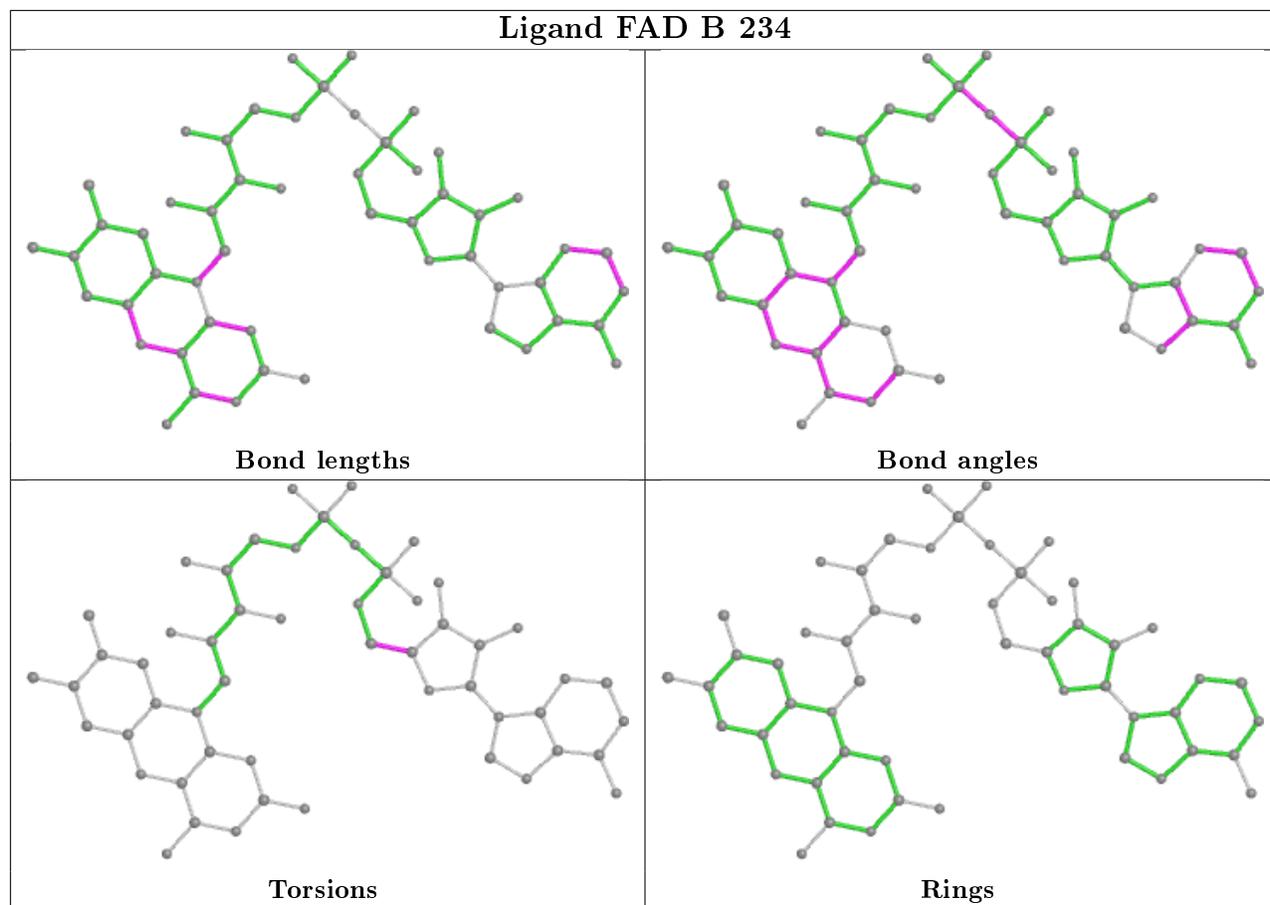
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	232	FAD	1	0
4	B	235	CB1	1	0
4	A	233	CB1	1	0
3	B	234	FAD	3	0

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, 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	227/231 (98%)	0.76	23 (10%) 7 10	33, 39, 46, 51	0
1	B	229/231 (99%)	0.48	12 (5%) 27 35	33, 39, 46, 64	0
All	All	456/462 (98%)	0.62	35 (7%) 13 18	33, 39, 46, 64	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	196	SER	4.9
1	B	62	PRO	4.8
1	A	58	THR	4.5
1	A	1	ALA	4.5
1	A	194	ILE	4.4
1	B	128	ILE	4.2
1	B	158	THR	3.8
1	A	59	LEU	3.8
1	A	60	SER	3.7
1	B	230	GLN	3.4
1	A	150	GLY	3.3
1	B	229	GLY	3.3
1	B	200	ARG	3.2
1	B	156	THR	3.1
1	A	201	LYS	3.1
1	A	70	GLU	3.0
1	B	60	SER	2.9
1	A	57	GLY	2.9
1	B	64	VAL	2.8
1	A	226	TRP	2.7
1	A	191	ALA	2.7
1	A	197	GLU	2.7
1	A	52	ASP	2.7
1	A	78	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	56	THR	2.5
1	A	62	PRO	2.5
1	A	224	ALA	2.5
1	B	61	ASN	2.4
1	A	199	GLU	2.3
1	B	129	PRO	2.2
1	A	61	ASN	2.2
1	A	195	ALA	2.1
1	B	195	ALA	2.1
1	A	143	LEU	2.0
1	A	3	LYS	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 carbohydrates in this entry.

6.4 Ligands [i](#)

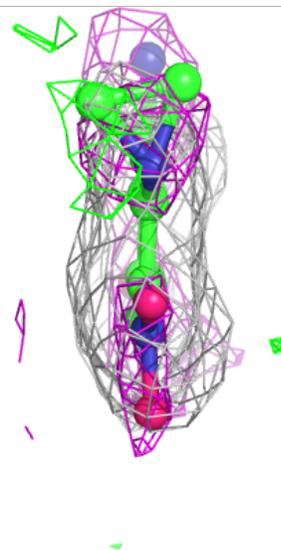
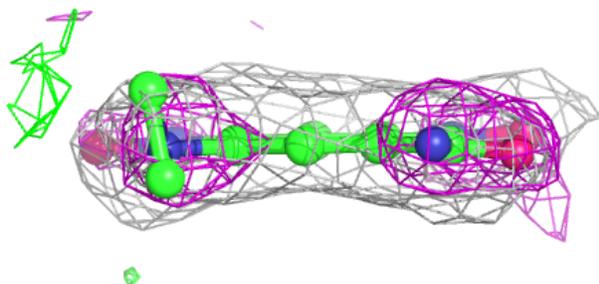
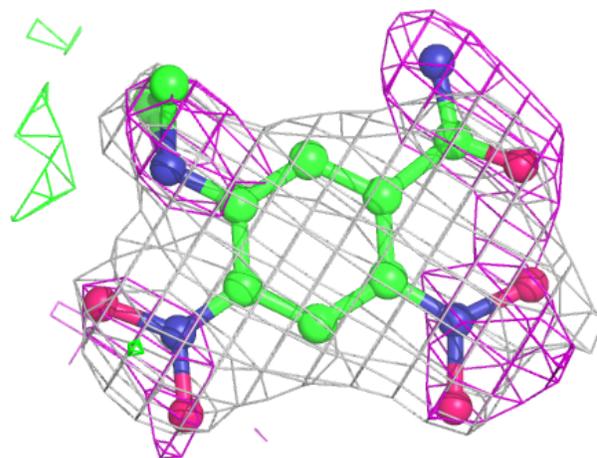
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CB1	A	233	18/18	0.87	0.32	50,50,54,54	0
3	FAD	A	232	53/53	0.88	0.28	28,36,72,73	0
4	CB1	B	235	18/18	0.90	0.32	48,50,54,55	0
3	FAD	B	234	53/53	0.91	0.22	30,38,70,71	0
2	ZN	A	231	1/1	0.94	0.14	53,53,53,53	0
2	ZN	B	231	1/1	0.99	0.14	41,41,41,41	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.

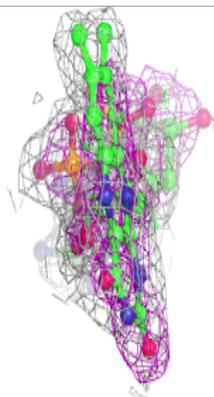
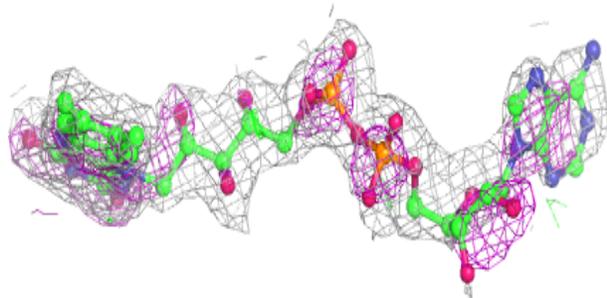
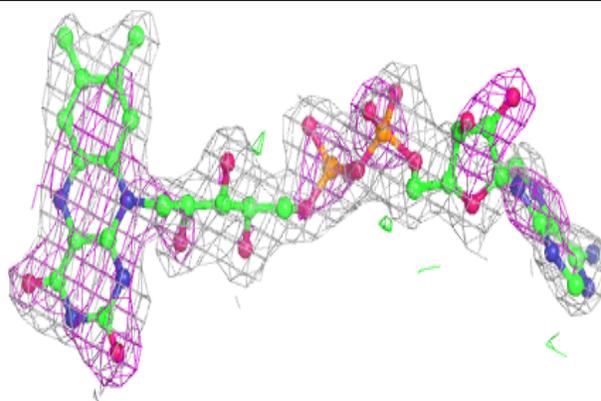
Electron density around CB1 A 233:

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



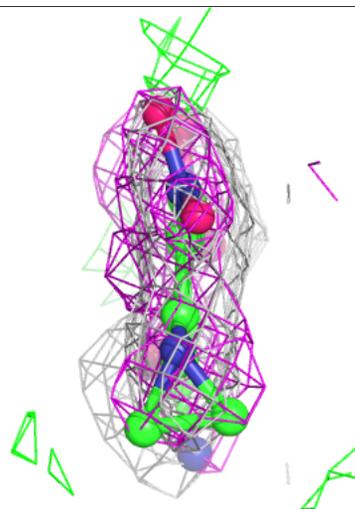
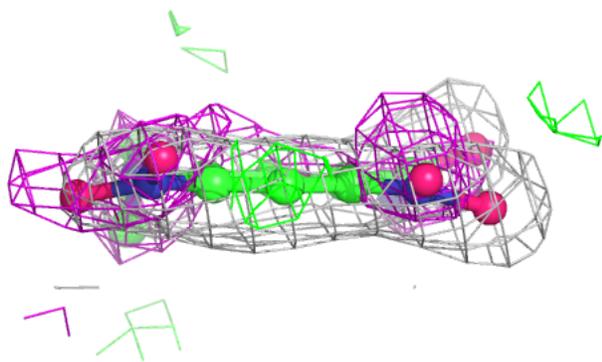
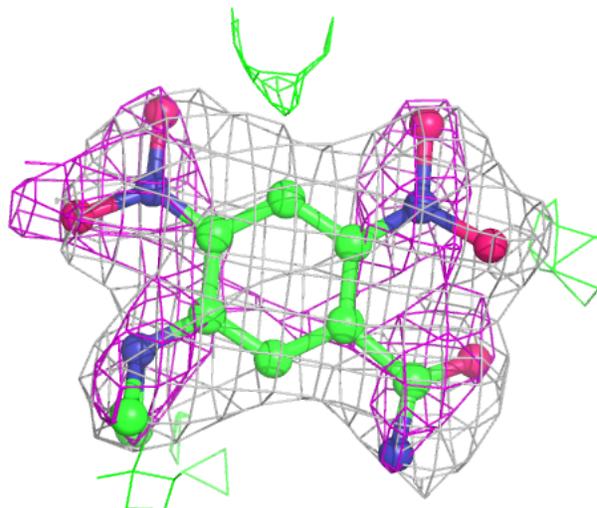
Electron density around FAD A 232:

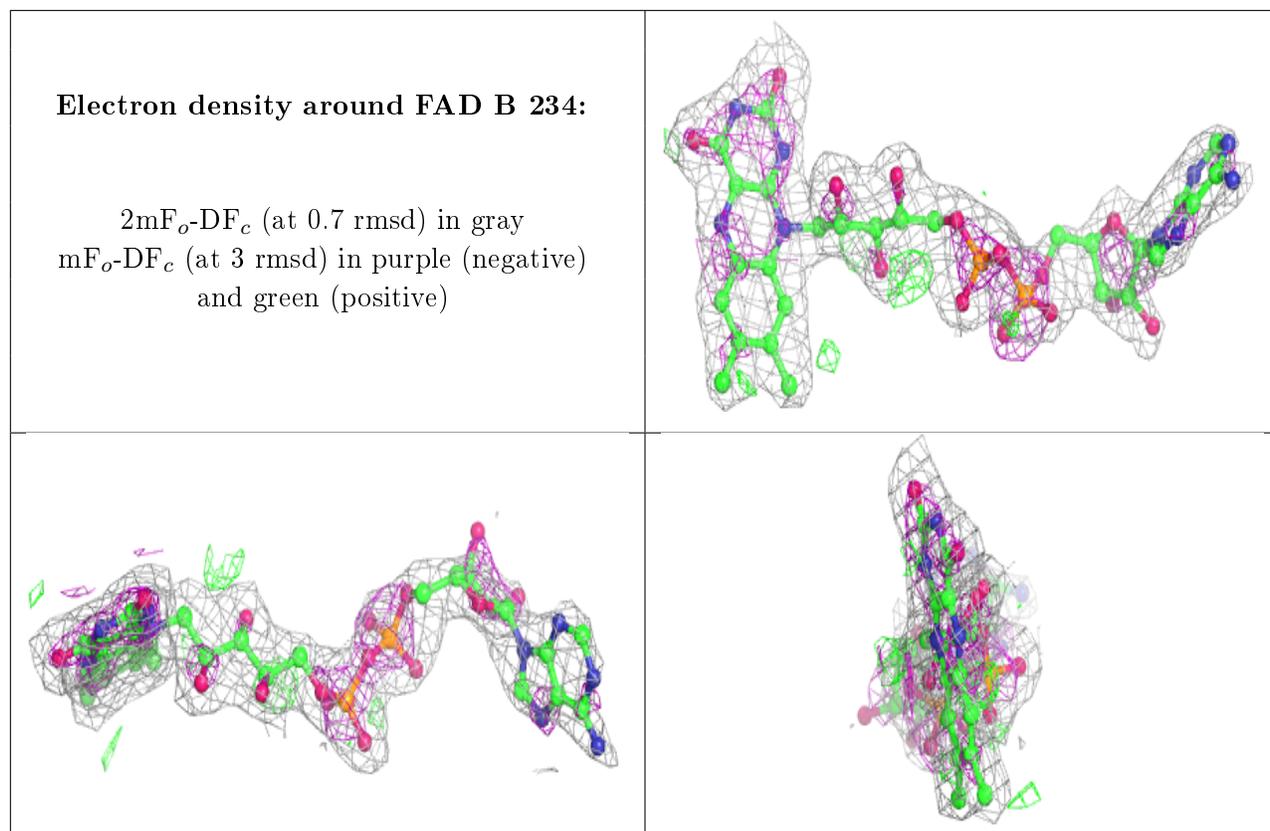
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CB1 B 235:

$2mF_o-DF_c$ (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.