



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2018 – 01:49 pm GMT

PDB ID : 2QAE
Title : Crystal Structure Analysis of Trypanosoma cruzi Lipoamide dehydrogenase
Authors : Werner, C.; Krauth-Siegel, R.L.; Stubbs, M.T.; Klebe, G.
Deposited on : 2007-06-15
Resolution : 1.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk31020
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk31020

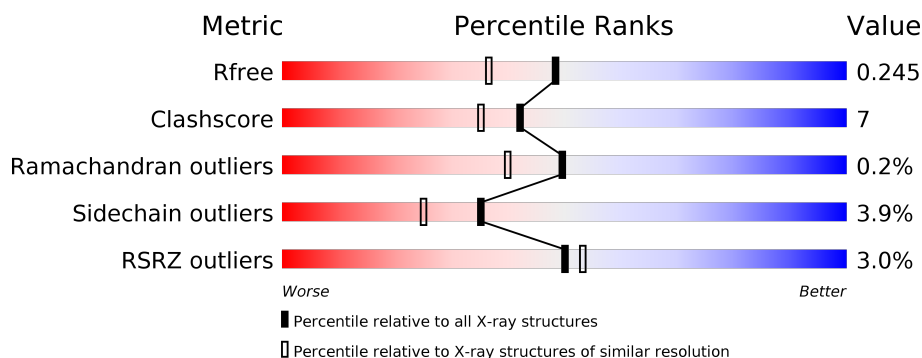
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	111664	5502 (1.90-1.90)
Clashscore	122126	6115 (1.90-1.90)
Ramachandran outliers	120053	6048 (1.90-1.90)
Sidechain outliers	120020	6048 (1.90-1.90)
RSRZ outliers	108989	5379 (1.90-1.90)

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 ≥ 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	468	<div> <div>3%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>..</div> </div> </div>
1	B	468	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>12%</div> <div>..</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7134 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 Dihydrolipoyl dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3440	2170	587	660	23			
1	B	465	Total	C	N	O	S	0	0	0
			3440	2170	587	660	23			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	58	LEU	VAL	VARIANT	UNP P90597
A	135	LEU	PHE	VARIANT	UNP P90597
A	199	GLU	LYS	VARIANT	UNP P90597
A	367	ASP	GLU	VARIANT	UNP P90597
A	462	PHE	VAL	VARIANT	UNP P90597
B	58	LEU	VAL	VARIANT	UNP P90597
B	135	LEU	PHE	VARIANT	UNP P90597
B	199	GLU	LYS	VARIANT	UNP P90597
B	367	ASP	GLU	VARIANT	UNP P90597
B	462	PHE	VAL	VARIANT	UNP P90597

- 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		

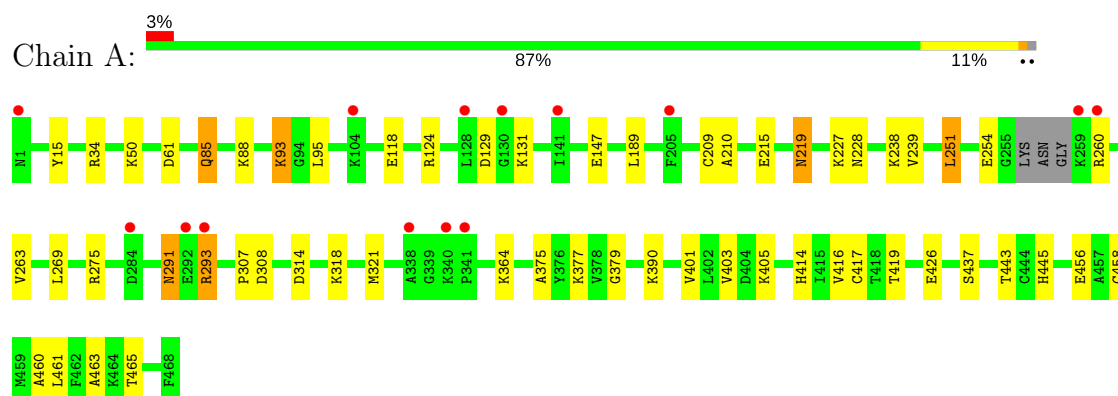
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	72	Total	O	0	0
			72	72		

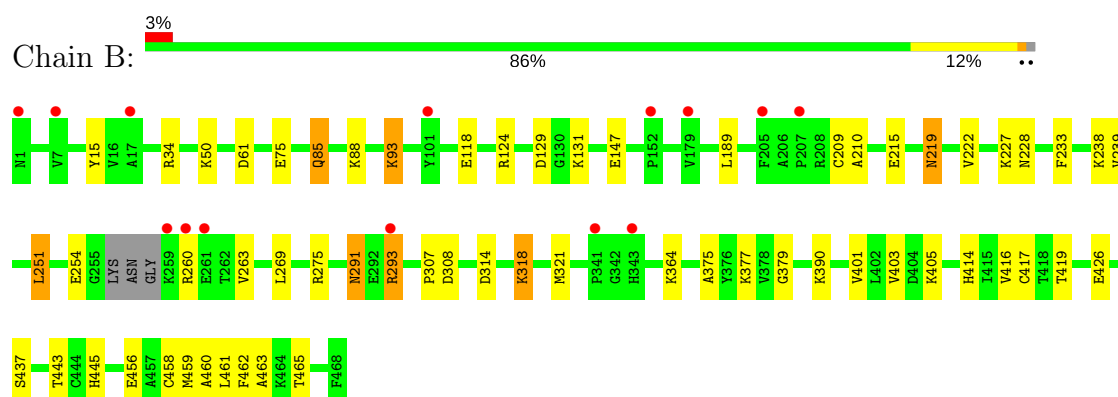
3 Residue-property plots [i](#)

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: Dihydrolipoyl dehydrogenase



• Molecule 1: Dihydrolipoyl dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.70Å 104.90Å 147.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.90 28.02 – 1.90	Depositor EDS
% Data completeness (in resolution range)	89.5 (30.00-1.90) 89.4 (28.02-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 1.91Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.187 , 0.217 0.225 , 0.245	Depositor DCC
R_{free} test set	7225 reflections (10.12%)	wwPDB-VP
Wilson B-factor (Å ²)	23.2	Xtriage
Anisotropy	0.251	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7134	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.76% 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

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	0.73	0/3496	0.82	1/4720 (0.0%)
1	B	0.73	0/3496	0.83	1/4720 (0.0%)
All	All	0.73	0/6992	0.82	2/9440 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	321	MET	CG-SD-CE	-5.45	91.48	100.20
1	A	321	MET	CG-SD-CE	-5.12	92.01	100.20

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	3440	0	3484	50	0
1	B	3440	0	3484	56	0
2	A	53	0	30	4	0
2	B	53	0	31	3	0
3	A	76	0	0	0	0
3	B	72	0	0	4	0
All	All	7134	0	7029	105	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 7.

All (105) 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:456:GLU:O	1:B:465:THR:HG22	1.68	0.94
1:A:260:ARG:HH11	1:A:260:ARG:HG2	1.36	0.91
1:A:456:GLU:O	1:A:465:THR:HG22	1.71	0.91
1:B:260:ARG:HG2	1:B:260:ARG:HH11	1.37	0.88
1:A:460:ALA:HB2	1:A:465:THR:HG23	1.58	0.83
1:B:460:ALA:HB2	1:B:465:THR:HG23	1.60	0.83
1:B:129:ASP:OD1	1:B:131:LYS:HG2	1.84	0.78
1:A:129:ASP:OD1	1:A:131:LYS:HG2	1.84	0.77
1:B:93:LYS:HB2	1:B:93:LYS:NZ	2.01	0.75
1:A:93:LYS:HB2	1:A:93:LYS:NZ	2.02	0.73
1:B:75:GLU:HB2	3:B:547:HOH:O	1.89	0.72
1:A:426:GLU:OE1	1:A:445:HIS:HE1	1.74	0.69
1:B:254:GLU:OE1	1:B:260:ARG:NH1	2.27	0.68
1:A:460:ALA:HB2	1:A:465:THR:CG2	2.24	0.67
1:A:254:GLU:OE1	1:A:260:ARG:NH1	2.26	0.66
1:B:426:GLU:OE1	1:B:445:HIS:HE1	1.79	0.65
1:B:291:ASN:C	1:B:291:ASN:HD22	1.99	0.65
1:B:34:ARG:HH21	2:B:480:FAD:H3B	1.62	0.64
1:A:260:ARG:HH11	1:A:260:ARG:CG	2.09	0.64
1:B:460:ALA:HB2	1:B:465:THR:CG2	2.28	0.64
1:A:34:ARG:HH21	2:A:480:FAD:H3B	1.64	0.63
1:B:85:GLN:HA	1:B:85:GLN:HE21	1.65	0.62
1:A:291:ASN:HD22	1:A:291:ASN:C	2.03	0.61
1:A:260:ARG:NH1	1:A:260:ARG:HG2	2.13	0.60
1:A:93:LYS:HZ3	1:A:93:LYS:HB2	1.66	0.60
1:A:85:GLN:HE21	1:A:85:GLN:HA	1.67	0.59
1:B:260:ARG:NH1	1:B:260:ARG:HG2	2.13	0.59
1:B:260:ARG:CG	1:B:260:ARG:HH11	2.10	0.59
1:A:293:ARG:HG3	1:A:293:ARG:NH1	2.18	0.58
1:B:293:ARG:HG3	1:B:293:ARG:NH1	2.18	0.58
1:B:93:LYS:HB2	1:B:93:LYS:HZ3	1.71	0.56
1:A:426:GLU:OE1	1:A:445:HIS:CE1	2.57	0.55
1:A:314:ASP:OD1	2:A:480:FAD:H5'1	2.07	0.55
1:B:426:GLU:OE1	1:B:445:HIS:CE1	2.60	0.54
1:B:85:GLN:HE22	1:B:88:LYS:NZ	2.06	0.54
1:B:227:LYS:HE3	1:B:228:ASN:OD1	2.08	0.54
1:B:314:ASP:OD1	2:B:480:FAD:H5'1	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:GLN:HE22	1:A:88:LYS:NZ	2.08	0.51
1:A:260:ARG:NH1	1:A:260:ARG:CG	2.70	0.50
1:A:443:THR:O	1:A:445:HIS:HD2	1.94	0.50
1:A:227:LYS:HE3	1:A:228:ASN:OD1	2.11	0.50
1:B:307:PRO:O	1:B:308:ASP:HB2	2.12	0.50
1:B:403:VAL:HG13	1:B:461:LEU:HD13	1.94	0.49
1:B:403:VAL:CG1	1:B:461:LEU:HD13	2.42	0.49
1:B:85:GLN:HE22	1:B:88:LYS:HZ2	1.60	0.49
1:B:379:GLY:N	1:B:461:LEU:HD21	2.28	0.49
1:B:85:GLN:NE2	1:B:88:LYS:NZ	2.60	0.49
1:B:293:ARG:HG3	1:B:293:ARG:HH11	1.77	0.49
1:A:239:VAL:HG13	1:A:251:LEU:HD22	1.95	0.49
1:A:293:ARG:HG3	1:A:293:ARG:HH11	1.77	0.49
1:B:443:THR:O	1:B:445:HIS:HD2	1.95	0.49
1:A:215:GLU:O	1:A:219:ASN:HB2	2.13	0.49
1:B:85:GLN:HA	1:B:85:GLN:NE2	2.28	0.48
1:A:379:GLY:N	1:A:461:LEU:HD21	2.27	0.48
1:A:403:VAL:CG1	1:A:461:LEU:HD13	2.43	0.48
1:A:403:VAL:HG22	1:A:461:LEU:HD22	1.94	0.48
1:B:416:VAL:O	1:B:417:CYS:HB3	2.13	0.48
1:A:307:PRO:O	1:A:308:ASP:HB2	2.14	0.48
1:B:403:VAL:HG22	1:B:461:LEU:HD22	1.94	0.48
1:B:419:THR:HG22	1:B:419:THR:O	2.14	0.48
1:B:147:GLU:OE1	1:B:275:ARG:NH1	2.47	0.47
1:A:403:VAL:HG13	1:A:461:LEU:HD13	1.97	0.47
1:B:215:GLU:O	1:B:219:ASN:HB2	2.13	0.47
1:B:291:ASN:C	1:B:291:ASN:ND2	2.68	0.47
1:B:318:LYS:CE	3:B:537:HOH:O	2.62	0.46
1:A:85:GLN:NE2	1:A:85:GLN:HA	2.30	0.46
1:B:85:GLN:CA	1:B:85:GLN:HE21	2.24	0.46
1:A:85:GLN:CA	1:A:85:GLN:HE21	2.28	0.46
1:B:239:VAL:HG13	1:B:251:LEU:HD22	1.96	0.46
1:A:85:GLN:NE2	1:A:88:LYS:NZ	2.63	0.46
1:B:375:ALA:HB1	1:B:405:LYS:HG2	1.98	0.46
1:A:416:VAL:O	1:A:417:CYS:HB3	2.16	0.45
1:A:419:THR:O	1:A:419:THR:HG22	2.17	0.45
1:B:458:CYS:HA	1:B:461:LEU:HD12	1.99	0.45
1:B:34:ARG:HH21	2:B:480:FAD:C3B	2.30	0.44
1:A:364:LYS:O	1:A:414:HIS:HE1	2.00	0.44
1:A:209:CYS:O	1:A:210:ALA:C	2.56	0.44
1:B:209:CYS:O	1:B:210:ALA:C	2.55	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:ARG:CG	1:B:260:ARG:NH1	2.71	0.44
1:B:293:ARG:CG	1:B:293:ARG:HH11	2.31	0.44
1:B:364:LYS:O	1:B:414:HIS:HE1	2.01	0.44
1:A:147:GLU:OE1	1:A:275:ARG:NH1	2.49	0.44
1:B:403:VAL:HG11	1:B:437:SER:HB3	2.00	0.44
1:B:222:VAL:HG13	1:B:233:PHE:CE1	2.53	0.43
1:A:95:LEU:CD2	1:B:390:LYS:HD3	2.48	0.43
1:A:458:CYS:HA	1:A:461:LEU:HD12	2.00	0.43
1:A:291:ASN:ND2	1:A:293:ARG:H	2.17	0.43
1:A:293:ARG:HH11	1:A:293:ARG:CG	2.32	0.43
1:A:34:ARG:HH21	2:A:480:FAD:C3B	2.31	0.43
1:A:401:VAL:HB	1:A:461:LEU:HD11	2.00	0.42
1:A:85:GLN:HE22	1:A:88:LYS:HZ2	1.66	0.42
1:A:238:LYS:NZ	1:A:239:VAL:H	2.17	0.42
1:A:390:LYS:HE3	3:B:532:HOH:O	2.18	0.42
1:B:401:VAL:HB	1:B:461:LEU:HD11	2.01	0.42
1:A:375:ALA:HB1	1:A:405:LYS:HG2	2.02	0.42
1:B:318:LYS:HE2	3:B:537:HOH:O	2.18	0.41
1:B:238:LYS:HZ1	1:B:239:VAL:H	1.69	0.41
1:B:93:LYS:CB	1:B:93:LYS:NZ	2.77	0.41
1:A:403:VAL:HG11	1:A:437:SER:HB3	2.03	0.41
1:A:118:GLU:OE2	1:A:124:ARG:NH2	2.54	0.40
1:A:314:ASP:CG	2:A:480:FAD:H5'1	2.41	0.40
1:B:118:GLU:OE2	1:B:124:ARG:NH2	2.54	0.40
1:B:291:ASN:ND2	1:B:293:ARG:H	2.19	0.40
1:B:459:MET:HA	1:B:462:PHE:CE1	2.57	0.40
1:B:93:LYS:HB2	1:B:93:LYS:HZ2	1.82	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	461/468 (98%)	445 (96%)	15 (3%)	1 (0%)	49	40
1	B	461/468 (98%)	446 (97%)	14 (3%)	1 (0%)	49	40
All	All	922/936 (98%)	891 (97%)	29 (3%)	2 (0%)	49	40

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	463	ALA
1	B	463	ALA

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	363/365 (100%)	349 (96%)	14 (4%)	35	25
1	B	363/365 (100%)	349 (96%)	14 (4%)	35	25
All	All	726/730 (100%)	698 (96%)	28 (4%)	35	25

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	TYR
1	A	50	LYS
1	A	61	ASP
1	A	85	GLN
1	A	93	LYS
1	A	189	LEU
1	A	219	ASN
1	A	251	LEU
1	A	263	VAL
1	A	269	LEU
1	A	291	ASN
1	A	293	ARG
1	A	318	LYS
1	A	377	LYS

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Mol	Chain	Res	Type
1	B	15	TYR
1	B	50	LYS
1	B	61	ASP
1	B	85	GLN
1	B	93	LYS
1	B	189	LEU
1	B	219	ASN
1	B	251	LEU
1	B	263	VAL
1	B	269	LEU
1	B	291	ASN
1	B	293	ARG
1	B	318	LYS
1	B	377	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	43	ASN
1	A	85	GLN
1	A	86	GLN
1	A	126	ASN
1	A	287	ASN
1	A	291	ASN
1	A	414	HIS
1	A	445	HIS
1	B	23	GLN
1	B	43	ASN
1	B	85	GLN
1	B	86	GLN
1	B	126	ASN
1	B	287	ASN
1	B	291	ASN
1	B	414	HIS
1	B	445	HIS

5.3.3 RNA ⓘ

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](#)

2 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	480	-	51,58,58	2.22	16 (31%)	57,89,89	3.59	28 (49%)
2	FAD	B	480	-	51,58,58	2.02	12 (23%)	57,89,89	3.60	27 (47%)

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	480	-	-	0/28/50/50	0/6/6/6
2	FAD	B	480	-	-	0/28/50/50	0/6/6/6

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	480	FAD	O5B-C5B	-4.78	1.26	1.44
2	A	480	FAD	O5B-C5B	-4.75	1.26	1.44
2	B	480	FAD	O5'-C5'	-4.34	1.27	1.44
2	A	480	FAD	O5'-C5'	-4.20	1.28	1.44
2	A	480	FAD	C3B-C4B	-3.54	1.43	1.53
2	B	480	FAD	C3B-C4B	-2.92	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	480	FAD	C8A-N7A	-2.83	1.29	1.34
2	A	480	FAD	C4'-C3'	-2.70	1.48	1.53
2	A	480	FAD	C8A-N7A	-2.59	1.29	1.34
2	A	480	FAD	O3B-C3B	-2.58	1.36	1.43
2	B	480	FAD	O2B-C2B	-2.12	1.37	1.43
2	A	480	FAD	O2B-C2B	-2.07	1.37	1.43
2	B	480	FAD	C4'-C3'	-2.07	1.49	1.53
2	B	480	FAD	O3B-C3B	-2.06	1.37	1.43
2	A	480	FAD	O3'-C3'	-2.04	1.38	1.43
2	A	480	FAD	C1'-N10	-2.00	1.46	1.48
2	A	480	FAD	C9A-C5X	2.22	1.47	1.42
2	A	480	FAD	C9A-N10	2.24	1.41	1.38
2	A	480	FAD	C2A-N3A	2.36	1.36	1.32
2	A	480	FAD	O4B-C1B	2.75	1.45	1.41
2	B	480	FAD	C10-N1	3.00	1.37	1.33
2	B	480	FAD	O4B-C1B	3.15	1.45	1.41
2	B	480	FAD	C4-C4X	3.87	1.48	1.41
2	A	480	FAD	C4-C4X	4.17	1.48	1.41
2	B	480	FAD	C5'-C4'	5.05	1.59	1.51
2	B	480	FAD	C4X-C10	5.83	1.50	1.41
2	A	480	FAD	C4X-C10	6.94	1.52	1.41
2	A	480	FAD	C5'-C4'	7.03	1.62	1.51

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	480	FAD	N3A-C2A-N1A	-9.88	120.40	128.86
2	B	480	FAD	N3A-C2A-N1A	-9.38	120.84	128.86
2	B	480	FAD	C4B-O4B-C1B	-8.44	101.03	109.83
2	A	480	FAD	C4B-O4B-C1B	-8.39	101.08	109.83
2	B	480	FAD	O2A-PA-O5B	-5.03	84.40	107.75
2	B	480	FAD	C4-C4X-C10	-4.98	116.23	119.95
2	A	480	FAD	C5B-C4B-C3B	-4.87	96.97	115.29
2	B	480	FAD	C5B-C4B-C3B	-4.61	97.92	115.29
2	A	480	FAD	O2A-PA-O5B	-4.54	86.64	107.75
2	B	480	FAD	O2A-PA-O1A	-4.21	90.73	112.14
2	A	480	FAD	C4X-C4-N3	-4.20	117.50	123.47
2	A	480	FAD	C4-C4X-C10	-3.97	116.98	119.95
2	B	480	FAD	C4X-C4-N3	-3.72	118.18	123.47
2	A	480	FAD	O5'-C5'-C4'	-3.71	99.46	109.36
2	A	480	FAD	P-O3P-PA	-3.60	120.52	132.63
2	B	480	FAD	P-O3P-PA	-3.52	120.81	132.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	480	FAD	O2A-PA-O1A	-3.48	94.45	112.14
2	B	480	FAD	O5'-C5'-C4'	-3.30	100.56	109.36
2	A	480	FAD	PA-O5B-C5B	-2.45	107.33	121.68
2	B	480	FAD	PA-O5B-C5B	-2.43	107.41	121.68
2	A	480	FAD	C4X-C10-N10	-2.35	118.21	120.40
2	A	480	FAD	C4A-C5A-N7A	-2.31	107.18	109.41
2	B	480	FAD	C1'-N10-C10	-2.18	116.28	118.46
2	A	480	FAD	C1'-N10-C10	-2.17	116.29	118.46
2	B	480	FAD	C2B-C3B-C4B	-2.11	98.57	102.62
2	B	480	FAD	C4X-C10-N10	-2.06	118.49	120.40
2	A	480	FAD	C2A-N1A-C6A	2.15	122.40	118.75
2	A	480	FAD	O2'-C2'-C3'	2.47	115.21	109.10
2	B	480	FAD	O3B-C3B-C2B	2.55	119.98	111.83
2	A	480	FAD	O3B-C3B-C2B	2.60	120.17	111.83
2	B	480	FAD	O2'-C2'-C3'	2.79	116.00	109.10
2	A	480	FAD	O4'-C4'-C5'	3.39	117.57	110.02
2	A	480	FAD	C4-C4X-N5	3.59	122.73	118.70
2	B	480	FAD	O4'-C4'-C5'	3.83	118.57	110.02
2	B	480	FAD	O3'-C3'-C2'	3.91	118.46	108.82
2	A	480	FAD	O3'-C3'-C2'	4.07	118.86	108.82
2	B	480	FAD	C4-C4X-N5	4.21	123.43	118.70
2	B	480	FAD	P-O5'-C5'	4.36	147.24	121.68
2	A	480	FAD	P-O5'-C5'	4.39	147.45	121.68
2	B	480	FAD	O4'-C4'-C3'	4.89	121.18	109.10
2	B	480	FAD	C4X-N5-C5X	5.06	122.06	116.76
2	A	480	FAD	C4X-N5-C5X	5.08	122.08	116.76
2	A	480	FAD	O4'-C4'-C3'	5.36	122.34	109.10
2	A	480	FAD	O5B-PA-O1A	5.48	130.48	109.07
2	B	480	FAD	O5'-P-O1P	5.93	132.24	109.07
2	B	480	FAD	O5B-PA-O1A	5.94	132.28	109.07
2	B	480	FAD	O4B-C4B-C3B	6.07	117.16	105.15
2	A	480	FAD	O5'-P-O1P	6.10	132.90	109.07
2	B	480	FAD	C1'-N10-C9A	6.23	123.85	118.31
2	A	480	FAD	O4B-C4B-C3B	6.33	117.67	105.15
2	A	480	FAD	C1'-N10-C9A	6.35	123.95	118.31
2	B	480	FAD	O4B-C4B-C5B	6.52	131.11	109.39
2	A	480	FAD	O4B-C4B-C5B	6.63	131.46	109.39
2	A	480	FAD	C4-N3-C2	8.43	122.32	115.14
2	B	480	FAD	C4-N3-C2	9.05	122.85	115.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	480	FAD	4	0
2	B	480	FAD	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	465/468 (99%)	0.37	14 (3%) 50 53	15, 26, 50, 75	0
1	B	465/468 (99%)	0.43	14 (3%) 50 53	15, 26, 50, 76	0
All	All	930/936 (99%)	0.40	28 (3%) 50 53	15, 26, 50, 76	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1	ASN	4.5
1	A	341	PRO	4.0
1	B	259	LYS	3.6
1	A	259	LYS	3.3
1	B	293	ARG	3.1
1	B	205	PHE	2.8
1	A	205	PHE	2.8
1	A	130	GLY	2.8
1	B	152	PRO	2.7
1	B	101	TYR	2.7
1	B	343	HIS	2.6
1	A	284	ASP	2.5
1	B	261	GLU	2.5
1	B	179	VAL	2.5
1	B	260	ARG	2.4
1	A	338	ALA	2.3
1	A	260	ARG	2.3
1	B	207	PRO	2.3
1	A	292	GLU	2.2
1	A	293	ARG	2.2
1	A	340	LYS	2.2
1	A	128	LEU	2.2
1	B	17	ALA	2.2
1	A	104	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	341	PRO	2.1
1	B	7	VAL	2.1
1	A	1	ASN	2.1
1	A	141	ILE	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, 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(\AA^2)	Q<0.9
2	FAD	A	480	53/53	0.91	0.15	16,21,26,34	0
2	FAD	B	480	53/53	0.92	0.13	15,20,26,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.