



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:28 PM UTC

PDB ID : 1EBD / pdb_00001ebd
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

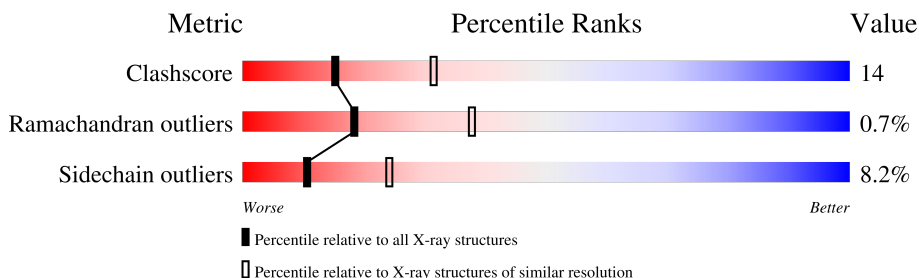
1 Overall quality at a glance

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	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$

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.

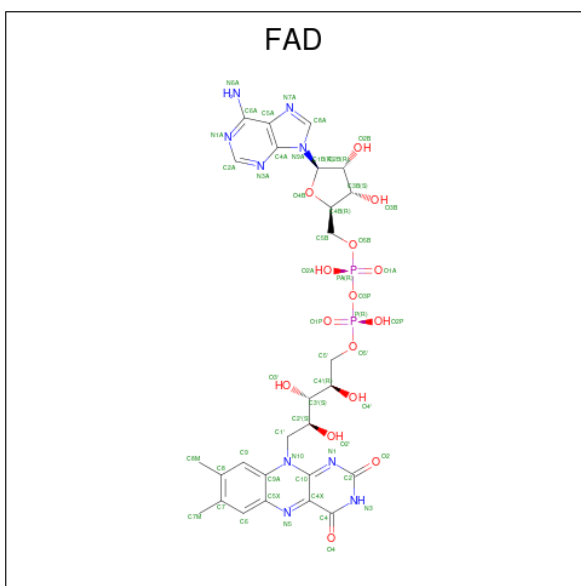
- Molecule 1 is a protein called DIHYDROLIPOAMIDE DEHYDROGENASE.

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

- Molecule 2 is a protein called DIHYDROLIPOAMIDE ACETYLTRANSFERASE.

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

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

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

- 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		

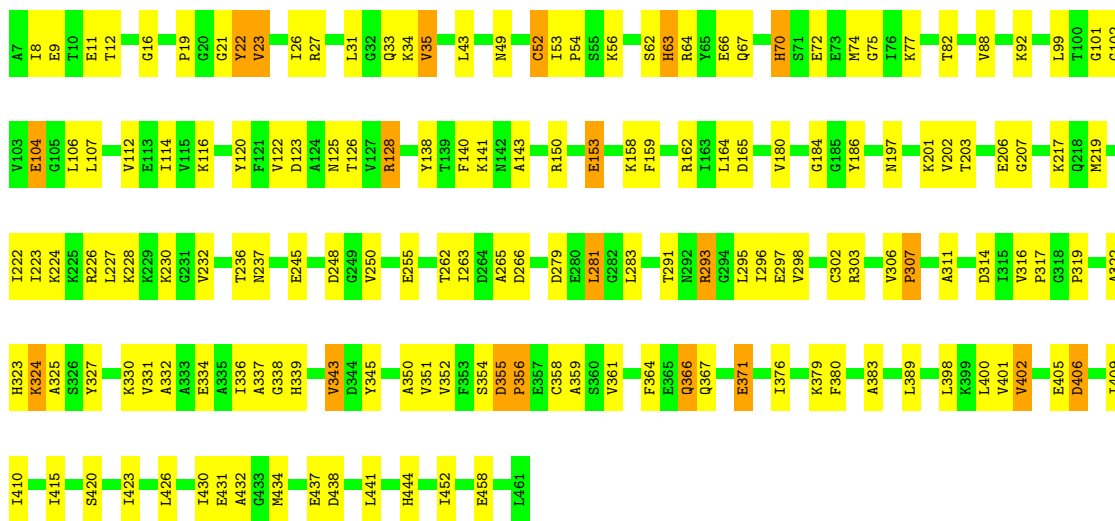
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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. 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. 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.

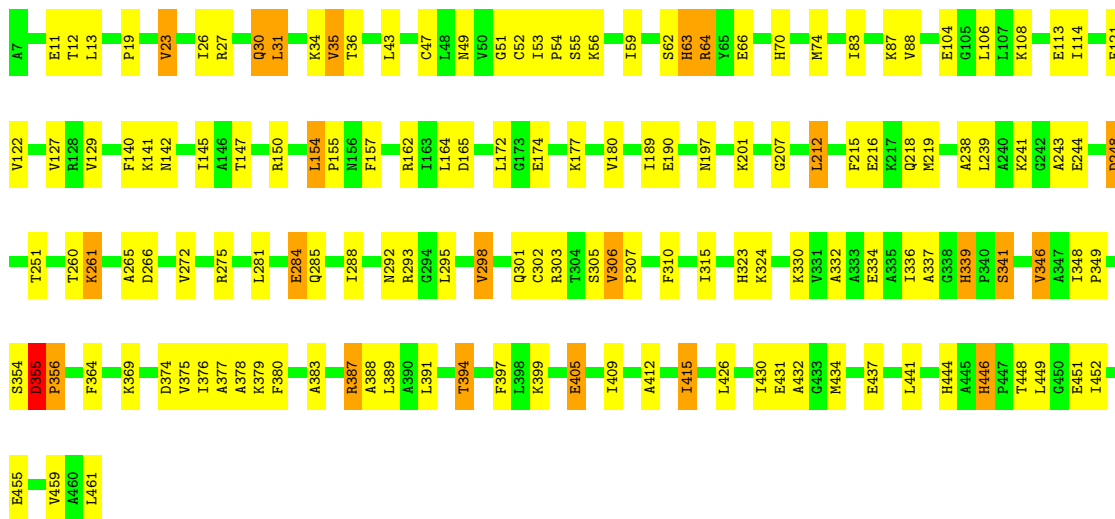
- Molecule 1: DIHYDROLIPOAMIDE DEHYDROGENASE

Chain A:  64% 31%



- Molecule 1: DIHYDROLIPOAMIDE DEHYDROGENASE

Chain B:  68% 27% 5%



● Molecule 2: DIHYDROLIPOAMIDE ACETYLTRANSFERASE

Chain C:  49% 44% 5%

4 Data and refinement statistics

Xtrriage (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.	Xtrriage
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	0.50	0/3397	1.01	21/4600 (0.5%)
1	B	0.48	0/3394	1.01	19/4596 (0.4%)
2	C	0.43	0/306	0.98	4/406 (1.0%)
All	All	0.49	0/7097	1.01	44/9602 (0.5%)

There are no bond length outliers.

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	355	ASP	CA-C-N	-8.97	109.12	120.79
1	B	355	ASP	C-N-CA	-8.97	109.12	120.79
1	A	355	ASP	CA-C-N	-8.11	109.71	119.84
1	A	355	ASP	C-N-CA	-8.11	109.71	119.84
1	B	356	PRO	N-CA-C	-7.86	97.69	110.74
1	A	364	PHE	N-CA-C	-6.90	100.31	110.52
1	B	49	ASN	N-CA-C	6.81	118.50	111.14
1	B	70	HIS	N-CA-C	6.79	119.77	110.06
1	B	165	ASP	N-CA-C	-6.75	99.49	109.81
1	B	306	VAL	CA-C-N	6.61	128.10	119.84
1	B	306	VAL	C-N-CA	6.61	128.10	119.84
1	A	306	VAL	CA-C-N	6.60	128.09	119.84
1	A	306	VAL	C-N-CA	6.60	128.09	119.84
1	A	324	LYS	N-CA-C	-6.50	104.20	111.28
1	A	70	HIS	N-CA-C	6.49	119.19	110.88
1	A	371	GLU	N-CA-C	-6.37	105.38	113.02
1	B	364	PHE	N-CA-C	-6.29	100.33	110.32
1	A	281	LEU	N-CA-C	-6.22	105.19	114.64
1	A	67	GLN	N-CA-C	6.06	117.68	111.14
1	A	153	GLU	N-CA-C	-6.05	100.00	109.25
1	A	35	VAL	N-CA-C	6.04	117.70	108.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	35	VAL	N-CA-C	5.98	117.20	108.53
1	A	164	LEU	N-CA-C	5.94	119.20	110.59
1	B	324	LYS	N-CA-C	-5.88	104.95	111.36
1	A	141	LYS	N-CA-C	-5.79	104.41	112.45
1	B	64	ARG	N-CA-C	-5.76	105.00	111.28
1	A	356	PRO	N-CA-C	-5.74	100.64	112.47
1	A	49	ASN	N-CA-C	5.71	118.36	111.40
1	B	164	LEU	N-CA-C	5.69	118.94	110.52
1	B	405	GLU	N-CA-C	5.50	119.61	113.01
1	B	306	VAL	N-CA-C	-5.39	102.65	107.56
1	A	64	ARG	N-CA-C	-5.37	105.53	111.71
2	C	155	ASN	N-CA-C	-5.35	106.69	114.12
2	C	163	ILE	N-CA-C	-5.34	105.51	110.53
1	A	101	GLY	N-CA-C	-5.32	106.34	112.73
1	B	154	LEU	CA-C-N	5.27	126.43	119.84
1	B	154	LEU	C-N-CA	5.27	126.43	119.84
1	A	314	ASP	N-CA-C	5.26	118.80	112.38
2	C	162	ASP	N-CA-C	-5.21	105.71	111.71
1	A	165	ASP	N-CA-C	-5.21	102.19	109.69
2	C	156	GLY	N-CA-C	-5.20	107.88	115.72
1	B	380	PHE	CA-C-N	5.16	125.57	119.99
1	B	380	PHE	C-N-CA	5.16	125.57	119.99
1	A	22	TYR	N-CA-C	5.04	116.85	111.36

There are no chirality outliers.

There are no planarity outliers.

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	3351	0	3411	99	0
1	B	3348	0	3404	91	0
2	C	304	0	321	14	0
3	A	53	0	31	2	0
3	B	53	0	31	0	0
4	A	21	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	16	0	0	0	0
All	All	7146	0	7198	194	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 14.

All (194) 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:HD2	1:A:217:LYS:H	1.47	0.79
1:B:303:ARG:HG2	1:B:310:PHE:CE2	2.22	0.76
1:B:47:CYS:SG	1:B:52:CYS:SG	2.64	0.75
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
1:A:351:VAL:HG11	1:A:420:SER:HB3	1.83	0.61
1:B:284:GLU:H	1:B:284:GLU:CD	2.08	0.60
1:A:222:ILE:HD12	1:A:415:ILE:HD13	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:33:GLN:HE21	1:A:337:ALA:HB2	1.69	0.58
1:A:207:GLY:O	1:A:237:ASN:HA	2.04	0.58
1:B:218:GLN:HE22	1:B:379:LYS:HE3	1.67	0.58
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:177:LYS:HB3	1:B:266:ASP:OD2	2.04	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:HE3	1:B:261:LYS:HA	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:241:LYS:HZ1	1:B:260:THR:HG23	1.74	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
2:C:154:LYS:HZ2	2:C:155:ASN:HB2	1.77	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:N	1:B:56:LYS:HD3	2.29	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:H	1:A:217:LYS:CD	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:63:HIS:CD2	1:B:66:GLU:OE1	2.67	0.47
1:B:346:VAL:CG2	1:B:430:ILE:HG22	2.44	0.47
1:B:177:LYS:HD3	1:B:177:LYS:C	2.38	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:HD3	1:A:201:LYS:HA	1.72	0.46
1:A:434:MET:HE2	1:A:438:ASP:HB3	1.96	0.46
3:A:462:FAD:N1	3:A:462:FAD:H2'	2.30	0.46
1:B:215:PHE:HB3	1:B:219:MET:HE3	1.97	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:A:22:TYR:O	1:A:26:ILE:HG12	2.17	0.45
1:B:36:THR:HA	1:B:113:GLU:HB3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
1:B:376:ILE:HG12	1:B:377:ALA:N	2.32	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:323:HIS:CE1	1:B:446:HIS:HD1	2.35	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: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:A:338:GLY:O	1:A:339:HIS:C	2.62	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: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
3:A:462:FAD:HM81	3:A:462:FAD:HM73	1.78	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:LEU:HD23	1:A:337:ALA:HB3	2.00	0.42
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:A:63:HIS:CE1	1:A:66:GLU:CD	2.97	0.42
1:A:219:MET:HE2	1:A:358:CYS:SG	2.60	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:HG2	2:C:157:ARG:HH11	1.84	0.42
1:A:88:VAL:HG22	1:B:74:MET:O	2.19	0.42
1:B:43:LEU:HD21	1:B:114:ILE:HD13	2.02	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: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:415:ILE:O	1:A:415:ILE:HG23	2.20	0.41
1:A:31:LEU:HD23	1:A:337:ALA:CB	2.51	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:327:TYR:CE2	1:A:345:TYR:HD2	2.39	0.41
1:A:401:VAL:HG23	1:A:401:VAL:O	2.20	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: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
1:B:354:SER:O	1:B:355:ASP:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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%)	14	30
1	B	453/455 (100%)	430 (95%)	22 (5%)	1 (0%)	43	66
2	C	39/41 (95%)	30 (77%)	7 (18%)	2 (5%)	1	2
All	All	945/951 (99%)	878 (93%)	60 (6%)	7 (1%)	18	38

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 [i](#)

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%)	318 (92%)	26 (8%)	12	27
1	B	343/346 (99%)	312 (91%)	31 (9%)	9	20
2	C	29/31 (94%)	27 (93%)	2 (7%)	14	32
All	All	716/723 (99%)	657 (92%)	59 (8%)	10	24

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLU
1	A	23	VAL
1	A	52	CYS
1	A	63	HIS
1	A	77	LYS
1	A	82	THR
1	A	104	GLU
1	A	116	LYS
1	A	123	ASP
1	A	125	ASN
1	A	128	ARG
1	A	150	ARG
1	A	158	LYS
1	A	236	THR
1	A	255	GLU
1	A	262	THR
1	A	279	ASP
1	A	281	LEU
1	A	293	ARG
1	A	297	GLU
1	A	307	PRO
1	A	316	VAL
1	A	354	SER
1	A	402	VAL
1	A	431	GLU
1	A	458	GLU
1	B	13	LEU
1	B	23	VAL
1	B	30	GLN
1	B	31	LEU
1	B	34	LYS
1	B	63	HIS
1	B	122	VAL
1	B	129	VAL
1	B	142	ASN
1	B	155	PRO
1	B	162	ARG
1	B	172	LEU
1	B	174	GLU
1	B	201	LYS
1	B	212	LEU
1	B	239	LEU

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Mol	Chain	Res	Type
1	B	248	ASP
1	B	261	LYS
1	B	284	GLU
1	B	295	LEU
1	B	298	VAL
1	B	339	HIS
1	B	341	SER
1	B	346	VAL
1	B	355	ASP
1	B	369	LYS
1	B	375	VAL
1	B	387	ARG
1	B	394	THR
1	B	415	ILE
1	B	446	HIS
2	C	146	ILE
2	C	155	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	63	HIS
1	A	70	HIS
1	A	136	GLN
1	A	161	ASN
1	A	237	ASN
1	A	444	HIS
1	B	49	ASN
1	B	70	HIS
1	B	131	ASN
1	B	171	ASN
1	B	218	GLN
1	B	237	ASN
1	B	366	GLN
1	B	367	GLN
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 [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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
3	FAD	A	462	-	58,58,58	1.09	5 (8%)	85,89,89	1.58	16 (18%)
3	FAD	B	462	-	58,58,58	1.04	3 (5%)	85,89,89	1.57	14 (16%)

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	-	-	2/34/50/50	0/6/6/6
3	FAD	B	462	-	-	3/34/50/50	0/6/6/6

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	462	FAD	C5A-N7A	-3.73	1.32	1.39
3	B	462	FAD	C5A-N7A	-3.67	1.32	1.39
3	B	462	FAD	C5X-N5	-3.06	1.33	1.39
3	A	462	FAD	C4A-N9A	-2.49	1.32	1.37
3	A	462	FAD	C5A-C6A	-2.36	1.34	1.41
3	B	462	FAD	C5A-C6A	-2.12	1.35	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	462	FAD	C5X-N5	-2.11	1.35	1.39
3	A	462	FAD	C10-N10	2.04	1.41	1.37

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	462	FAD	N3A-C2A-N1A	-5.50	120.26	128.58
3	A	462	FAD	N3A-C2A-N1A	-5.44	120.35	128.58
3	A	462	FAD	C4A-N9A-C8A	4.04	109.98	105.74
3	B	462	FAD	N3A-C4A-N9A	3.94	133.86	127.17
3	B	462	FAD	C5A-C4A-N3A	-3.88	121.38	126.72
3	A	462	FAD	C4B-O4B-C1B	-3.66	101.38	109.47
3	B	462	FAD	N9A-C8A-N7A	-3.60	108.83	113.94
3	B	462	FAD	C4A-N9A-C8A	3.58	109.50	105.74
3	B	462	FAD	C4-N3-C2	-3.44	119.54	125.64
3	A	462	FAD	N3A-C4A-N9A	3.37	132.90	127.17
3	B	462	FAD	C2A-N3A-C4A	3.23	119.71	111.83
3	A	462	FAD	C4-N3-C2	-3.21	119.95	125.64
3	A	462	FAD	N9A-C8A-N7A	-2.96	109.73	113.94
3	A	462	FAD	C5X-N5-C4X	2.89	122.76	118.09
3	B	462	FAD	C5X-N5-C4X	2.61	122.31	118.09
3	A	462	FAD	C2A-N3A-C4A	2.61	118.20	111.83
3	A	462	FAD	C10-N1-C2	2.45	122.16	116.85
3	B	462	FAD	O4-C4-C4X	-2.44	120.09	126.53
3	A	462	FAD	C5A-C4A-N3A	-2.38	123.44	126.72
3	B	462	FAD	C6-C5X-C9A	2.37	122.30	119.05
3	B	462	FAD	C9A-C5X-N5	-2.36	119.95	122.45
3	A	462	FAD	O4-C4-C4X	-2.36	120.31	126.53
3	B	462	FAD	C5A-N7A-C8A	2.20	106.90	103.45
3	B	462	FAD	C4X-C4-N3	2.19	118.82	113.25
3	A	462	FAD	C4X-C4-N3	2.18	118.81	113.25
3	B	462	FAD	C4'-C3'-C2'	-2.16	109.98	113.57
3	A	462	FAD	C10-C4X-N5	-2.14	120.44	124.81
3	A	462	FAD	C9A-C5X-N5	-2.14	120.19	122.45
3	A	462	FAD	C5A-C4A-N9A	-2.01	103.62	105.81
3	A	462	FAD	O3'-C3'-C2'	-2.01	104.36	108.93

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	462	FAD	PA-O3P-P-O5'

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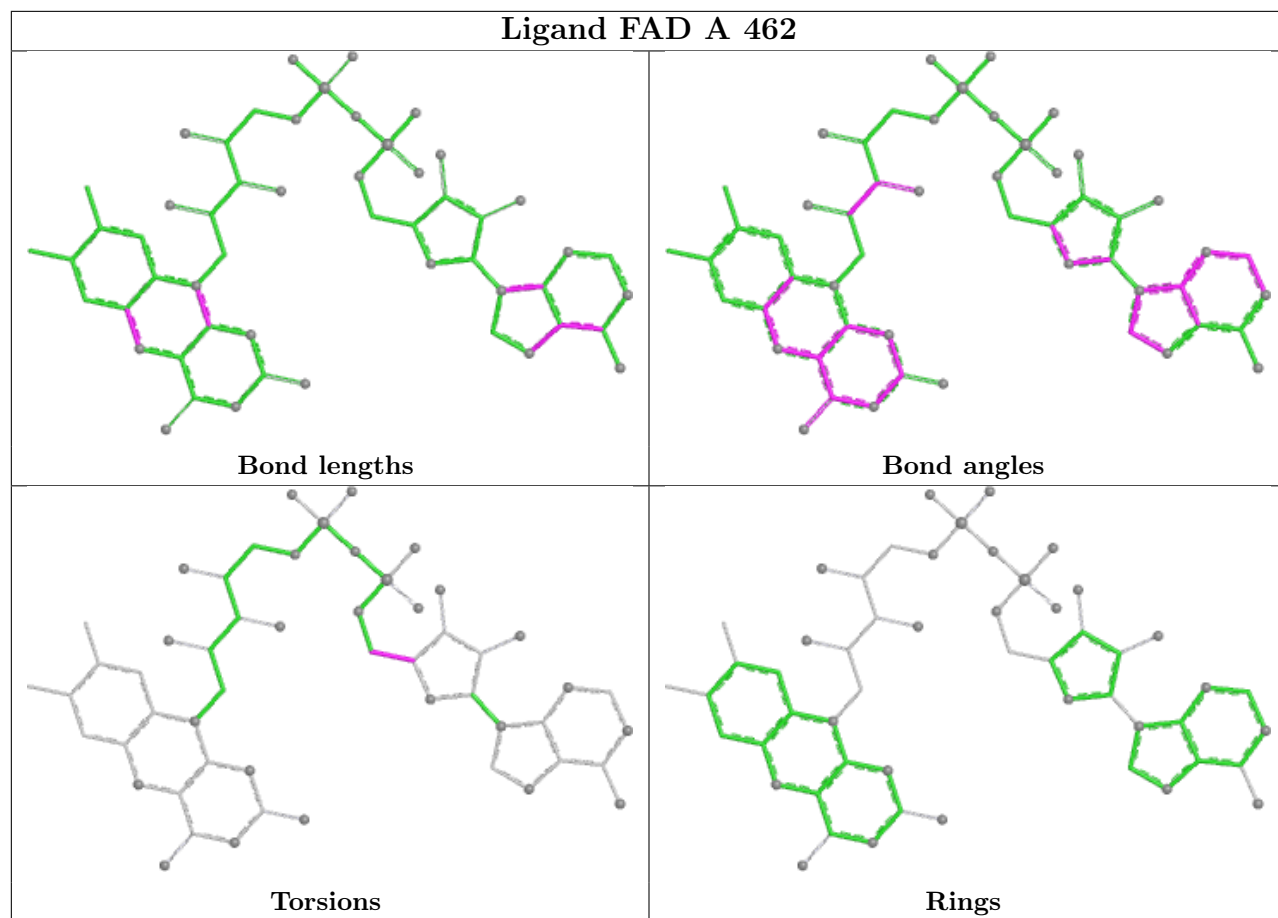
Mol	Chain	Res	Type	Atoms
3	A	462	FAD	C3B-C4B-C5B-O5B
3	A	462	FAD	O4B-C4B-C5B-O5B
3	B	462	FAD	P-O3P-PA-O1A
3	B	462	FAD	P-O3P-PA-O2A

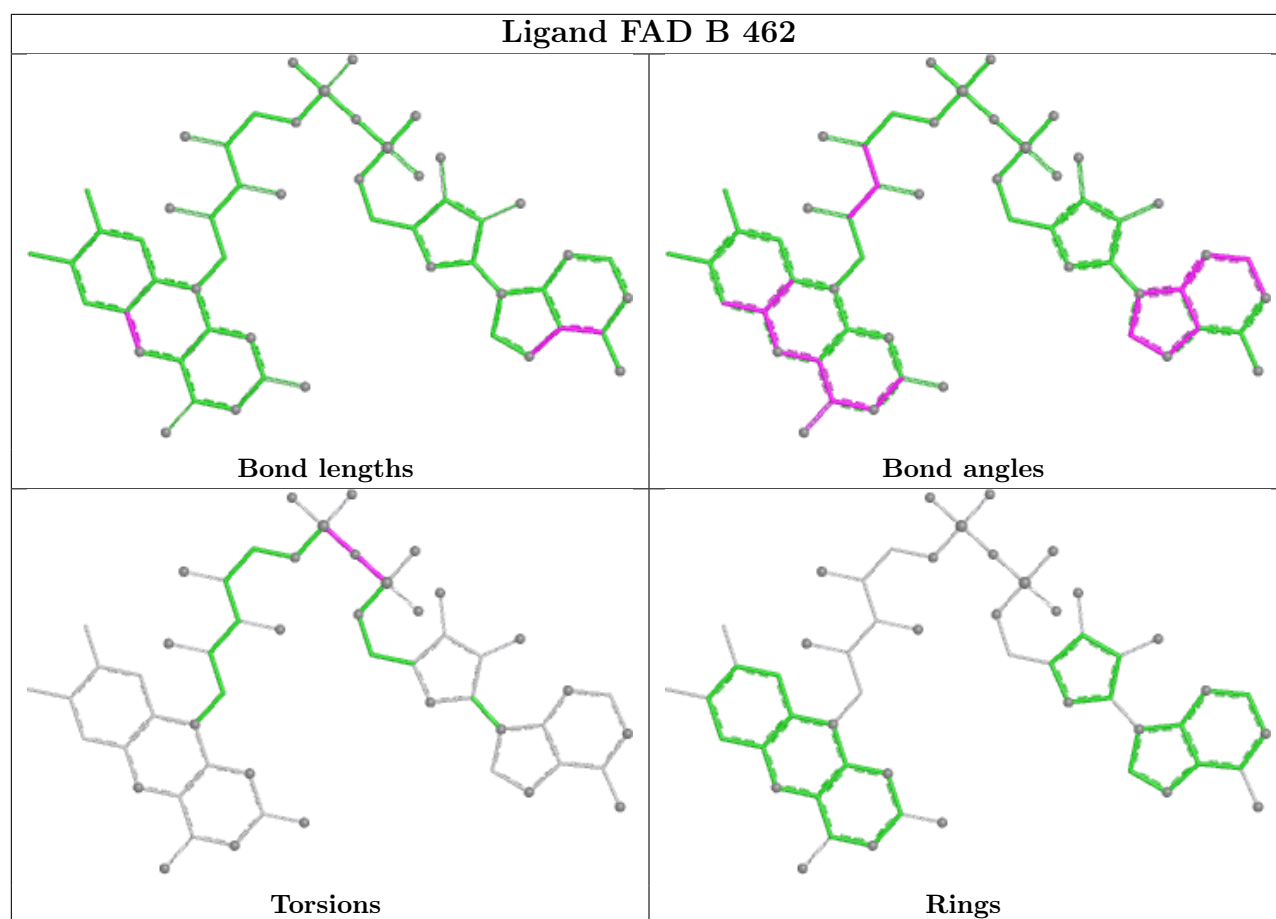
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	462	FAD	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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.