



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:00 am GMT

PDB ID : 1EBD
Title : DIHYDROLIPOAMIDE DEHYDROGENASE COMPLEXED WITH THE
BINDING DOMAIN OF THE DIHYDROLIPOAMIDE ACETYLASE
Authors : Mande, S.S.; Sarfaty, S.; Allen, M.D.; Perham, R.N.; Hol, W.G.J.
Deposited on : 1996-02-03
Resolution : 2.60 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

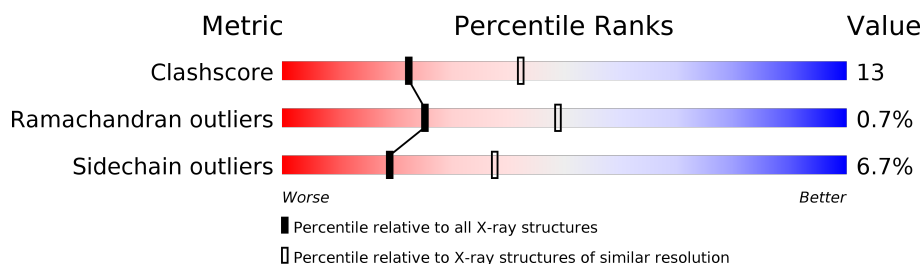
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	455	
1	B	455	
2	C	41	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7146 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.

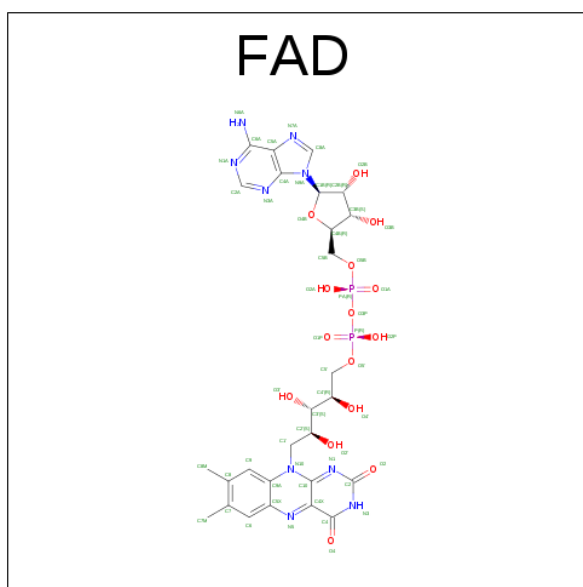
- Molecule 1 is a protein called DIHYDROLIPOAMIDE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total	C	N	O	S	0	0	0
			3351	2121	568	652	10			
1	B	455	Total	C	N	O	S	0	0	0
			3348	2118	568	652	10			

- Molecule 2 is a protein called DIHYDROLIPOAMIDE ACETYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	41	Total	C	N	O	S	0	0	0
			304	191	59	53	1			

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

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

- Molecule 4 is water.

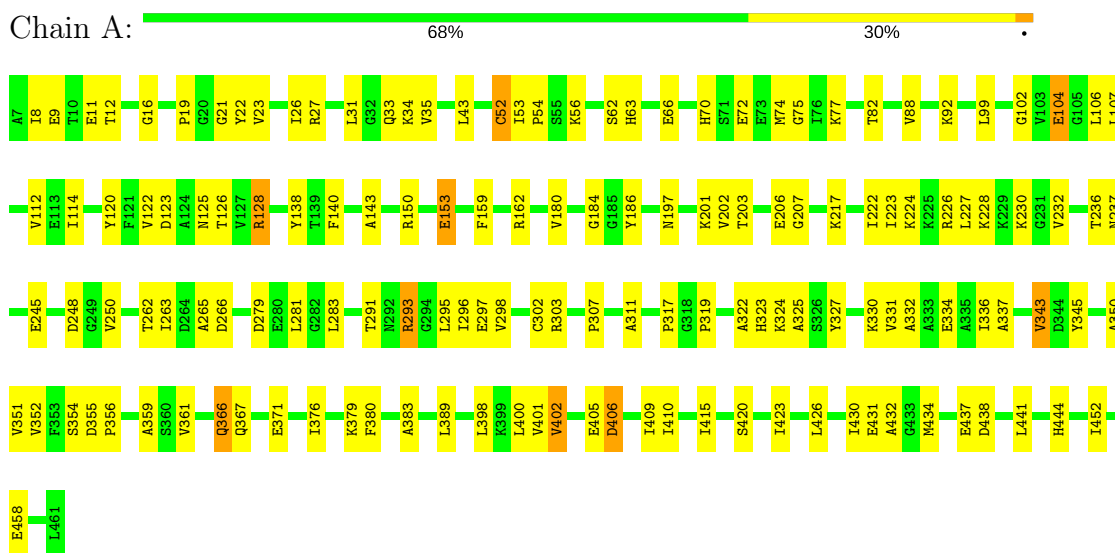
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total	O	0	0
			21	21		
4	B	16	Total	O	0	0
			16	16		

3 Residue-property plots

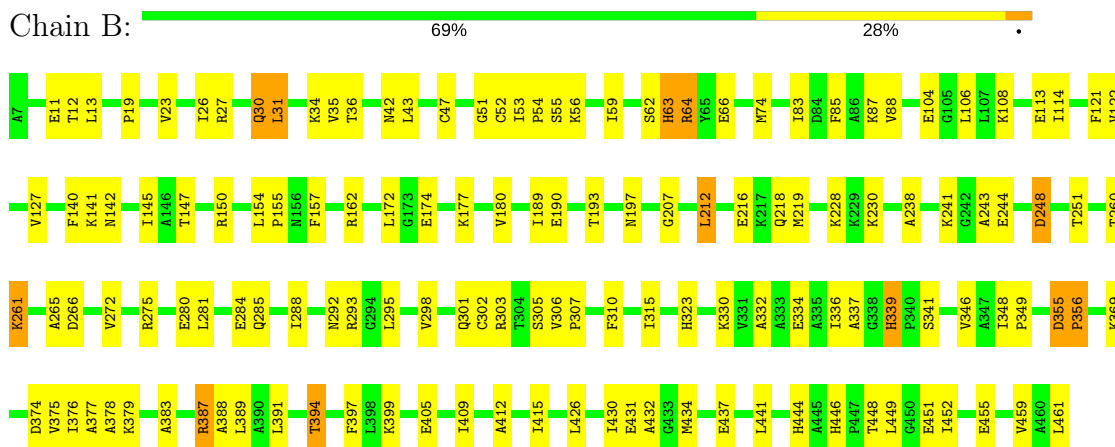
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

• Molecule 1: DIHYDROLIPOAMIDE DEHYDROGENASE



• Molecule 1: DIHYDROLIPOAMIDE DEHYDROGENASE



• Molecule 2: DIHYDROLIPOAMIDE ACETYLTRANSFERASE



I130	A131	M132	P133	A139	K142	D145	I146	R147	L148	V149	Q150	G151	T152	G153	K154	N155	G156	R157	V158	L159	K160	E161	D162	I163	D164	G170
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	106.60Å 106.60Å 204.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.60)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.215 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7146	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

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	0.44	2/3397 (0.1%)	0.68	0/4600
1	B	0.40	0/3394	0.66	1/4596 (0.0%)
2	C	0.40	0/306	0.60	0/406
All	All	0.42	2/7097 (0.0%)	0.67	1/9602 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	153	GLU	CB-CG	7.04	1.65	1.52
1	A	153	GLU	CG-CD	5.87	1.60	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	356	PRO	N-CA-C	-5.54	97.69	112.10

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	3351	0	3411	95	0
1	B	3348	0	3404	89	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	304	0	321	14	0
3	A	53	0	31	1	0
3	B	53	0	31	0	0
4	A	21	0	0	0	0
4	B	16	0	0	0	0
All	All	7146	0	7198	187	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 13.

All (187) 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:A:184:GLY:H	1:A:206:GLU:HG2	1.41	0.85
1:B:11:GLU:HG3	1:B:141:LYS:HD3	1.59	0.84
1:B:383:ALA:HA	1:B:394:THR:OG1	1.78	0.83
1:A:12:THR:HB	1:A:35:VAL:HG12	1.63	0.81
1:A:355:ASP:HB3	1:A:356:PRO:HD3	1.63	0.81
1:A:217:LYS:H	1:A:217:LYS:HD2	1.47	0.79
1:B:303:ARG:HG2	1:B:310:PHE:CE2	2.22	0.76
1:A:120:TYR:HB3	1:A:128:ARG:HG3	1.68	0.73
1:A:376:ILE:HG12	1:A:402:VAL:HG22	1.72	0.71
1:A:153:GLU:HG3	1:A:159:PHE:CD2	2.25	0.71
1:A:54:PRO:HB3	1:A:92:LYS:HE2	1.73	0.70
1:A:99:LEU:HD21	1:B:389:LEU:HD13	1.72	0.70
1:A:291:THR:HB	1:A:293:ARG:NH2	2.07	0.69
1:A:323:HIS:NE2	1:B:446:HIS:ND1	2.39	0.69
1:B:47:CYS:HG	1:B:52:CYS:CB	2.05	0.69
1:A:53:ILE:HB	1:A:54:PRO:HD3	1.75	0.68
1:B:154:LEU:HD13	1:B:272:VAL:HG21	1.74	0.68
1:A:298:VAL:HB	1:A:302:CYS:HA	1.76	0.67
1:A:75:GLY:HA2	1:B:87:LYS:HD3	1.77	0.66
1:B:127:VAL:HG13	1:B:140:PHE:HE2	1.62	0.65
1:A:303:ARG:HD3	1:A:307:PRO:HA	1.79	0.65
1:B:298:VAL:HG13	1:B:302:CYS:HA	1.77	0.65
1:A:224:LYS:O	1:A:228:LYS:HG3	1.96	0.65
1:A:226:ARG:O	1:A:230:LYS:HG3	1.98	0.63
1:B:292:ASN:HD21	1:B:293:ARG:HH11	1.46	0.63
1:A:35:VAL:HG23	1:A:112:VAL:HG13	1.80	0.63
1:A:52:CYS:O	1:A:56:LYS:HG2	1.99	0.62
1:B:455:GLU:O	1:B:459:VAL:HG23	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:351:VAL:HG11	1:A:420:SER:HB3	1.83	0.61
1:A:222:ILE:HD12	1:A:415:ILE:HD13	1.84	0.60
1:B:244:GLU:CG	1:B:251:THR:HB	2.32	0.60
1:A:366:GLN:HG3	1:A:367:GLN:H	1.67	0.60
1:A:120:TYR:HB3	1:A:128:ARG:CG	2.33	0.58
1:B:426:LEU:O	1:B:430:ILE:HG12	2.03	0.58
1:A:207:GLY:O	1:A:237:ASN:HA	2.04	0.58
1:A:33:GLN:HE21	1:A:337:ALA:HB2	1.69	0.58
1:B:218:GLN:HE22	1:B:379:LYS:HE3	1.67	0.58
1:B:177:LYS:HB3	1:B:266:ASP:OD2	2.04	0.57
1:B:51:GLY:O	1:B:55:SER:HB2	2.04	0.57
1:B:63:HIS:CE1	1:B:66:GLU:OE1	2.56	0.57
1:B:284:GLU:H	1:B:284:GLU:CD	2.08	0.57
2:C:154:LYS:HZ2	2:C:155:ASN:HB2	1.69	0.57
1:A:322:ALA:O	1:A:325:ALA:HB3	2.06	0.56
1:A:355:ASP:HB3	1:A:356:PRO:CD	2.36	0.56
1:B:19:PRO:O	1:B:23:VAL:HG13	2.06	0.55
1:B:261:LYS:HA	1:B:261:LYS:HE3	1.88	0.55
1:B:180:VAL:HG23	1:B:265:ALA:HB2	1.89	0.55
1:A:27:ARG:HD3	1:A:31:LEU:HD13	1.88	0.54
1:B:355:ASP:HB3	1:B:356:PRO:HD3	1.88	0.54
1:B:412:ALA:HB1	1:B:426:LEU:HD13	1.89	0.54
1:A:444:HIS:O	1:B:323:HIS:ND1	2.41	0.54
1:B:53:ILE:HB	1:B:54:PRO:HD3	1.90	0.54
1:B:180:VAL:HG23	1:B:265:ALA:CB	2.38	0.54
1:A:16:GLY:O	1:A:21:GLY:HA3	2.09	0.53
1:A:367:GLN:O	1:A:371:GLU:HB2	2.07	0.53
1:B:446:HIS:ND1	1:B:451:GLU:OE1	2.42	0.53
1:B:437:GLU:HG2	1:B:441:LEU:HD22	1.91	0.53
2:C:145:ASP:O	2:C:149:VAL:HG23	2.08	0.53
1:A:43:LEU:HD13	1:A:107:LEU:HD12	1.91	0.52
1:A:43:LEU:HD12	1:A:104:GLU:HG2	1.90	0.52
1:B:284:GLU:HG2	1:B:285:GLN:OE1	2.10	0.52
1:B:207:GLY:HA2	1:B:238:ALA:O	2.10	0.51
1:B:216:GLU:HB2	1:B:219:MET:HG3	1.91	0.51
2:C:154:LYS:NZ	2:C:155:ASN:HB2	2.25	0.51
1:B:241:LYS:NZ	1:B:260:THR:HG23	2.25	0.51
1:A:319:PRO:O	1:A:324:LYS:HD3	2.11	0.51
1:B:332:ALA:O	1:B:336:ILE:HG13	2.10	0.50
1:B:298:VAL:CG1	1:B:302:CYS:HA	2.41	0.50
1:A:303:ARG:CD	1:A:307:PRO:HA	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:HIS:ND1	1:B:444:HIS:O	2.43	0.50
1:A:331:VAL:HG11	1:A:343:VAL:HG13	1.92	0.50
1:A:62:SER:HB2	1:A:197:ASN:HB2	1.93	0.50
1:A:43:LEU:HD11	1:A:114:ILE:HD13	1.93	0.49
1:B:292:ASN:HD21	1:B:293:ARG:NH1	2.08	0.49
1:A:184:GLY:N	1:A:206:GLU:HG2	2.19	0.49
1:B:62:SER:HB2	1:B:197:ASN:HB2	1.94	0.49
1:B:431:GLU:HG3	2:C:133:PRO:HG3	1.95	0.49
1:A:227:LEU:HD12	1:A:230:LYS:HD3	1.93	0.49
1:A:11:GLU:HG3	1:A:34:LYS:HB2	1.95	0.48
1:A:186:TYR:HB2	1:A:352:VAL:HG13	1.96	0.48
1:A:437:GLU:O	1:A:441:LEU:HG	2.13	0.48
1:B:56:LYS:HD3	1:B:56:LYS:N	2.29	0.48
1:B:193:THR:HG23	1:B:230:LYS:HZ2	1.78	0.48
1:A:203:THR:HG21	1:A:263:ILE:CD1	2.44	0.48
1:A:245:GLU:HG3	1:A:250:VAL:HG22	1.96	0.48
1:A:217:LYS:CD	1:A:217:LYS:H	2.19	0.48
2:C:160:LYS:O	2:C:163:ILE:HB	2.13	0.48
1:A:380:PHE:CD2	1:A:452:ILE:HG23	2.48	0.48
1:B:12:THR:HB	1:B:35:VAL:HG22	1.95	0.47
1:B:346:VAL:CG2	1:B:430:ILE:HG22	2.44	0.47
1:B:63:HIS:CD2	1:B:66:GLU:OE1	2.67	0.47
1:B:63:HIS:NE2	1:B:66:GLU:OE1	2.47	0.47
1:A:376:ILE:CG1	1:A:402:VAL:HG22	2.42	0.47
1:B:83:ILE:HD11	1:B:88:VAL:HG21	1.95	0.47
1:A:332:ALA:O	1:A:336:ILE:HG12	2.13	0.47
1:B:27:ARG:HH21	1:B:330:LYS:HE2	1.78	0.47
1:B:437:GLU:HA	1:B:461:LEU:HD11	1.95	0.47
2:C:130:ILE:HG22	2:C:157:ARG:HB3	1.97	0.47
1:A:122:VAL:HG22	1:A:126:THR:O	2.14	0.47
1:A:361:VAL:HG12	1:A:423:ILE:HD11	1.97	0.46
1:A:201:LYS:HA	1:A:201:LYS:HD3	1.72	0.46
3:A:462:FAD:H2'	3:A:462:FAD:N1	2.30	0.46
1:A:74:MET:SD	1:B:64:ARG:HG3	2.56	0.46
1:A:227:LEU:HG	1:A:232:VAL:HB	1.97	0.46
1:B:244:GLU:HG2	1:B:251:THR:HB	1.97	0.46
2:C:152:THR:HB	2:C:159:LEU:HD13	1.97	0.46
1:A:409:ILE:HG22	1:A:430:ILE:HG12	1.98	0.46
1:A:379:LYS:HA	1:A:398:LEU:O	2.15	0.46
1:B:376:ILE:HG12	1:B:377:ALA:N	2.32	0.45
1:A:22:TYR:O	1:A:26:ILE:HG12	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:THR:HA	1:B:113:GLU:HB3	1.97	0.45
1:A:380:PHE:HD2	1:A:452:ILE:HG23	1.82	0.45
2:C:145:ASP:OD2	2:C:148:LEU:HB2	2.16	0.45
2:C:161:GLU:HA	2:C:164:ASP:HB2	1.99	0.45
1:A:351:VAL:HG11	1:A:420:SER:CB	2.47	0.45
1:A:438:ASP:OD1	2:C:132:MET:HB3	2.17	0.45
1:A:405:GLU:O	1:A:406:ASP:HB3	2.17	0.45
1:B:27:ARG:O	1:B:31:LEU:HD13	2.16	0.45
1:A:19:PRO:O	1:A:23:VAL:HG13	2.17	0.45
1:A:70:HIS:HD2	1:A:72:GLU:OE2	2.00	0.44
1:B:150:ARG:CZ	1:B:275:ARG:HD3	2.47	0.44
1:B:334:GLU:HB3	1:B:339:HIS:HB3	1.99	0.44
1:B:157:PHE:CE2	1:B:243:ALA:HB2	2.52	0.44
1:A:401:VAL:HG23	1:A:410:ILE:HG13	1.98	0.44
1:B:244:GLU:HG3	1:B:251:THR:HB	1.99	0.44
1:A:330:LYS:O	1:A:334:GLU:HG2	2.18	0.44
1:A:383:ALA:O	1:A:389:LEU:HD11	2.17	0.44
1:B:189:ILE:HG13	1:B:212:LEU:HD13	2.00	0.44
1:A:323:HIS:CE1	1:B:446:HIS:HD1	2.35	0.44
1:A:8:ILE:O	1:A:138:TYR:HA	2.18	0.44
1:A:202:VAL:HB	1:A:232:VAL:HG22	1.99	0.44
1:B:288:ILE:HD13	1:B:306:VAL:HG23	2.00	0.44
1:A:11:GLU:HB2	1:A:34:LYS:O	2.18	0.44
1:B:31:LEU:HD23	1:B:337:ALA:CB	2.48	0.44
1:A:63:HIS:CE1	1:A:66:GLU:OE1	2.70	0.44
1:B:47:CYS:HG	1:B:52:CYS:HB2	1.81	0.44
1:A:283:LEU:HD13	1:A:296:ILE:HD11	2.00	0.43
1:B:448:THR:O	1:B:451:GLU:HG2	2.18	0.43
1:B:388:ALA:HA	1:B:391:LEU:HD12	1.99	0.43
1:A:361:VAL:HG13	1:A:430:ILE:HD12	2.01	0.43
1:B:59:ILE:HD11	1:B:190:GLU:HB3	2.01	0.43
1:B:177:LYS:HD3	1:B:177:LYS:C	2.38	0.43
1:B:409:ILE:HG22	1:B:430:ILE:CD1	2.48	0.43
1:A:432:ALA:HB3	1:A:434:MET:HG3	2.00	0.43
1:B:26:ILE:HD11	1:B:106:LEU:HB3	2.00	0.43
1:B:145:ILE:O	1:B:315:ILE:HD11	2.19	0.43
1:B:374:ASP:OD2	1:B:405:GLU:HG3	2.19	0.43
1:B:432:ALA:O	2:C:157:ARG:NH2	2.52	0.43
1:A:223:ILE:O	1:A:227:LEU:HB2	2.18	0.43
1:B:154:LEU:HD12	1:B:154:LEU:HA	1.91	0.42
1:A:31:LEU:HD23	1:A:337:ALA:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:GLN:HG2	1:B:341:SER:O	2.19	0.42
1:A:120:TYR:HB2	1:A:128:ARG:CZ	2.50	0.42
1:B:288:ILE:HA	1:B:305:SER:OG	2.19	0.42
1:B:378:ALA:O	1:B:399:LYS:HA	2.19	0.42
2:C:157:ARG:HH11	2:C:157:ARG:HG2	1.84	0.42
1:B:43:LEU:HD21	1:B:114:ILE:HD13	2.02	0.42
1:A:88:VAL:HG22	1:B:74:MET:O	2.19	0.42
1:B:104:GLU:HG2	1:B:108:LYS:HE2	2.02	0.42
1:B:121:PHE:CD2	1:B:127:VAL:HG12	2.55	0.42
2:C:139:ALA:HB2	2:C:163:ILE:HD13	2.02	0.42
1:A:415:ILE:O	1:A:415:ILE:HG23	2.20	0.42
1:A:102:GLY:O	1:A:106:LEU:HD13	2.20	0.41
1:A:302:CYS:HB3	1:A:311:ALA:O	2.20	0.41
1:B:397:PHE:H	1:B:415:ILE:HG23	1.85	0.41
1:A:400:LEU:HD11	1:A:426:LEU:HD22	2.01	0.41
1:A:434:MET:SD	1:B:434:MET:CE	3.09	0.41
1:B:449:LEU:O	1:B:452:ILE:HG13	2.20	0.41
1:A:140:PHE:CE1	1:A:143:ALA:HB2	2.56	0.41
1:B:27:ARG:NH1	1:B:30:GLN:HG2	2.36	0.41
1:B:147:THR:HG21	1:B:281:LEU:HD11	2.01	0.41
1:B:305:SER:O	1:B:307:PRO:HD3	2.21	0.41
1:A:350:ALA:O	1:A:359:ALA:HA	2.21	0.41
1:A:203:THR:HG21	1:A:263:ILE:HD13	2.01	0.41
1:A:295:LEU:HD13	1:A:317:PRO:HA	2.03	0.41
1:A:401:VAL:HG23	1:A:401:VAL:O	2.20	0.41
1:A:31:LEU:HD23	1:A:337:ALA:CB	2.51	0.41
1:A:327:TYR:CE2	1:A:345:TYR:HD2	2.39	0.41
2:C:153:GLY:O	2:C:154:LYS:HB3	2.21	0.41
1:A:180:VAL:HG23	1:A:265:ALA:CB	2.50	0.41
1:B:241:LYS:HZ1	1:B:260:THR:HG23	1.86	0.41
1:B:387:ARG:NH1	1:B:391:LEU:HD21	2.36	0.40
1:B:348:ILE:HA	1:B:349:PRO:HD3	1.88	0.40
1:B:446:HIS:CE1	1:B:451:GLU:OE1	2.74	0.40
1:A:226:ARG:NH2	1:A:356:PRO:HG3	2.37	0.40
1:A:162:ARG:NH1	1:A:266:ASP:O	2.54	0.40
1:A:405:GLU:O	1:A:406:ASP:CB	2.70	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	453/455 (100%)	418 (92%)	31 (7%)	4 (1%)	20	40
1	B	453/455 (100%)	430 (95%)	22 (5%)	1 (0%)	51	76
2	C	39/41 (95%)	30 (77%)	7 (18%)	2 (5%)	2	3
All	All	945/951 (99%)	878 (93%)	60 (6%)	7 (1%)	25	49

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	ASP
1	A	248	ASP
1	B	248	ASP
2	C	152	THR
1	A	366	GLN
2	C	150	GLN
1	A	343	VAL

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	344/346 (99%)	325 (94%)	19 (6%)	25	49
1	B	343/346 (99%)	317 (92%)	26 (8%)	15	30
2	C	29/31 (94%)	26 (90%)	3 (10%)	8	15
All	All	716/723 (99%)	668 (93%)	48 (7%)	19	38

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLU
1	A	52	CYS
1	A	77	LYS
1	A	82	THR
1	A	104	GLU
1	A	123	ASP
1	A	125	ASN
1	A	128	ARG
1	A	150	ARG
1	A	236	THR
1	A	262	THR
1	A	279	ASP
1	A	281	LEU
1	A	293	ARG
1	A	297	GLU
1	A	354	SER
1	A	402	VAL
1	A	431	GLU
1	A	458	GLU
1	B	13	LEU
1	B	30	GLN
1	B	31	LEU
1	B	34	LYS
1	B	42	ASN
1	B	63	HIS
1	B	64	ARG
1	B	85	PHE
1	B	122	VAL
1	B	142	ASN
1	B	155	PRO
1	B	162	ARG
1	B	172	LEU
1	B	174	GLU
1	B	212	LEU
1	B	228	LYS
1	B	248	ASP
1	B	261	LYS
1	B	280	GLU
1	B	295	LEU
1	B	339	HIS
1	B	355	ASP
1	B	369	LYS

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Mol	Chain	Res	Type
1	B	375	VAL
1	B	387	ARG
1	B	394	THR
2	C	142	LYS
2	C	146	ILE
2	C	155	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	63	HIS
1	A	70	HIS
1	A	125	ASN
1	A	161	ASN
1	A	444	HIS
1	B	33	GLN
1	B	49	ASN
1	B	125	ASN
1	B	131	ASN
1	B	142	ASN
1	B	171	ASN
1	B	218	GLN
1	B	237	ASN
2	C	150	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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

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$
3	FAD	A	462	-	51,58,58	1.11	4 (7%)	54,89,89	2.42	9 (16%)
3	FAD	B	462	-	51,58,58	1.03	2 (3%)	54,89,89	2.13	8 (14%)

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
3	FAD	A	462	-	-	0/28/50/50	0/6/6/6
3	FAD	B	462	-	-	0/28/50/50	0/6/6/6

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	462	FAD	C5A-N7A	-2.10	1.32	1.39
3	A	462	FAD	C2-N1	-2.10	1.34	1.38
3	B	462	FAD	C5A-N7A	-2.07	1.32	1.39
3	A	462	FAD	C9A-N10	2.09	1.41	1.38
3	B	462	FAD	C4-N3	3.72	1.39	1.33
3	A	462	FAD	C4-N3	4.33	1.40	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	462	FAD	N3A-C2A-N1A	-9.88	120.26	128.86
3	A	462	FAD	N3A-C2A-N1A	-9.77	120.35	128.86
3	A	462	FAD	C4B-O4B-C1B	-7.88	101.38	109.77
3	A	462	FAD	C4X-C10-N10	-3.99	117.75	120.52
3	B	462	FAD	C4X-C10-N10	-3.59	118.03	120.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	462	FAD	C4X-C4-N3	-3.29	118.81	123.48
3	B	462	FAD	C4X-C4-N3	-3.28	118.82	123.48
3	B	462	FAD	C4B-O4B-C1B	-2.33	107.29	109.77
3	A	462	FAD	C4-C4X-N5	2.08	120.96	118.68
3	A	462	FAD	C4A-C5A-N7A	2.23	111.56	109.41
3	B	462	FAD	C6-C5X-C9A	2.54	122.30	119.00
3	A	462	FAD	C1'-N10-C9A	4.08	122.09	118.35
3	B	462	FAD	C1'-N10-C9A	4.72	122.67	118.35
3	B	462	FAD	C4-N3-C2	5.01	119.54	115.16
3	B	462	FAD	C4X-N5-C5X	5.26	122.31	116.76
3	A	462	FAD	C4-N3-C2	5.48	119.95	115.16
3	A	462	FAD	C4X-N5-C5X	5.68	122.76	116.76

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
3	A	462	FAD	1	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 ⓘ

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.