



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:09 PM GMT

PDB ID : 4NWZ
Title : Structure of bacterial type II NADH dehydrogenase from *Caldalkalibacillus thermarum* at 2.5Å resolution
Authors : Nakatani, Y.; Heikal, A.; Lott, J.S.; Sazanov, L.A.; Baker, E.N.; Cook, G.M.
Deposited on : 2013-12-07
Resolution : 2.50 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
<http://wwpdb.org/validation/2016/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 (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

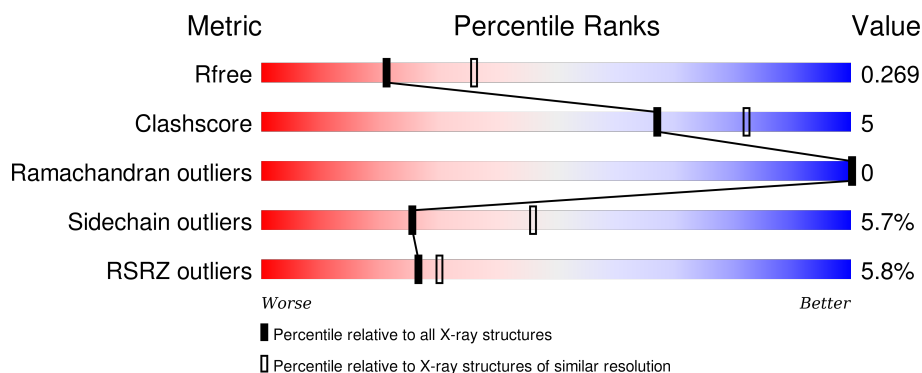
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	405	<div> <div>2%</div> <div>84%</div> <div>12%</div> <div>••</div> </div>
1	B	405	<div> <div>%</div> <div>84%</div> <div>12%</div> <div>••</div> </div>
1	C	405	<div> <div>2%</div> <div>84%</div> <div>11%</div> <div>••</div> </div>
1	D	405	<div> <div>17%</div> <div>79%</div> <div>12%</div> <div>• 8%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12454 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 FAD-dependent pyridine nucleotide-disulfide oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	395	Total	C	N	O	S	84	1	0
			3049	1944	529	567	9			
1	B	395	Total	C	N	O	S	40	2	0
			3055	1948	529	569	9			
1	C	394	Total	C	N	O	S	80	0	0
			3039	1938	528	564	9			
1	D	373	Total	C	N	O	S	106	1	0
			2886	1846	493	538	9			

There are 24 discrepancies between the modelled and reference sequences:

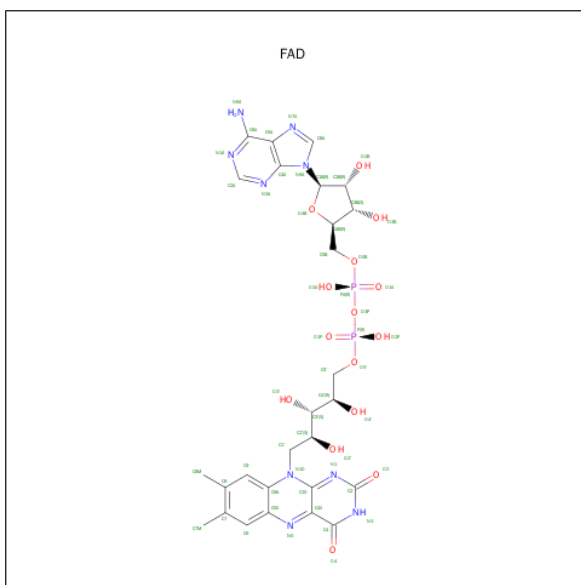
Chain	Residue	Modelled	Actual	Comment	Reference
A	400	HIS	-	EXPRESSION TAG	UNP F5L3B8
A	401	HIS	-	EXPRESSION TAG	UNP F5L3B8
A	402	HIS	-	EXPRESSION TAG	UNP F5L3B8
A	403	HIS	-	EXPRESSION TAG	UNP F5L3B8
A	404	HIS	-	EXPRESSION TAG	UNP F5L3B8
A	405	HIS	-	EXPRESSION TAG	UNP F5L3B8
B	400	HIS	-	EXPRESSION TAG	UNP F5L3B8
B	401	HIS	-	EXPRESSION TAG	UNP F5L3B8
B	402	HIS	-	EXPRESSION TAG	UNP F5L3B8
B	403	HIS	-	EXPRESSION TAG	UNP F5L3B8
B	404	HIS	-	EXPRESSION TAG	UNP F5L3B8
B	405	HIS	-	EXPRESSION TAG	UNP F5L3B8
C	400	HIS	-	EXPRESSION TAG	UNP F5L3B8
C	401	HIS	-	EXPRESSION TAG	UNP F5L3B8
C	402	HIS	-	EXPRESSION TAG	UNP F5L3B8
C	403	HIS	-	EXPRESSION TAG	UNP F5L3B8
C	404	HIS	-	EXPRESSION TAG	UNP F5L3B8
C	405	HIS	-	EXPRESSION TAG	UNP F5L3B8
D	400	HIS	-	EXPRESSION TAG	UNP F5L3B8
D	401	HIS	-	EXPRESSION TAG	UNP F5L3B8
D	402	HIS	-	EXPRESSION TAG	UNP F5L3B8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	403	HIS	-	EXPRESSION TAG	UNP F5L3B8
D	404	HIS	-	EXPRESSION TAG	UNP F5L3B8
D	405	HIS	-	EXPRESSION TAG	UNP F5L3B8

- 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		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	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	65	Total	O	0	0
			65	65		
3	B	77	Total	O	0	0
			77	77		
3	C	53	Total	O	0	0
			53	53		

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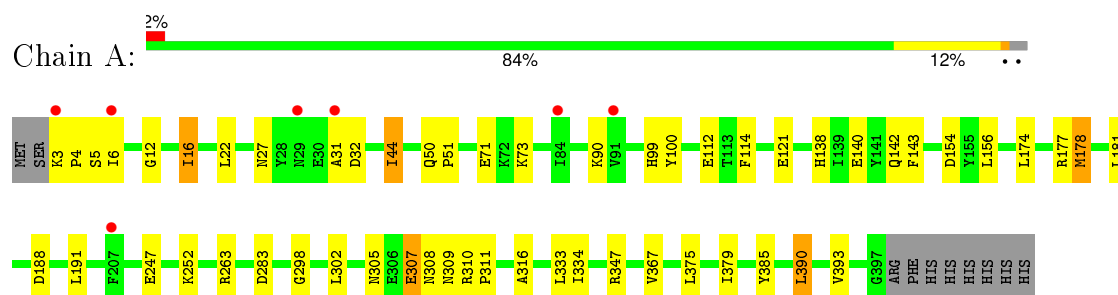
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	18	Total	O	0	0
			18	18		

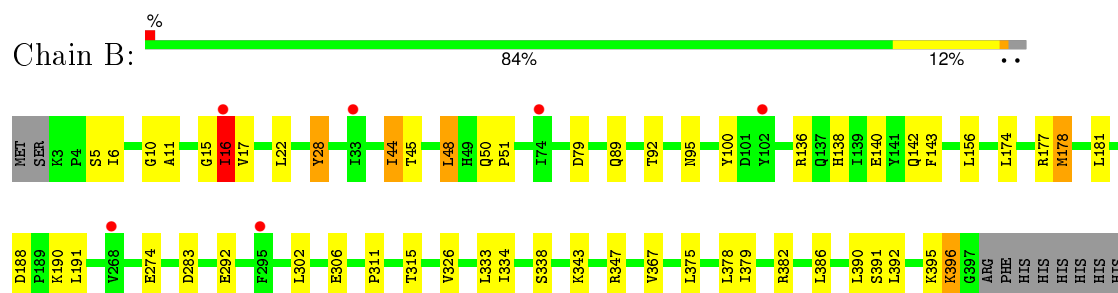
3 Residue-property plots [i](#)

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

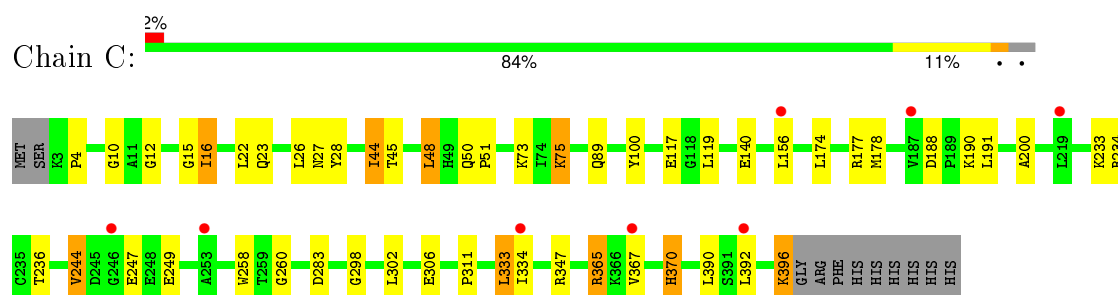
- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



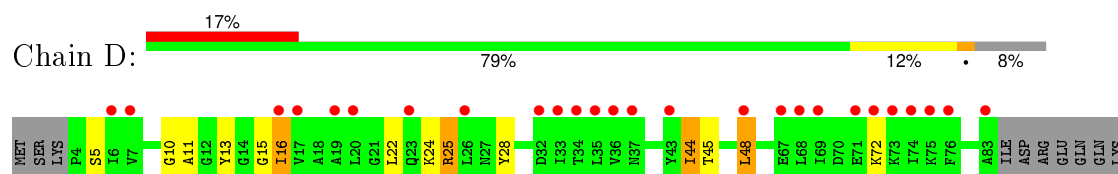
- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase

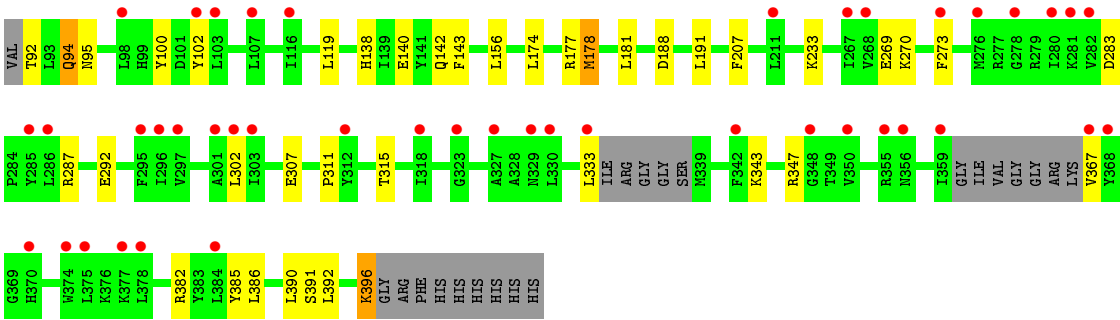


- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase



- Molecule 1: FAD-dependent pyridine nucleotide-disulfide oxidoreductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.79 Å 114.48 Å 130.59 Å 90.00° 92.04° 90.00°	Depositor
Resolution (Å)	64.61 – 2.50 64.53 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (64.61-2.50) 99.5 (64.53-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.51 Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.217 , 0.269 0.219 , 0.269	Depositor DCC
R_{free} test set	3717 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	69.2	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.2	EDS
Estimated twinning fraction	0.026 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 73728 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12454	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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, 0.375 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

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	0.61	2/3111 (0.1%)	0.79	7/4214 (0.2%)
1	B	0.61	3/3120 (0.1%)	0.73	6/4226 (0.1%)
1	C	0.57	4/3098 (0.1%)	0.63	4/4197 (0.1%)
1	D	0.57	6/2945 (0.2%)	0.74	11/3991 (0.3%)
All	All	0.59	15/12274 (0.1%)	0.72	28/16628 (0.2%)

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

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

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	28	TYR	CB-CG	11.55	1.69	1.51
1	C	28	TYR	CA-CB	-9.54	1.32	1.53
1	D	102	TYR	CA-CB	9.51	1.74	1.53
1	D	269	GLU	CA-CB	-8.21	1.35	1.53
1	D	307	GLU	CB-CG	8.02	1.67	1.52
1	A	347	ARG	CB-CG	-7.79	1.31	1.52
1	C	370	HIS	CB-CG	-6.80	1.37	1.50
1	D	347	ARG	CB-CG	-6.59	1.34	1.52
1	B	190	LYS	CG-CD	6.05	1.73	1.52
1	C	75	LYS	CB-CG	-5.97	1.36	1.52
1	A	308	ASN	N-CA	5.82	1.57	1.46
1	D	273	PHE	CB-CG	5.33	1.60	1.51
1	C	365	ARG	CB-CG	5.15	1.66	1.52
1	B	395	LYS	CB-CG	-5.07	1.38	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	25	ARG	CA-CB	-5.01	1.43	1.53

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	307	GLU	O-C-N	-21.69	88.00	122.70
1	B	28	TYR	CB-CG-CD2	16.30	130.78	121.00
1	B	28	TYR	CB-CG-CD1	-16.14	111.31	121.00
1	A	307	GLU	CB-CA-C	-13.97	82.46	110.40
1	D	28	TYR	CB-CA-C	12.70	135.81	110.40
1	D	28	TYR	CA-CB-CG	10.23	132.83	113.40
1	D	28	TYR	N-CA-CB	-9.74	93.07	110.60
1	A	307	GLU	CA-C-N	9.43	137.94	117.20
1	A	310	ARG	CB-CG-CD	-8.99	88.24	111.60
1	D	207	PHE	CB-CG-CD2	-8.95	114.53	120.80
1	D	207	PHE	CB-CG-CD1	8.14	126.50	120.80
1	D	343	LYS	CA-CB-CG	8.14	131.30	113.40
1	A	247	GLU	CB-CA-C	7.62	125.64	110.40
1	B	347	ARG	CG-CD-NE	-7.12	96.85	111.80
1	A	307	GLU	C-N-CA	6.68	138.41	121.70
1	D	270	LYS	CB-CA-C	-6.47	97.45	110.40
1	C	247	GLU	CB-CA-C	6.37	123.15	110.40
1	D	307	GLU	CA-CB-CG	-6.31	99.51	113.40
1	B	190	LYS	CB-CG-CD	-6.25	95.36	111.60
1	D	95	ASN	CA-CB-CG	-6.14	99.89	113.40
1	C	347	ARG	CB-CG-CD	5.90	126.95	111.60
1	B	396	LYS	CB-CG-CD	5.75	126.55	111.60
1	D	72	LYS	CB-CA-C	5.57	121.53	110.40
1	C	306	GLU	N-CA-CB	-5.47	100.75	110.60
1	C	28	TYR	CB-CA-C	5.45	121.31	110.40
1	B	16	ILE	CG1-CB-CG2	-5.14	100.09	111.40
1	A	252	LYS	CB-CG-CD	-5.11	98.32	111.60
1	D	95	ASN	CB-CA-C	-5.02	100.36	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	307	GLU	Mainchain

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	3049	0	3084	34	0
1	B	3055	0	3090	33	0
1	C	3039	0	3075	23	0
1	D	2886	0	2904	26	0
2	A	53	0	31	2	0
2	B	53	0	31	1	0
2	C	53	0	31	1	0
2	D	53	0	31	2	0
3	A	65	0	0	0	0
3	B	77	0	0	3	0
3	C	53	0	0	0	0
3	D	18	0	0	0	0
All	All	12454	0	12277	114	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 (114) 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:C:117:GLU:HB3	1:C:234:ARG:HG2	1.57	0.86
1:B:178:MET:HA	1:B:178:MET:HE3	1.70	0.74
1:A:178:MET:HA	1:A:178:MET:HE3	1.71	0.73
1:A:390:LEU:HA	1:A:393:VAL:HG22	1.71	0.72
1:C:26:LEU:O	1:C:73:LYS:NZ	2.22	0.72
1:A:302:LEU:HD11	1:A:311:PRO:HB3	1.72	0.70
1:D:178:MET:HA	1:D:178:MET:HE3	1.74	0.68
1:D:386:LEU:HD23	1:D:386:LEU:O	1.93	0.67
1:B:386:LEU:HD23	1:B:386:LEU:O	1.95	0.66
1:A:4:PRO:HD2	1:A:31:ALA:HA	1.77	0.66
1:D:5:SER:OG	1:D:100:TYR:HA	2.01	0.61
1:A:5:SER:OG	1:A:100:TYR:HA	2.01	0.60
1:D:315:THR:HB	2:D:601:FAD:O2	2.01	0.60
1:D:13:TYR:CD2	2:D:601:FAD:O4'	2.54	0.60
1:D:94:GLN:HA	1:D:94:GLN:NE2	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:GLU:CD	1:A:71:GLU:H	2.06	0.57
1:B:5:SER:OG	1:B:100:TYR:HA	2.05	0.57
1:C:45:THR:HA	1:C:48:LEU:HD22	1.86	0.57
1:B:338:SER:HB3	1:C:200:ALA:HB2	1.88	0.56
1:B:45:THR:HA	1:B:48:LEU:HD22	1.88	0.56
1:A:188:ASP:HB3	1:A:191:LEU:HD12	1.88	0.56
1:A:71:GLU:N	1:A:71:GLU:OE2	2.22	0.55
1:D:45:THR:HA	1:D:48:LEU:HD22	1.87	0.55
1:B:178:MET:HE1	1:B:181:LEU:HD23	1.90	0.54
1:C:396:LYS:C	1:C:396:LYS:HD2	2.29	0.53
1:D:302:LEU:HD11	1:D:311:PRO:HB3	1.90	0.53
1:B:188:ASP:HB3	1:B:191:LEU:HD12	1.90	0.53
1:D:143:PHE:CZ	1:D:178:MET:HE1	2.44	0.53
1:C:302:LEU:HD11	1:C:311:PRO:HB3	1.90	0.53
1:C:188:ASP:HB3	1:C:191:LEU:HD12	1.90	0.53
1:A:44:ILE:N	1:A:44:ILE:CD1	2.72	0.52
1:D:188:ASP:HB3	1:D:191:LEU:HD12	1.90	0.52
1:A:112:GLU:OE2	1:A:263:ARG:HD2	2.10	0.51
1:C:283:ASP:C	1:C:283:ASP:OD1	2.49	0.51
1:B:89:GLN:HA	1:B:100:TYR:CZ	2.46	0.51
1:A:143:PHE:CZ	1:A:178:MET:HE1	2.46	0.51
1:D:386:LEU:HD23	1:D:386:LEU:C	2.29	0.51
1:D:283:ASP:C	1:D:283:ASP:OD1	2.50	0.50
1:B:283:ASP:OD1	1:B:283:ASP:C	2.50	0.49
1:B:306:GLU:HG3	3:B:762:HOH:O	2.10	0.49
1:B:178:MET:HA	1:B:178:MET:CE	2.39	0.49
1:D:44:ILE:CD1	1:D:44:ILE:N	2.75	0.49
1:A:178:MET:HE3	1:A:181:LEU:HB3	1.95	0.49
1:A:6:ILE:HD11	1:A:334:ILE:HD11	1.93	0.49
1:C:117:GLU:OE1	1:C:234:ARG:HD3	2.12	0.49
1:C:244:VAL:HG21	1:C:249:GLU:OE1	2.13	0.49
1:D:178:MET:HE3	1:D:181:LEU:HB3	1.93	0.49
1:B:315:THR:HB	2:B:601:FAD:O2	2.12	0.49
1:B:326:VAL:HA	3:B:730:HOH:O	2.12	0.49
1:B:302:LEU:HD11	1:B:311:PRO:HB3	1.95	0.48
1:B:386:LEU:C	1:B:386:LEU:HD23	2.34	0.48
1:A:112:GLU:HG3	1:A:114:PHE:CD1	2.49	0.47
1:C:258:TRP:CZ2	1:C:260:GLY:HA3	2.49	0.47
1:C:298:GLY:HA3	2:C:601:FAD:O2P	2.14	0.47
1:D:287:ARG:NH1	1:D:292:GLU:HG3	2.30	0.47
1:A:178:MET:HE1	1:A:181:LEU:HD23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:GLU:HG3	1:A:114:PHE:CE1	2.49	0.47
1:A:283:ASP:OD1	1:A:283:ASP:C	2.53	0.47
1:D:178:MET:HA	1:D:178:MET:CE	2.40	0.46
1:B:338:SER:CB	1:C:200:ALA:HB2	2.45	0.46
1:A:44:ILE:N	1:A:44:ILE:HD13	2.31	0.45
1:C:89:GLN:HA	1:C:100:TYR:CZ	2.51	0.45
1:A:178:MET:HA	1:A:178:MET:CE	2.42	0.45
1:B:178:MET:HE3	1:B:181:LEU:HB3	1.99	0.45
1:C:4:PRO:HG2	1:C:334:ILE:HD12	1.98	0.44
1:C:44:ILE:CD1	1:C:44:ILE:N	2.80	0.44
1:A:114:PHE:CD1	1:A:114:PHE:N	2.85	0.44
1:A:143:PHE:CE2	1:A:178:MET:CE	3.00	0.44
1:D:396:LYS:HD3	1:D:396:LYS:O	2.17	0.44
1:B:140:GLU:OE2	1:B:177:ARG:NH2	2.51	0.44
1:C:12:GLY:O	1:C:16:ILE:HB	2.17	0.44
1:A:316:ALA:HB2	2:A:601:FAD:H2'	2.00	0.43
1:D:138:HIS:O	1:D:142:GLN:HG2	2.18	0.43
1:B:44:ILE:CD1	1:B:44:ILE:N	2.81	0.43
1:C:50:GLN:HB2	1:C:51:PRO:CD	2.48	0.43
1:C:140:GLU:OE2	1:C:177:ARG:NH2	2.52	0.43
1:B:382:ARG:HG2	1:B:382:ARG:HH11	1.83	0.43
1:A:298:GLY:HA3	2:A:601:FAD:O2P	2.18	0.43
1:A:12:GLY:O	1:A:16:ILE:HB	2.18	0.43
1:B:6:ILE:HD11	1:B:334:ILE:HD11	2.00	0.43
1:A:138:HIS:O	1:A:142:GLN:HG2	2.20	0.42
1:B:138:HIS:O	1:B:142:GLN:HG2	2.20	0.42
1:D:143:PHE:CE2	1:D:178:MET:HE1	2.54	0.42
1:A:140:GLU:OE2	1:A:177:ARG:NH2	2.52	0.42
1:A:3:LYS:HA	1:A:3:LYS:HD2	1.89	0.42
1:B:143:PHE:CE2	1:B:178:MET:CE	3.03	0.42
1:A:143:PHE:CZ	1:A:178:MET:CE	3.03	0.42
1:C:10:GLY:O	1:C:15:GLY:HA3	2.19	0.42
1:D:385:TYR:CD1	1:D:385:TYR:C	2.93	0.42
1:B:89:GLN:HA	1:B:100:TYR:CE1	2.55	0.42
1:B:79:ASP:OD2	1:B:95:ASN:HB2	2.20	0.42
1:C:119:LEU:HA	1:C:119:LEU:HD12	1.93	0.42
1:B:375:LEU:O	1:B:379:ILE:HG13	2.20	0.41
1:A:5:SER:HA	1:A:32:ASP:HB2	2.02	0.41
1:A:375:LEU:O	1:A:379:ILE:HG13	2.21	0.41
1:D:140:GLU:OE2	1:D:177:ARG:NH2	2.53	0.41
1:B:143:PHE:CE2	1:B:178:MET:HE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:ARG:HD2	3:B:712:HOH:O	2.20	0.41
1:D:11:ALA:HB1	1:D:16:ILE:HD13	2.02	0.41
1:A:50:GLN:HB2	1:A:51:PRO:CD	2.50	0.41
1:D:10:GLY:O	1:D:15:GLY:HA3	2.21	0.41
1:A:385:TYR:C	1:A:385:TYR:CD1	2.94	0.41
1:D:44:ILE:HD13	1:D:44:ILE:N	2.35	0.41
1:B:16:ILE:CG2	1:B:17:VAL:N	2.84	0.41
1:C:333:LEU:HA	1:C:333:LEU:HD12	1.97	0.41
1:B:10:GLY:O	1:B:15:GLY:HA3	2.21	0.40
1:A:154:ASP:OD1	1:A:154:ASP:N	2.50	0.40
1:D:119:LEU:HA	1:D:119:LEU:HD12	1.94	0.40
1:C:23:GLN:HA	1:C:73:LYS:HE2	2.03	0.40
1:B:50:GLN:HB2	1:B:51:PRO:CD	2.52	0.40
1:A:305:ASN:O	1:A:309:ASN:N	2.53	0.40
1:D:386:LEU:CD2	1:D:386:LEU:C	2.90	0.40
1:B:11:ALA:HB1	1:B:16:ILE:HD13	2.03	0.40
1:B:16:ILE:HG22	1:B:17:VAL:N	2.35	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	394/405 (97%)	380 (96%)	14 (4%)	0	100	100
1	B	395/405 (98%)	384 (97%)	11 (3%)	0	100	100
1	C	392/405 (97%)	381 (97%)	11 (3%)	0	100	100
1	D	366/405 (90%)	353 (96%)	13 (4%)	0	100	100
All	All	1547/1620 (96%)	1498 (97%)	49 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	320/329 (97%)	306 (96%)	14 (4%)	35	60
1	B	321/329 (98%)	302 (94%)	19 (6%)	24	44
1	C	319/329 (97%)	299 (94%)	20 (6%)	22	40
1	D	304/329 (92%)	285 (94%)	19 (6%)	22	40
All	All	1264/1316 (96%)	1192 (94%)	72 (6%)	25	46

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	22	LEU
1	A	27	ASN
1	A	44	ILE
1	A	73	LYS
1	A	90	LYS
1	A	99	HIS
1	A	121	GLU
1	A	156	LEU
1	A	174	LEU
1	A	178	MET
1	A	333	LEU
1	A	367	VAL
1	A	390	LEU
1	B	16	ILE
1	B	22	LEU
1	B	28	TYR
1	B	44	ILE
1	B	48	LEU
1	B	92	THR
1	B	156	LEU
1	B	174	LEU
1	B	178	MET
1	B	274	GLU

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Mol	Chain	Res	Type
1	B	292	GLU
1	B	333	LEU
1	B	343	LYS
1	B	367	VAL
1	B	378	LEU
1	B	390	LEU
1	B	391	SER
1	B	392	LEU
1	B	396	LYS
1	C	16	ILE
1	C	22	LEU
1	C	27	ASN
1	C	44	ILE
1	C	48	LEU
1	C	75	LYS
1	C	156	LEU
1	C	174	LEU
1	C	178	MET
1	C	190	LYS
1	C	233	LYS
1	C	236	THR
1	C	244	VAL
1	C	333	LEU
1	C	365	ARG
1	C	367	VAL
1	C	370	HIS
1	C	390	LEU
1	C	392	LEU
1	C	396	LYS
1	D	16	ILE
1	D	22	LEU
1	D	24	LYS
1	D	25	ARG
1	D	44	ILE
1	D	48	LEU
1	D	92	THR
1	D	94	GLN
1	D	156	LEU
1	D	174	LEU
1	D	178	MET
1	D	233	LYS
1	D	333	LEU

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Mol	Chain	Res	Type
1	D	367	VAL
1	D	382	ARG
1	D	390	LEU
1	D	391	SER
1	D	392	LEU
1	D	396	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	291	HIS
1	B	89	GLN
1	D	94	GLN

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	601	-	48,58,58	1.50	8 (16%)	54,89,89	2.33	11 (20%)
2	FAD	B	601	-	48,58,58	1.41	8 (16%)	54,89,89	2.07	11 (20%)
2	FAD	C	601	-	48,58,58	1.25	9 (18%)	54,89,89	2.19	16 (29%)
2	FAD	D	601	-	48,58,58	1.46	6 (12%)	54,89,89	2.12	11 (20%)

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	601	-	-	0/30/50/50	0/6/6/6
2	FAD	B	601	-	-	0/30/50/50	0/6/6/6
2	FAD	C	601	-	-	0/30/50/50	0/6/6/6
2	FAD	D	601	-	-	0/30/50/50	0/6/6/6

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	FAD	C4'-C3'	-3.55	1.46	1.53
2	A	601	FAD	C6-C5X	-3.17	1.36	1.41
2	B	601	FAD	C6-C5X	-3.17	1.36	1.41
2	A	601	FAD	C4'-C3'	-2.27	1.48	1.53
2	C	601	FAD	C6-C5X	-2.20	1.38	1.41
2	A	601	FAD	C1'-N10	-2.12	1.46	1.48
2	C	601	FAD	C2A-N3A	2.03	1.35	1.32
2	B	601	FAD	C4-C4X	2.04	1.45	1.41
2	B	601	FAD	C4X-C10	2.16	1.45	1.41
2	C	601	FAD	O4B-C1B	2.17	1.43	1.41
2	D	601	FAD	C8A-N7A	2.18	1.38	1.34
2	A	601	FAD	C8-C7	2.20	1.46	1.41
2	C	601	FAD	C5A-C4A	2.21	1.45	1.40
2	B	601	FAD	O4B-C1B	2.35	1.44	1.41
2	B	601	FAD	C5A-C4A	2.46	1.46	1.40
2	C	601	FAD	C4-C4X	2.47	1.46	1.41
2	C	601	FAD	C9A-C5X	2.59	1.47	1.42
2	C	601	FAD	C8-C7	2.63	1.48	1.41
2	B	601	FAD	C8-C7	2.68	1.48	1.41
2	C	601	FAD	C4X-C10	2.71	1.46	1.41
2	B	601	FAD	C9A-C5X	3.07	1.48	1.42
2	C	601	FAD	C9A-N10	3.07	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	C9A-C5X	3.18	1.49	1.42
2	D	601	FAD	C8-C7	3.40	1.50	1.41
2	D	601	FAD	C5A-C4A	3.46	1.48	1.40
2	A	601	FAD	C9A-N10	3.53	1.43	1.38
2	D	601	FAD	C4-C4X	3.56	1.48	1.41
2	D	601	FAD	C9A-C5X	3.60	1.49	1.42
2	A	601	FAD	C4X-C10	3.75	1.48	1.41
2	D	601	FAD	C4X-C10	4.20	1.48	1.41
2	A	601	FAD	O4B-C1B	4.67	1.47	1.41

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	N3A-C2A-N1A	-9.98	121.25	128.89
2	C	601	FAD	N3A-C2A-N1A	-7.55	123.11	128.89
2	D	601	FAD	N3A-C2A-N1A	-7.49	123.16	128.89
2	B	601	FAD	N3A-C2A-N1A	-5.92	124.36	128.89
2	A	601	FAD	C4-C4X-C10	-5.29	116.55	119.94
2	D	601	FAD	C4X-C4-N3	-4.44	117.52	123.59
2	B	601	FAD	C4X-C4-N3	-4.31	117.70	123.59
2	A	601	FAD	C1B-N9A-C4A	-3.86	121.12	126.94
2	D	601	FAD	C1B-N9A-C4A	-3.79	121.22	126.94
2	D	601	FAD	C4-C4X-C10	-3.78	117.52	119.94
2	C	601	FAD	C4X-C4-N3	-3.73	118.48	123.59
2	C	601	FAD	C4A-C5A-N7A	-3.72	106.06	109.48
2	A	601	FAD	C4X-C4-N3	-3.45	118.88	123.59
2	B	601	FAD	C4-C4X-C10	-3.26	117.86	119.94
2	C	601	FAD	P-O3P-PA	-3.11	124.00	132.73
2	C	601	FAD	C4-C4X-C10	-3.11	117.95	119.94
2	C	601	FAD	C2B-C1B-N9A	-2.84	109.95	114.29
2	B	601	FAD	C1B-N9A-C4A	-2.70	122.87	126.94
2	B	601	FAD	P-O3P-PA	-2.52	125.65	132.73
2	C	601	FAD	O3B-C3B-C2B	-2.52	103.63	111.83
2	B	601	FAD	C4A-C5A-N7A	-2.51	107.17	109.48
2	D	601	FAD	P-O3P-PA	-2.30	126.26	132.73
2	D	601	FAD	C4A-C5A-N7A	-2.26	107.40	109.48
2	B	601	FAD	O5'-P-O1P	-2.14	101.31	109.62
2	A	601	FAD	P-O3P-PA	-2.08	126.88	132.73
2	C	601	FAD	C1B-N9A-C4A	-2.08	123.80	126.94
2	C	601	FAD	C4X-C10-N10	-2.01	119.33	120.52
2	C	601	FAD	O4B-C1B-N9A	2.07	112.42	108.10
2	C	601	FAD	C5X-C9A-N10	2.08	119.20	117.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	601	FAD	C4-C4X-N5	2.09	121.26	118.72
2	B	601	FAD	O2P-P-O1P	2.11	123.97	112.53
2	A	601	FAD	C2B-C3B-C4B	2.13	106.99	102.61
2	B	601	FAD	C4X-N5-C5X	2.14	119.23	116.76
2	C	601	FAD	C2B-C3B-C4B	2.25	107.23	102.61
2	A	601	FAD	C5X-C9A-N10	2.29	119.36	117.62
2	C	601	FAD	O3P-PA-O5B	2.31	109.07	102.94
2	A	601	FAD	C2A-N1A-C6A	2.34	122.94	118.77
2	D	601	FAD	C2A-N1A-C6A	2.72	123.64	118.77
2	A	601	FAD	O2P-P-O1P	2.75	127.44	112.53
2	D	601	FAD	C5X-C9A-N10	3.04	119.93	117.62
2	D	601	FAD	C4X-N5-C5X	3.30	120.56	116.76
2	C	601	FAD	C4X-N5-C5X	3.49	120.78	116.76
2	B	601	FAD	C1'-N10-C9A	3.52	122.81	118.86
2	C	601	FAD	C1'-N10-C9A	3.87	123.21	118.86
2	A	601	FAD	C1'-N10-C9A	4.82	124.27	118.86
2	A	601	FAD	C4-N3-C2	7.42	121.66	115.25
2	C	601	FAD	C4-N3-C2	7.73	121.93	115.25
2	D	601	FAD	C4-N3-C2	7.87	122.05	115.25
2	B	601	FAD	C4-N3-C2	8.54	122.63	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	2	0
2	B	601	FAD	1	0
2	C	601	FAD	1	0
2	D	601	FAD	2	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	395/405 (97%)	0.27	7 (1%) 71 75	36, 67, 104, 130	19 (4%)
1	B	395/405 (97%)	0.36	6 (1%) 76 79	39, 70, 103, 126	9 (2%)
1	C	394/405 (97%)	0.36	8 (2%) 68 72	45, 76, 112, 135	17 (4%)
1	D	373/405 (92%)	1.00	69 (18%) 2 2	53, 101, 139, 164	20 (5%)
All	All	1557/1620 (96%)	0.49	90 (5%) 26 30	36, 76, 122, 164	65 (4%)

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	302	LEU	7.4
1	D	282	VAL	6.3
1	D	295	PHE	6.2
1	D	34	THR	6.1
1	D	280	ILE	5.3
1	D	76	PHE	5.2
1	D	7	VAL	5.1
1	A	84	ILE	5.1
1	D	36	VAL	4.9
1	D	267	ILE	4.9
1	D	98	LEU	4.8
1	A	3	LYS	4.7
1	D	303	ILE	4.7
1	D	367	VAL	4.5
1	D	35	LEU	4.5
1	D	327	ALA	4.2
1	D	350	VAL	4.1
1	D	374	TRP	4.0
1	D	285	TYR	3.8
1	D	103	LEU	3.6
1	D	83	ALA	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	31	ALA	3.5
1	D	330	LEU	3.4
1	D	296	ILE	3.4
1	D	23	GLN	3.4
1	D	301	ALA	3.3
1	D	348	GLY	3.2
1	D	375	LEU	3.2
1	D	342	PHE	3.1
1	D	378	LEU	3.1
1	D	73	LYS	3.1
1	D	278	GLY	3.1
1	D	72	LYS	3.1
1	D	276	MET	3.0
1	C	219	LEU	3.0
1	D	116	ILE	3.0
1	D	368	TYR	3.0
1	D	355	ARG	2.9
1	D	102	TYR	2.9
1	D	17	VAL	2.9
1	D	211	LEU	2.8
1	D	32	ASP	2.8
1	A	207	PHE	2.8
1	D	67	GLU	2.8
1	D	359	ILE	2.7
1	D	71	GLU	2.7
1	D	356	ASN	2.7
1	D	19	ALA	2.7
1	D	74	ILE	2.7
1	D	16	ILE	2.7
1	D	6	ILE	2.6
1	D	286	LEU	2.6
1	D	20	LEU	2.6
1	A	29	ASN	2.6
1	B	16	ILE	2.6
1	C	367	VAL	2.5
1	D	297	VAL	2.5
1	D	312	TYR	2.5
1	C	253	ALA	2.5
1	D	37	ASN	2.5
1	D	68	LEU	2.4
1	D	333	LEU	2.4
1	D	26	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	268	VAL	2.3
1	D	48	LEU	2.3
1	B	268	VAL	2.3
1	D	75	LYS	2.3
1	C	334	ILE	2.3
1	D	318	ILE	2.3
1	C	392	LEU	2.3
1	D	281	LYS	2.2
1	B	295	PHE	2.2
1	B	102	TYR	2.2
1	D	69	ILE	2.2
1	D	323	GLY	2.2
1	A	91	VAL	2.2
1	B	74	ILE	2.2
1	D	370	HIS	2.1
1	C	187	VAL	2.1
1	D	329	ASN	2.1
1	D	107	LEU	2.1
1	D	384	LEU	2.1
1	D	33	ILE	2.0
1	D	377	LYS	2.0
1	B	33	ILE	2.0
1	C	156	LEU	2.0
1	D	43	TYR	2.0
1	D	273	PHE	2.0
1	C	246	GLY	2.0
1	A	6	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. 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, 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	LLDF	B-factors(Å ²)	Q<0.9
2	FAD	C	601	53/53	0.98	0.18	0.22	44,54,75,88	0
2	FAD	B	601	53/53	0.98	0.17	-0.31	28,46,58,63	0
2	FAD	A	601	53/53	0.98	0.16	-0.41	39,49,59,64	0
2	FAD	D	601	53/53	0.94	0.17	-0.72	63,80,106,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.