



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:36 PM GMT

PDB ID : 1GEU  
Title : ANATOMY OF AN ENGINEERED NAD-BINDING SITE  
Authors : Mittl, P.R.E.; Schulz, G.E.  
Deposited on : 1994-01-18  
Resolution : 2.20 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

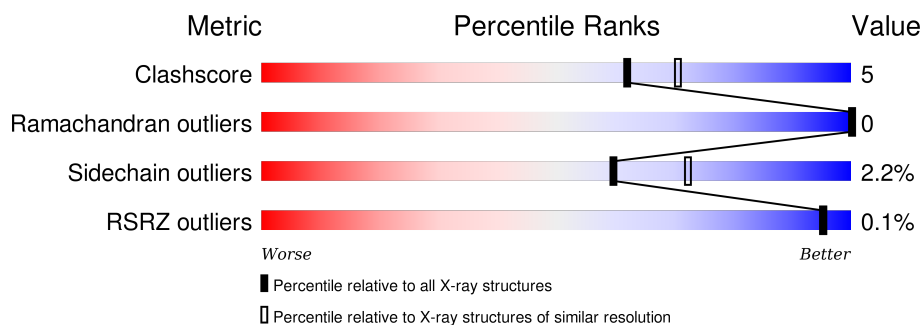
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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. 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	450	
1	B	450	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7425 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.

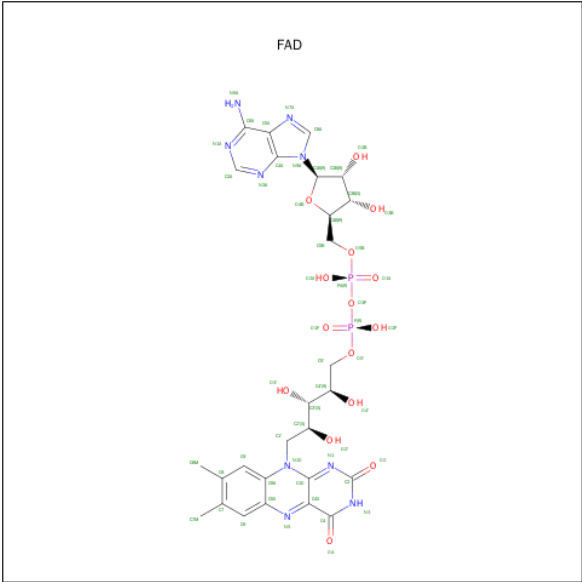
- Molecule 1 is a protein called GLUTATHIONE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3407	2153	580	656	18			
1	B	449	Total	C	N	O	S	0	0	0
			3414	2157	581	658	18			

There are 14 discrepancies between the modelled and reference sequences:

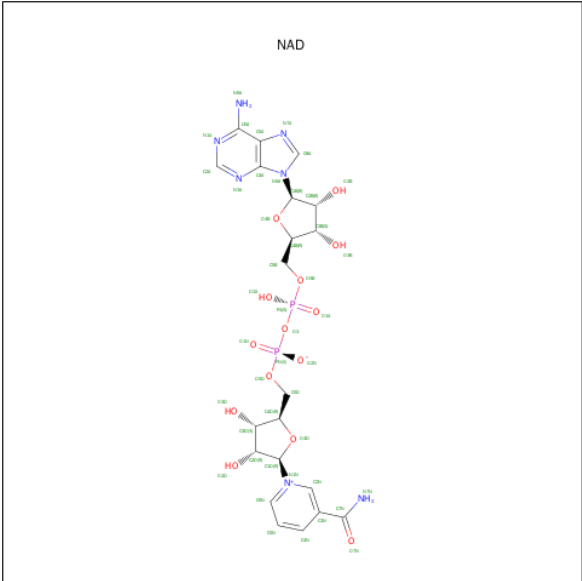
Chain	Residue	Modelled	Actual	Comment	Reference
A	179	GLY	ALA	ENGINEERED	UNP P06715
A	183	GLY	ALA	ENGINEERED	UNP P06715
A	197	GLU	VAL	ENGINEERED	UNP P06715
A	198	MET	ARG	ENGINEERED	UNP P06715
A	199	PHE	LYS	ENGINEERED	UNP P06715
A	200	ASP	HIS	ENGINEERED	UNP P06715
A	204	PRO	ARG	ENGINEERED	UNP P06715
B	179	GLY	ALA	ENGINEERED	UNP P06715
B	183	GLY	ALA	ENGINEERED	UNP P06715
B	197	GLU	VAL	ENGINEERED	UNP P06715
B	198	MET	ARG	ENGINEERED	UNP P06715
B	199	PHE	LYS	ENGINEERED	UNP P06715
B	200	ASP	HIS	ENGINEERED	UNP P06715
B	204	PRO	ARG	ENGINEERED	UNP P06715

- 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
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C<sub>21</sub>H<sub>27</sub>N<sub>7</sub>O<sub>14</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

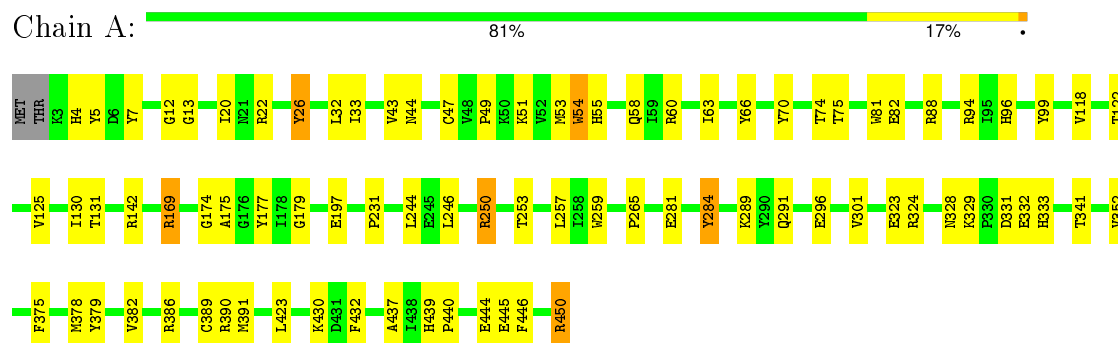
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	192	Total	O	0	0
			192	192		
4	B	218	Total	O	0	0
			218	218		

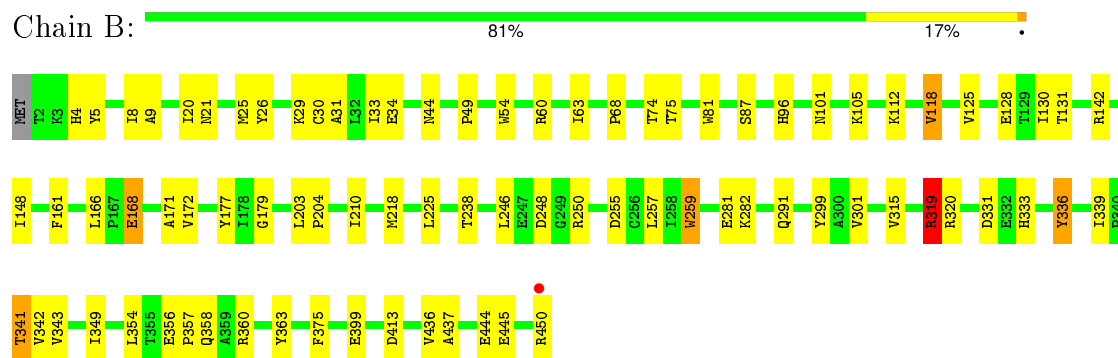
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: GLUTATHIONE REDUCTASE



#### • Molecule 1: GLUTATHIONE REDUCTASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 1 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	120.50 Å   72.50 Å   60.80 Å 90.00°   90.00°   82.80°	Depositor
Resolution (Å)	7.00 – 2.20 19.51 – 2.15	Depositor EDS
% Data completeness (in resolution range)	(Not available) (7.00-2.20) 79.2 (19.51-2.15)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.169 , (Not available) 0.159 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	14.6	Xtriage
Anisotropy	0.431	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 69.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>1</sup>	$\langle  L  \rangle = 0.37$ , $\langle L^2 \rangle = 0.20$	Xtriage
Outliers	2 of 45272 reflections (0.004%)	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7425	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.98% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: FAD, NAD

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	0.82	0/3478	1.45	31/4722 (0.7%)
1	B	0.84	0/3485	1.44	28/4732 (0.6%)
All	All	0.83	0/6963	1.45	59/9454 (0.6%)

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	A	0	1

There are no bond length outliers.

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	ARG	NE-CZ-NH2	-11.26	114.67	120.30
1	A	88	ARG	NE-CZ-NH2	-9.39	115.61	120.30
1	A	54	TRP	CD1-CG-CD2	8.79	113.33	106.30
1	A	88	ARG	NE-CZ-NH1	8.55	124.58	120.30
1	A	390	ARG	NE-CZ-NH2	-7.94	116.33	120.30
1	B	118	VAL	CG1-CB-CG2	-7.94	98.20	110.90
1	B	259	TRP	CD1-CG-CD2	7.92	112.64	106.30
1	A	169	ARG	NE-CZ-NH1	7.78	124.19	120.30
1	B	259	TRP	CE2-CD2-CG	-7.61	101.22	107.30
1	B	320	ARG	NE-CZ-NH1	7.58	124.09	120.30
1	A	60	ARG	NE-CZ-NH1	7.49	124.04	120.30
1	A	250	ARG	CA-CB-CG	7.25	129.34	113.40
1	B	81	TRP	CD1-CG-CD2	7.24	112.09	106.30
1	B	81	TRP	CE2-CD2-CG	-7.21	101.53	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	81	TRP	CE2-CD2-CG	-7.09	101.62	107.30
1	B	60	ARG	NE-CZ-NH1	7.05	123.83	120.30
1	A	94	ARG	NE-CZ-NH1	7.03	123.81	120.30
1	A	81	TRP	CD1-CG-CD2	6.88	111.81	106.30
1	A	70	TYR	CB-CG-CD2	-6.87	116.88	121.00
1	A	324	ARG	NE-CZ-NH1	6.83	123.71	120.30
1	B	142	ARG	NE-CZ-NH2	-6.67	116.96	120.30
1	A	259	TRP	CD1-CG-CD2	6.62	111.60	106.30
1	A	60	ARG	NE-CZ-NH2	-6.61	117.00	120.30
1	A	54	TRP	CE2-CD2-CG	-6.58	102.04	107.30
1	A	250	ARG	CG-CD-NE	6.53	125.50	111.80
1	A	142	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	A	250	ARG	CB-CG-CD	-6.42	94.90	111.60
1	A	259	TRP	CE2-CD2-CG	-6.37	102.21	107.30
1	A	257	LEU	CA-CB-CG	6.24	129.66	115.30
1	B	26	TYR	CB-CG-CD1	-6.16	117.31	121.00
1	A	379	TYR	CB-CG-CD2	-6.01	117.39	121.00
1	B	5	TYR	CB-CG-CD1	-5.99	117.41	121.00
1	B	81	TRP	CB-CG-CD1	-5.98	119.22	127.00
1	A	54	TRP	CG-CD1-NE1	-5.95	104.15	110.10
1	A	66	TYR	CB-CG-CD2	-5.92	117.45	121.00
1	B	81	TRP	CG-CD2-CE3	5.92	139.23	133.90
1	B	142	ARG	NE-CZ-NH1	5.91	123.25	120.30
1	B	54	TRP	CE2-CD2-CG	-5.89	102.58	107.30
1	A	390	ARG	CA-CB-CG	-5.89	100.44	113.40
1	B	319	ARG	NE-CZ-NH1	5.85	123.22	120.30
1	A	284	TYR	CB-CG-CD2	-5.82	117.50	121.00
1	A	391	MET	CG-SD-CE	-5.77	90.97	100.20
1	B	363	TYR	CB-CG-CD1	-5.77	117.54	121.00
1	A	26	TYR	CB-CG-CD1	-5.77	117.54	121.00
1	B	54	TRP	CD1-CG-CD2	5.76	110.91	106.30
1	B	413	ASP	CB-CG-OD1	5.65	123.39	118.30
1	B	336	TYR	CB-CG-CD1	-5.57	117.66	121.00
1	B	248	ASP	CB-CG-OD1	5.55	123.30	118.30
1	B	299	TYR	CB-CG-CD2	-5.43	117.74	121.00
1	B	436	VAL	CG1-CB-CG2	-5.43	102.21	110.90
1	A	430	LYS	CA-CB-CG	5.41	125.31	113.40
1	B	30	CYS	CA-CB-SG	5.35	123.64	114.00
1	B	342	VAL	CG1-CB-CG2	-5.28	102.45	110.90
1	B	125	VAL	CB-CA-C	-5.24	101.45	111.40
1	B	60	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	B	360	ARG	NE-CZ-NH1	5.21	122.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	81	TRP	CB-CG-CD1	-5.17	120.28	127.00
1	A	5	TYR	CD1-CG-CD2	5.04	123.44	117.90
1	B	450	ARG	NE-CZ-NH2	-5.03	117.78	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	26	TYR	Sidechain

## 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	3407	0	3352	43	0
1	B	3414	0	3359	39	0
2	A	53	0	31	1	0
2	B	53	0	31	0	0
3	A	44	0	26	0	0
3	B	44	0	26	0	0
4	A	192	0	0	1	0
4	B	218	0	0	1	0
All	All	7425	0	6825	75	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 5.

All (75) 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:4:HIS:HD2	1:B:131:THR:HG23	1.48	0.78
1:B:4:HIS:CD2	1:B:131:THR:HG23	2.29	0.68
1:B:246:LEU:HD12	1:B:250:ARG:HB2	1.77	0.66
1:A:177:TYR:OH	1:A:341:THR:HG21	1.98	0.63
1:B:44:ASN:OD1	1:B:96:HIS:HE1	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:HIS:HD2	1:A:131:THR:HG23	1.64	0.61
1:A:437:ALA:HB1	1:A:444:GLU:HB2	1.83	0.60
1:A:231:PRO:HB3	1:A:244:LEU:HD11	1.83	0.59
1:B:177:TYR:OH	1:B:341:THR:HG21	2.06	0.56
1:B:49:PRO:HB2	1:B:161:PHE:CE1	2.40	0.56
1:B:437:ALA:HB1	1:B:444:GLU:HB2	1.89	0.55
1:B:8:ILE:HG12	1:B:31:ALA:HB3	1.89	0.54
1:A:33:ILE:HD11	1:A:130:ILE:HD11	1.89	0.54
1:B:204:PRO:HD2	4:B:595:HOH:O	2.08	0.54
1:A:331:ASP:O	1:A:333:HIS:HD2	1.91	0.54
1:A:174:GLY:O	1:A:179:GLY:HA3	2.06	0.54
1:B:168:GLU:HG3	1:B:255:ASP:OD2	2.08	0.54
1:A:63:ILE:HG12	1:B:63:ILE:HG12	1.89	0.53
1:A:246:LEU:HD12	1:A:250:ARG:HB2	1.91	0.53
1:A:44:ASN:OD1	1:A:96:HIS:HE1	1.92	0.53
1:B:171:ALA:HB3	1:B:257:LEU:HD12	1.91	0.52
1:A:74:THR:HG21	1:B:63:ILE:HD13	1.92	0.52
1:B:218:MET:HE3	1:B:225:LEU:HD22	1.92	0.52
1:A:43:VAL:HG21	1:A:99:TYR:HE2	1.76	0.51
1:A:12:GLY:HA2	1:A:32:LEU:HD11	1.92	0.51
1:A:450:ARG:HH12	1:B:319:ARG:NH1	2.11	0.49
1:B:21:ASN:O	1:B:25:MET:HB2	2.12	0.49
1:A:296:GLU:H	1:A:296:GLU:CD	2.16	0.48
1:B:315:VAL:O	1:B:319:ARG:HB2	2.14	0.47
1:B:354:LEU:HD22	1:B:358:GLN:HB3	1.96	0.47
1:A:49:PRO:O	1:A:53:MET:HG2	2.14	0.47
1:A:291:GLN:HG3	1:A:301:VAL:HG12	1.96	0.47
1:A:169:ARG:NH2	1:A:253:THR:HG22	2.29	0.47
1:B:336:TYR:HA	1:B:339:ILE:HD11	1.97	0.47
1:A:175:ALA:HB3	1:A:197:GLU:HB2	1.97	0.47
1:B:101:ASN:HD21	1:B:105:LYS:HE2	1.79	0.46
1:B:34:GLU:O	1:B:112:LYS:HA	2.15	0.46
1:A:7:TYR:CD2	1:A:20:ILE:HD12	2.50	0.46
1:A:323:GLU:HB2	1:A:332:GLU:HG2	1.98	0.46
1:B:33:ILE:HD11	1:B:130:ILE:HD11	1.98	0.46
1:B:218:MET:CE	1:B:225:LEU:HD22	2.46	0.45
1:B:210:ILE:HD13	1:B:349:ILE:HG23	1.99	0.45
1:A:328:ASN:HB2	4:A:543:HOH:O	2.16	0.45
1:B:9:ALA:HB2	1:B:20:ILE:HD13	1.99	0.45
1:A:4:HIS:CD2	1:A:131:THR:HG23	2.48	0.44
1:B:148:ILE:HD11	1:B:259:TRP:CH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:GLU:HB3	1:A:329:LYS:HD2	1.99	0.44
1:A:375:PHE:CZ	1:A:389:CYS:SG	3.09	0.44
1:A:13:GLY:HA3	2:A:451:FAD:O5B	2.17	0.44
1:B:356:GLU:HB3	1:B:357:PRO:HD3	1.99	0.44
1:A:51:LYS:HE2	1:A:55:HIS:CE1	2.53	0.44
1:A:75:THR:O	1:B:74:THR:HA	2.18	0.44
1:B:29:LYS:HA	1:B:29:LYS:HD2	1.83	0.44
1:A:375:PHE:CE1	1:A:445:GLU:HG3	2.54	0.43
1:B:172:VAL:HG11	1:B:179:GLY:HA2	1.99	0.43
1:A:74:THR:HA	1:B:75:THR:O	2.19	0.42
1:A:118:VAL:HB	1:A:122:THR:HB	2.00	0.42
1:A:54:TRP:O	1:A:58:GLN:HG2	2.19	0.42
1:B:375:PHE:CE1	1:B:445:GLU:HG3	2.55	0.42
1:A:22:ARG:HD2	1:A:323:GLU:OE2	2.19	0.42
1:B:44:ASN:O	1:B:49:PRO:HG3	2.19	0.42
1:A:33:ILE:HD13	1:A:125:VAL:HG21	2.02	0.42
1:A:439:HIS:HA	1:A:440:PRO:HA	1.84	0.41
1:B:331:ASP:O	1:B:333:HIS:HD2	2.04	0.41
1:A:352:VAL:HG22	1:A:423:LEU:HD23	2.03	0.41
1:A:289:LYS:HZ1	1:A:333:HIS:CE1	2.38	0.41
1:A:265:PRO:HB3	1:A:284:TYR:CD1	2.55	0.41
1:A:450:ARG:HH22	1:B:319:ARG:HH12	1.68	0.41
1:A:382:VAL:HB	1:B:87:SER:HB3	2.02	0.41
1:B:291:GLN:HG3	1:B:301:VAL:HG12	2.03	0.41
1:B:177:TYR:CE2	1:B:343:VAL:HG22	2.56	0.41
1:B:148:ILE:HD11	1:B:259:TRP:HH2	1.87	0.40
1:A:378:MET:O	1:A:378:MET:HG2	2.21	0.40
1:A:386:ARG:HH11	1:A:386:ARG:HD2	1.78	0.40
1:A:432:PHE:HD2	1:A:446:PHE:O	2.04	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	446/450 (99%)	432 (97%)	14 (3%)	0	100	100
1	B	447/450 (99%)	429 (96%)	18 (4%)	0	100	100
All	All	893/900 (99%)	861 (96%)	32 (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	360/362 (99%)	356 (99%)	4 (1%)	80	89
1	B	361/362 (100%)	349 (97%)	12 (3%)	45	56
All	All	721/724 (100%)	705 (98%)	16 (2%)	60	72

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	CYS
1	A	82	GLU
1	A	281	GLU
1	A	450	ARG
1	B	68	PRO
1	B	118	VAL
1	B	128	GLU
1	B	166	LEU
1	B	168	GLU
1	B	203	LEU
1	B	238	THR
1	B	281	GLU
1	B	282	LYS
1	B	319	ARG
1	B	341	THR
1	B	399	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	HIS
1	A	55	HIS
1	A	96	HIS
1	A	237	ASN
1	A	327	ASN
1	A	333	HIS
1	A	362	GLN
1	A	434	ASN
1	B	4	HIS
1	B	55	HIS
1	B	96	HIS
1	B	237	ASN
1	B	271	ASN
1	B	327	ASN
1	B	333	HIS
1	B	434	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

4 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	451	-	48,58,58	1.12	4 (8%)	54,89,89	1.77	7 (12%)
3	NAD	A	452	-	38,48,48	1.04	3 (7%)	47,73,73	2.14	6 (12%)
2	FAD	B	451	-	48,58,58	1.23	4 (8%)	54,89,89	1.87	6 (11%)
3	NAD	B	452	-	38,48,48	0.97	2 (5%)	47,73,73	2.08	6 (12%)

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	451	-	-	0/30/50/50	0/6/6/6
3	NAD	A	452	-	-	0/22/62/62	0/5/5/5
2	FAD	B	451	-	-	0/30/50/50	0/6/6/6
3	NAD	B	452	-	-	0/22/62/62	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	451	FAD	C8A-N7A	-2.54	1.29	1.34
2	B	451	FAD	C8A-N7A	-2.24	1.30	1.34
3	A	452	NAD	C5A-N7A	-2.02	1.32	1.39
2	A	451	FAD	C4-C4X	2.06	1.45	1.41
3	A	452	NAD	C3N-C7N	2.17	1.53	1.50
3	B	452	NAD	O4B-C1B	2.18	1.44	1.41
2	A	451	FAD	C4X-N5	2.39	1.37	1.33
2	B	451	FAD	C4-C4X	2.74	1.46	1.41
3	B	452	NAD	O4D-C1D	2.93	1.44	1.41
2	A	451	FAD	C4-N3	3.17	1.39	1.33
2	B	451	FAD	C4X-N5	3.18	1.38	1.33
3	A	452	NAD	O4D-C1D	3.63	1.45	1.41
2	B	451	FAD	C4-N3	4.51	1.41	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	452	NAD	N3A-C2A-N1A	-9.35	121.73	128.89
3	B	452	NAD	N3A-C2A-N1A	-9.04	121.97	128.89
2	A	451	FAD	C4X-C10-N10	-6.50	116.69	120.52
2	B	451	FAD	C4X-C10-N10	-5.56	117.24	120.52
2	B	451	FAD	C4X-C4-N3	-5.07	116.66	123.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	451	FAD	C4X-C4-N3	-4.94	116.83	123.59
2	B	451	FAD	C4B-O4B-C1B	-4.70	104.55	109.72
3	B	452	NAD	C4D-O4D-C1D	-3.63	105.73	109.72
2	A	451	FAD	C4B-O4B-C1B	-3.45	105.93	109.72
2	B	451	FAD	C4-C4X-C10	-3.04	118.00	119.94
3	A	452	NAD	C4B-O4B-C1B	-2.87	106.56	109.72
3	A	452	NAD	C2B-C1B-N9A	-2.16	110.99	114.29
3	A	452	NAD	O5D-PN-O1N	-2.03	101.73	109.62
3	B	452	NAD	O2A-PA-O1A	2.05	123.66	112.53
2	A	451	FAD	C6-C5X-C9A	2.06	121.69	118.98
2	A	451	FAD	O2A-PA-O3P	2.11	114.68	105.09
2	B	451	FAD	N6A-C6A-N1A	2.12	123.75	119.20
3	B	452	NAD	O4D-C4D-C5D	2.23	117.29	109.32
2	A	451	FAD	C4X-N5-C5X	2.29	119.40	116.76
3	A	452	NAD	C4A-C5A-N7A	4.79	113.89	109.48
3	B	452	NAD	C4A-C5A-N7A	4.82	113.92	109.48
3	B	452	NAD	O4D-C1D-N1N	6.25	115.00	108.13
3	A	452	NAD	O4D-C1D-N1N	6.96	115.77	108.13
2	A	451	FAD	C4-N3-C2	7.11	121.40	115.25
2	B	451	FAD	C4-N3-C2	7.53	121.75	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	451	FAD	1	0

## 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 [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, 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	448/450 (99%)	-0.86	0 <b>100</b> <b>100</b>	3, 13, 36, 55	0
1	B	449/450 (99%)	-0.92	1 (0%) <b>95</b> <b>95</b>	3, 12, 32, 63	0
All	All	897/900 (99%)	-0.89	1 (0%) <b>95</b> <b>95</b>	3, 13, 34, 63	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	450	ARG	2.3

### 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. 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	FAD	A	451	53/53	0.98	0.07	-0.16	2,11,21,28	0
2	FAD	B	451	53/53	0.98	0.07	-0.28	2,7,14,15	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAD	A	452	44/44	0.98	0.07	-0.77	2,17,30,32	0
3	NAD	B	452	44/44	0.98	0.07	-0.83	2,15,21,25	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.