



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 11:14 PM GMT

PDB ID : 3DA1  
Title : X-Ray structure of the glycerol-3-phosphate dehydrogenase from *Bacillus halodurans* complexed with FAD. Northeast Structural Genomics Consortium target BhR167.  
Authors : Kuzin, A.P.; Abashidze, M.; Seetharaman, J.; Wang, D.; Janjua, H.; Owens, L.; Xiao, R.; Nair, R.; Baran, M.C.; Acton, T.B.; Rost, B.; Montelione, G.T.; Hunt, J.F.; Tong, L.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2008-05-28  
Resolution : 2.70 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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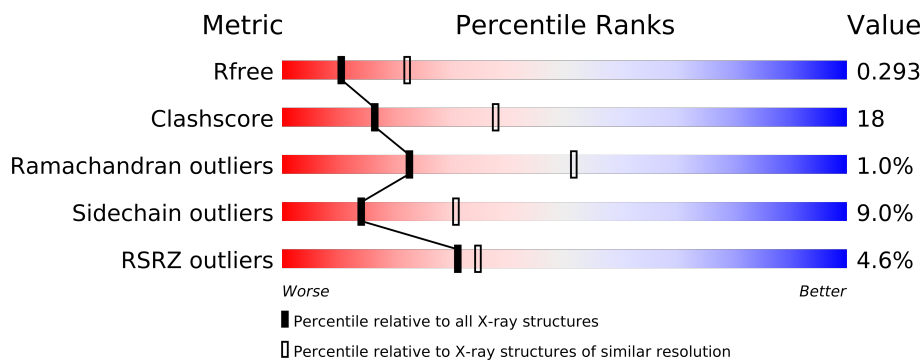
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.70 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))
$R_{free}$	66092	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	561	

## 2 Entry composition i

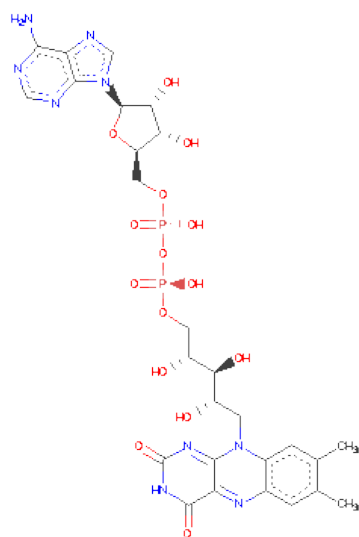
There are 3 unique types of molecules in this entry. The entry contains 4005 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Glycerol-3-phosphate dehydrogenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	496	3916	2486	674	740	2	14	0	0	0

- 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	53	27	9	15	2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	36	Total 36 O 36	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.77Å 96.95Å 226.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.74 – 2.70 29.75 – 2.67	Depositor EDS
% Data completeness (in resolution range)	83.0 (19.74-2.70) 93.3 (29.75-2.67)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.85 (at 2.68Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, $R_{free}$	0.222 , 0.282 0.240 , 0.293	Depositor DCC
$R_{free}$ test set	760 reflections (4.83%)	DCC
Wilson B-factor (Å <sup>2</sup> )	39.6	Xtriage
Anisotropy	0.758	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 31.6	EDS
Estimated twinning fraction	0.029 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.046 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30266 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	4005	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.44	0/3975	0.67	0/5338

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3916	0	3907	143	0
2	A	53	0	31	2	0
3	A	36	0	0	2	0
All	All	4005	0	3938	143	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 18.

All (143) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:61:THR:HG22	1:A:254:HIS:NE2	1.86	0.90

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:488:LEU:O	1:A:492:VAL:HG12	1.71	0.88
1:A:106:ILE:HG12	1:A:159:GLY:HA2	1.56	0.88
1:A:107:PHE:HZ	1:A:321:MSE:HE1	1.44	0.83
1:A:306:MSE:HE3	1:A:544:PRO:HB3	1.61	0.82
1:A:151:LEU:HD22	1:A:321:MSE:HE2	1.64	0.79
1:A:171:ARG:O	1:A:175:GLU:HG2	1.83	0.78
1:A:279:ILE:HG23	1:A:291:GLY:HA3	1.65	0.78
1:A:486:THR:HG22	1:A:488:LEU:H	1.49	0.76
1:A:374:LEU:HB2	2:A:609:FAD:O2'	1.85	0.76
1:A:1:MSE:H2	1:A:6:LYS:HD3	1.52	0.74
1:A:3:SER:HB2	1:A:480:GLU:HA	1.69	0.73
1:A:478:GLY:HA2	1:A:482:GLU:HB2	1.70	0.71
1:A:85:GLU:O	1:A:89:VAL:HG23	1.92	0.69
1:A:544:PRO:HG2	1:A:545:LEU:HD22	1.78	0.66
1:A:290:ILE:O	1:A:291:GLY:O	2.14	0.66
1:A:24:ILE:HD12	1:A:193:VAL:HG21	1.78	0.66
1:A:486:THR:HG22	1:A:488:LEU:N	2.14	0.61
1:A:488:LEU:HD21	1:A:534:LEU:HA	1.81	0.61
1:A:213:THR:OG1	1:A:215:THR:HG22	2.00	0.61
1:A:13:GLU:O	1:A:17:LYS:HG2	2.01	0.61
1:A:340:ARG:HD3	3:A:616:HOH:O	2.00	0.60
1:A:311:ARG:HG2	1:A:311:ARG:HH11	1.67	0.60
1:A:103:LEU:HB3	1:A:268:VAL:HG12	1.84	0.60
1:A:55:GLY:O	1:A:59:ARG:HD2	2.02	0.60
1:A:524:GLU:O	1:A:528:THR:HG22	2.02	0.58
1:A:357:ASP:HB2	1:A:380:MSE:HE1	1.85	0.58
1:A:536:THR:O	1:A:540:MSE:HG3	2.03	0.58
1:A:33:ILE:HD12	1:A:384:THR:HB	1.86	0.58
1:A:523:THR:OG1	1:A:525:GLU:HG2	2.02	0.57
1:A:548:VAL:HG12	1:A:549:GLU:H	1.69	0.57
1:A:357:ASP:HA	1:A:371:GLY:HA2	1.85	0.57
1:A:156:LEU:HD22	1:A:158:GLY:H	1.71	0.56
1:A:487:PRO:HB2	1:A:534:LEU:HD12	1.87	0.56
1:A:198:TYR:OH	1:A:239:ASP:HA	2.05	0.56
1:A:148:GLU:OE2	1:A:150:LEU:HB2	2.06	0.56
1:A:193:VAL:HG13	1:A:206:VAL:HB	1.88	0.55
1:A:9:LYS:O	1:A:13:GLU:HB2	2.06	0.55
1:A:450:HIS:HA	1:A:453:ASN:HD22	1.71	0.54
1:A:21:LEU:HD12	1:A:22:LEU:N	2.23	0.54
1:A:372:GLY:HA3	3:A:635:HOH:O	2.08	0.54
1:A:50:ASN:O	1:A:189:ASN:HB2	2.07	0.53
1:A:35:LEU:O	1:A:39:VAL:HG13	2.08	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:383:ARG:O	1:A:386:ASP:HB2	2.09	0.53
1:A:29:THR:O	1:A:33:ILE:HG12	2.08	0.52
1:A:433:ASP:HB3	1:A:436:GLU:HB2	1.90	0.52
1:A:378:ARG:NH1	1:A:401:ALA:HB2	2.25	0.52
1:A:95:HIS:HB2	1:A:442:LYS:O	2.09	0.52
1:A:20:ASP:OD2	1:A:222:LYS:HD2	2.10	0.52
1:A:378:ARG:CZ	1:A:401:ALA:HB2	2.39	0.52
1:A:504:LEU:C	1:A:504:LEU:HD23	2.30	0.52
1:A:1:MSE:HE2	1:A:481:GLN:HG2	1.92	0.51
1:A:325:LEU:N	1:A:325:LEU:HD12	2.26	0.51
1:A:292:THR:HG23	1:A:293:THR:N	2.24	0.51
1:A:104:LEU:HD12	1:A:269:TYR:O	2.11	0.51
1:A:192:LYS:O	1:A:208:ALA:HA	2.11	0.51
1:A:539:LYS:HD3	1:A:546:PHE:CE1	2.46	0.50
1:A:4:ALA:HA	1:A:483:MSE:HE3	1.94	0.49
1:A:240:ARG:HG3	1:A:240:ARG:HH11	1.77	0.49
1:A:479:VAL:HG23	1:A:480:GLU:N	2.27	0.49
1:A:33:ILE:HD13	1:A:381:ALA:HA	1.93	0.49
1:A:516:MSE:HE3	1:A:520:PHE:HE2	1.78	0.49
1:A:445:GLY:O	1:A:448:VAL:HG22	2.13	0.49
1:A:14:MSE:HG3	1:A:219:ILE:HD12	1.94	0.49
1:A:374:LEU:HB3	2:A:609:FAD:O2	2.12	0.49
1:A:359:ILE:HD13	1:A:359:ILE:N	2.28	0.49
1:A:156:LEU:HD22	1:A:158:GLY:N	2.28	0.49
1:A:105:PRO:O	1:A:106:ILE:HD13	2.12	0.48
1:A:343:ILE:HD12	1:A:343:ILE:N	2.28	0.48
1:A:142:LYS:O	1:A:146:GLU:HG3	2.12	0.48
1:A:292:THR:HG23	1:A:293:THR:O	2.13	0.48
1:A:360:PHE:HB2	1:A:368:SER:HB2	1.95	0.48
1:A:515:TRP:CZ2	1:A:519:GLU:HG3	2.49	0.47
1:A:21:LEU:HD12	1:A:22:LEU:H	1.80	0.47
1:A:246:TYR:HE2	1:A:248:LYS:HE3	1.80	0.47
1:A:17:LYS:HG3	1:A:17:LYS:O	2.15	0.47
1:A:450:HIS:HA	1:A:453:ASN:ND2	2.29	0.47
1:A:232:VAL:O	1:A:236:ARG:HG3	2.15	0.47
1:A:215:THR:HG23	1:A:217:HIS:CE1	2.51	0.46
1:A:49:MSE:O	1:A:190:TYR:HA	2.14	0.46
1:A:280:PHE:O	1:A:291:GLY:N	2.48	0.46
1:A:11:ILE:HG21	1:A:217:HIS:CD2	2.51	0.46
1:A:448:VAL:HA	1:A:451:VAL:HG12	1.96	0.46
1:A:51:ASP:HB3	1:A:54:SER:HB3	1.98	0.46
1:A:1:MSE:HG3	1:A:2:PHE:H	1.79	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:103:LEU:O	1:A:105:PRO:HD3	2.16	0.46
1:A:268:VAL:HG21	1:A:322:PHE:HZ	1.81	0.46
1:A:504:LEU:HD23	1:A:504:LEU:O	2.16	0.46
1:A:416:PHE:N	1:A:417:PRO:CD	2.78	0.46
1:A:465:TYR:O	1:A:467:LEU:HG	2.14	0.45
1:A:1:MSE:N	1:A:6:LYS:HD3	2.26	0.45
1:A:216:THR:HG22	1:A:217:HIS:N	2.31	0.45
1:A:311:ARG:HG2	1:A:311:ARG:NH1	2.31	0.45
1:A:492:VAL:HG13	1:A:493:ARG:HG3	1.99	0.45
1:A:341:PRO:O	1:A:342:LEU:HD12	2.17	0.45
1:A:215:THR:HG23	1:A:217:HIS:HE1	1.82	0.45
1:A:28:ILE:CD1	1:A:172:LEU:HG	2.47	0.45
1:A:484:VAL:HG21	1:A:494:ARG:HG3	1.99	0.45
1:A:292:THR:CG2	1:A:293:THR:N	2.79	0.45
1:A:461:GLU:CD	1:A:461:GLU:H	2.20	0.45
1:A:25:GLY:O	1:A:30:GLY:HA3	2.17	0.44
1:A:461:GLU:CD	1:A:461:GLU:N	2.71	0.44
1:A:325:LEU:HB3	1:A:327:LEU:HD13	2.00	0.44
1:A:96:VAL:HG11	1:A:172:LEU:HB2	2.00	0.44
1:A:325:LEU:N	1:A:325:LEU:CD1	2.81	0.44
1:A:44:THR:HG22	1:A:45:GLY:N	2.32	0.44
1:A:229:GLY:O	1:A:232:VAL:HG23	2.18	0.44
1:A:33:ILE:CD1	1:A:381:ALA:HA	2.48	0.44
1:A:19:LEU:O	1:A:221:ALA:HA	2.18	0.44
1:A:175:GLU:OE1	1:A:178:LYS:HD3	2.18	0.44
1:A:141:GLU:HG3	1:A:156:LEU:HD13	2.00	0.44
1:A:476:GLN:HA	1:A:479:VAL:HG22	2.00	0.43
1:A:341:PRO:C	1:A:342:LEU:HD12	2.39	0.43
1:A:424:SER:HA	1:A:437:VAL:HG23	2.00	0.43
1:A:175:GLU:HA	1:A:175:GLU:OE1	2.18	0.43
1:A:301:ILE:HG23	1:A:339:LEU:HD23	1.99	0.43
1:A:369:ILE:N	1:A:369:ILE:HD12	2.33	0.43
1:A:175:GLU:HA	1:A:406:SER:HB3	2.00	0.43
1:A:380:MSE:HE3	1:A:383:ARG:NH2	2.33	0.43
1:A:1:MSE:H2	1:A:6:LYS:CD	2.25	0.43
1:A:462:ALA:HB1	1:A:467:LEU:O	2.18	0.43
1:A:194:GLU:O	1:A:238:LYS:HD3	2.19	0.43
1:A:344:HIS:HD2	1:A:345:GLU:HG3	1.83	0.43
1:A:107:PHE:CZ	1:A:321:MSE:HE1	2.36	0.42
1:A:305:ARG:HA	1:A:337:ALA:CB	2.49	0.42
1:A:305:ARG:NH2	1:A:546:PHE:O	2.53	0.42
1:A:84:LYS:O	1:A:88:ILE:HG12	2.20	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:274:SER:OG	1:A:275:ASP:N	2.50	0.42
1:A:95:HIS:HE1	1:A:175:GLU:OE2	2.02	0.42
1:A:140:ASN:OD1	1:A:143:GLN:HG3	2.20	0.42
1:A:516:MSE:HE2	1:A:522:TRP:CZ2	2.55	0.41
1:A:293:THR:HG23	1:A:314:ILE:HD11	2.02	0.41
1:A:48:GLU:O	1:A:190:TYR:N	2.53	0.41
1:A:222:LYS:HB3	1:A:392:LEU:HD21	2.02	0.41
1:A:178:LYS:HE2	1:A:406:SER:OG	2.21	0.41
1:A:24:ILE:HG12	1:A:225:VAL:O	2.20	0.41
1:A:36:ASP:O	1:A:39:VAL:HG22	2.21	0.41
1:A:487:PRO:CB	1:A:534:LEU:HD12	2.50	0.40
1:A:517:ALA:HA	1:A:522:TRP:CE3	2.55	0.40
1:A:97:THR:CG2	1:A:165:TYR:HB3	2.51	0.40
1:A:270:PHE:N	1:A:270:PHE:CD1	2.89	0.40
1:A:154:GLU:O	1:A:155:ASN:HB2	2.22	0.40
1:A:7:ARG:HG2	1:A:7:ARG:HH11	1.87	0.40

There are no symmetry-related clashes.

## 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	488/561 (87%)	446 (91%)	37 (8%)	5 (1%)	22 51

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	275	ASP
1	A	291	GLY
1	A	273	GLU
1	A	393	ASN
1	A	372	GLY

### 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	411/454 (90%)	374 (91%)	37 (9%)	14 31

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	ARG
1	A	156	LEU
1	A	161	ILE
1	A	209	LYS
1	A	214	ASP
1	A	215	THR
1	A	240	ARG
1	A	245	LYS
1	A	249	LEU
1	A	250	SER
1	A	262	PHE
1	A	266	GLN
1	A	271	ASP
1	A	278	MSE
1	A	292	THR
1	A	326	ARG
1	A	327	LEU
1	A	339	LEU
1	A	359	ILE
1	A	380	MSE
1	A	418	ARG
1	A	426	LYS
1	A	443	LEU
1	A	448	VAL
1	A	452	LEU
1	A	459	LYS
1	A	467	LEU
1	A	470	LEU
1	A	472	LEU
1	A	483	MSE
1	A	501	ASN

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Mol	Chain	Res	Type
1	A	510	GLU
1	A	514	ARG
1	A	515	TRP
1	A	518	GLU
1	A	524	GLU
1	A	548	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	HIS
1	A	155	ASN
1	A	217	HIS
1	A	453	ASN
1	A	501	ASN
1	A	532	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

1 ligand is 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	609	-	58,58,58	1.76	11 (18%)	85,89,89	3.19	30 (35%)

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	609	-	-	2/34/50/50	0/1/6/6

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	609	FAD	C4X-C10	6.10	1.51	1.40
2	A	609	FAD	C4-C4X	5.15	1.49	1.41
2	A	609	FAD	C9A-N10	4.18	1.45	1.38
2	A	609	FAD	O4B-C1B	3.53	1.46	1.41
2	A	609	FAD	PA-O3P	3.17	1.65	1.59
2	A	609	FAD	C1'-C2'	2.58	1.54	1.51
2	A	609	FAD	C2-N3	-2.54	1.32	1.37
2	A	609	FAD	C4X-N5	-2.36	1.31	1.36
2	A	609	FAD	C8A-N7A	-2.23	1.30	1.34
2	A	609	FAD	P-O3P	-2.04	1.56	1.59
2	A	609	FAD	C10-N10	2.01	1.43	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	609	FAD	P-O3P-PA	-9.87	102.74	131.68
2	A	609	FAD	O3P-P-O5'	9.41	145.50	103.41
2	A	609	FAD	C2'-C1'-N10	8.57	123.82	112.45
2	A	609	FAD	N3A-C2A-N1A	-8.15	121.89	128.71
2	A	609	FAD	O5'-P-O1P	-7.26	80.93	109.37
2	A	609	FAD	C2-N1-C10	6.61	121.64	114.98
2	A	609	FAD	O4B-C1B-N9A	6.24	114.25	108.44
2	A	609	FAD	C4B-O4B-C1B	-5.91	103.33	109.75
2	A	609	FAD	C4X-C10-N1	-5.72	117.01	122.73
2	A	609	FAD	N3A-C4A-N9A	5.27	134.94	125.43
2	A	609	FAD	C1'-N10-C10	5.23	126.58	119.17
2	A	609	FAD	C4X-N5-C5X	4.92	122.21	116.69
2	A	609	FAD	O2P-P-O5'	-4.71	84.75	108.51
2	A	609	FAD	O2P-P-O3P	-4.58	83.40	105.14

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	609	FAD	O3P-P-O1P	-4.51	78.92	111.28
2	A	609	FAD	C9A-N10-C10	-4.45	117.40	121.77
2	A	609	FAD	C4X-C10-N10	-4.35	118.34	120.51
2	A	609	FAD	C5'-C4'-C3'	-3.67	105.14	112.06
2	A	609	FAD	P-O5'-C5'	-3.51	96.76	122.03
2	A	609	FAD	C4A-C5A-N7A	-3.37	106.64	109.52
2	A	609	FAD	N1-C10-N10	3.22	124.43	115.97
2	A	609	FAD	C5A-C4A-N3A	-3.07	119.02	125.70
2	A	609	FAD	C5X-C9A-N10	2.98	119.73	116.80
2	A	609	FAD	C1'-N10-C9A	-2.93	116.02	118.87
2	A	609	FAD	C2A-N3A-C4A	2.76	121.86	114.01
2	A	609	FAD	O4B-C1B-C2B	-2.54	102.87	106.77
2	A	609	FAD	O4B-C4B-C5B	2.24	117.34	109.36
2	A	609	FAD	C8A-N7A-C5A	2.09	110.06	103.58
2	A	609	FAD	O3'-C3'-C4'	-2.04	103.58	108.74
2	A	609	FAD	O2P-P-O1P	2.01	123.46	112.21

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	609	FAD	C2'-C1'-N10-C9A
2	A	609	FAD	C2'-C1'-N10-C10

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	496/561 (88%)	0.33	22 (4%)	33 37	23, 45, 70, 93	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	137	TYR	3.7
1	A	136	ARG	3.6
1	A	16	GLU	3.6
1	A	394	VAL	3.4
1	A	138	MSE	3.3
1	A	1	MSE	3.2
1	A	395	ASN	3.2
1	A	108	LYS	2.9
1	A	547	GLN	2.8
1	A	23	VAL	2.7
1	A	357	ASP	2.5
1	A	17	LYS	2.5
1	A	525	GLU	2.5
1	A	393	ASN	2.5
1	A	66	GLY	2.4
1	A	298	ASP	2.3
1	A	363	ASP	2.3
1	A	414	GLN	2.2
1	A	397	PRO	2.2
1	A	344	HIS	2.1
1	A	548	VAL	2.1
1	A	154	GLU	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

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAD	A	609	53/53	0.29	1.90	47,60,73,77	0

### 6.5 Other polymers

There are no such residues in this entry.