



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:14 AM UTC

PDB ID : 4DQK / pdb_00004dqk
Title : Crystal structure of the FAD binding domain of cytochrome P450 BM3
Authors : Joyce, M.G.; Leys, D.
Deposited on : 2012-02-16
Resolution : 2.40 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

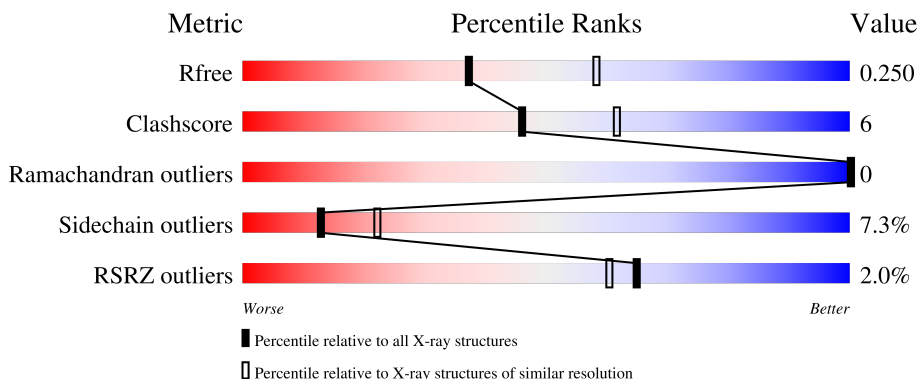
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

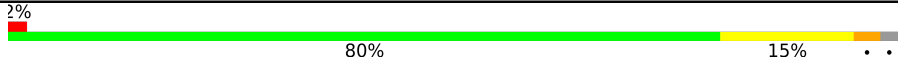
The reported resolution of this entry is 2.40 Å.

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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	391	
1	B	391	

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	PG4	A	1106	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6310 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.

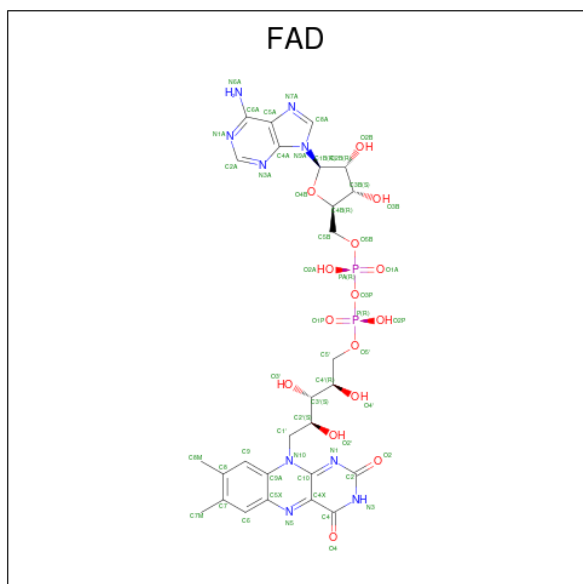
- Molecule 1 is a protein called Bifunctional P-450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	385	Total	C	N	O	S	0	0	0
			2933	1863	504	555	11			
1	B	377	Total	C	N	O	S	0	0	0
			2922	1854	500	557	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	774	ALA	CYS	engineered mutation	UNP P14779
B	774	ALA	CYS	engineered mutation	UNP P14779

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



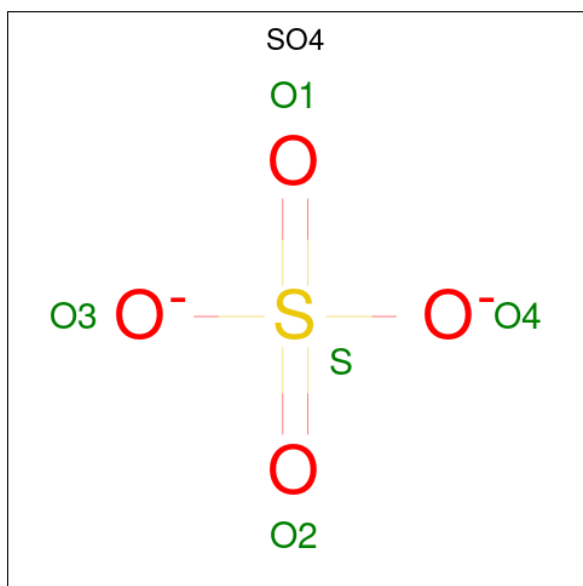
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

Continued on next page...

Continued from previous page...

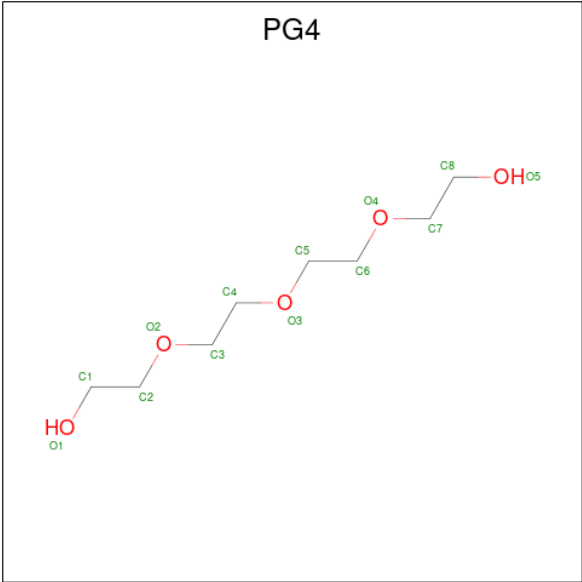
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			13	8	5		
4	B	1	Total	C	O	0	0
			12	8	4		

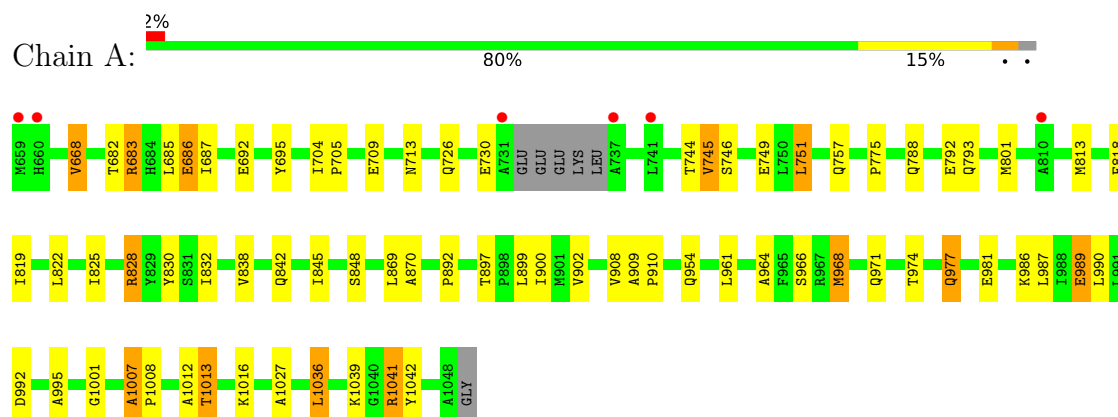
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	137	Total	O	0	0
			137	137		
5	B	137	Total	O	0	0
			137	137		

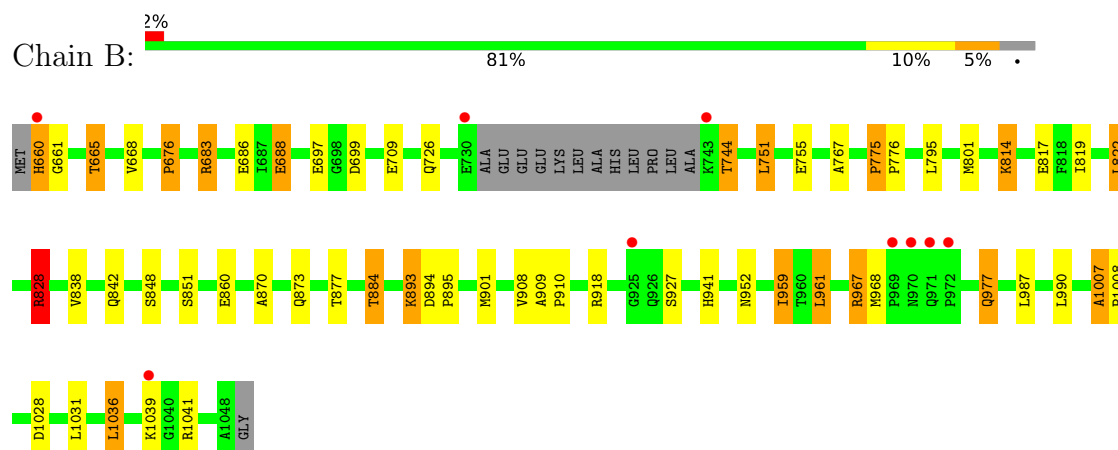
3 Residue-property plots [i](#)

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: Bifunctional P-450/NADPH-P450 reductase



- Molecule 1: Bifunctional P-450/NADPH-P450 reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	191.39Å 191.39Å 74.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.36 – 2.40 29.36 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.36-2.40) 99.8 (29.36-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.213 , 0.256 0.205 , 0.250	Depositor DCC
R_{free} test set	3086 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	44.2	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 29.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6310	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% 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: FAD, SO4, PG4

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	1.56	22/2994 (0.7%)	1.17	8/4067 (0.2%)
1	B	1.54	9/2983 (0.3%)	1.20	8/4044 (0.2%)
All	All	1.55	31/5977 (0.5%)	1.18	16/8111 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1001	GLY	C-O	-10.18	1.18	1.24
1	B	1007	ALA	C-O	-7.36	1.17	1.24
1	A	1012	ALA	CA-CB	7.14	1.64	1.53
1	B	795	LEU	N-CA	6.98	1.54	1.46
1	B	967	ARG	CD-NE	6.61	1.55	1.46
1	A	992	ASP	C-O	-6.24	1.16	1.24
1	A	845	ILE	C-O	-6.15	1.17	1.23
1	A	1007	ALA	CA-CB	-6.03	1.46	1.53
1	B	683	ARG	CD-NE	-5.97	1.37	1.46
1	A	900	ILE	CA-CB	-5.92	1.47	1.54
1	B	1039	LYS	CD-CE	5.92	1.70	1.52
1	A	1027	ALA	CA-CB	5.70	1.62	1.53
1	A	819	ILE	N-CA	-5.66	1.39	1.46
1	A	1013	THR	CA-C	5.58	1.59	1.52
1	A	954	GLN	N-CA	-5.49	1.39	1.46
1	B	901	MET	CA-C	-5.40	1.46	1.52

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	848	SER	N-CA	-5.33	1.39	1.46
1	A	751	LEU	CA-C	5.33	1.60	1.52
1	A	989	GLU	CG-CD	5.33	1.65	1.52
1	A	989	GLU	C-O	-5.30	1.18	1.24
1	A	713	ASN	C-O	-5.30	1.17	1.24
1	B	661	GLY	N-CA	5.25	1.52	1.45
1	A	682	THR	C-O	-5.20	1.18	1.23
1	A	995	ALA	CA-CB	-5.17	1.44	1.53
1	A	902	VAL	CA-CB	-5.17	1.47	1.54
1	A	685	LEU	CA-C	-5.16	1.46	1.52
1	A	961	LEU	CA-C	-5.16	1.46	1.52
1	A	830	TYR	CA-C	5.13	1.58	1.52
1	B	709	GLU	C-O	-5.09	1.18	1.24
1	A	705	PRO	N-CA	-5.06	1.42	1.46
1	A	869	LEU	C-O	-5.03	1.18	1.24

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	676	PRO	CA-C-N	-8.87	109.90	122.86
1	B	676	PRO	C-N-CA	-8.87	109.90	122.86
1	B	959	ILE	CB-CA-C	-7.60	100.68	111.40
1	A	897	THR	CA-C-N	7.38	127.36	119.76
1	A	897	THR	C-N-CA	7.38	127.36	119.76
1	A	1041	ARG	NE-CZ-NH2	6.02	124.62	119.20
1	B	775	PRO	N-CA-C	5.88	117.88	110.70
1	A	813	MET	N-CA-C	-5.76	102.37	110.50
1	A	704	ILE	CB-CA-C	-5.69	104.35	110.95
1	A	989	GLU	N-CA-CB	5.64	118.42	110.12
1	A	775	PRO	N-CA-C	5.62	117.56	110.70
1	A	757	GLN	N-CA-C	5.55	117.55	110.61
1	B	660	HIS	CB-CA-C	-5.46	99.74	110.10
1	B	767	ALA	N-CA-C	-5.45	105.50	111.82
1	B	967	ARG	CB-CG-CD	5.21	123.29	111.30
1	B	828	ARG	CB-CG-CD	5.16	123.18	111.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	660	HIS	Peptide

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	2933	0	2835	38	1
1	B	2922	0	2858	30	1
2	A	53	0	31	1	0
2	B	53	0	31	2	0
3	A	15	0	0	1	0
3	B	15	0	0	0	0
4	A	20	0	27	8	0
4	B	25	0	33	1	0
5	A	137	0	0	4	1
5	B	137	0	0	4	1
All	All	6310	0	5815	69	2

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 (69) 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:981:GLU:OE2	1:A:981:GLU:CB	2.12	0.95
1:A:977:GLN:HE21	1:A:977:GLN:H	1.26	0.82
1:B:726:GLN:OE1	1:B:744:THR:HG22	1.79	0.82
1:B:977:GLN:HE21	1:B:977:GLN:H	1.28	0.79
1:A:745:VAL:HG23	1:A:749:GLU:HG3	1.67	0.75
1:A:683:ARG:HD2	1:A:870:ALA:HA	1.71	0.72
1:A:668:VAL:HG22	1:A:686:GLU:HB3	1.73	0.70
1:A:1036:LEU:HD13	1:A:1042:TYR:HB2	1.74	0.70
1:A:964:ALA:HB1	1:A:974:THR:HG22	1.73	0.70
1:B:726:GLN:HB3	1:B:744:THR:CG2	2.21	0.70
1:B:860:GLU:H	4:B:1105:PG4:H81	1.57	0.70
1:B:941:HIS:HD2	5:B:1284:HOH:O	1.78	0.66
1:A:1016:LYS:HD3	4:A:1106:PG4:H81	1.75	0.66
1:B:918:ARG:NH2	5:B:1320:HOH:O	2.28	0.65
1:A:788:GLN:HE21	1:A:792:GLU:CD	2.04	0.65
1:B:1036:LEU:HD23	1:B:1041:ARG:HB2	1.77	0.65
1:A:1016:LYS:CD	4:A:1106:PG4:H81	2.27	0.65

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1036:LEU:HD22	1:A:1041:ARG:HB2	1.81	0.62
1:B:959:ILE:HD12	1:B:961:LEU:HD13	1.82	0.61
1:A:1013:THR:HA	4:A:1106:PG4:H82	1.85	0.59
1:B:726:GLN:HB3	1:B:744:THR:HG21	1.83	0.59
1:A:801:MET:HE3	1:A:818:PHE:CD1	2.38	0.58
1:B:665:THR:HG21	1:B:688:GLU:O	2.04	0.57
1:A:977:GLN:H	1:A:977:GLN:NE2	2.01	0.56
3:A:1104:SO4:O1	5:A:1329:HOH:O	2.17	0.56
1:A:892:PRO:CD	1:A:899:LEU:HD21	2.36	0.56
1:A:788:GLN:NE2	1:A:792:GLU:OE1	2.38	0.55
1:B:683:ARG:HD3	5:B:1306:HOH:O	2.07	0.55
1:B:683:ARG:HD2	1:B:870:ALA:HA	1.88	0.54
1:A:730:GLU:CB	5:A:1255:HOH:O	2.56	0.54
1:B:683:ARG:CD	5:B:1306:HOH:O	2.56	0.53
1:B:699:ASP:OD1	1:B:884:THR:HB	2.09	0.53
1:A:745:VAL:CG2	1:A:749:GLU:HG3	2.39	0.53
1:A:1016:LYS:HB3	4:A:1106:PG4:H81	1.91	0.53
1:B:814:LYS:HG2	1:B:817:GLU:HB2	1.90	0.52
1:A:683:ARG:HD3	5:A:1298:HOH:O	2.10	0.52
1:B:1028:ASP:HA	1:B:1031:LEU:HD12	1.91	0.52
1:A:1007:ALA:HB3	1:A:1008:PRO:HD3	1.92	0.52
1:B:775:PRO:HB2	1:B:776:PRO:HD3	1.94	0.50
1:B:751:LEU:HD13	1:B:819:ILE:HD11	1.92	0.50
1:B:828:ARG:HG2	2:B:1101:FAD:H3'	1.95	0.49
1:B:977:GLN:H	1:B:977:GLN:NE2	2.05	0.49
1:A:1013:THR:HA	4:A:1106:PG4:H61	1.93	0.49
1:A:726:GLN:OE1	1:A:744:THR:HG22	2.13	0.49
1:A:968:MET:HE3	1:A:971:GLN:HG3	1.94	0.48
1:B:801:MET:CE	1:B:822:LEU:HD13	2.43	0.48
1:B:1036:LEU:CD2	1:B:1041:ARG:HB2	2.43	0.48
1:A:1016:LYS:HB3	4:A:1106:PG4:C8	2.45	0.47
1:A:746:SER:OG	1:A:749:GLU:HG2	2.15	0.46
1:A:828:ARG:HD2	2:A:1101:FAD:O1P	2.16	0.46
1:B:1007:ALA:HB3	1:B:1008:PRO:HD3	1.96	0.46
1:A:801:MET:HE3	1:A:818:PHE:CE1	2.50	0.46
1:B:909:ALA:HB3	1:B:910:PRO:HD3	1.98	0.46
1:A:683:ARG:CD	5:A:1298:HOH:O	2.63	0.45
1:A:692:GLU:H	1:A:692:GLU:CD	2.25	0.44
1:A:1039:LYS:HE3	4:A:1105:PG4:H11	1.98	0.44
1:A:686:GLU:HG2	1:A:842:GLN:OE1	2.19	0.43
1:A:1016:LYS:CG	4:A:1106:PG4:H81	2.48	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:908:VAL:HG22	1:A:908:VAL:O	2.20	0.42
1:A:909:ALA:HB3	1:A:910:PRO:HD3	2.01	0.42
1:B:755:GLU:OE2	1:B:828:ARG:NH2	2.52	0.42
1:B:828:ARG:HD2	2:B:1101:FAD:O1P	2.20	0.41
1:B:894:ASP:HA	1:B:895:PRO:HD3	1.87	0.41
1:B:908:VAL:O	1:B:909:ALA:C	2.63	0.41
1:B:893:LYS:HD2	1:B:893:LYS:H	1.85	0.41
1:A:695:TYR:CE1	1:A:832:ILE:HG21	2.56	0.41
1:A:964:ALA:CB	1:A:974:THR:HG22	2.45	0.41
1:A:971:GLN:HA	1:A:971:GLN:OE1	2.20	0.40
1:B:686:GLU:CG	1:B:842:GLN:NE2	2.85	0.40

All (2) 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)
5:A:1326:HOH:O	5:B:1329:HOH:O[3_665]	2.03	0.17
1:A:793:GLN:OE1	1:B:676:PRO:O[6_665]	2.14	0.06

5.3 Torsion angles

5.3.1 Protein backbone

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	381/391 (97%)	366 (96%)	15 (4%)	0	100	100
1	B	373/391 (95%)	362 (97%)	11 (3%)	0	100	100
All	All	754/782 (96%)	728 (97%)	26 (3%)	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	296/330 (90%)	276 (93%)	20 (7%)	14	25
1	B	305/330 (92%)	281 (92%)	24 (8%)	11	19
All	All	601/660 (91%)	557 (93%)	44 (7%)	13	22

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	668	VAL
1	A	683	ARG
1	A	686	GLU
1	A	687	ILE
1	A	709	GLU
1	A	745	VAL
1	A	751	LEU
1	A	822	LEU
1	A	825	ILE
1	A	828	ARG
1	A	838	VAL
1	A	848	SER
1	A	966	SER
1	A	968	MET
1	A	977	GLN
1	A	986	LYS
1	A	987	LEU
1	A	989	GLU
1	A	990	LEU
1	A	1036	LEU
1	B	665	THR
1	B	668	VAL
1	B	688	GLU
1	B	697	GLU
1	B	744	THR
1	B	751	LEU
1	B	814	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	822	LEU
1	B	828	ARG
1	B	838	VAL
1	B	851	SER
1	B	873	GLN
1	B	877	THR
1	B	884	THR
1	B	893	LYS
1	B	927	SER
1	B	952	ASN
1	B	961	LEU
1	B	967	ARG
1	B	968	MET
1	B	977	GLN
1	B	987	LEU
1	B	990	LEU
1	B	1036	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	666	ASN
1	A	777	HIS
1	A	788	GLN
1	A	977	GLN
1	A	993	GLN
1	B	675	GLN
1	B	842	GLN
1	B	920	GLN
1	B	924	GLN
1	B	977	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

12 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$
4	PG4	B	1105	-	12,12,12	0.76	0	11,11,11	0.80	0
2	FAD	A	1101	-	58,58,58	1.43	7 (12%)	85,89,89	1.79	21 (24%)
3	SO4	B	1102	-	4,4,4	0.44	0	6,6,6	1.27	0
4	PG4	B	1106	-	11,11,12	0.72	0	10,10,11	0.82	0
3	SO4	A	1102	-	4,4,4	0.24	0	6,6,6	1.15	1 (16%)
3	SO4	B	1104	-	4,4,4	0.34	0	6,6,6	0.43	0
4	PG4	A	1106	-	6,6,12	1.48	1 (16%)	5,5,11	1.57	1 (20%)
3	SO4	A	1103	-	4,4,4	0.42	0	6,6,6	0.99	1 (16%)
3	SO4	A	1104	-	4,4,4	0.46	0	6,6,6	0.41	0
2	FAD	B	1101	-	58,58,58	1.29	5 (8%)	85,89,89	1.54	16 (18%)
4	PG4	A	1105	-	12,12,12	0.79	0	11,11,11	1.04	0
3	SO4	B	1103	-	4,4,4	0.30	0	6,6,6	0.37	0

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	PG4	B	1105	-	-	7/10/10/10	-
2	FAD	A	1101	-	-	3/34/50/50	0/6/6/6
4	PG4	B	1106	-	-	4/9/9/10	-
4	PG4	A	1106	-	-	3/4/4/10	-
2	FAD	B	1101	-	-	3/34/50/50	0/6/6/6

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PG4	A	1105	-	-	8/10/10/10	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1101	FAD	C4X-N5	4.16	1.39	1.30
2	A	1101	FAD	C5A-N7A	-4.04	1.31	1.39
2	B	1101	FAD	C10-N1	3.48	1.40	1.33
2	B	1101	FAD	C1'-C2'	3.45	1.57	1.52
2	B	1101	FAD	C4X-N5	3.44	1.38	1.30
2	A	1101	FAD	C1'-C2'	3.11	1.57	1.52
2	A	1101	FAD	C8M-C8	3.10	1.56	1.51
2	B	1101	FAD	C9A-C5X	-2.52	1.37	1.41
2	A	1101	FAD	PA-O3P	2.39	1.62	1.59
2	A	1101	FAD	C10-N1	2.34	1.38	1.33
2	B	1101	FAD	O4B-C4B	-2.21	1.40	1.45
2	A	1101	FAD	C2A-N3A	2.20	1.37	1.33
4	A	1106	PG4	C7-C8	2.10	1.60	1.49

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	FAD	C5A-C4A-N3A	-5.81	118.72	126.72
2	B	1101	FAD	N3A-C2A-N1A	-5.34	120.50	128.58
2	A	1101	FAD	C5A-N7A-C8A	4.10	109.89	103.45
2	A	1101	FAD	C4A-C5A-N7A	-3.73	106.32	110.58
2	A	1101	FAD	C2A-N3A-C4A	3.39	120.11	111.83
2	B	1101	FAD	C5A-C4A-N3A	-3.30	122.17	126.72
2	A	1101	FAD	C5X-C9A-N10	3.23	120.89	117.97
2	B	1101	FAD	C2B-C3B-C4B	3.23	108.85	102.61
2	A	1101	FAD	N3A-C4A-N9A	3.19	132.60	127.17
2	A	1101	FAD	N3A-C2A-N1A	-3.16	123.79	128.58
2	B	1101	FAD	N3A-C4A-N9A	3.02	132.31	127.17
2	A	1101	FAD	C9A-C5X-N5	-3.00	119.27	122.45
2	A	1101	FAD	C4X-C4-N3	2.99	120.87	113.25
2	A	1101	FAD	C4-N3-C2	-2.99	120.34	125.64
2	A	1101	FAD	O2A-PA-O3P	2.93	115.19	107.27
2	B	1101	FAD	C2A-N3A-C4A	2.84	118.77	111.83
2	B	1101	FAD	O3B-C3B-C4B	-2.83	102.95	111.08
2	A	1101	FAD	N9A-C8A-N7A	-2.80	109.96	113.94
2	B	1101	FAD	N9A-C8A-N7A	-2.74	110.05	113.94
2	B	1101	FAD	C5A-N7A-C8A	2.71	107.72	103.45

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	FAD	C6A-C5A-C4A	2.69	120.85	117.18
2	A	1101	FAD	C5A-C4A-N9A	2.61	108.65	105.81
2	A	1101	FAD	O2'-C2'-C3'	-2.60	103.16	109.25
2	A	1101	FAD	O3B-C3B-C4B	-2.54	103.78	111.08
3	A	1102	SO4	O4-S-O3	2.46	122.12	108.54
2	B	1101	FAD	C4X-C4-N3	2.45	119.48	113.25
2	B	1101	FAD	C5'-C4'-C3'	-2.37	107.74	112.22
2	B	1101	FAD	C4X-C10-N10	2.31	119.78	116.48
2	A	1101	FAD	C9A-C9-C8	2.28	123.80	119.22
2	A	1101	FAD	C1'-C2'-C3'	2.24	115.73	109.66
2	B	1101	FAD	C9-C8-C7	-2.24	116.40	119.69
2	A	1101	FAD	O4-C4-C4X	-2.18	120.78	126.53
2	A	1101	FAD	C7M-C7-C8	-2.17	116.32	120.76
2	B	1101	FAD	C2A-N1A-C6A	2.16	122.28	118.73
3	A	1103	SO4	O4-S-O1	2.11	120.58	109.56
2	B	1101	FAD	C9A-C9-C8	2.07	123.39	119.22
2	A	1101	FAD	O3P-PA-O1A	-2.07	104.49	110.70
2	B	1101	FAD	C4A-N9A-C8A	2.06	107.90	105.74
2	B	1101	FAD	C10-C4X-N5	-2.02	120.68	124.81
4	A	1106	PG4	O4-C6-C5	2.01	118.97	110.11

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1105	PG4	C1-C2-O2-C3
4	B	1105	PG4	C6-C5-O3-C4
4	A	1105	PG4	O1-C1-C2-O2
4	A	1106	PG4	O3-C5-C6-O4
4	B	1105	PG4	O4-C7-C8-O5
4	B	1105	PG4	O2-C3-C4-O3
4	B	1105	PG4	O3-C5-C6-O4
4	B	1106	PG4	O3-C5-C6-O4
4	A	1105	PG4	O4-C7-C8-O5
4	B	1106	PG4	O4-C7-C8-O5
4	B	1105	PG4	C4-C3-O2-C2
4	A	1106	PG4	O4-C7-C8-O5
4	A	1106	PG4	C8-C7-O4-C6
2	A	1101	FAD	C2B-C1B-N9A-C8A
4	A	1105	PG4	C8-C7-O4-C6
4	A	1105	PG4	C6-C5-O3-C4
4	A	1105	PG4	C5-C6-O4-C7

Continued on next page...

Continued from previous page...

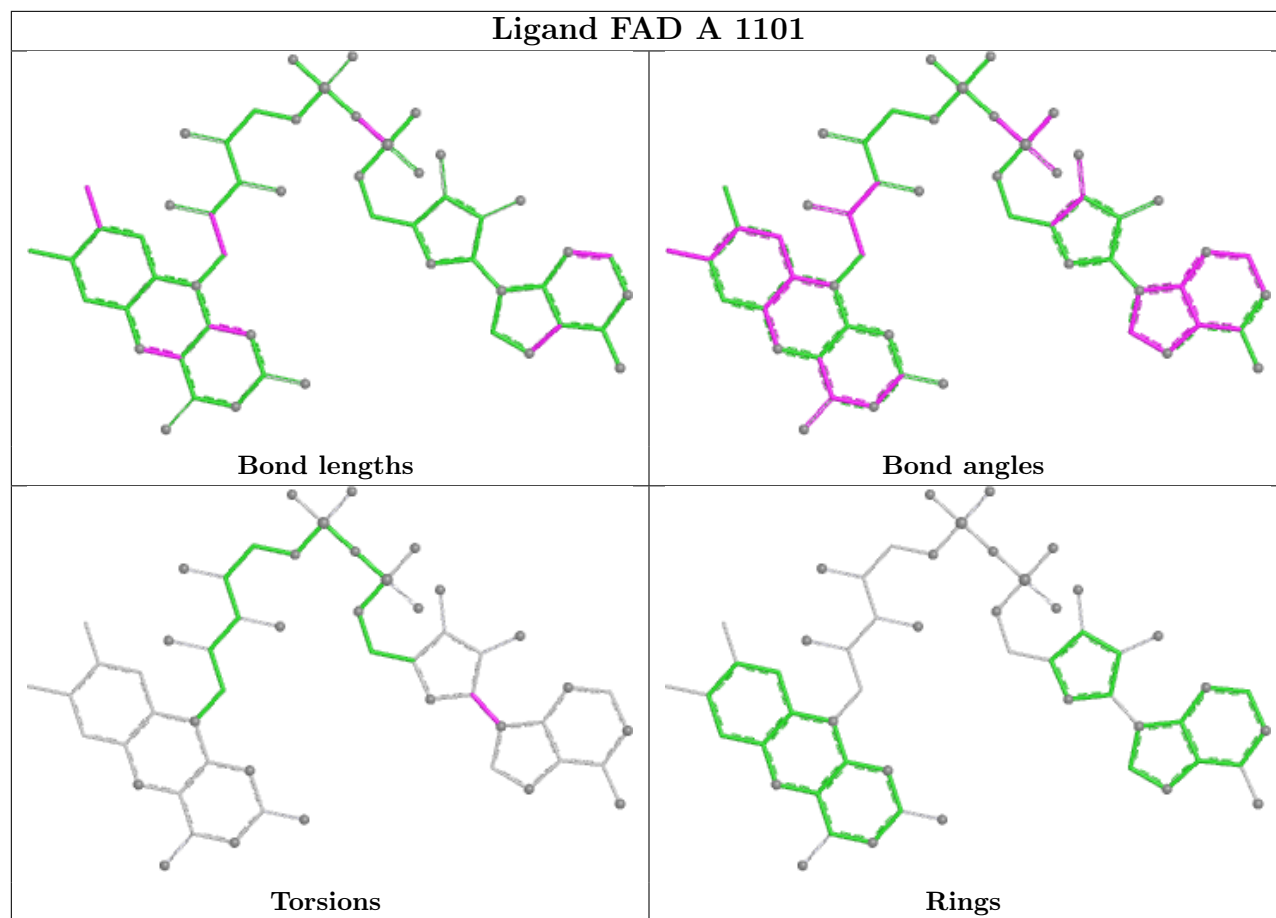
Mol	Chain	Res	Type	Atoms
2	B	1101	FAD	C2B-C1B-N9A-C8A
4	A	1105	PG4	C3-C4-O3-C5
4	B	1106	PG4	C8-C7-O4-C6
2	A	1101	FAD	C2B-C1B-N9A-C4A
4	A	1105	PG4	C4-C3-O2-C2
4	B	1105	PG4	C3-C4-O3-C5
2	B	1101	FAD	O4B-C1B-N9A-C8A
2	A	1101	FAD	O4B-C1B-N9A-C8A
2	B	1101	FAD	C2B-C1B-N9A-C4A
4	B	1106	PG4	C5-C6-O4-C7
4	A	1105	PG4	O3-C5-C6-O4

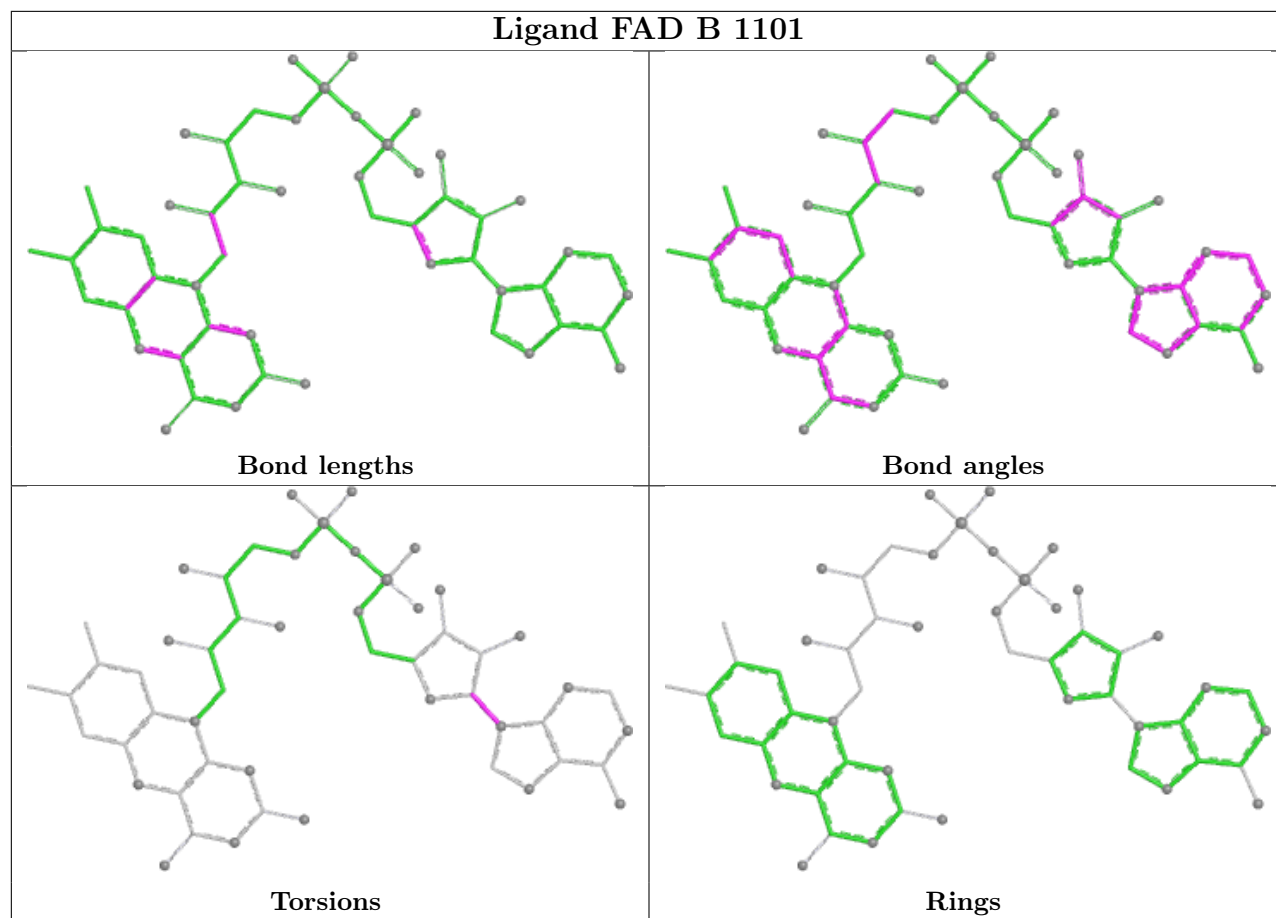
There are no ring outliers.

6 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1105	PG4	1	0
2	A	1101	FAD	1	0
4	A	1106	PG4	7	0
3	A	1104	SO4	1	0
2	B	1101	FAD	2	0
4	A	1105	PG4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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	385/391 (98%)	-0.16	6 (1%) 70 66	22, 35, 59, 78	0
1	B	377/391 (96%)	-0.14	9 (2%) 59 55	20, 36, 59, 79	0
All	All	762/782 (97%)	-0.15	15 (1%) 65 60	20, 36, 59, 79	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	659	MET	4.1
1	A	660	HIS	3.9
1	B	660	HIS	3.8
1	A	731	ALA	3.8
1	A	737	ALA	3.7
1	B	970	ASN	3.7
1	B	969	PRO	3.0
1	B	743	LYS	2.9
1	B	971	GLN	2.8
1	B	972	PRO	2.7
1	A	741	LEU	2.6
1	A	810	ALA	2.5
1	B	925	GLY	2.2
1	B	1039	LYS	2.2
1	B	730	GLU	2.1

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 oligosaccharides in this entry.

6.4 Ligands

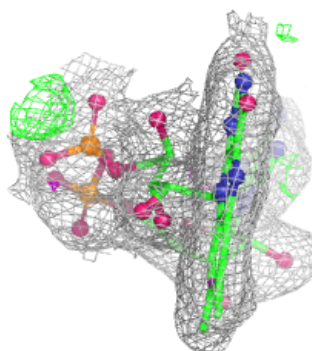
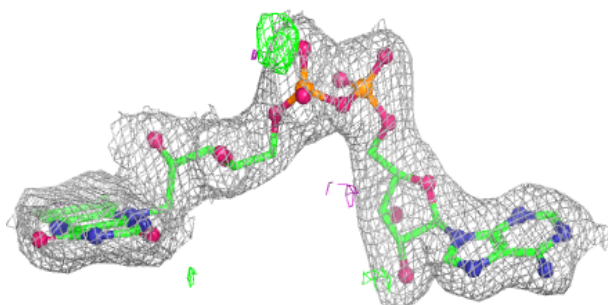
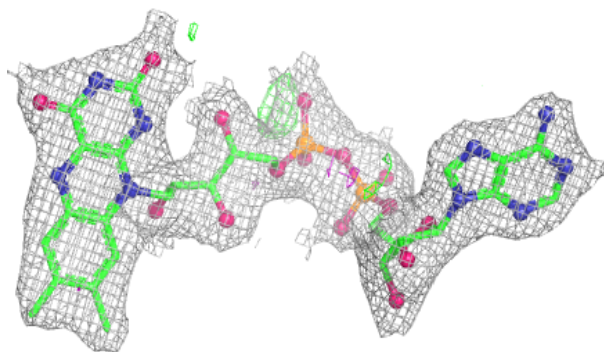
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	SO4	B	1103	5/5	0.75	0.14	90,91,94,95	0
3	SO4	A	1104	5/5	0.79	0.21	80,82,83,84	0
3	SO4	A	1102	5/5	0.79	0.31	78,81,88,89	0
3	SO4	B	1102	5/5	0.82	0.21	72,73,73,76	0
4	PG4	A	1106	7/13	0.82	0.29	37,48,54,56	0
4	PG4	B	1105	13/13	0.84	0.20	76,77,81,81	0
3	SO4	B	1104	5/5	0.86	0.19	84,84,86,87	0
4	PG4	B	1106	12/13	0.86	0.18	52,60,63,65	0
4	PG4	A	1105	13/13	0.88	0.20	49,60,71,75	0
3	SO4	A	1103	5/5	0.89	0.13	57,58,69,69	0
2	FAD	A	1101	53/53	0.98	0.05	18,26,31,33	0
2	FAD	B	1101	53/53	0.98	0.05	13,22,29,31	0

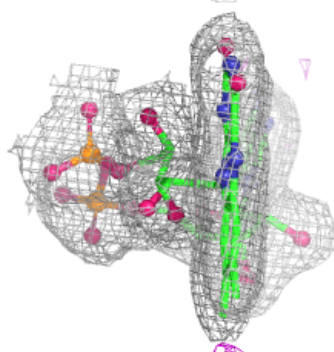
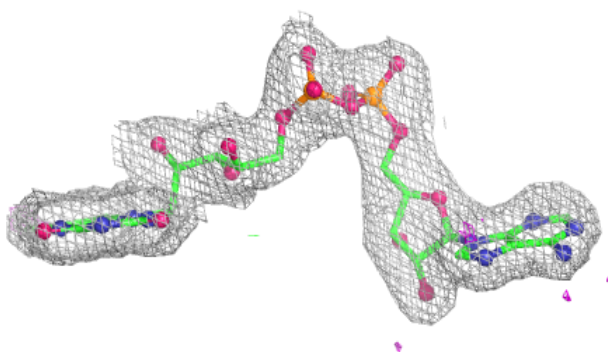
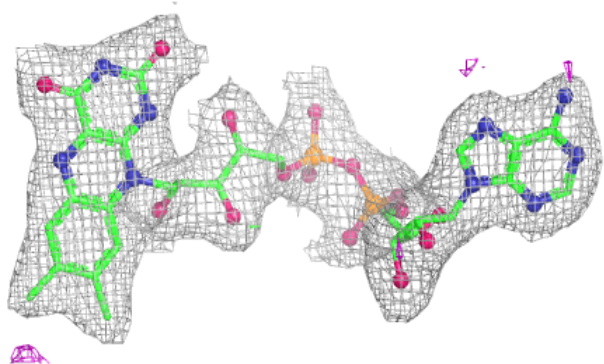
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FAD A 1101:

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

**Electron density around FAD B 1101:**

$2mF_o-DF_c$ (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.