



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 12:57 am BST

PDB ID : 3H8I  
Title : The first X-ray structure of a sulfide:quinone oxidoreductase: Insights into sulfide oxidation mechanism  
Authors : Brito, J.A.; Sousa, F.L.; Stelter, M.; Bandeiras, T.M.; Vonrhein, C.; Teixeira, M.; Pereira, M.M.; Archer, M.  
Deposited on : 2009-04-29  
Resolution : 2.65 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.

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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

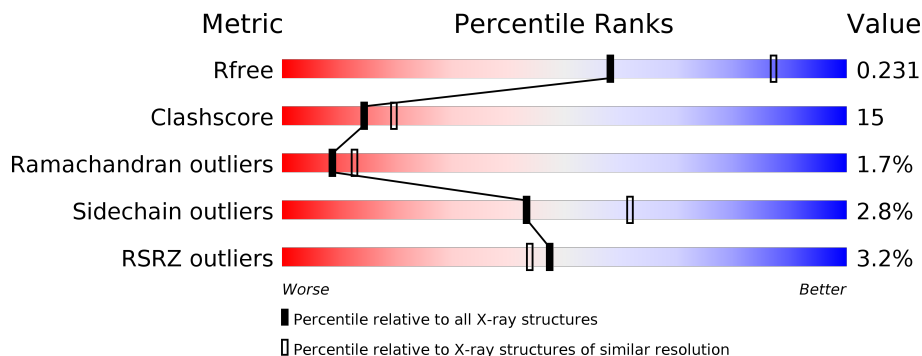
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.65 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	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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 on 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	409	<p>2% 65% 20% 13%</p>
1	B	409	<p>4% 62% 22% 13%</p>

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
3	S3H	A	1002	-	-	X	-

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5916 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.

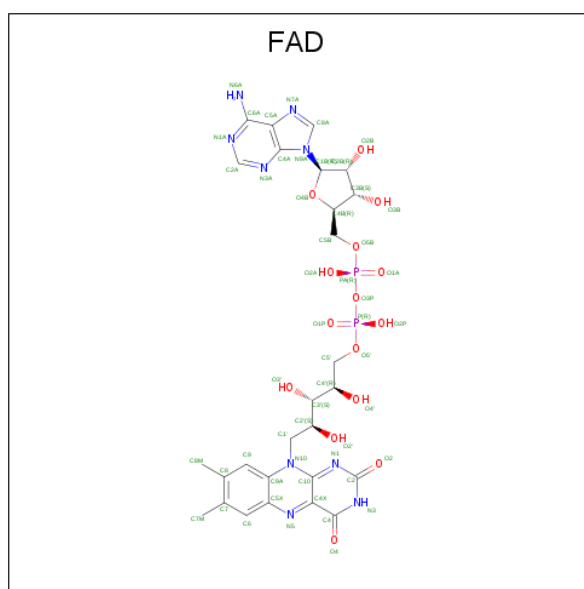
- Molecule 1 is a protein called NADH oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	356	Total 2785	C 1785	N 460	O 527	S 13	0	4	0
1	B	356	Total 2803	C 1796	N 464	O 529	S 14	0	6	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	349	VAL	MET	SEE REMARK 999	UNP Q7ZAG8
B	349	VAL	MET	SEE REMARK 999	UNP Q7ZAG8

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



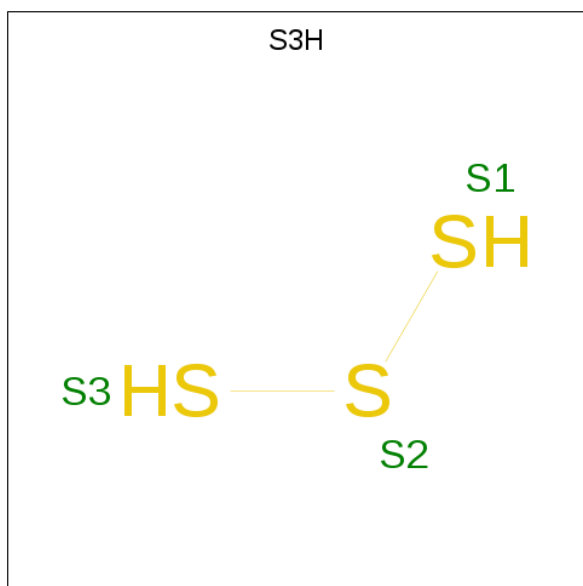
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 53	C 27	N 9	O 15	P 2	0	0

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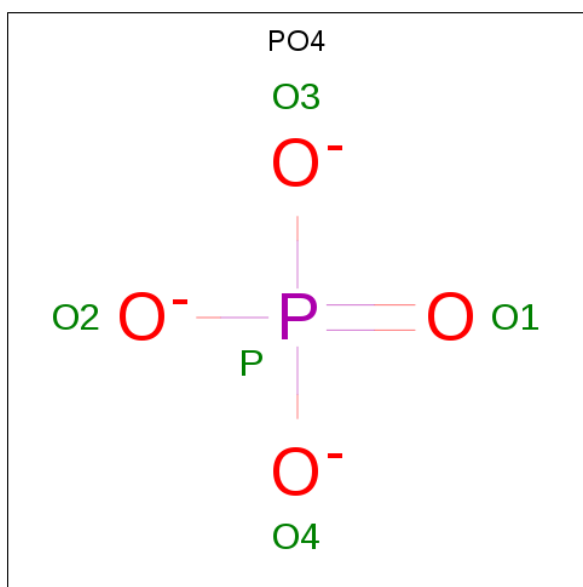
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
2	B	1	53	27	9	15	2	0	0

- Molecule 3 is trisulfane (three-letter code: S3H) (formula: H<sub>2</sub>S<sub>3</sub>).



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

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



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

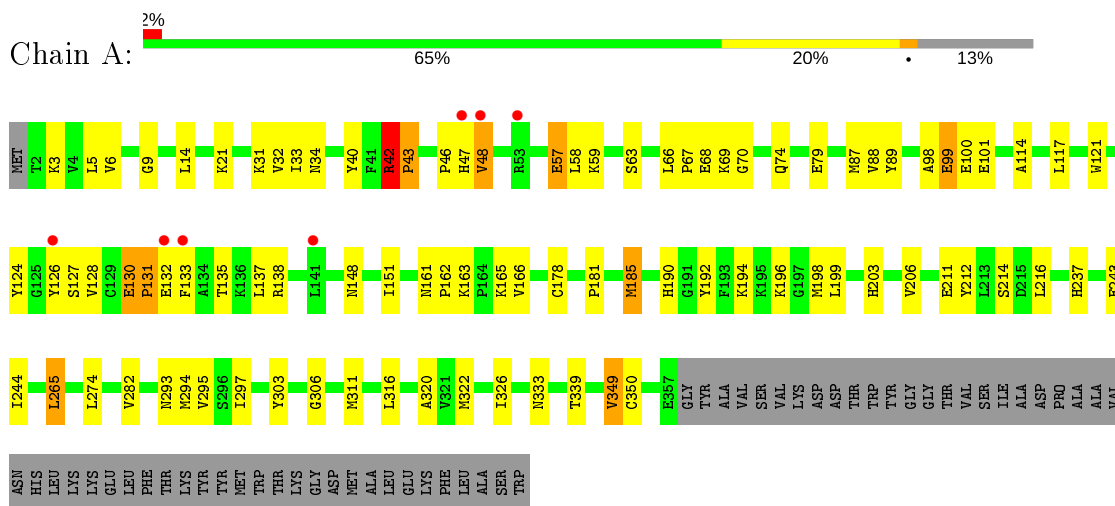
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	104	Total O 104 104	0	0
5	B	67	Total O 67 67	0	0

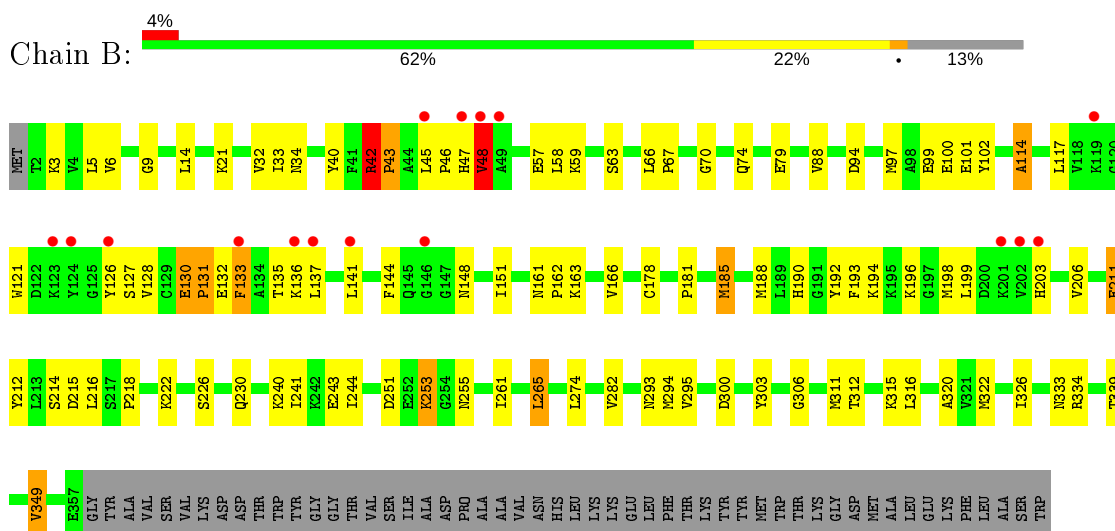
### 3 Residue-property plots [i](#)

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: NADH oxidase



- Molecule 1: NADH oxidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	178.55Å 178.55Å 162.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.66 – 2.65 47.43 – 2.65	Depositor EDS
% Data completeness (in resolution range)	99.0 (26.66-2.65) 99.4 (47.43-2.65)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.98 (at 2.65Å)	Xtrriage
Refinement program	BUSTER-TNT 2.7.0	Depositor
R, $R_{free}$	0.196 , 0.225 0.202 , 0.231	Depositor DCC
$R_{free}$ test set	2236 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	55.7	Xtrriage
Anisotropy	0.053	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 62.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5916	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.78% 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: PO4, S3H, 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	0.38	0/2849	0.60	2/3864 (0.1%)
1	B	0.35	0/2868	0.65	4/3889 (0.1%)
All	All	0.37	0/5717	0.63	6/7753 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	212	TYR	N-CA-CB	-12.15	88.73	110.60
1	A	42	ARG	C-N-CD	-11.55	95.18	120.60
1	B	42	ARG	C-N-CD	-11.46	95.38	120.60
1	B	211	GLU	N-CA-C	-8.98	86.76	111.00
1	B	130	GLU	C-N-CD	-8.55	101.80	120.60
1	A	130	GLU	C-N-CD	-7.63	103.82	120.60

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	2785	0	2782	86	0
1	B	2803	0	2795	92	0
2	A	53	0	28	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	53	0	28	0	0
3	A	3	0	0	2	0
3	B	3	0	0	1	0
4	A	20	0	0	1	0
4	B	25	0	0	2	0
5	A	104	0	0	1	0
5	B	67	0	0	2	0
All	All	5916	0	5633	176	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 15.

All (176) 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:130:GLU:HB2	1:B:133:PHE:HB2	1.35	1.07
1:B:130:GLU:HB2	1:B:133:PHE:CB	2.01	0.90
1:A:126:TYR:HB3	1:A:137:LEU:HD13	1.55	0.89
1:A:130:GLU:HB2	1:A:133:PHE:CB	2.02	0.88
1:A:121:TRP:CH2	1:A:127:SER:HB2	2.11	0.86
1:B:196:LYS:CE	1:B:198:MET:HE2	2.09	0.82
1:B:166:VAL:HG11	1:B:311:MET:HE2	1.61	0.81
1:A:196:LYS:HE2	1:A:198:MET:HE2	1.60	0.80
1:B:21:LYS:HE2	1:B:70:GLY:O	1.84	0.78
1:B:127:SER:HB3	1:B:130:GLU:HG2	1.64	0.78
1:B:121:TRP:CH2	1:B:127:SER:HB2	2.19	0.77
1:B:253:LYS:HB2	1:B:255:ASN:ND2	1.99	0.77
1:A:121:TRP:HH2	1:A:127:SER:HB2	1.49	0.77
1:A:166:VAL:HG11	1:A:311:MET:HE2	1.68	0.76
1:A:89:TYR:CD2	1:A:99:GLU:HG2	2.21	0.76
1:B:166:VAL:HG11	1:B:311:MET:CE	2.16	0.75
1:A:127:SER:HB3	1:A:130:GLU:HG2	1.69	0.75
1:A:166:VAL:HG11	1:A:311:MET:CE	2.16	0.74
1:B:194:LYS:HG3	1:B:199:LEU:HD11	1.67	0.74
1:A:130:GLU:HB2	1:A:133:PHE:HB2	1.70	0.73
1:A:194:LYS:HG3	1:A:199:LEU:HD11	1.68	0.73
1:A:127:SER:HB3	1:A:130:GLU:CG	2.22	0.70
1:A:21:LYS:HE2	1:A:70:GLY:O	1.91	0.69
1:B:196:LYS:HE2	1:B:198:MET:HE2	1.74	0.68
1:A:128:VAL:HG11	1:A:265:LEU:HD12	1.76	0.68
1:B:126:TYR:CE2	1:B:136:LYS:HE2	2.30	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:127:SER:CB	1:B:130:GLU:HG2	2.25	0.66
1:B:121:TRP:HH2	1:B:127:SER:HB2	1.58	0.65
1:A:130:GLU:HB2	1:A:133:PHE:HB3	1.78	0.65
1:B:211:GLU:O	1:B:211:GLU:HG2	1.96	0.64
1:A:127:SER:CB	1:A:130:GLU:HG2	2.28	0.64
1:A:40:TYR:HE2	1:A:59:LYS:HD2	1.63	0.64
1:A:126:TYR:HB3	1:A:137:LEU:CD1	2.25	0.63
1:B:196:LYS:HD3	1:B:198:MET:CE	2.28	0.63
1:B:130:GLU:CB	1:B:133:PHE:HB2	2.22	0.62
1:A:98:ALA:O	1:A:99:GLU:HG3	2.00	0.62
1:B:211:GLU:O	1:B:211:GLU:CG	2.48	0.62
1:A:196:LYS:CE	1:A:198:MET:HE2	2.30	0.61
1:B:181:PRO:O	1:B:185[B]:MET:HG3	1.99	0.61
1:A:63:SER:O	1:A:67:PRO:HG2	1.99	0.61
1:B:42:ARG:HB3	1:B:43:PRO:HD2	1.83	0.61
1:A:98:ALA:C	1:A:99:GLU:HG3	2.20	0.61
1:B:333:ASN:ND2	1:B:339:THR:O	2.28	0.60
1:A:121:TRP:CZ2	1:A:127:SER:HB2	2.36	0.60
1:B:40:TYR:HE2	1:B:59:LYS:HD2	1.66	0.59
1:A:181:PRO:O	1:A:185[B]:MET:HG3	2.02	0.59
1:A:333:ASN:ND2	1:A:339:THR:O	2.29	0.59
1:B:66:LEU:HB2	1:B:67:PRO:HD3	1.84	0.58
1:A:42:ARG:HB3	1:A:43:PRO:HD2	1.86	0.57
1:B:196:LYS:CD	1:B:198:MET:HE2	2.34	0.57
1:B:46:PRO:HG2	1:B:185[A]:MET:HE3	1.85	0.57
1:B:48:VAL:HG23	1:B:188:MET:HE3	1.85	0.57
1:B:63:SER:O	1:B:67:PRO:HG2	2.04	0.57
1:B:99:GLU:HG2	5:B:428:HOH:O	2.06	0.56
1:A:350[A]:CYS:SG	3:A:1002:S3H:S3	2.39	0.56
1:B:300:ASP:O	1:B:334:ARG:NH2	2.38	0.55
1:A:66:LEU:HB2	1:A:67:PRO:HD3	1.89	0.55
1:A:46:PRO:HG2	1:A:185[A]:MET:HE3	1.88	0.55
1:A:128:VAL:HG21	1:A:185[B]:MET:SD	2.47	0.55
1:A:196:LYS:HB3	1:A:198:MET:HE2	1.89	0.55
1:B:46:PRO:HG2	1:B:185[A]:MET:CE	2.37	0.55
1:B:132:GLU:OE1	1:B:132:GLU:HA	2.07	0.55
1:A:40:TYR:CE2	1:A:59:LYS:HD2	2.42	0.54
1:B:130:GLU:HB2	1:B:133:PHE:CG	2.42	0.54
1:B:45:LEU:HD22	1:B:185[B]:MET:CG	2.37	0.54
1:B:46:PRO:CG	1:B:185[A]:MET:HE3	2.38	0.54
1:B:114:ALA:HB1	1:B:117:LEU:HG	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:TRP:CZ2	1:B:127:SER:HB2	2.42	0.54
1:B:128:VAL:HG21	1:B:185[B]:MET:SD	2.47	0.54
1:B:33:ILE:HA	1:B:74:GLN:O	2.08	0.54
1:A:89:TYR:CE2	1:A:99:GLU:HG2	2.43	0.53
1:B:193:PHE:HD1	1:B:198:MET:HE3	1.74	0.53
1:A:161:ASN:HD21	1:A:163:LYS:HE3	1.73	0.53
1:A:33:ILE:HA	1:A:74:GLN:O	2.09	0.53
1:B:196:LYS:HD3	1:B:198:MET:HE2	1.90	0.53
1:A:128:VAL:HG11	1:A:265:LEU:CD1	2.38	0.53
1:A:166:VAL:HG11	1:A:311:MET:HE1	1.89	0.52
1:A:46:PRO:HG2	1:A:185[A]:MET:CE	2.39	0.52
1:A:130:GLU:HB3	1:A:131:PRO:HD2	1.91	0.52
1:A:306:GLY:HA2	1:A:320:ALA:HA	1.92	0.52
1:A:148:ASN:HD22	1:A:203:HIS:HB2	1.74	0.52
1:B:58:LEU:HD12	1:B:59:LYS:N	2.25	0.52
1:A:46:PRO:CG	1:A:185[A]:MET:HE3	2.40	0.51
1:A:58:LEU:HD12	1:A:59:LYS:N	2.25	0.51
1:B:40:TYR:CE2	1:B:59:LYS:HD2	2.44	0.51
1:B:192:TYR:CZ	1:B:196:LYS:HD2	2.46	0.51
1:B:151:ILE:HG23	1:B:206:VAL:HG22	1.93	0.51
1:B:306:GLY:HA2	1:B:320:ALA:HA	1.93	0.51
1:B:42:ARG:HB3	1:B:43:PRO:CD	2.41	0.50
1:B:3:LYS:NZ	1:B:100:GLU:OE1	2.27	0.50
1:B:127:SER:HB3	1:B:130:GLU:CG	2.39	0.50
1:A:190:HIS:HE1	4:A:2002:PO4:O2	1.94	0.50
1:B:148:ASN:HD22	1:B:203:HIS:HB2	1.75	0.50
1:B:293:ASN:O	1:B:294:MET:HB2	2.11	0.50
1:A:132:GLU:OE1	1:A:132:GLU:HA	2.11	0.50
1:A:151:ILE:HG23	1:A:206:VAL:HG22	1.94	0.50
1:A:293:ASN:O	1:A:294:MET:HB2	2.12	0.50
1:A:161:ASN:HA	1:A:162:PRO:C	2.32	0.49
1:B:243:GLU:HG2	1:B:244:ILE:N	2.27	0.49
1:A:48:VAL:O	1:A:48:VAL:HG13	2.12	0.49
1:B:137:LEU:HG	1:B:141:LEU:CD1	2.42	0.49
1:A:178:CYS:O	1:A:265:LEU:HD22	2.12	0.49
1:B:240:LYS:O	1:B:251:ASP:HB2	2.11	0.49
1:B:130:GLU:HB3	1:B:131:PRO:HD2	1.95	0.49
1:B:161:ASN:HA	1:B:162:PRO:C	2.33	0.49
1:A:130:GLU:CB	1:A:131:PRO:HD2	2.43	0.48
1:B:3:LYS:HE2	1:B:5:LEU:HD21	1.95	0.48
1:A:127:SER:HB3	1:A:130:GLU:HG3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:THR:HG23	1:A:138:ARG:HH22	1.78	0.48
1:A:42:ARG:HB3	1:A:43:PRO:CD	2.42	0.48
1:B:48:VAL:HG23	1:B:188:MET:CE	2.44	0.48
1:B:128:VAL:CG1	1:B:265:LEU:HG	2.44	0.48
1:A:42:ARG:HG2	1:A:57:GLU:HB2	1.96	0.48
1:B:132:GLU:O	1:B:135:THR:HB	2.14	0.48
1:A:114:ALA:HB1	1:A:117:LEU:HG	1.96	0.47
1:B:94:ASP:OD2	4:B:411:PO4:O3	2.33	0.47
1:A:322:MET:O	1:A:326:ILE:HD12	2.14	0.47
1:A:126:TYR:HA	1:A:133:PHE:HE1	1.80	0.47
1:A:79:GLU:HG3	5:A:451:HOH:O	2.14	0.47
1:B:166:VAL:HG11	1:B:311:MET:HE1	1.96	0.47
1:B:253:LYS:HB2	1:B:255:ASN:HD22	1.75	0.47
1:B:322:MET:O	1:B:326:ILE:HD12	2.14	0.46
1:B:130:GLU:HB2	1:B:133:PHE:CD2	2.51	0.46
1:A:88:VAL:O	1:A:99:GLU:HA	2.16	0.46
1:B:45:LEU:HD22	1:B:185[B]:MET:HG2	1.97	0.46
1:B:144:PHE:CD1	1:B:261:ILE:HD11	2.51	0.45
1:B:316:LEU:HD21	3:B:1002:S3H:S3	2.55	0.45
1:B:45:LEU:HD22	1:B:185[B]:MET:SD	2.55	0.45
1:A:349:VAL:HG13	1:B:216:LEU:CD2	2.45	0.45
1:A:135:THR:HG23	1:A:138:ARG:NH2	2.31	0.45
1:A:3:LYS:HD3	1:A:101:GLU:O	2.16	0.45
1:B:211:GLU:HG2	1:B:214:SER:HB3	1.98	0.44
1:A:161:ASN:ND2	1:A:163:LYS:HE3	2.33	0.44
1:B:196:LYS:NZ	1:B:198:MET:HE2	2.33	0.44
1:A:274:LEU:O	1:A:282:VAL:HG11	2.18	0.43
1:A:316:LEU:HD21	3:A:1002:S3H:S3	2.59	0.43
1:B:215:ASP:N	1:B:215:ASP:OD1	2.46	0.43
1:A:67:PRO:O	1:A:69:LYS:N	2.52	0.43
1:B:57:GLU:O	1:B:57:GLU:HG3	2.19	0.43
1:B:190:HIS:HE1	4:B:2002:PO4:O2	2.02	0.43
1:B:312:THR:O	1:B:315:LYS:HD3	2.19	0.43
1:A:151:ILE:O	1:A:206:VAL:HA	2.19	0.42
1:B:88:VAL:HG23	1:B:102:TYR:CD2	2.54	0.42
1:B:42:ARG:CB	1:B:43:PRO:CD	2.96	0.42
1:A:196:LYS:CE	1:A:198:MET:CE	2.97	0.42
1:A:216:LEU:CD2	1:B:349:VAL:HG13	2.49	0.42
1:A:42:ARG:CB	1:A:43:PRO:CD	2.97	0.42
1:A:130:GLU:CB	1:A:133:PHE:HB2	2.47	0.42
1:A:165:LYS:HD2	1:A:297:ILE:HG22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:GLY:N	1:A:34:ASN:OD1	2.52	0.42
1:A:3:LYS:HE2	1:A:5:LEU:HD21	2.02	0.42
1:B:3:LYS:HD3	1:B:101:GLU:O	2.20	0.42
1:A:6:VAL:HB	1:A:32:VAL:HG22	2.01	0.42
1:B:194:LYS:HG3	1:B:199:LEU:CD1	2.43	0.42
1:A:211:GLU:HG2	1:A:214:SER:HB3	2.02	0.42
1:A:295:VAL:HG22	1:A:303:TYR:CE1	2.55	0.42
1:A:127:SER:OG	1:A:130:GLU:HG2	2.19	0.42
1:A:194:LYS:HG3	1:A:199:LEU:CD1	2.45	0.42
1:B:130:GLU:O	1:B:133:PHE:HB2	2.20	0.41
1:B:161:ASN:HD21	1:B:163:LYS:HE3	1.86	0.41
1:B:194:LYS:HA	1:B:199:LEU:HG	2.02	0.41
1:A:31:LYS:NZ	1:A:100:GLU:OE2	2.54	0.41
1:A:211:GLU:O	1:A:237:HIS:HD2	2.03	0.41
1:B:311:MET:HE3	5:B:432:HOH:O	2.19	0.41
1:B:226:SER:O	1:B:230:GLN:HG2	2.21	0.41
1:A:48:VAL:HG22	1:A:192:TYR:CD1	2.56	0.41
1:B:274:LEU:O	1:B:282:VAL:HG11	2.20	0.41
1:B:9:GLY:N	1:B:34:ASN:OD1	2.54	0.41
1:A:243:GLU:HG2	1:A:244:ILE:N	2.35	0.41
1:B:218:PRO:O	1:B:222:LYS:HG3	2.21	0.41
1:B:295:VAL:HG22	1:B:303:TYR:CE1	2.56	0.41
1:A:196:LYS:HB2	1:A:198:MET:HE3	2.02	0.41
1:A:87:MET:CE	1:A:99:GLU:HB2	2.51	0.41
1:B:6:VAL:HB	1:B:32:VAL:HG22	2.02	0.40
1:B:79:GLU:HG3	1:B:97[B]:MET:HE1	2.02	0.40
1:B:151:ILE:O	1:B:206:VAL:HA	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	358/409 (88%)	335 (94%)	17 (5%)	6 (2%)	9	13
1	B	360/409 (88%)	338 (94%)	16 (4%)	6 (2%)	9	13
All	All	718/818 (88%)	673 (94%)	33 (5%)	12 (2%)	9	13

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	ARG
1	A	131	PRO
1	B	42	ARG
1	B	131	PRO
1	A	68	GLU
1	B	114	ALA
1	B	253	LYS
1	A	212	TYR
1	A	43	PRO
1	B	48	VAL
1	A	48	VAL
1	B	43	PRO

### 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	304/343 (89%)	295 (97%)	9 (3%)	41	59
1	B	306/343 (89%)	296 (97%)	10 (3%)	38	54
All	All	610/686 (89%)	591 (97%)	19 (3%)	43	57

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	47	HIS
1	A	57	GLU
1	A	99	GLU

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Mol	Chain	Res	Type
1	A	124	TYR
1	A	185[A]	MET
1	A	185[B]	MET
1	A	265	LEU
1	A	349	VAL
1	B	14	LEU
1	B	47	HIS
1	B	48	VAL
1	B	133	PHE
1	B	178	CYS
1	B	185[A]	MET
1	B	185[B]	MET
1	B	241	ILE
1	B	265	LEU
1	B	349	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	ASN
1	A	161	ASN
1	A	190	HIS
1	A	237	HIS
1	B	148	ASN
1	B	161	ASN
1	B	190	HIS
1	B	255	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

13 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	FAD	A	1001	1	51,58,58	2.28	8 (15%)	60,89,89	2.72	15 (25%)
4	PO4	A	2001	-	4,4,4	2.06	1 (25%)	6,6,6	0.50	0
4	PO4	B	411	-	4,4,4	2.37	2 (50%)	6,6,6	0.78	0
4	PO4	A	410	-	4,4,4	2.20	2 (50%)	6,6,6	1.00	0
3	S3H	B	1002	1	0,2,2	0.00	-	0,1,1	0.00	-
3	S3H	A	1002	1	0,2,2	0.00	-	0,1,1	0.00	-
4	PO4	B	2002	-	4,4,4	2.35	2 (50%)	6,6,6	1.11	1 (16%)
4	PO4	B	2001	-	4,4,4	2.17	1 (25%)	6,6,6	0.56	0
4	PO4	B	410	-	4,4,4	2.29	1 (25%)	6,6,6	1.36	1 (16%)
2	FAD	B	1001	1	51,58,58	2.21	7 (13%)	60,89,89	2.90	15 (25%)
4	PO4	B	2003	-	4,4,4	2.09	1 (25%)	6,6,6	0.55	0
4	PO4	A	2002	-	4,4,4	2.31	1 (25%)	6,6,6	0.78	0
4	PO4	A	2003	-	4,4,4	2.35	1 (25%)	6,6,6	0.81	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
2	FAD	A	1001	1	-	6/30/50/50	0/6/6/6
2	FAD	B	1001	1	-	6/30/50/50	0/6/6/6

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	FAD	C4X-C10	12.40	1.51	1.38
2	B	1001	FAD	C4X-C10	11.78	1.50	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1001	FAD	C4-C4X	5.36	1.50	1.41
2	B	1001	FAD	C4-N3	5.08	1.41	1.33
2	A	1001	FAD	C4-N3	4.76	1.41	1.33
2	A	1001	FAD	C4-C4X	4.65	1.49	1.41
4	B	411	PO4	P-O1	3.91	1.60	1.50
4	A	2002	PO4	P-O1	3.76	1.59	1.50
4	B	410	PO4	P-O1	3.72	1.59	1.50
4	A	2003	PO4	P-O1	3.68	1.59	1.50
4	A	410	PO4	P-O1	3.67	1.59	1.50
4	B	2001	PO4	P-O1	3.54	1.59	1.50
4	B	2002	PO4	P-O1	3.48	1.59	1.50
4	A	2001	PO4	P-O1	3.45	1.58	1.50
4	B	2003	PO4	P-O1	3.41	1.58	1.50
2	B	1001	FAD	C10-N1	-3.20	1.29	1.33
2	B	1001	FAD	C9A-N10	2.99	1.42	1.38
2	A	1001	FAD	C9A-N10	2.73	1.42	1.38
2	A	1001	FAD	C5X-N5	2.72	1.39	1.35
2	A	1001	FAD	C9-C8	2.40	1.43	1.37
2	A	1001	FAD	C10-N1	-2.38	1.30	1.33
2	B	1001	FAD	C9-C8	2.37	1.43	1.37
2	B	1001	FAD	C8A-N7A	-2.25	1.30	1.34
2	A	1001	FAD	O4B-C1B	2.22	1.44	1.41
4	B	2002	PO4	P-O4	2.15	1.61	1.54
4	B	411	PO4	P-O2	2.13	1.61	1.54
4	A	410	PO4	P-O3	2.12	1.61	1.54

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1001	FAD	C4-N3-C2	14.66	127.52	115.14
2	A	1001	FAD	C4-N3-C2	12.94	126.06	115.14
2	B	1001	FAD	C4X-C4-N3	-7.69	112.91	123.43
2	A	1001	FAD	C10-C4X-N5	6.52	125.77	121.26
2	A	1001	FAD	C4X-C4-N3	-6.46	114.59	123.43
2	B	1001	FAD	C10-C4X-N5	6.12	125.49	121.26
2	A	1001	FAD	C4-C4X-C10	-6.06	115.94	119.95
2	A	1001	FAD	C1'-N10-C9A	5.83	122.88	118.29
2	B	1001	FAD	C4-C4X-C10	-5.54	116.28	119.95
2	B	1001	FAD	C1'-N10-C9A	4.91	122.16	118.29
2	B	1001	FAD	C5X-C9A-N10	4.81	121.20	117.72
2	A	1001	FAD	C5X-C9A-N10	4.21	120.77	117.72
2	B	1001	FAD	C4X-C10-N10	-4.17	116.01	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	FAD	C4X-C10-N10	-3.82	116.38	120.30
2	B	1001	FAD	C6-C5X-C9A	3.46	123.59	119.05
2	B	1001	FAD	C6-C5X-N5	-3.09	115.64	119.05
2	B	1001	FAD	C4'-C3'-C2'	-2.74	107.67	113.36
2	A	1001	FAD	C6-C5X-N5	-2.71	116.06	119.05
2	A	1001	FAD	O3'-C3'-C4'	2.69	115.30	108.81
2	A	1001	FAD	C4'-C3'-C2'	-2.62	107.91	113.36
2	B	1001	FAD	C4A-C5A-N7A	2.61	112.12	109.40
2	B	1001	FAD	C1'-C2'-C3'	-2.58	102.58	109.79
4	B	410	PO4	O2-P-O1	-2.54	101.58	110.89
2	A	1001	FAD	C6-C5X-C9A	2.42	122.23	119.05
2	A	1001	FAD	C1'-C2'-C3'	-2.40	103.07	109.79
2	B	1001	FAD	O3'-C3'-C4'	2.38	114.56	108.81
2	A	1001	FAD	C5A-C6A-N6A	2.38	123.96	120.35
4	B	2002	PO4	O4-P-O2	2.33	115.44	107.97
2	A	1001	FAD	C9A-N10-C10	-2.32	118.88	121.91
2	B	1001	FAD	C9-C9A-C5X	-2.14	116.23	119.88
2	B	1001	FAD	C9A-N10-C10	-2.06	119.22	121.91
2	A	1001	FAD	O2'-C2'-C3'	-2.01	104.21	109.10

There are no chirality outliers.

All (12) torsion outliers are listed below:

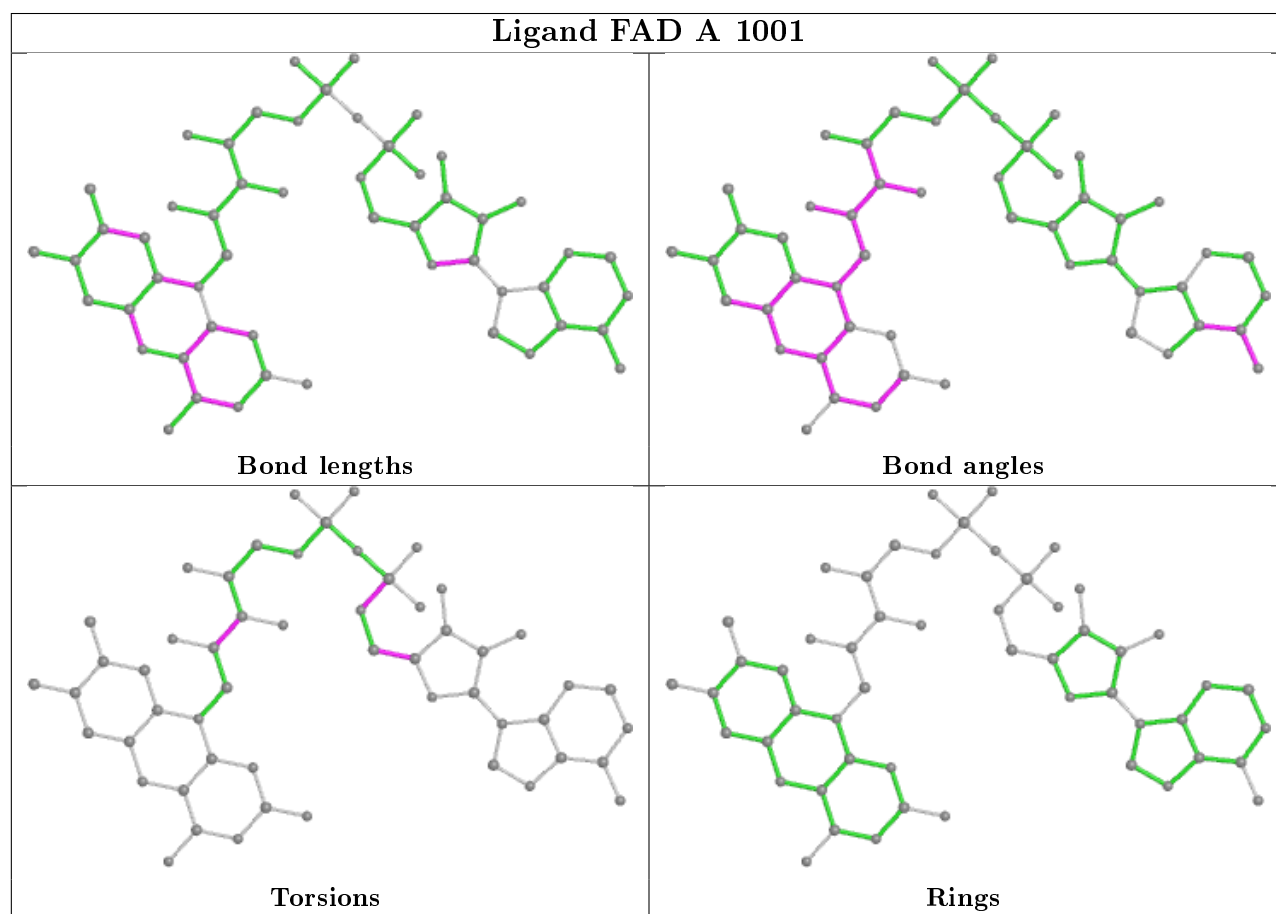
Mol	Chain	Res	Type	Atoms
2	A	1001	FAD	O2'-C2'-C3'-O3'
2	B	1001	FAD	O2'-C2'-C3'-O3'
2	B	1001	FAD	O4B-C4B-C5B-O5B
2	B	1001	FAD	C3B-C4B-C5B-O5B
2	A	1001	FAD	C3B-C4B-C5B-O5B
2	A	1001	FAD	O4B-C4B-C5B-O5B
2	B	1001	FAD	O2'-C2'-C3'-C4'
2	A	1001	FAD	O2'-C2'-C3'-C4'
2	B	1001	FAD	C5B-O5B-PA-O3P
2	A	1001	FAD	C5B-O5B-PA-O3P
2	A	1001	FAD	C1'-C2'-C3'-O3'
2	B	1001	FAD	C1'-C2'-C3'-O3'

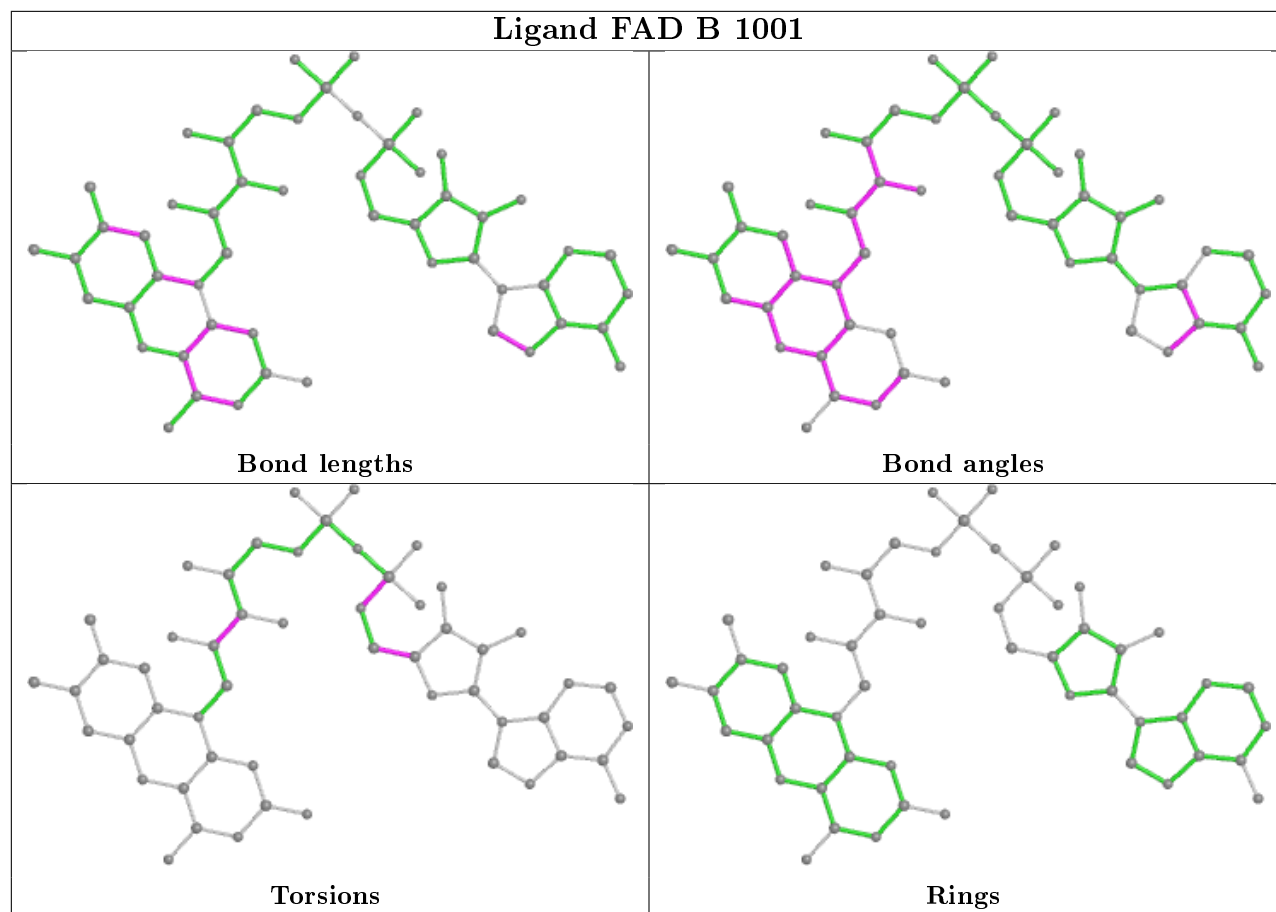
There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	411	PO4	1	0
3	B	1002	S3H	1	0
3	A	1002	S3H	2	0
4	B	2002	PO4	1	0
4	A	2002	PO4	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

### 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	356/409 (87%)	0.00	7 (1%) 65 60	30, 48, 90, 120	0
1	B	356/409 (87%)	0.12	16 (4%) 33 30	32, 54, 96, 125	0
All	All	712/818 (87%)	0.06	23 (3%) 47 44	30, 51, 94, 125	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	48	VAL	3.8
1	A	133	PHE	3.8
1	B	48	VAL	3.2
1	A	141	LEU	3.1
1	B	141	LEU	3.0
1	A	47	HIS	2.9
1	B	133	PHE	2.8
1	B	201	LYS	2.8
1	B	137	LEU	2.8
1	B	124	TYR	2.7
1	B	49	ALA	2.5
1	B	123	LYS	2.5
1	B	146	GLY	2.4
1	B	136	LYS	2.3
1	B	47	HIS	2.3
1	B	126	TYR	2.2
1	B	203	HIS	2.2
1	A	132	GLU	2.2
1	B	202	VAL	2.2
1	B	119	LYS	2.1
1	B	45	LEU	2.1
1	A	126	TYR	2.0
1	A	53	ARG	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](#)

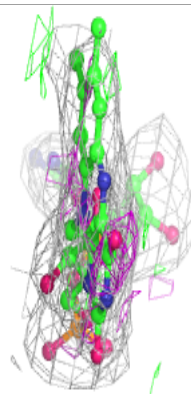
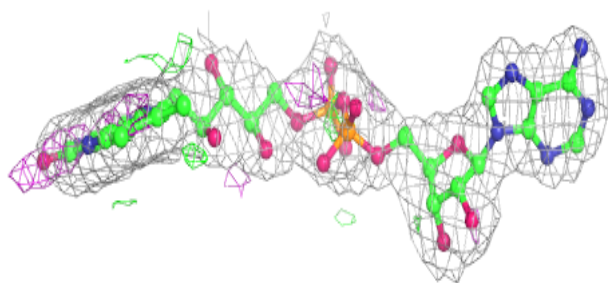
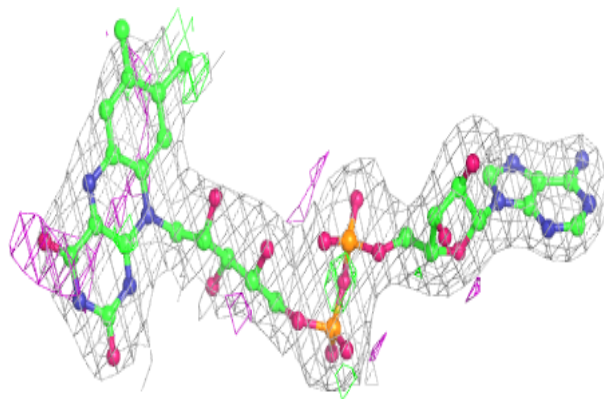
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	PO4	A	410	5/5	0.89	0.27	73,76,78,79	5
4	PO4	B	2003	5/5	0.89	0.25	127,131,133,134	0
4	PO4	B	2002	5/5	0.90	0.25	90,94,95,97	0
3	S3H	B	1002	3/3	0.90	0.15	60,60,63,64	2
4	PO4	A	2002	5/5	0.90	0.13	106,110,112,112	0
4	PO4	A	2001	5/5	0.91	0.18	92,96,99,100	0
4	PO4	A	2003	5/5	0.91	0.20	67,70,74,75	5
4	PO4	B	2001	5/5	0.94	0.12	98,101,104,105	0
4	PO4	B	411	5/5	0.94	0.14	73,75,78,80	5
3	S3H	A	1002	3/3	0.95	0.16	57,57,59,62	2
2	FAD	A	1001	53/53	0.97	0.18	46,49,52,54	0
4	PO4	B	410	5/5	0.97	0.17	63,66,68,69	0
2	FAD	B	1001	53/53	0.97	0.17	46,48,51,53	0

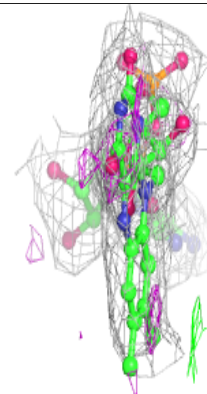
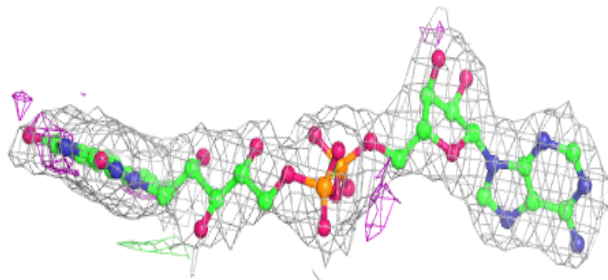
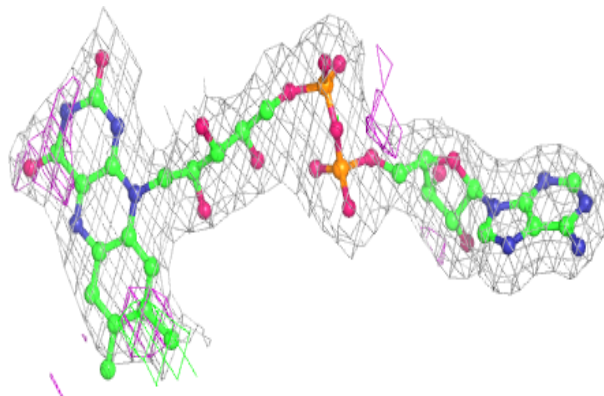
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around FAD A 1001:**

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

**Electron density around FAD B 1001:**

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





## 6.5 Other polymers

There are no such residues in this entry.