



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 04:01 AM UTC

PDB ID : 3NFR / pdb_00003nfr
Title : Casimiroin analog inhibitor of quinone reductase 2
Authors : Sturdy, M.; Mesecar, A.D.; Jermihov, K.; Cushman, M.; Maiti, A.
Deposited on : 2010-06-10
Resolution : 1.57 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

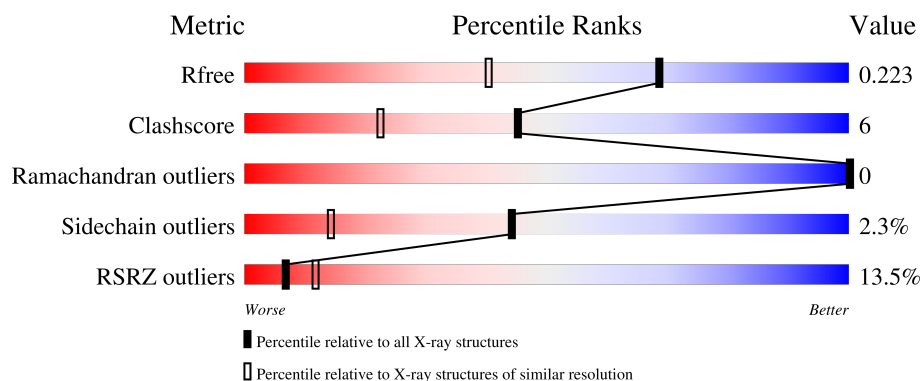
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.57 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}	180053	1094 (1.58-1.58)
Clashscore	190562	1105 (1.58-1.58)
Ramachandran outliers	187476	1082 (1.58-1.58)
Sidechain outliers	187428	1081 (1.58-1.58)
RSRZ outliers	180081	1094 (1.58-1.58)

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	230	20% 73% 24% .
1	B	230	7% 83% 17%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4086 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.

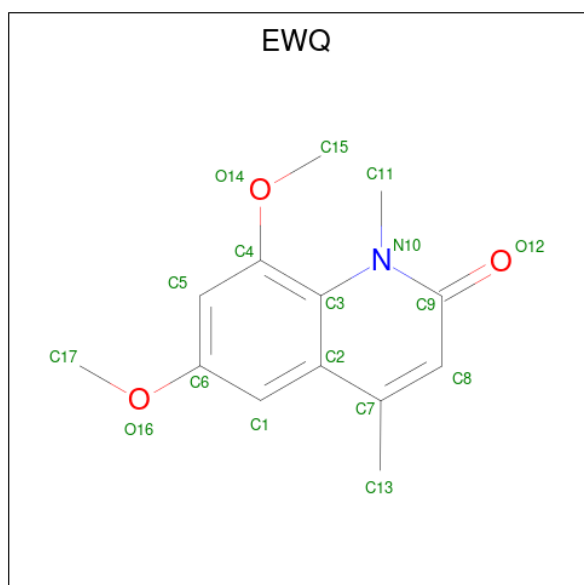
- Molecule 1 is a protein called Ribosyldihydronicotinamide dehydrogenase [quinone].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	0	0
			1824	1174	304	338	8			
1	B	230	Total	C	N	O	S	0	0	0
			1824	1174	304	338	8			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

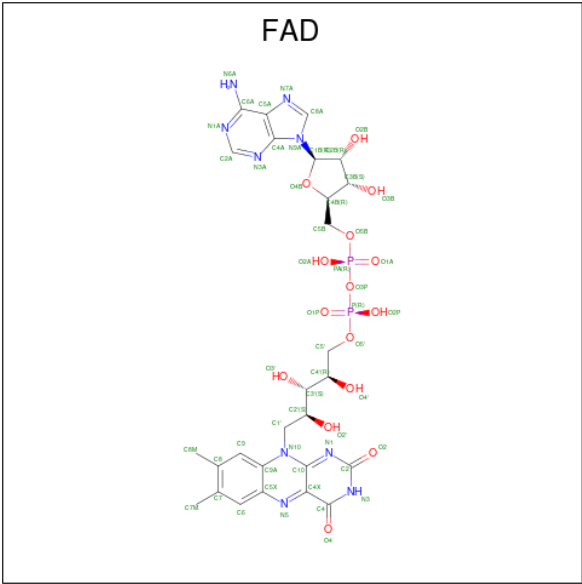
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

- Molecule 3 is 6,8-dimethoxy-1,4-dimethylquinolin-2(1H)-one (CCD ID: EWQ) (formula: C₁₃H₁₅NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	13	1	3		
3	B	1	Total	C	N	O	0	0
			17	13	1	3		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

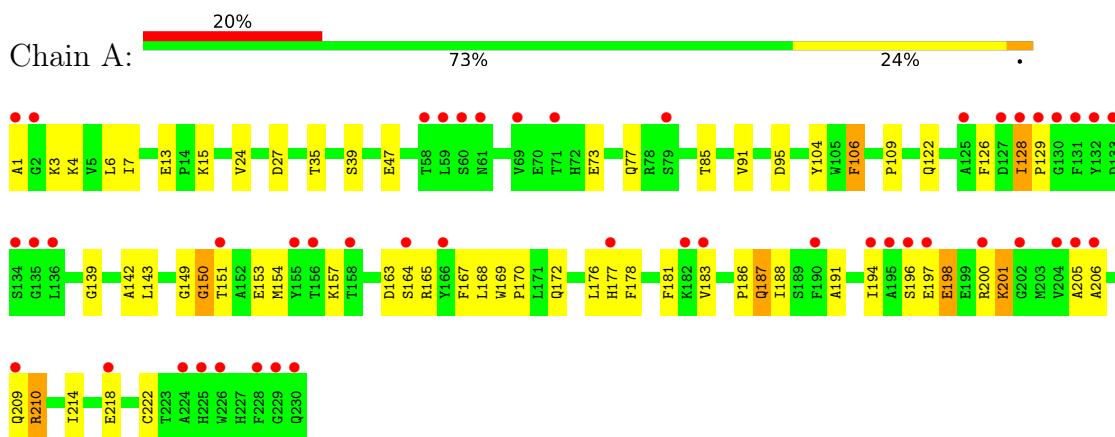
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	120	Total	O	0	0
			120	120		
5	B	176	Total	O	0	0
			176	176		

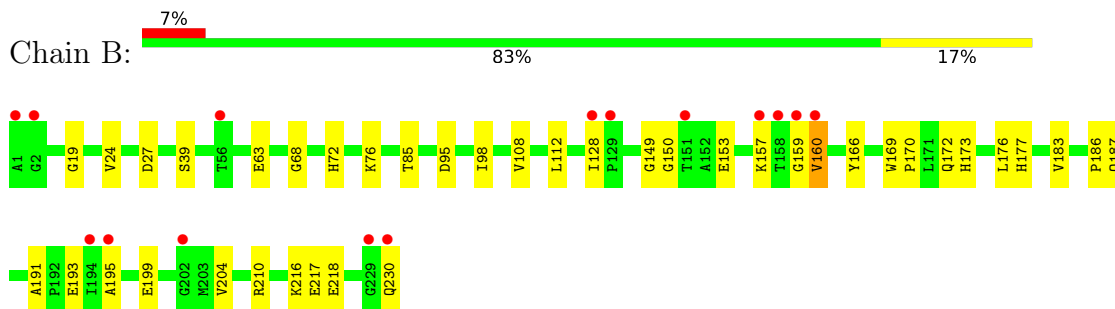
3 Residue-property plots

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: Ribosyldihydronicotinamide dehydrogenase [quinone]



- Molecule 1: Ribosyldihydronicotinamide dehydrogenase [quinone]



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.06Å 83.21Å 106.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.56 – 1.57 65.56 – 1.57	Depositor EDS
% Data completeness (in resolution range)	90.7 (65.56-1.57) 90.8 (65.56-1.57)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.15 (at 1.57Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.194 , 0.219 (Not available) , 0.223	Depositor DCC
R_{free} test set	3228 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.075	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4086	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.29% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EWQ, ZN, 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	1.65	21/1874 (1.1%)	1.43	12/2542 (0.5%)
1	B	1.58	15/1874 (0.8%)	1.39	9/2542 (0.4%)
All	All	1.62	36/3748 (1.0%)	1.41	21/5084 (0.4%)

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	176	LEU	N-CA	9.10	1.57	1.46
1	A	178	PHE	N-CA	-7.90	1.36	1.46
1	B	191	ALA	C-O	7.78	1.27	1.23
1	B	98	ILE	CA-CB	7.17	1.62	1.54
1	A	7	ILE	CA-CB	-6.78	1.45	1.53
1	A	191	ALA	CA-CB	6.70	1.58	1.52
1	B	27	ASP	N-CA	6.61	1.54	1.46
1	A	188	ILE	C-O	6.52	1.31	1.24
1	B	177	HIS	C-O	-6.38	1.16	1.24
1	A	187	GLN	N-CA	-6.30	1.38	1.46
1	B	39	SER	C-O	6.18	1.31	1.23
1	B	216	LYS	N-CA	6.01	1.54	1.46
1	A	191	ALA	C-O	5.97	1.26	1.23
1	A	39	SER	N-CA	5.73	1.54	1.46
1	A	109	PRO	N-CA	5.69	1.54	1.47
1	B	112	LEU	N-CA	5.67	1.53	1.46
1	B	108	VAL	CA-C	5.66	1.58	1.52
1	A	169	TRP	N-CA	-5.59	1.41	1.46
1	A	206	ALA	CA-C	5.53	1.59	1.52
1	A	13	GLU	C-O	-5.50	1.18	1.24
1	A	188	ILE	CA-C	-5.47	1.45	1.52
1	B	95	ASP	C-O	5.47	1.31	1.24
1	A	143	LEU	N-CA	-5.35	1.40	1.46
1	A	106	PHE	C-N	5.33	1.38	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	214	ILE	CA-CB	5.31	1.62	1.54
1	B	173	HIS	N-CA	-5.28	1.40	1.46
1	B	183	VAL	CA-CB	5.27	1.60	1.54
1	A	6	LEU	N-CA	5.21	1.52	1.46
1	A	183	VAL	CA-CB	5.18	1.60	1.54
1	B	217	GLU	CA-C	-5.18	1.46	1.52
1	A	27	ASP	N-CA	5.17	1.52	1.46
1	A	150	GLY	N-CA	5.15	1.52	1.45
1	B	76	LYS	N-CA	5.11	1.52	1.46
1	B	68	GLY	CA-C	5.10	1.57	1.52
1	A	178	PHE	C-O	-5.10	1.18	1.24
1	B	176	LEU	N-CA	5.00	1.52	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	47	GLU	CA-C-N	-7.52	112.42	120.47
1	A	47	GLU	C-N-CA	-7.52	112.42	120.47
1	B	160	VAL	CB-CA-C	-6.30	103.91	111.97
1	A	206	ALA	N-CA-C	6.21	118.12	111.36
1	A	95	ASP	N-CA-C	-6.16	105.93	113.50
1	A	149	GLY	CA-C-N	6.10	126.14	120.34
1	A	149	GLY	C-N-CA	6.10	126.14	120.34
1	B	72	HIS	O-C-N	5.91	128.16	122.07
1	A	128	ILE	CB-CA-C	5.91	122.11	111.36
1	A	210	ARG	NE-CZ-NH1	-5.85	115.65	121.50
1	A	139	GLY	N-CA-C	-5.78	107.03	113.79
1	B	218	GLU	CB-CA-C	-5.61	100.03	108.61
1	B	169	TRP	CA-C-N	-5.50	114.01	119.56
1	B	169	TRP	C-N-CA	-5.50	114.01	119.56
1	A	218	GLU	N-CA-C	-5.31	103.38	110.07
1	B	19	GLY	CA-C-O	-5.13	115.46	120.75
1	A	91	VAL	O-C-N	5.13	127.11	121.83
1	B	159	GLY	CA-C-N	5.10	126.99	120.56
1	B	159	GLY	C-N-CA	5.10	126.99	120.56
1	B	85	THR	OG1-CB-CG2	-5.05	99.20	109.30
1	A	164	SER	N-CA-C	5.01	117.39	111.33

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1824	0	1779	28	0
1	B	1824	0	1779	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	17	0	15	3	0
3	B	17	0	15	5	0
4	A	53	0	31	7	0
4	B	53	0	31	0	0
5	A	120	0	0	2	0
5	B	176	0	0	2	0
All	All	4086	0	3650	44	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 6.

All (44) 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:A:232:EWQ:H11B	3:A:232:EWQ:O14	1.78	0.82
3:B:231:EWQ:H11	3:B:231:EWQ:O14	1.85	0.77
1:A:200:ARG:HD2	4:A:233:FAD:N3A	2.04	0.73
1:A:200:ARG:NH1	4:A:233:FAD:N3A	2.39	0.69
1:A:205:ALA:O	1:A:209:GLN:HG3	1.93	0.69
1:A:200:ARG:HB3	4:A:233:FAD:C2A	2.28	0.64
1:A:172:GLN:HE22	1:A:186:PRO:HD3	1.64	0.63
3:A:232:EWQ:O14	3:A:232:EWQ:C11	2.48	0.62
1:B:149:GLY:C	3:B:231:EWQ:H11A	2.26	0.61
1:B:24:VAL:HG11	1:B:204:VAL:CG1	2.31	0.60
1:A:200:ARG:HH11	4:A:233:FAD:C2A	2.15	0.60
3:B:231:EWQ:O14	3:B:231:EWQ:C11	2.48	0.60
1:B:187:GLN:HE21	1:B:210:ARG:HH11	1.50	0.59
1:A:196:SER:O	1:A:198:GLU:O	2.20	0.59
1:A:157:LYS:HG2	1:A:163:ASP:HB2	1.86	0.58
1:B:150:GLY:N	3:B:231:EWQ:H11A	2.19	0.57
1:A:165:ARG:HA	1:A:168:LEU:HD12	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:GLU:O	1:A:200:ARG:N	2.38	0.55
1:A:85:THR:HG21	5:A:288:HOH:O	2.07	0.53
1:B:24:VAL:HG13	5:B:328:HOH:O	2.12	0.49
1:A:200:ARG:HB3	4:A:233:FAD:H2A	1.94	0.49
1:A:187:GLN:HE21	1:A:210:ARG:HH11	1.59	0.49
1:A:4:LYS:HG2	1:A:35:THR:HB	1.94	0.48
1:A:24:VAL:HG13	5:A:350:HOH:O	2.16	0.46
1:A:128:ILE:HA	1:A:129:PRO:HA	1.75	0.46
1:B:157:LYS:HG3	1:B:166:TYR:OH	2.16	0.45
1:A:1:ALA:HA	1:A:3:LYS:H	1.81	0.45
1:B:187:GLN:NE2	1:B:210:ARG:HH11	2.15	0.45
1:A:142:ALA:HB2	1:A:181:PHE:CD1	2.52	0.44
1:A:73:GLU:O	1:A:77:GLN:HG2	2.17	0.44
3:B:231:EWQ:H17B	3:B:231:EWQ:H1	1.62	0.44
1:A:106:PHE:HB3	1:B:170:PRO:HB3	2.00	0.43
1:B:24:VAL:HG11	1:B:204:VAL:HG12	2.00	0.43
1:A:201:LYS:HD3	4:A:233:FAD:N1A	2.34	0.43
1:A:151:THR:OG1	1:A:154:MET:HG3	2.18	0.42
1:B:193:GLU:HB2	5:B:398:HOH:O	2.18	0.42
1:A:177:HIS:ND1	1:A:222:CYS:HB3	2.35	0.42
1:A:104:TYR:HA	4:A:233:FAD:C5X	2.49	0.41
1:B:172:GLN:HE22	1:B:186:PRO:HD3	1.85	0.41
1:A:15:LYS:NZ	1:B:63:GLU:HG2	2.36	0.41
1:B:195:ALA:HB1	1:B:199:GLU:HB2	2.03	0.41
1:A:150:GLY:HA2	3:A:232:EWQ:H11	2.02	0.41
1:A:122:GLN:HB2	1:A:126:PHE:CZ	2.56	0.40
1:A:167:PHE:O	1:A:170:PRO:HD2	2.21	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	228/230 (99%)	215 (94%)	13 (6%)	0	100	100
1	B	228/230 (99%)	221 (97%)	7 (3%)	0	100	100
All	All	456/460 (99%)	436 (96%)	20 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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	194/194 (100%)	189 (97%)	5 (3%)	40	11
1	B	194/194 (100%)	190 (98%)	4 (2%)	47	17
All	All	388/388 (100%)	379 (98%)	9 (2%)	44	14

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	153	GLU
1	A	194	ILE
1	A	197	GLU
1	A	198	GLU
1	A	201	LYS
1	B	128	ILE
1	B	153	GLU
1	B	160	VAL
1	B	230	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	172	GLN
1	A	187	GLN
1	A	212	GLN
1	B	100	GLN

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Mol	Chain	Res	Type
1	B	138	GLN
1	B	172	GLN
1	B	187	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 oligosaccharides 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	EWQ	B	231	-	17,18,18	2.91	6 (35%)	20,26,26	2.03	6 (30%)
4	FAD	B	233	-	58,58,58	1.74	12 (20%)	85,89,89	2.04	25 (29%)
3	EWQ	A	232	-	17,18,18	1.98	3 (17%)	20,26,26	2.31	4 (20%)
4	FAD	A	233	-	58,58,58	1.96	16 (27%)	85,89,89	2.08	23 (27%)

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	EWQ	B	231	-	-	0/4/4/4	0/2/2/2
4	FAD	B	233	-	-	9/34/50/50	0/6/6/6
3	EWQ	A	232	-	-	0/4/4/4	0/2/2/2
4	FAD	A	233	-	-	9/34/50/50	0/6/6/6

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	231	EWQ	C8-C7	7.01	1.44	1.35
3	B	231	EWQ	C2-C3	6.61	1.49	1.40
4	A	233	FAD	P-O3P	6.17	1.66	1.59
4	A	233	FAD	C4X-N5	5.61	1.42	1.30
4	B	233	FAD	P-O3P	5.52	1.65	1.59
4	B	233	FAD	C4X-N5	5.39	1.42	1.30
3	A	232	EWQ	C2-C3	5.21	1.47	1.40
4	A	233	FAD	C10-N1	5.10	1.43	1.33
3	B	231	EWQ	C3-C4	4.82	1.46	1.40
3	A	232	EWQ	C3-C4	4.24	1.45	1.40
4	B	233	FAD	C10-N1	3.87	1.41	1.33
4	B	233	FAD	C2A-N3A	3.48	1.40	1.33
4	A	233	FAD	C6-C7	-3.37	1.35	1.39
3	A	232	EWQ	C8-C7	3.35	1.39	1.35
4	A	233	FAD	C1'-C2'	3.26	1.57	1.52
4	B	233	FAD	C2A-N1A	3.23	1.39	1.33
4	B	233	FAD	PA-O3P	3.12	1.62	1.59
4	A	233	FAD	O2'-C2'	-3.09	1.36	1.43
4	B	233	FAD	C1'-C2'	2.93	1.56	1.52
3	B	231	EWQ	C2-C7	2.80	1.49	1.45
4	A	233	FAD	C2A-N3A	2.76	1.38	1.33
4	B	233	FAD	C9-C8	2.71	1.43	1.39
4	A	233	FAD	P-O2P	-2.70	1.42	1.55
3	B	231	EWQ	O12-C9	2.57	1.29	1.23
4	A	233	FAD	PA-O3P	2.53	1.62	1.59
4	A	233	FAD	C4X-C4	-2.47	1.35	1.44
4	A	233	FAD	C2A-N1A	2.28	1.38	1.33
4	B	233	FAD	P-O1P	-2.24	1.43	1.50
4	A	233	FAD	P-O1P	2.23	1.58	1.50
4	A	233	FAD	C5X-N5	-2.22	1.35	1.39
4	B	233	FAD	C5X-N5	-2.17	1.35	1.39
4	B	233	FAD	C4A-N3A	2.13	1.38	1.34
4	A	233	FAD	C1'-N10	2.13	1.53	1.47
4	A	233	FAD	O4-C4	-2.12	1.19	1.23
3	B	231	EWQ	C13-C7	2.10	1.54	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	233	FAD	C5'-C4'	2.06	1.54	1.51
4	B	233	FAD	C6-C5X	2.02	1.43	1.40

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	232	EWQ	C8-C9-N10	8.88	124.11	115.44
4	B	233	FAD	N3A-C2A-N1A	-6.08	119.38	128.58
3	B	231	EWQ	C8-C9-N10	5.72	121.03	115.44
4	A	233	FAD	N3A-C2A-N1A	-5.58	120.13	128.58
4	A	233	FAD	C5A-C4A-N3A	-5.41	119.27	126.72
4	A	233	FAD	C7M-C7-C6	-4.83	111.06	119.57
4	B	233	FAD	C4X-C10-N10	4.80	123.36	116.48
4	B	233	FAD	N9A-C8A-N7A	-4.51	107.54	113.94
4	B	233	FAD	C5A-N7A-C8A	4.46	110.46	103.45
4	A	233	FAD	C5A-N7A-C8A	4.28	110.17	103.45
4	A	233	FAD	N9A-C8A-N7A	-4.11	108.10	113.94
4	A	233	FAD	C4X-C10-N10	3.97	122.17	116.48
4	A	233	FAD	C2A-N3A-C4A	3.92	121.41	111.83
4	B	233	FAD	C5A-C4A-N3A	-3.91	121.33	126.72
4	B	233	FAD	C8M-C8-C7	3.88	128.68	120.76
4	A	233	FAD	C6-C7-C8	3.56	124.91	119.69
4	A	233	FAD	O4'-C4'-C3'	-3.52	101.02	109.25
4	B	233	FAD	C4-N3-C2	-3.51	119.41	125.64
4	B	233	FAD	C4A-C5A-N7A	-3.49	106.59	110.58
4	A	233	FAD	C4A-C5A-N7A	-3.49	106.60	110.58
4	B	233	FAD	C7M-C7-C6	-3.48	113.45	119.57
4	B	233	FAD	C4-C4X-N5	3.46	122.98	118.21
4	A	233	FAD	N3A-C4A-N9A	3.36	132.87	127.17
4	A	233	FAD	C2B-C1B-N9A	3.30	121.51	113.30
4	B	233	FAD	O4'-C4'-C5'	-3.30	102.72	109.99
4	B	233	FAD	C10-C4X-N5	-3.28	118.12	124.81
3	B	231	EWQ	C5-C6-C1	3.27	125.94	120.98
4	A	233	FAD	C9A-C5X-N5	3.25	125.90	122.45
4	B	233	FAD	C2A-N3A-C4A	3.24	119.75	111.83
4	B	233	FAD	O4B-C1B-N9A	3.14	114.13	108.09
3	B	231	EWQ	C17-O16-C6	-3.13	110.78	117.50
4	B	233	FAD	O4-C4-N3	-3.00	114.47	120.11
4	A	233	FAD	O3B-C3B-C4B	-2.97	102.55	111.08
4	A	233	FAD	C8M-C8-C7	2.92	126.72	120.76
4	B	233	FAD	C8M-C8-C9	-2.90	114.47	119.57
4	B	233	FAD	O4'-C4'-C3'	-2.83	102.63	109.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	233	FAD	N3A-C4A-N9A	2.81	131.94	127.17
4	B	233	FAD	C4A-N9A-C8A	2.81	108.69	105.74
4	A	233	FAD	O2B-C2B-C1B	2.79	119.72	110.10
3	B	231	EWQ	C15-O14-C4	2.78	121.59	117.51
4	B	233	FAD	C4X-C4-N3	2.74	120.24	113.25
3	A	232	EWQ	O12-C9-C8	-2.69	118.98	125.61
3	A	232	EWQ	C9-C8-C7	-2.60	120.23	123.48
4	B	233	FAD	O2A-PA-O3P	2.58	114.24	107.27
4	A	233	FAD	O2-C2-N1	-2.52	117.62	121.80
4	B	233	FAD	O5'-P-O1P	2.49	118.79	108.94
3	A	232	EWQ	C5-C6-C1	2.40	124.61	120.98
4	A	233	FAD	O2B-C2B-C3B	2.39	119.47	111.82
4	A	233	FAD	C9-C8-C7	-2.31	116.30	119.69
3	B	231	EWQ	C13-C7-C2	2.30	122.54	119.94
4	A	233	FAD	O4B-C1B-C2B	-2.29	101.72	106.62
4	B	233	FAD	C6-C7-C8	2.28	123.04	119.69
4	A	233	FAD	O3P-PA-O1A	-2.27	103.88	110.70
4	A	233	FAD	C10-C4X-N5	-2.26	120.19	124.81
3	B	231	EWQ	C2-C7-C8	-2.12	116.61	118.65
4	B	233	FAD	C9-C8-C7	-2.10	116.60	119.69
4	B	233	FAD	C4X-C10-N1	-2.09	119.47	124.59
4	A	233	FAD	C4-N3-C2	-2.00	122.09	125.64

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	233	FAD	C5B-O5B-PA-O1A
4	A	233	FAD	C5B-O5B-PA-O3P
4	A	233	FAD	C2B-C1B-N9A-C8A
4	B	233	FAD	C5B-O5B-PA-O1A
4	B	233	FAD	C5B-O5B-PA-O2A
4	B	233	FAD	C5B-O5B-PA-O3P
4	A	233	FAD	C2B-C1B-N9A-C4A
4	B	233	FAD	C2B-C1B-N9A-C8A
4	A	233	FAD	C5B-O5B-PA-O2A
4	B	233	FAD	C2B-C1B-N9A-C4A
4	B	233	FAD	C4'-C5'-O5'-P
4	A	233	FAD	C4B-C5B-O5B-PA
4	A	233	FAD	C4'-C5'-O5'-P
4	B	233	FAD	O4'-C4'-C5'-O5'
4	A	233	FAD	P-O3P-PA-O1A

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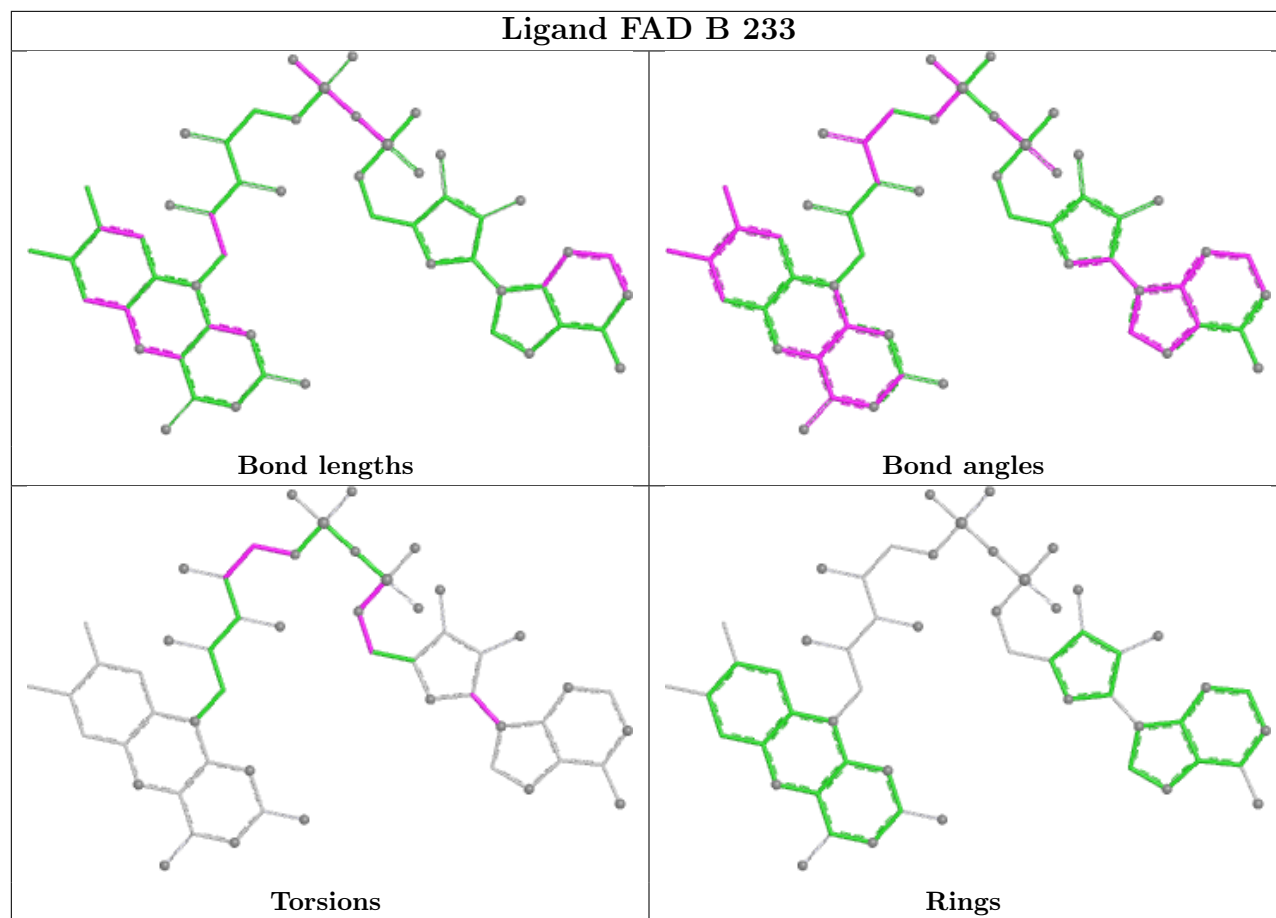
Mol	Chain	Res	Type	Atoms
4	A	233	FAD	P-O3P-PA-O2A
4	B	233	FAD	O4B-C1B-N9A-C8A
4	B	233	FAD	C4B-C5B-O5B-PA

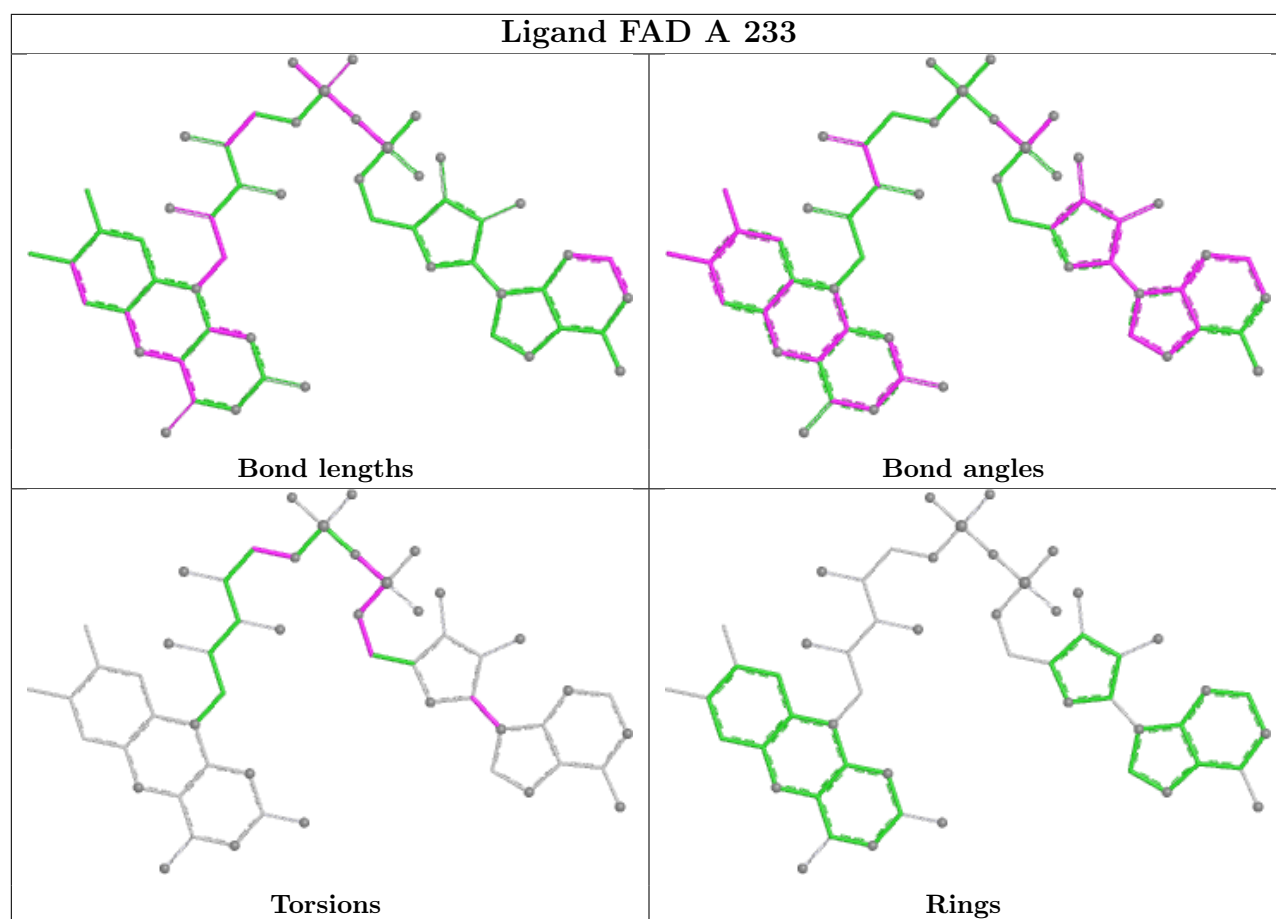
There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	231	EWQ	5	0
3	A	232	EWQ	3	0
4	A	233	FAD	7	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	230/230 (100%)	0.93	47 (20%) 3 4	10, 25, 48, 62	0
1	B	230/230 (100%)	0.22	15 (6%) 25 34	9, 18, 41, 54	0
All	All	460/460 (100%)	0.58	62 (13%) 7 12	9, 20, 45, 62	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	128	ILE	6.1
1	B	1	ALA	5.1
1	A	1	ALA	5.0
1	A	131	PHE	4.8
1	A	194	ILE	4.8
1	B	2	GLY	4.5
1	A	79	SER	4.5
1	A	130	GLY	4.4
1	A	58	THR	4.0
1	A	226	TRP	3.8
1	A	129	PRO	3.6
1	A	132	TYR	3.6
1	B	229	GLY	3.5
1	B	230	GLN	3.4
1	B	194	ILE	3.3
1	A	2	GLY	3.3
1	A	196	SER	3.2
1	A	133	ASP	3.2
1	B	195	ALA	3.2
1	A	59	LEU	3.1
1	B	151	THR	3.0
1	A	61	ASN	2.9
1	A	202	GLY	2.8
1	A	230	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	177	HIS	2.7
1	A	224	ALA	2.7
1	A	155	TYR	2.7
1	A	166	TYR	2.7
1	A	71	THR	2.7
1	A	134	SER	2.7
1	A	197	GLU	2.7
1	A	228	PHE	2.6
1	A	182	LYS	2.6
1	B	160	VAL	2.6
1	A	209	GLN	2.5
1	A	151	THR	2.5
1	A	60	SER	2.4
1	A	135	GLY	2.4
1	A	204	VAL	2.4
1	B	128	ILE	2.4
1	B	157	LYS	2.4
1	A	195	ALA	2.3
1	A	205	ALA	2.3
1	A	206	ALA	2.3
1	B	158	THR	2.3
1	A	190	PHE	2.3
1	A	69	VAL	2.3
1	A	229	GLY	2.2
1	B	159	GLY	2.2
1	A	156	THR	2.2
1	A	125	ALA	2.2
1	B	56	THR	2.2
1	A	183	VAL	2.2
1	B	202	GLY	2.1
1	A	158	THR	2.1
1	A	200	ARG	2.1
1	A	136	LEU	2.1
1	A	127	ASP	2.1
1	B	129	PRO	2.1
1	A	225	HIS	2.0
1	A	218	GLU	2.0
1	A	164	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

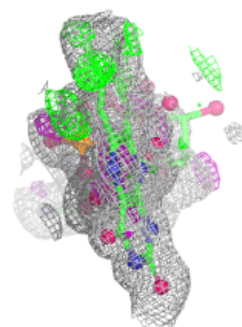
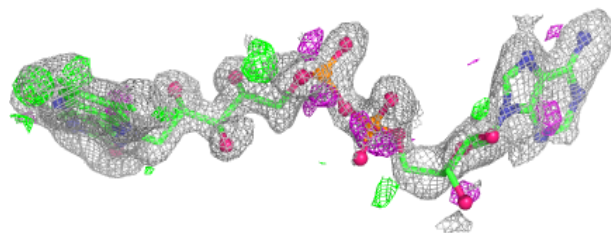
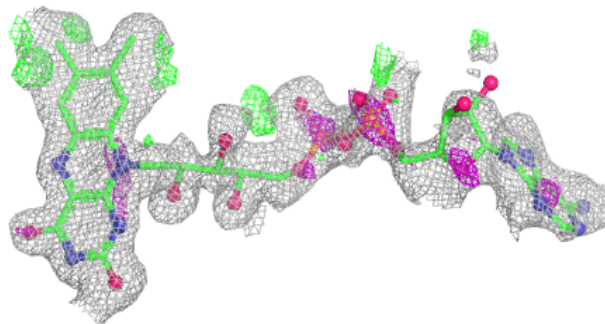
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

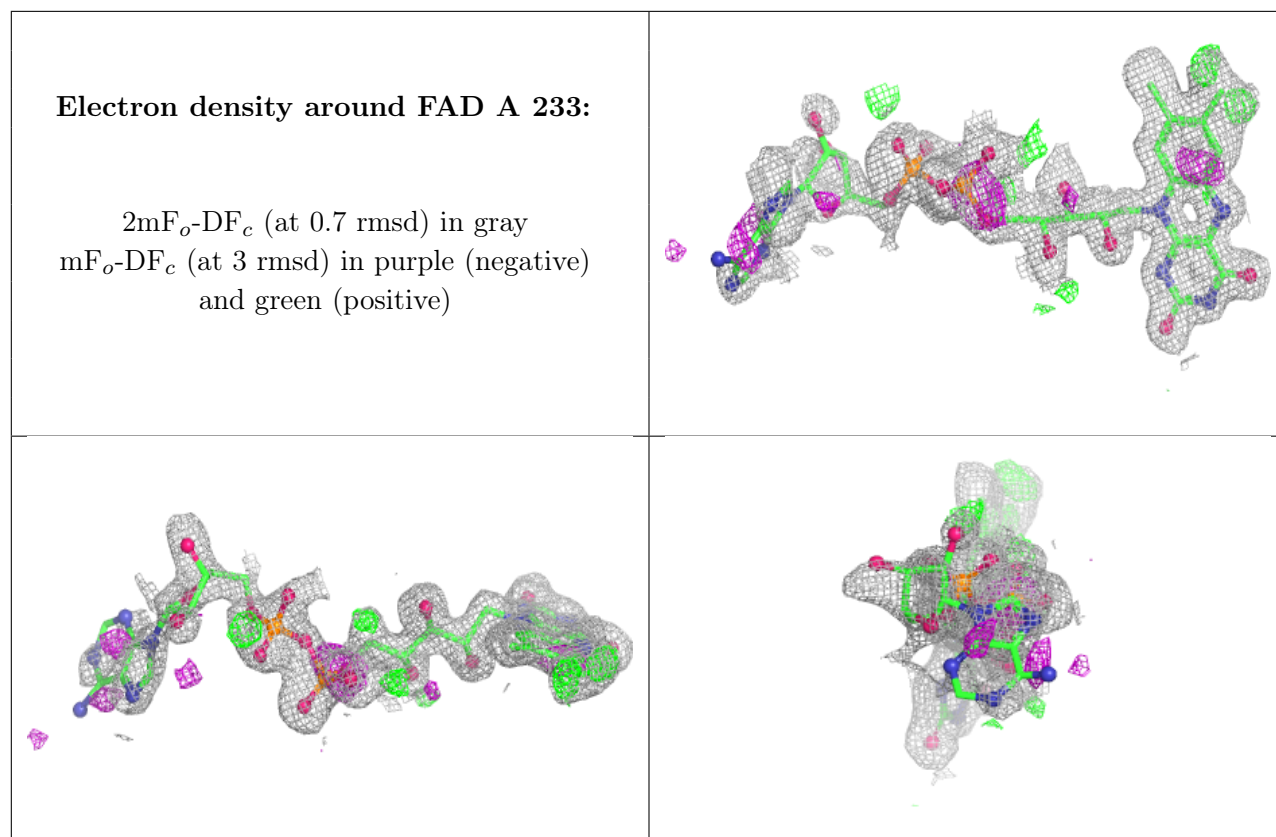
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EWQ	B	231	17/17	0.86	0.12	23,30,36,39	0
2	ZN	A	231	1/1	0.91	0.12	29,29,29,29	0
4	FAD	B	233	53/53	0.91	0.11	16,26,53,57	0
3	EWQ	A	232	17/17	0.92	0.10	18,22,29,33	0
4	FAD	A	233	53/53	0.93	0.11	15,25,58,59	0
2	ZN	B	232	1/1	0.98	0.04	19,19,19,19	0

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

Electron density around FAD B 233:

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.