



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 07:55 pm BST

PDB ID : 4Y9L
Title : Crystal Structure of Caenorhabditis elegans ACDH-11
Authors : Li, Z.J.; Zhai, Y.J.; Zhang, K.; Sun, F.
Deposited on : 2015-02-17
Resolution : 2.27 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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

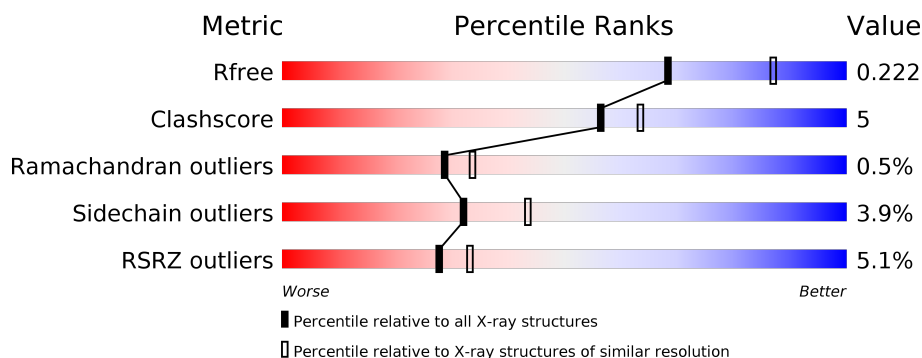
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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}	130704	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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 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 $\leq 5\%$. 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	593	<div> <div>5%</div> <div> <div></div> <div>88%</div> <div>10%</div> </div> <div></div> </div>
1	B	593	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>12%</div> </div> <div></div> </div>

2 Entry composition [i](#)

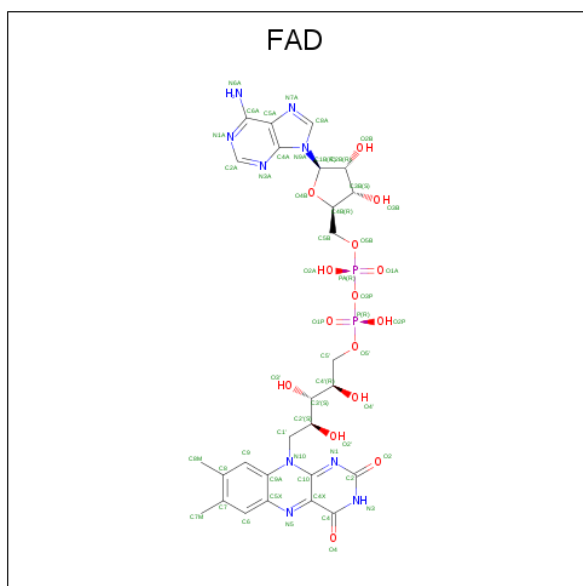
There are 3 unique types of molecules in this entry. The entry contains 9915 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein ACDH-11, isoform b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	592	Total	C	N	O	S	0	0	0
			4615	2900	819	876	20			
1	B	593	Total	C	N	O	S	0	2	0
			4632	2912	822	877	21			

- 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		

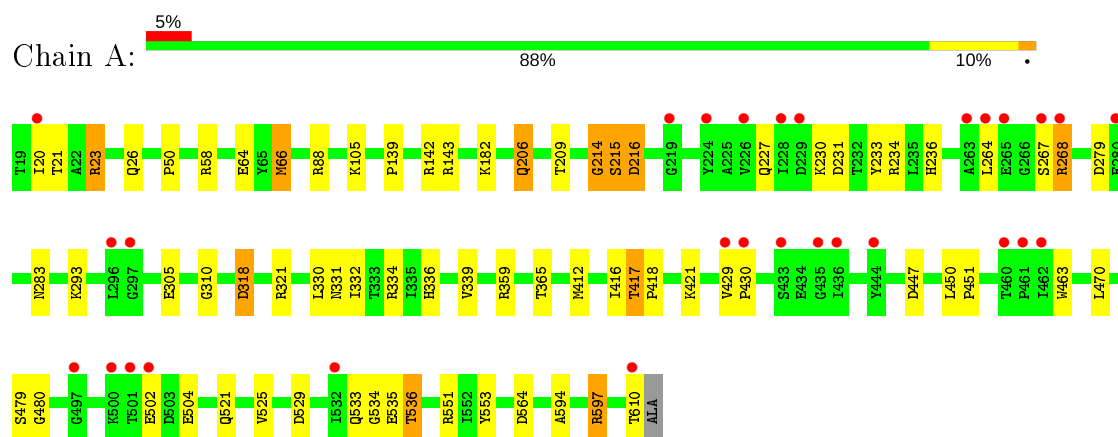
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	265	Total 265	O 265	0	0
3	B	297	Total 297	O 297	0	0

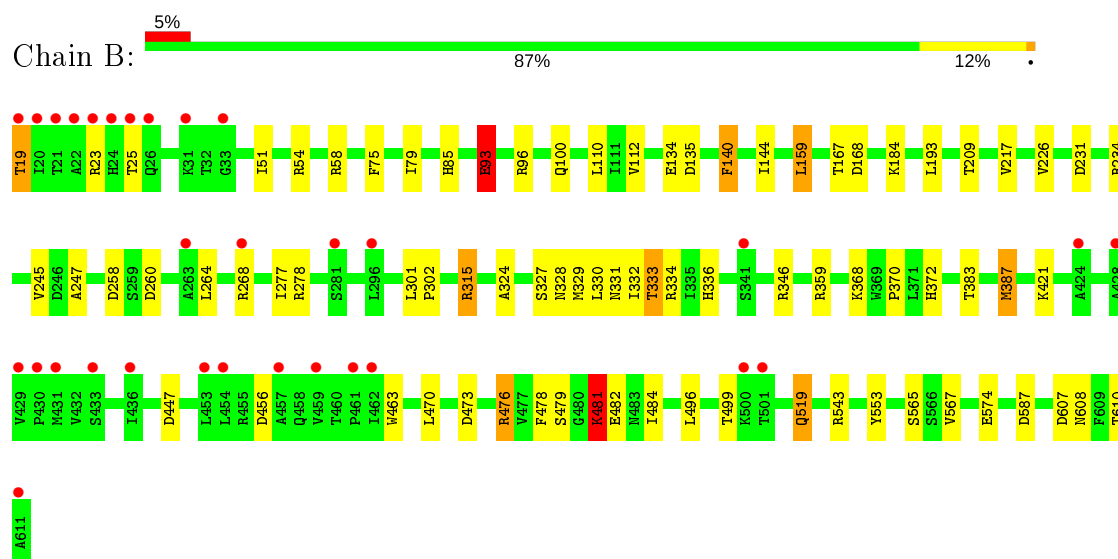
3 Residue-property plots [i](#)

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: Protein ACDH-11, isoform b



- Molecule 1: Protein ACDH-11, isoform b



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	137.93Å 113.40Å 113.90Å 90.00° 123.98° 90.00°	Depositor
Resolution (Å)	50.00 – 2.27 48.61 – 2.27	Depositor EDS
% Data completeness (in resolution range)	98.7 (50.00-2.27) 98.7 (48.61-2.27)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.09 (at 2.27Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.167 , 0.223 0.169 , 0.222	Depositor DCC
R_{free} test set	3260 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	40.5	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9915	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.43% 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	0.85	0/4692	0.94	9/6336 (0.1%)
1	B	0.87	1/4716 (0.0%)	0.94	11/6368 (0.2%)
All	All	0.86	1/9408 (0.0%)	0.94	20/12704 (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
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	93	GLU	CD-OE2	5.49	1.31	1.25

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	159	LEU	CA-CB-CG	-8.51	95.74	115.30
1	A	551	ARG	NE-CZ-NH2	-8.08	116.26	120.30
1	B	387	MET	CG-SD-CE	-6.90	89.17	100.20
1	A	214	GLY	N-CA-C	-6.36	97.19	113.10
1	A	480	GLY	N-CA-C	-6.18	97.66	113.10
1	B	23	ARG	NE-CZ-NH2	6.01	123.31	120.30
1	A	551	ARG	NE-CZ-NH1	5.96	123.28	120.30
1	B	334	ARG	NE-CZ-NH2	-5.91	117.35	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	23	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	143	ARG	NE-CZ-NH2	-5.71	117.44	120.30
1	B	346	ARG	NE-CZ-NH2	-5.64	117.48	120.30
1	B	96	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	B	447	ASP	CB-CG-OD2	-5.45	113.39	118.30
1	A	88	ARG	NE-CZ-NH1	5.35	122.97	120.30
1	B	543	ARG	NE-CZ-NH2	-5.35	117.63	120.30
1	B	456	ASP	CB-CG-OD1	5.25	123.03	118.30
1	B	447	ASP	CB-CG-OD1	5.23	123.01	118.30
1	B	278	ARG	NE-CZ-NH1	5.21	122.90	120.30
1	A	334	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	A	447	ASP	CB-CG-OD1	5.03	122.83	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	479	SER	Peptide
1	B	481	LYS	Peptide

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	4615	0	4665	42	0
1	B	4632	0	4686	53	1
2	A	53	0	31	3	0
2	B	53	0	31	1	0
3	A	265	0	0	12	0
3	B	297	0	0	23	0
All	All	9915	0	9413	94	1

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 (94) 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:332:ILE:HD12	3:B:1032:HOH:O	1.75	0.85
1:B:519:GLN:HG3	3:B:1055:HOH:O	1.78	0.84
1:B:85[B]:HIS:CD2	3:B:824:HOH:O	2.31	0.82
1:B:167:THR:HG23	1:B:193:LEU:HD22	1.61	0.82
1:A:610:THR:HA	3:A:1014:HOH:O	1.79	0.81
1:A:610:THR:HB	3:A:1001:HOH:O	1.80	0.80
1:B:85[B]:HIS:CG	3:B:824:HOH:O	2.34	0.79
1:B:330:LEU:HD23	3:B:973:HOH:O	1.89	0.71
1:B:476:ARG:HB2	3:B:1063:HOH:O	1.92	0.69
1:A:504:GLU:HG3	3:A:1027:HOH:O	1.93	0.69
1:A:504:GLU:CG	3:A:1027:HOH:O	2.44	0.65
1:B:574:GLU:HG2	3:B:1012:HOH:O	1.96	0.65
1:A:234:ARG:HH11	1:A:236:HIS:CE1	2.15	0.64
1:B:333:THR:HB	3:B:815:HOH:O	1.96	0.64
1:A:214:GLY:O	1:A:216:ASP:N	2.31	0.63
1:A:318:ASP:HB2	1:A:321:ARG:HG3	1.79	0.63
1:A:534:GLY:HA3	3:A:1032:HOH:O	1.98	0.63
1:B:496:LEU:O	1:B:499:THR:OG1	2.14	0.62
1:A:421:LYS:HD2	1:A:470:LEU:CD1	2.32	0.59
1:B:19:THR:N	3:B:806:HOH:O	2.35	0.58
1:A:209:THR:OG1	2:A:701:FAD:H1'1	2.03	0.58
1:A:502:GLU:HG2	3:A:1041:HOH:O	2.03	0.58
1:B:519:GLN:HA	1:B:519:GLN:HE21	1.70	0.57
1:A:331:ASN:HA	1:A:417:THR:HG21	1.86	0.57
1:B:231:ASP:OD1	1:B:315:ARG:NH2	2.40	0.54
1:A:502:GLU:CG	3:A:1041:HOH:O	2.56	0.53
1:B:331:ASN:HB2	3:B:1028:HOH:O	2.07	0.53
1:B:93:GLU:HG2	3:B:1070:HOH:O	2.07	0.53
1:B:226:VAL:HG11	1:B:234:ARG:NH2	2.23	0.53
1:A:209:THR:HG1	2:A:701:FAD:H1'1	1.73	0.51
1:B:330:LEU:O	1:B:331:ASN:C	2.49	0.51
1:A:64:GLU:HA	1:A:64:GLU:OE1	2.11	0.51
1:B:330:LEU:HA	3:B:815:HOH:O	2.10	0.51
1:B:168:ASP:OD2	1:B:336:HIS:HE1	1.93	0.50
1:B:134:GLU:HG2	3:B:1075:HOH:O	2.11	0.50
1:B:209:THR:OG1	2:B:701:FAD:H1'1	2.12	0.50
1:B:333:THR:CB	3:B:815:HOH:O	2.58	0.50
1:B:58:ARG:NH1	3:B:803:HOH:O	2.27	0.50
1:A:234:ARG:HD2	1:A:236:HIS:HE1	1.76	0.50
1:A:139:PRO:HA	1:A:142:ARG:HG2	1.93	0.49
1:A:336:HIS:O	1:A:339:VAL:HB	2.12	0.49
1:A:594:ALA:O	1:A:597:ARG:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:PHE:O	1:B:144:ILE:HD12	2.14	0.48
1:A:535:GLU:O	1:A:535:GLU:HG3	2.13	0.48
1:B:258:ASP:OD1	1:B:260:ASP:N	2.42	0.48
1:B:51:ILE:HD11	1:B:387:MET:HG2	1.95	0.48
1:B:54:ARG:HD2	1:B:567:VAL:HG12	1.96	0.48
1:B:301:LEU:HD12	1:B:302:PRO:HD2	1.95	0.47
2:A:701:FAD:O2A	1:B:359:ARG:NE	2.38	0.47
1:B:85[B]:HIS:CE1	3:B:824:HOH:O	2.66	0.47
1:B:336:HIS:HD2	3:B:912:HOH:O	1.97	0.47
1:A:206:GLN:HB2	3:A:940:HOH:O	2.15	0.47
1:A:533:GLN:HA	1:A:536:THR:HB	1.96	0.47
1:B:478:PHE:O	1:B:481:LYS:HG2	2.15	0.47
1:A:521:GLN:O	1:A:525:VAL:HG23	2.14	0.47
1:B:247:ALA:O	1:B:277:ILE:HD12	2.15	0.46
1:A:264:LEU:HD23	1:A:268:ARG:HB3	1.97	0.46
1:A:504:GLU:HG2	3:A:1027:HOH:O	2.12	0.46
1:B:368:LYS:HE2	3:B:1072:HOH:O	2.15	0.45
1:A:234:ARG:HB2	1:A:310:GLY:HA2	1.98	0.45
1:A:58:ARG:HD3	3:A:810:HOH:O	2.17	0.45
1:B:75:PHE:O	1:B:79:ILE:HG12	2.17	0.45
1:B:519:GLN:HA	1:B:519:GLN:NE2	2.32	0.44
1:A:23:ARG:NH2	1:A:26:GLN:HG3	2.33	0.44
1:B:332:ILE:HA	1:B:332:ILE:HD13	1.72	0.44
1:B:473:ASP:OD1	1:B:476:ARG:NH1	2.43	0.44
1:B:482:GLU:O	1:B:484:ILE:N	2.51	0.44
1:A:450:LEU:HB2	1:A:451:PRO:HD3	1.99	0.44
1:A:529:ASP:HB3	3:A:921:HOH:O	2.17	0.44
1:B:324:ALA:O	1:B:327:SER:OG	2.28	0.44
3:A:866:HOH:O	1:B:372:HIS:HD2	2.00	0.43
1:B:110:LEU:HD22	1:B:245:VAL:HG23	2.01	0.43
1:A:50:PRO:HB3	1:B:607:ASP:HB3	2.00	0.43
1:A:330:LEU:O	1:A:331:ASN:C	2.55	0.43
1:A:564:ASP:OD1	1:A:564:ASP:C	2.57	0.43
1:B:258:ASP:C	1:B:258:ASP:OD1	2.57	0.43
1:B:330:LEU:HD12	3:B:1026:HOH:O	2.18	0.42
1:B:327:SER:O	1:B:329:MET:N	2.52	0.42
1:A:66:MET:HE3	1:A:66:MET:HA	2.02	0.42
1:B:93:GLU:HG3	3:B:1079:HOH:O	2.20	0.42
1:A:227:GLN:HA	1:A:233:TYR:CD2	2.55	0.41
1:A:359:ARG:O	1:A:365:THR:HA	2.19	0.41
1:B:383:THR:O	1:B:387:MET:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:MET:CE	1:A:66:MET:HA	2.50	0.41
1:B:368:LYS:CE	3:B:1072:HOH:O	2.67	0.41
1:B:329:MET:HG2	3:B:973:HOH:O	2.20	0.41
1:B:421:LYS:HD2	1:B:470:LEU:CD1	2.51	0.41
1:A:412:MET:O	1:A:416:ILE:HG22	2.20	0.41
1:A:429:VAL:HB	1:A:430:PRO:HD3	2.02	0.40
1:A:293:LYS:NZ	1:A:305:GLU:OE2	2.39	0.40
1:A:417:THR:HG22	1:A:418:PRO:HD3	2.03	0.40
1:B:329:MET:O	1:B:332:ILE:HB	2.20	0.40
1:B:587:ASP:HB3	3:B:836:HOH:O	2.20	0.40
1:A:279:ASP:OD1	1:A:283:ASN:HB2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:565:SER:OG	1:B:565:SER:OG[2_455]	2.01	0.19

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	590/593 (100%)	571 (97%)	15 (2%)	4 (1%)	22	25
1	B	593/593 (100%)	568 (96%)	23 (4%)	2 (0%)	41	49
All	All	1183/1186 (100%)	1139 (96%)	38 (3%)	6 (0%)	29	34

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	215	SER
1	A	230	LYS

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Mol	Chain	Res	Type
1	A	597	ARG
1	B	481	LYS
1	A	216	ASP
1	B	328	ASN

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	497/497 (100%)	481 (97%)	16 (3%)	39	52
1	B	499/497 (100%)	476 (95%)	23 (5%)	27	35
All	All	996/994 (100%)	957 (96%)	39 (4%)	32	43

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ILE
1	A	21	THR
1	A	66	MET
1	A	105	LYS
1	A	182	LYS
1	A	206	GLN
1	A	215	SER
1	A	231	ASP
1	A	267	SER
1	A	268	ARG
1	A	318	ASP
1	A	332	ILE
1	A	417	THR
1	A	463	TRP
1	A	536	THR
1	A	553	TYR
1	B	19	THR
1	B	25	THR
1	B	93	GLU

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Mol	Chain	Res	Type
1	B	100	GLN
1	B	112	VAL
1	B	135	ASP
1	B	140	PHE
1	B	159	LEU
1	B	184	LYS
1	B	217	VAL
1	B	264	LEU
1	B	268	ARG
1	B	315	ARG
1	B	333	THR
1	B	370	PRO
1	B	463	TRP
1	B	476	ARG
1	B	479	SER
1	B	481	LYS
1	B	519	GLN
1	B	553	TYR
1	B	608	ASN
1	B	610	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	236	HIS
1	A	336	HIS
1	A	458	GLN
1	A	533	GLN
1	B	336	HIS
1	B	372	HIS
1	B	458	GLN
1	B	498	ASN
1	B	519	GLN
1	B	608	ASN

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$
2	FAD	B	701	-	51,58,58	1.76	6 (11%)	60,89,89	2.16	13 (21%)
2	FAD	A	701	-	51,58,58	1.85	9 (17%)	60,89,89	2.52	14 (23%)

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	B	701	-	-	4/30/50/50	0/6/6/6
2	FAD	A	701	-	-	16/30/50/50	0/6/6/6

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	FAD	C4X-C10	9.30	1.48	1.38
2	B	701	FAD	C4X-C10	9.03	1.47	1.38
2	B	701	FAD	C4-C4X	4.59	1.49	1.41
2	A	701	FAD	C9A-C5X	3.62	1.49	1.42
2	A	701	FAD	C4-C4X	3.57	1.47	1.41
2	B	701	FAD	C9A-C5X	3.23	1.49	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	FAD	C5A-C4A	2.69	1.48	1.40
2	A	701	FAD	C8-C7	2.68	1.47	1.40
2	B	701	FAD	C5A-C4A	2.45	1.47	1.40
2	A	701	FAD	C2A-N3A	2.44	1.36	1.32
2	B	701	FAD	C5'-C4'	2.38	1.55	1.51
2	A	701	FAD	O4B-C1B	2.34	1.44	1.41
2	A	701	FAD	C6-C5X	-2.10	1.38	1.41
2	A	701	FAD	C2-N1	-2.07	1.34	1.38
2	B	701	FAD	C8-C7	2.05	1.46	1.40

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	FAD	C4-N3-C2	10.89	124.34	115.14
2	A	701	FAD	C1'-N10-C9A	9.33	125.64	118.29
2	B	701	FAD	C4-N3-C2	8.40	122.24	115.14
2	B	701	FAD	C1'-N10-C9A	6.53	123.43	118.29
2	A	701	FAD	C4-C4X-C10	-5.68	116.19	119.95
2	A	701	FAD	C4X-C4-N3	-4.98	116.63	123.43
2	B	701	FAD	N3A-C2A-N1A	-4.79	121.19	128.68
2	B	701	FAD	C4-C4X-C10	-4.49	116.98	119.95
2	B	701	FAD	C4X-C4-N3	-4.34	117.49	123.43
2	B	701	FAD	C4-C4X-N5	4.21	123.41	118.60
2	A	701	FAD	C4-C4X-N5	3.94	123.10	118.60
2	B	701	FAD	C4X-N5-C5X	3.63	120.40	116.77
2	A	701	FAD	C1'-N10-C10	-3.27	115.48	118.41
2	A	701	FAD	N6A-C6A-N1A	3.21	125.25	118.57
2	A	701	FAD	N3A-C2A-N1A	-2.79	124.32	128.68
2	B	701	FAD	O4B-C4B-C3B	2.71	110.48	105.11
2	B	701	FAD	N6A-C6A-N1A	2.71	124.19	118.57
2	B	701	FAD	O2A-PA-O1A	2.64	125.31	112.24
2	B	701	FAD	C9A-N10-C10	-2.50	118.64	121.91
2	A	701	FAD	C5A-C6A-N6A	-2.49	116.57	120.35
2	A	701	FAD	C9A-N10-C10	-2.48	118.65	121.91
2	A	701	FAD	O2A-PA-O1A	2.47	124.45	112.24
2	B	701	FAD	C2A-N1A-C6A	2.47	122.98	118.75
2	A	701	FAD	O3'-C3'-C4'	2.44	114.71	108.81
2	B	701	FAD	C10-C4X-N5	-2.39	119.61	121.26
2	A	701	FAD	C4X-N5-C5X	2.15	118.92	116.77
2	A	701	FAD	C6-C5X-N5	-2.14	116.69	119.05

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	FAD	C2'-C1'-N10-C9A
2	A	701	FAD	C2'-C1'-N10-C10
2	A	701	FAD	N10-C1'-C2'-O2'
2	A	701	FAD	N10-C1'-C2'-C3'
2	A	701	FAD	C1'-C2'-C3'-C4'
2	A	701	FAD	C2'-C3'-C4'-O4'
2	A	701	FAD	C2'-C3'-C4'-C5'
2	A	701	FAD	O3'-C3'-C4'-O4'
2	A	701	FAD	O3'-C3'-C4'-C5'
2	A	701	FAD	C3'-C4'-C5'-O5'
2	A	701	FAD	O4'-C4'-C5'-O5'
2	A	701	FAD	C5'-O5'-P-O1P
2	A	701	FAD	O2'-C2'-C3'-C4'
2	B	701	FAD	PA-O3P-P-O1P
2	A	701	FAD	C4'-C5'-O5'-P
2	B	701	FAD	PA-O3P-P-O2P
2	B	701	FAD	O2'-C2'-C3'-C4'
2	A	701	FAD	C5'-O5'-P-O2P
2	B	701	FAD	O3'-C3'-C4'-C5'
2	A	701	FAD	C5'-O5'-P-O3P

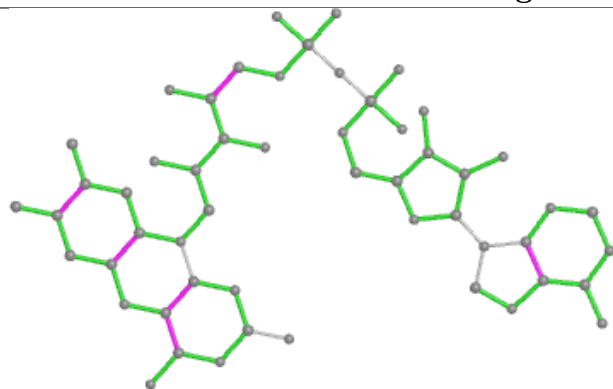
There are no ring outliers.

2 monomers are involved in 4 short contacts:

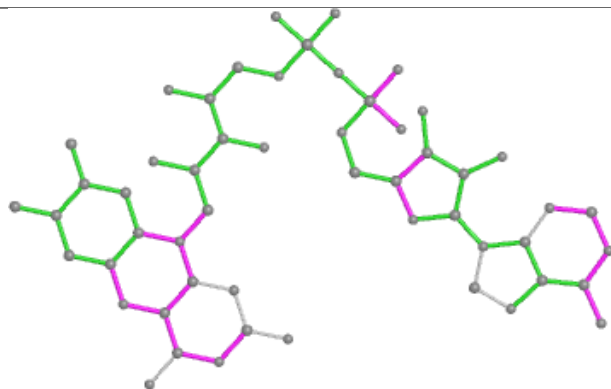
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	FAD	1	0
2	A	701	FAD	3	0

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.

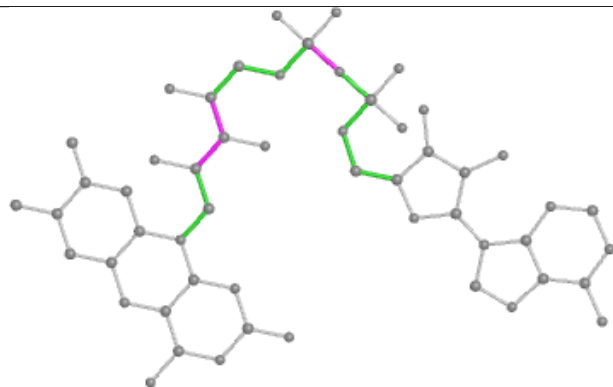
Ligand FAD B 701



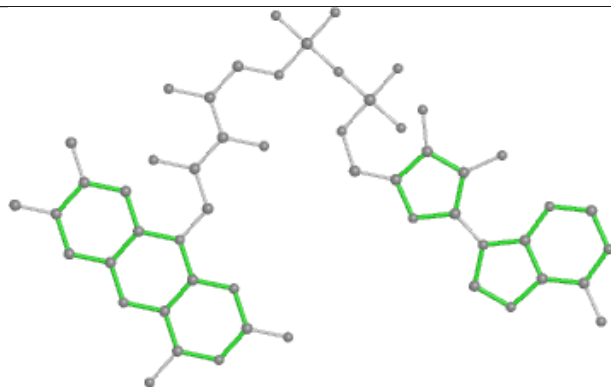
Bond lengths



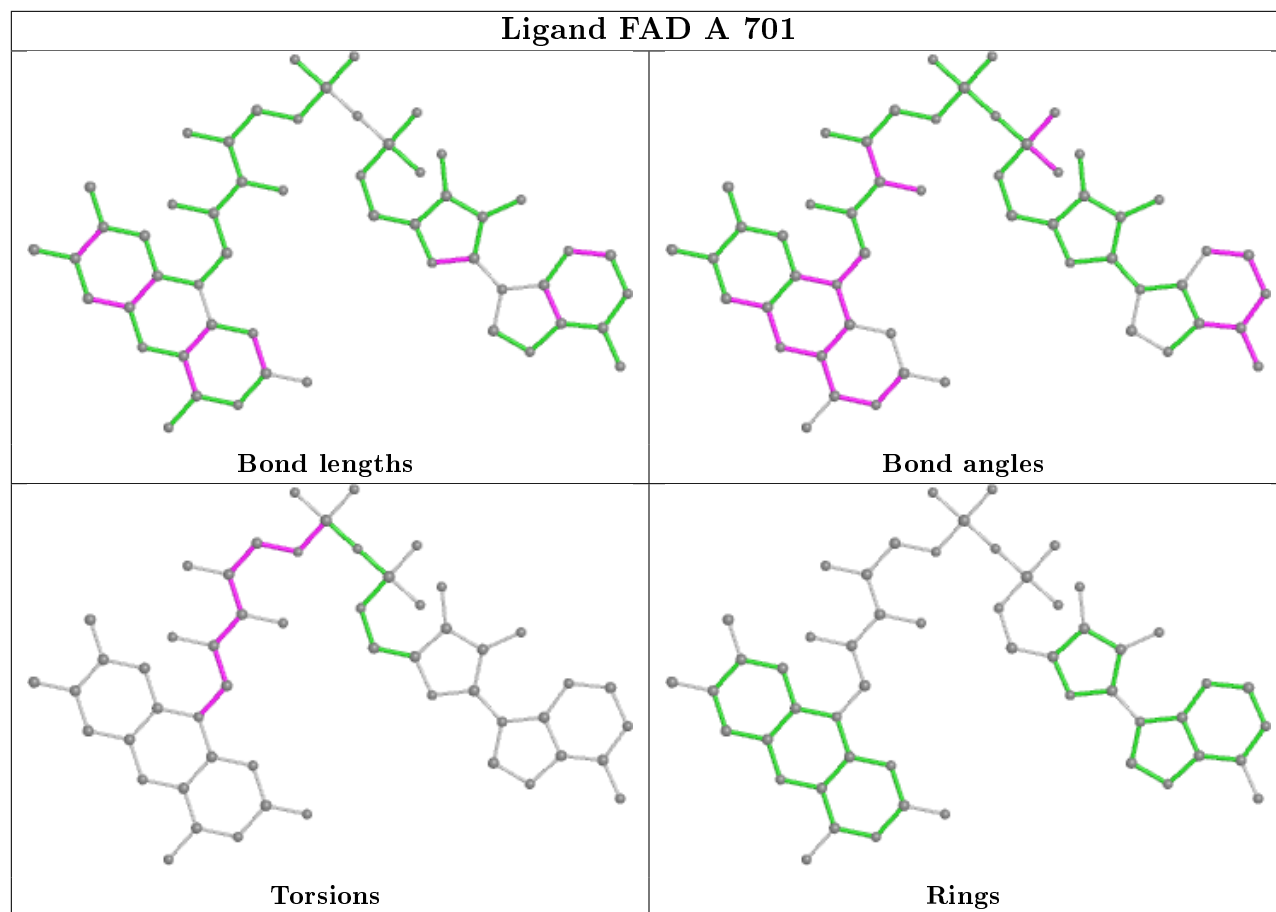
Bond angles



Torsions



Rings



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	592/593 (99%)	-0.06	29 (4%)	29 35	30, 51, 77, 105	0
1	B	593/593 (100%)	0.01	31 (5%)	27 32	29, 49, 80, 132	0
All	All	1185/1186 (99%)	-0.03	60 (5%)	28 33	29, 50, 78, 132	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	264	LEU	5.1
1	B	19	THR	4.5
1	A	265	GLU	4.3
1	B	611	ALA	4.1
1	B	21	THR	4.1
1	B	25	THR	3.9
1	A	429	VAL	3.7
1	A	263	ALA	3.6
1	B	20	ILE	3.6
1	B	500	LYS	3.3
1	B	430	PRO	3.2
1	B	429	VAL	3.1
1	A	20	ILE	3.0
1	B	461	PRO	2.9
1	A	296	LEU	2.8
1	B	428	ALA	2.8
1	B	22	ALA	2.8
1	B	263	ALA	2.8
1	A	462	ILE	2.8
1	A	430	PRO	2.7
1	B	454	LEU	2.7
1	A	228	ILE	2.7
1	A	610	THR	2.7
1	B	431	MET	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	433	SER	2.6
1	B	341	SER	2.6
1	A	436	ILE	2.6
1	B	296	LEU	2.6
1	A	268	ARG	2.5
1	A	502	GLU	2.5
1	A	532	ILE	2.5
1	B	457	ALA	2.4
1	A	229	ASP	2.4
1	B	24	HIS	2.4
1	B	462	ILE	2.4
1	A	461	PRO	2.4
1	A	501	THR	2.3
1	B	433	SER	2.3
1	A	224	TYR	2.3
1	B	268	ARG	2.3
1	B	436	ILE	2.3
1	A	297	GLY	2.3
1	A	497	GLY	2.3
1	B	281	SER	2.3
1	B	33	GLY	2.3
1	B	31	LYS	2.3
1	A	460	THR	2.2
1	B	26	GLN	2.2
1	A	280	GLU	2.2
1	A	226	VAL	2.2
1	B	453	LEU	2.2
1	B	501	THR	2.1
1	A	219	GLY	2.1
1	B	459	VAL	2.1
1	A	444	TYR	2.1
1	B	23	ARG	2.1
1	A	267	SER	2.1
1	A	435	GLY	2.0
1	A	500	LYS	2.0
1	B	424	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

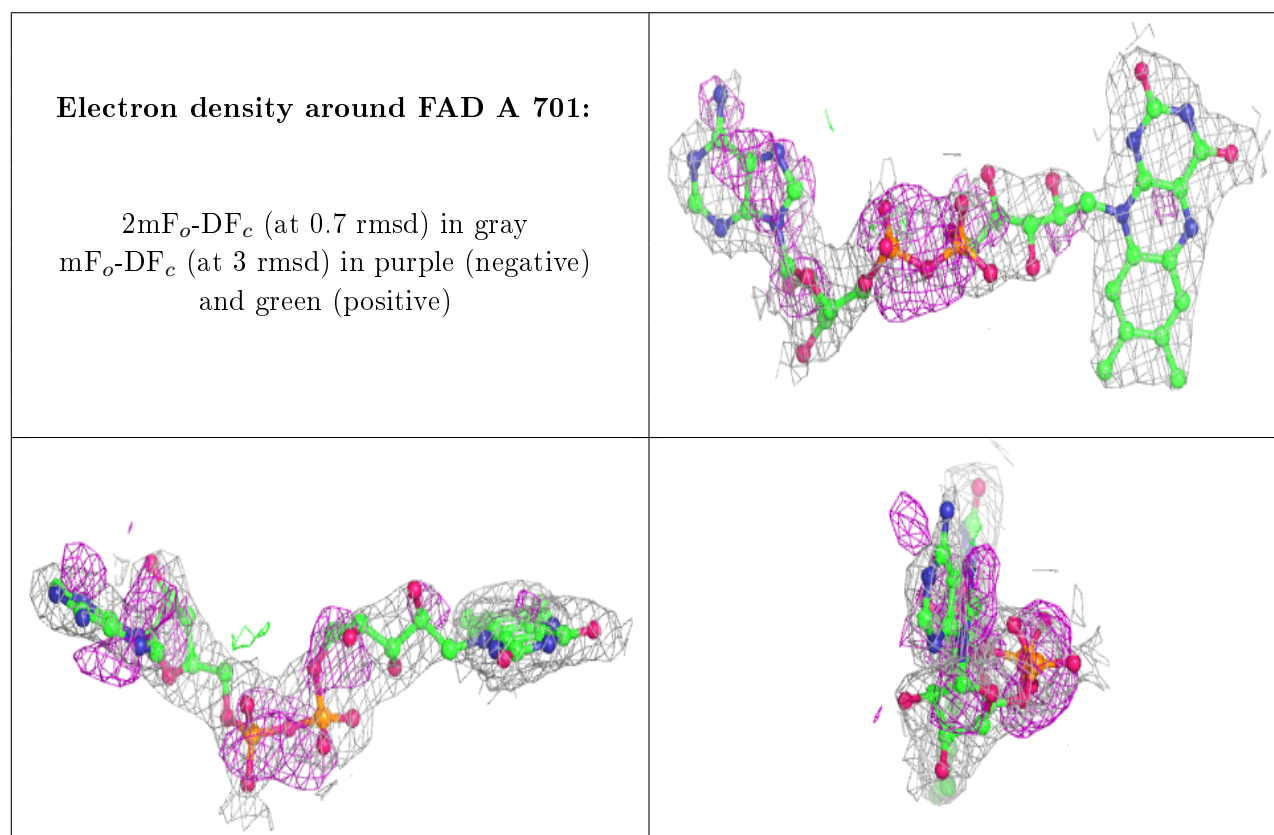
There are no carbohydrates in this entry.

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.

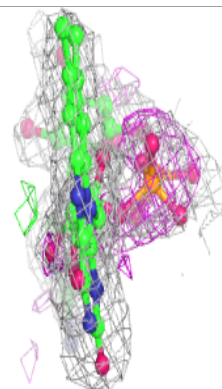
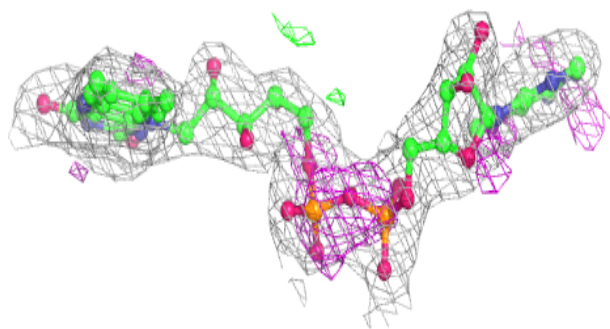
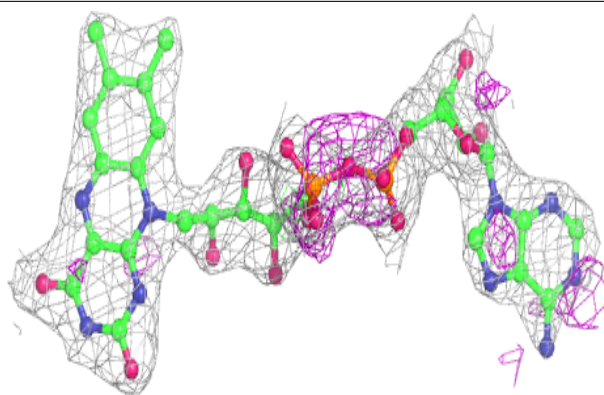
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FAD	A	701	53/53	0.90	0.15	50,66,81,87	0
2	FAD	B	701	53/53	0.93	0.12	41,55,62,65	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around FAD B 701:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.