Full wwPDB X-ray Structure Validation Report (i)

Jun 24, 2024 – 02:06 PM EDT

PDB ID	:	5Y77
Title	:	Crystal structure of Pseudomonas fluorescens Kynurenine 3-monooxygenase in complex with L-KYN (seMet derivative)
Authors	:	Xiang, Y.; Gao, J.J.; Zhu, D.Y.
Deposited on	:	2017-08-16
Resolution	:	1.60 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 (i) 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 (i)) were used in the production of this report:

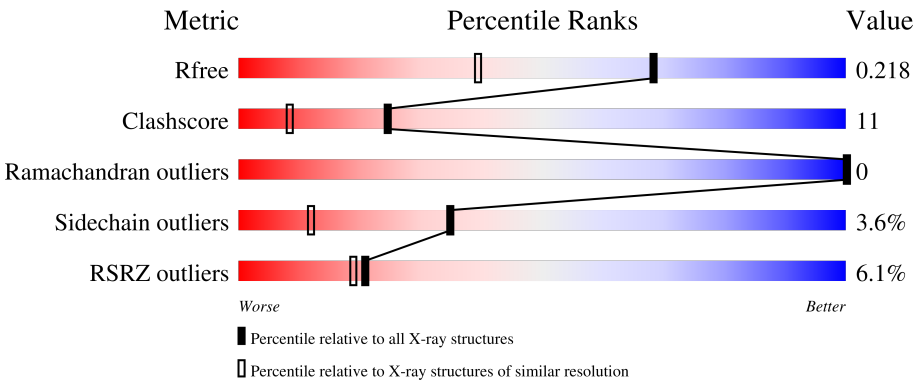
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 1.60 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))
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 <=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	473	4% 81% 13%
1	B	473	8% 74% 20% 5%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	KYN	A	502	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8388 atoms, of which 22 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 Kynurenine 3-monooxygenase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	455	3610	2267	668	653	7	15	0	16	0
1	B	451	3566	2239	658	649	6	14	0	14	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MSE	-	expression tag	UNP Q84HF5
A	1	GLY	-	expression tag	UNP Q84HF5
A	462	ALA	-	expression tag	UNP Q84HF5
A	463	ALA	-	expression tag	UNP Q84HF5
A	464	ALA	-	expression tag	UNP Q84HF5
A	465	LEU	-	expression tag	UNP Q84HF5
A	466	GLU	-	expression tag	UNP Q84HF5
A	467	HIS	-	expression tag	UNP Q84HF5
A	468	HIS	-	expression tag	UNP Q84HF5
A	469	HIS	-	expression tag	UNP Q84HF5
A	470	HIS	-	expression tag	UNP Q84HF5
A	471	HIS	-	expression tag	UNP Q84HF5
A	472	HIS	-	expression tag	UNP Q84HF5
B	0	MSE	-	expression tag	UNP Q84HF5
B	1	GLY	-	expression tag	UNP Q84HF5
B	462	ALA	-	expression tag	UNP Q84HF5
B	463	ALA	-	expression tag	UNP Q84HF5
B	464	ALA	-	expression tag	UNP Q84HF5
B	465	LEU	-	expression tag	UNP Q84HF5
B	466	GLU	-	expression tag	UNP Q84HF5
B	467	HIS	-	expression tag	UNP Q84HF5
B	468	HIS	-	expression tag	UNP Q84HF5
B	469	HIS	-	expression tag	UNP Q84HF5
B	470	HIS	-	expression tag	UNP Q84HF5
B	471	HIS	-	expression tag	UNP Q84HF5

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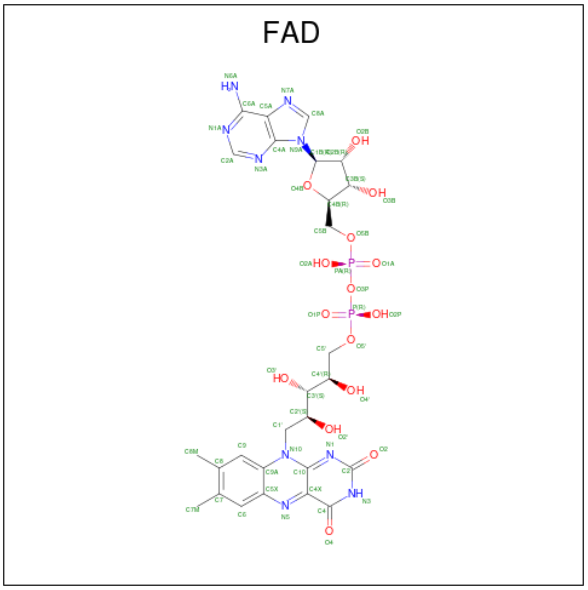
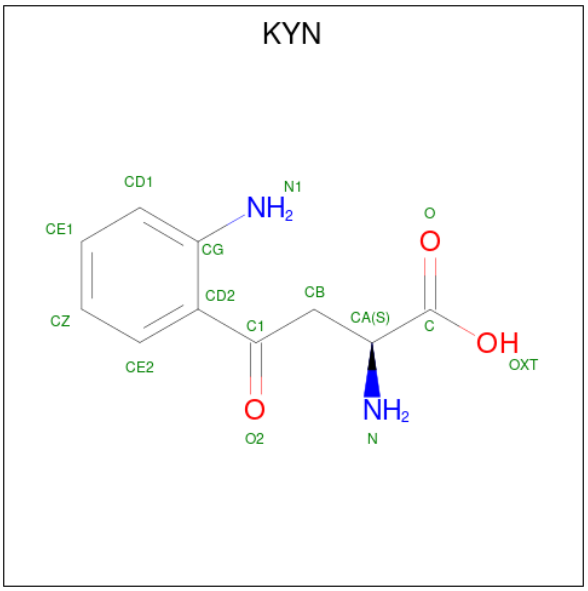
Chain	Residue	Modelled	Actual	Comment	Reference
B	472	HIS	-	expression tag	UNP Q84HF5

- 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
			Total	C	N	O	P		
2	A	1	53	27	9	15	2	0	0
			Total	C	N	O	P		
2	B	1	53	27	9	15	2	0	0

- Molecule 3 is (2S)-2-amino-4-(2-aminophenyl)-4-oxobutanoic acid (three-letter code: KYN) (formula: $C_{10}H_{12}N_2O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
3	A	1	26	10	11	2	3	0	0
			Total	C	H	N	O		
3	B	1	26	10	11	2	3	0	0

- Molecule 4 is water.

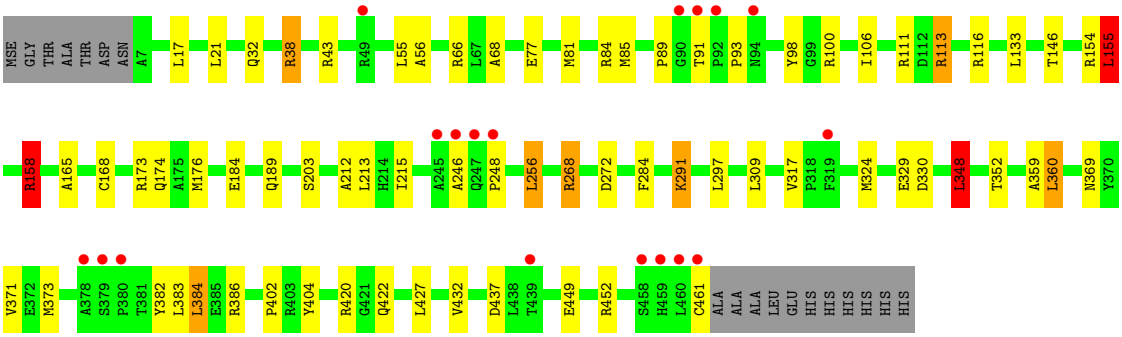
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	643	643	643	0	0
			Total	O		
4	B	411	411	411	0	0

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 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.

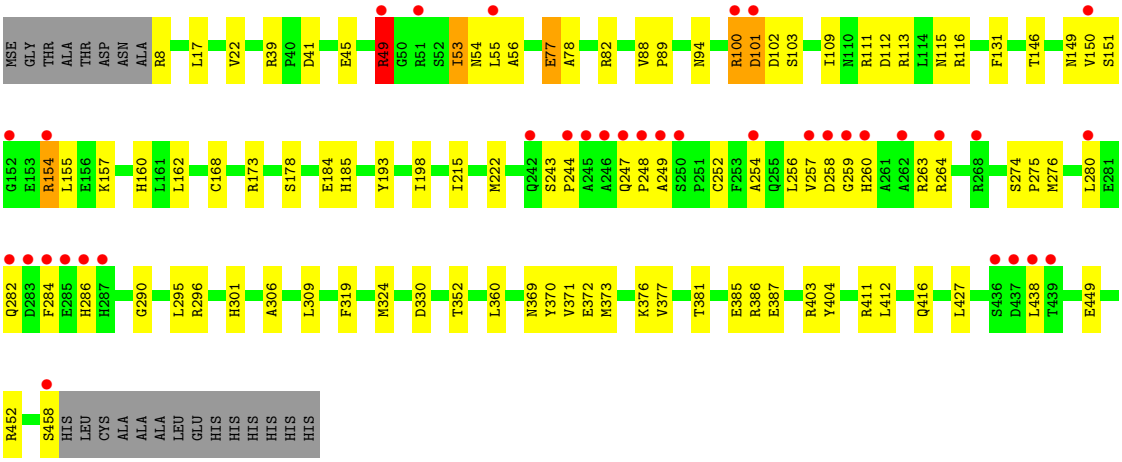
• Molecule 1: Kynurenine 3-monooxygenase

Chain A:



• Molecule 1: Kynurenine 3-monooxygenase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	69.56Å 52.45Å 135.71Å 90.00° 104.33° 90.00°	Depositor
Resolution (Å)	34.78 – 1.60 34.78 – 1.60	Depositor EDS
% Data completeness (in resolution range)	95.0 (34.78-1.60) 95.0 (34.78-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	6.96 (at 1.60Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.185 , 0.217 0.185 , 0.218	Depositor DCC
R_{free} test set	5969 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	17.7	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 56.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8388	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 42.61 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.9904e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: KYN, 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	0.81	3/3721 (0.1%)	0.99	20/5028 (0.4%)
1	B	0.69	0/3671	0.85	5/4963 (0.1%)
All	All	0.75	3/7392 (0.0%)	0.92	25/9991 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	176	MSE	SE-CE	-7.50	1.51	1.95
1	A	329	GLU	CG-CD	5.14	1.59	1.51
1	A	203	SER	CB-OG	-5.12	1.35	1.42

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	173	ARG	NE-CZ-NH1	11.64	126.12	120.30
1	B	173	ARG	NE-CZ-NH2	-10.25	115.17	120.30
1	A	173	ARG	NE-CZ-NH2	-10.10	115.25	120.30
1	A	38[A]	ARG	NE-CZ-NH2	-8.91	115.84	120.30
1	A	38[B]	ARG	NE-CZ-NH2	-8.91	115.84	120.30
1	B	173	ARG	NE-CZ-NH1	8.74	124.67	120.30
1	A	348	LEU	CB-CG-CD1	7.46	123.68	111.00
1	A	158	ARG	NE-CZ-NH1	7.36	123.98	120.30
1	A	427	LEU	CA-CB-CG	7.32	132.12	115.30
1	A	330	ASP	CB-CG-OD1	7.19	124.77	118.30
1	A	111	ARG	NE-CZ-NH2	-7.03	116.78	120.30
1	A	38[A]	ARG	NE-CZ-NH1	7.02	123.81	120.30
1	A	38[B]	ARG	NE-CZ-NH1	7.02	123.81	120.30
1	A	66	ARG	NE-CZ-NH1	6.83	123.72	120.30
1	A	111	ARG	CG-CD-NE	-6.53	98.09	111.80
1	A	66	ARG	NE-CZ-NH2	-6.01	117.29	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55	LEU	CA-CB-CG	5.76	128.56	115.30
1	B	330	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	A	360	LEU	CB-CG-CD1	-5.50	101.66	111.00
1	A	43	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	B	49	ARG	NE-CZ-NH1	-5.42	117.59	120.30
1	A	155	LEU	CA-CB-CG	5.40	127.72	115.30
1	B	330	ASP	CB-CG-OD1	5.22	123.00	118.30
1	A	158	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	420	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	3610	0	3616	64	0
1	B	3566	0	3569	88	1
2	A	53	0	29	1	0
2	B	53	0	31	3	0
3	A	15	11	11	9	0
3	B	15	11	11	3	0
4	A	643	0	0	22	3
4	B	411	0	0	27	0
All	All	8366	22	7267	159	4

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 11.

All (159) 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:452:ARG:NH1	4:B:601:HOH:O	1.96	0.96
1:B:101:ASP:OD2	1:B:103:SER:N	2.03	0.90
1:B:403[B]:ARG:NH1	4:B:602:HOH:O	2.04	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:ARG:HD3	1:B:49:ARG:H	1.38	0.87
1:B:39:ARG:NH2	4:B:603:HOH:O	2.08	0.85
1:B:100:ARG:HG3	1:B:100:ARG:HH11	1.45	0.81
1:B:49:ARG:HD3	1:B:49:ARG:N	1.96	0.80
1:A:404:TYR:OH	3:A:502:KYN:N	2.17	0.78
1:B:184:GLU:HG3	1:B:295:LEU:HD11	1.64	0.77
1:A:133:LEU:HD22	1:A:155:LEU:HD21	1.65	0.76
1:B:449:GLU:HA	4:B:601:HOH:O	1.84	0.76
1:A:77[A]:GLU:OE1	1:A:113:ARG:HG2	1.86	0.75
1:A:461:CYS:SG	4:A:795:HOH:O	2.43	0.75
1:A:360:LEU:HB3	4:A:780:HOH:O	1.86	0.75
1:B:352[A]:THR:HG21	4:B:855:HOH:O	1.87	0.74
3:A:502:KYN:HE1	4:A:966:HOH:O	1.89	0.72
1:B:77:GLU:OE1	1:B:113[B]:ARG:HG2	1.90	0.71
1:B:168[B]:CYS:SG	4:B:1000:HOH:O	2.49	0.71
1:A:84[B]:ARG:NH1	4:A:606:HOH:O	2.24	0.71
1:A:174:GLN:OE1	4:A:601:HOH:O	2.09	0.70
1:A:404:TYR:HE2	3:A:502:KYN:H	1.37	0.70
1:A:384:LEU:HD13	1:A:432:VAL:HG13	1.74	0.70
1:B:146:THR:CG2	1:B:154:ARG:HD3	2.22	0.70
1:B:263:ARG:HA	1:B:280:LEU:HD21	1.74	0.69
1:B:452:ARG:NH1	4:B:604:HOH:O	2.08	0.69
1:A:98:TYR:O	4:A:602:HOH:O	2.11	0.69
1:B:458:SER:OG	4:B:605:HOH:O	2.11	0.68
1:B:100:ARG:HH11	1:B:100:ARG:CG	2.07	0.68
1:B:100:ARG:HG3	1:B:100:ARG:NH1	2.05	0.68
1:B:215:ILE:CD1	1:B:373:MSE:HE2	2.24	0.68
1:B:53:ILE:HD12	1:B:54:ASN:N	2.08	0.67
1:B:371:VAL:HG23	4:B:722:HOH:O	1.93	0.67
1:B:78:ALA:HB2	1:B:109[A]:ILE:HG22	1.77	0.67
1:B:112[A]:ASP:OD1	4:B:606:HOH:O	2.12	0.67
1:A:81[A]:MSE:HE3	1:A:106:ILE:HG13	1.77	0.66
2:B:501:FAD:HM72	4:B:724:HOH:O	1.96	0.65
1:A:38[A]:ARG:HD2	4:A:761:HOH:O	1.96	0.65
1:B:411:ARG:HG2	4:B:671:HOH:O	1.96	0.65
1:B:215:ILE:HD11	1:B:373:MSE:HE2	1.79	0.65
1:B:17:LEU:HD13	1:B:324[B]:MSE:HG3	1.79	0.65
1:B:115[B]:ASN:ND2	4:B:611:HOH:O	2.20	0.64
1:B:215:ILE:HG23	1:B:222:MSE:CE	2.28	0.64
1:A:422[B]:GLN:HG3	4:A:668:HOH:O	1.96	0.64
1:A:113:ARG:NH2	4:A:610:HOH:O	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:HIS:O	1:B:264:ARG:HG3	1.98	0.63
1:B:404:TYR:OH	3:B:502:KYN:N	2.32	0.63
1:B:185:HIS:CE1	1:B:296[B]:ARG:HH21	2.16	0.63
3:A:502:KYN:C	3:A:502:KYN:O2	2.45	0.63
1:B:49:ARG:H	1:B:49:ARG:CD	1.96	0.63
1:B:263:ARG:HA	1:B:280:LEU:CD2	2.29	0.63
1:B:449:GLU:OE2	4:B:604:HOH:O	2.15	0.62
1:A:268:ARG:NE	4:A:611:HOH:O	2.31	0.62
1:A:449:GLU:OE1	1:A:452:ARG:NH2	2.33	0.62
4:A:918:HOH:O	1:B:360:LEU:HD11	1.98	0.62
3:B:502:KYN:C	3:B:502:KYN:O2	2.47	0.62
1:A:291:LYS:NZ	4:A:607:HOH:O	2.24	0.60
1:A:168[A]:CYS:SG	4:A:1065:HOH:O	2.56	0.60
1:B:372:GLU:OE2	4:B:608:HOH:O	2.16	0.59
1:A:309:LEU:N	1:A:309:LEU:HD23	2.16	0.58
1:B:78:ALA:HB2	1:B:109[A]:ILE:CG2	2.34	0.57
1:B:56:ALA:HB2	3:B:502:KYN:HN1	1.67	0.57
1:A:291:LYS:H	1:A:291:LYS:CE	2.18	0.57
1:A:184:GLU:HG2	1:A:297:LEU:CD2	2.35	0.57
1:B:381:THR:O	1:B:385:GLU:HG3	2.04	0.57
1:B:215:ILE:HG23	1:B:222:MSE:HE2	1.86	0.56
1:B:258:ASP:OD1	1:B:259:GLY:N	2.39	0.56
1:A:77[A]:GLU:OE1	1:A:113:ARG:CG	2.53	0.55
1:A:133:LEU:CD2	1:A:155:LEU:HD21	2.37	0.55
1:B:215:ILE:CG2	1:B:222:MSE:CE	2.85	0.55
1:B:416:GLN:OE1	4:B:609:HOH:O	2.17	0.54
1:A:84[A]:ARG:HD3	1:A:215:ILE:CD1	2.37	0.54
1:B:247:GLN:N	1:B:248:PRO:HD3	2.22	0.54
1:A:184:GLU:HG2	1:A:297:LEU:HD22	1.90	0.54
1:A:404:TYR:CE2	3:A:502:KYN:N	2.72	0.54
1:A:256:LEU:HD23	1:A:284:PHE:CZ	2.44	0.53
1:A:32:GLN:NE2	4:A:621:HOH:O	2.41	0.53
1:B:8:ARG:N	4:B:634:HOH:O	2.42	0.52
1:B:103:SER:HA	4:B:752:HOH:O	2.10	0.52
1:A:77[B]:GLU:OE2	4:A:603:HOH:O	2.19	0.51
1:A:352[A]:THR:HG21	4:A:978:HOH:O	2.11	0.51
1:B:49:ARG:N	1:B:49:ARG:CD	2.62	0.51
1:A:85:MSE:SE	1:A:93:PRO:HB2	2.61	0.51
1:B:372:GLU:HA	1:B:376:LYS:HB2	1.91	0.51
1:A:84[A]:ARG:HD3	1:A:215:ILE:HD12	1.92	0.50
1:A:449:GLU:CD	1:A:452:ARG:HH21	2.14	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:SER:OG	4:B:610:HOH:O	2.19	0.50
1:A:348:LEU:O	1:A:352[B]:THR:HG23	2.12	0.50
1:B:387:GLU:OE1	1:B:438:LEU:HD11	2.12	0.50
1:A:402:PRO:HG3	4:A:772:HOH:O	2.10	0.50
1:B:215:ILE:HD11	1:B:373:MSE:CE	2.41	0.50
1:A:291:LYS:H	1:A:291:LYS:CD	2.25	0.49
1:A:371:VAL:HG13	4:A:742:HOH:O	2.11	0.49
1:B:215:ILE:CG2	1:B:222:MSE:HE1	2.42	0.49
1:A:89:PRO:HA	1:A:272:ASP:OD2	2.12	0.49
1:A:158:ARG:HH11	1:A:158:ARG:CG	2.26	0.48
1:A:309:LEU:HD23	1:A:309:LEU:H	1.78	0.48
1:B:369:ASN:O	1:B:373:MSE:HG3	2.13	0.48
1:A:369:ASN:O	1:A:373:MSE:HG3	2.14	0.48
1:B:252:CYS:SG	1:B:254:ALA:HB3	2.53	0.48
1:A:146:THR:OG1	4:A:604:HOH:O	2.20	0.47
1:A:359:ALA:HB2	4:A:916:HOH:O	2.14	0.47
1:B:77:GLU:CD	1:B:113[B]:ARG:HE	2.18	0.47
1:B:411:ARG:NH1	4:B:641:HOH:O	2.46	0.47
1:A:158:ARG:HH11	1:A:158:ARG:HG2	1.80	0.47
1:B:146:THR:HG21	1:B:154:ARG:HD3	1.95	0.47
1:B:150:VAL:O	4:B:612:HOH:O	2.20	0.47
1:A:452:ARG:HH11	1:A:452:ARG:HG2	1.80	0.46
1:A:113:ARG:HH12	1:A:116:ARG:CZ	2.29	0.46
1:B:257:VAL:O	1:B:257:VAL:HG12	2.16	0.46
1:B:411:ARG:NH1	4:B:607:HOH:O	2.12	0.46
1:B:160:HIS:HE1	4:B:942:HOH:O	1.98	0.45
1:A:404:TYR:CZ	3:A:502:KYN:N	2.85	0.45
1:B:149[A]:ASN:CG	1:B:150:VAL:H	2.20	0.45
1:A:17:LEU:HD13	1:A:324[B]:MSE:HG3	1.98	0.45
1:B:248:PRO:O	1:B:249:ALA:HB3	2.16	0.45
1:B:372:GLU:HA	1:B:376:LYS:HD2	1.99	0.45
1:A:382:TYR:HE2	1:A:386:ARG:HH21	1.64	0.44
1:A:113:ARG:HH12	1:A:116:ARG:NE	2.14	0.44
1:B:113[B]:ARG:HD3	1:B:116:ARG:NH1	2.32	0.44
1:A:384:LEU:CD1	1:A:432:VAL:HG13	2.44	0.44
1:B:88:VAL:HA	1:B:89:PRO:HD3	1.89	0.44
1:B:162:LEU:O	1:B:306:ALA:HA	2.17	0.44
1:B:78:ALA:CB	1:B:109[A]:ILE:HG22	2.44	0.44
1:B:113[B]:ARG:HD3	1:B:116:ARG:HH12	1.82	0.44
1:B:243:SER:HA	1:B:244:PRO:HD2	1.86	0.43
1:A:56:ALA:HB2	3:A:502:KYN:HN1	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:372:GLU:O	1:B:377:VAL:HG13	2.19	0.43
1:A:89:PRO:C	1:A:91:THR:H	2.22	0.43
1:B:146:THR:HG23	1:B:154:ARG:HD3	1.98	0.43
1:B:198:ILE:HG21	1:B:276:MSE:HE2	2.01	0.43
1:A:189[A]:GLN:HG3	1:A:317:VAL:HG11	2.01	0.42
1:B:309:LEU:N	1:B:309:LEU:HD23	2.34	0.42
1:B:94:ASN:ND2	4:B:629:HOH:O	2.39	0.42
3:A:502:KYN:HN1A	3:A:502:KYN:HA	1.84	0.42
1:A:246:ALA:O	1:A:248:PRO:HD3	2.19	0.42
1:B:41:ASP:HB2	1:B:131:PHE:CD2	2.55	0.42
1:B:155:LEU:HD21	1:B:157:LYS:HE2	2.01	0.42
1:B:274:SER:HB3	1:B:275:PRO:HD3	2.01	0.42
1:A:146:THR:HG22	1:A:154:ARG:HD2	2.01	0.42
1:B:280:LEU:O	1:B:284:PHE:HB2	2.19	0.42
1:A:38[A]:ARG:HD3	4:A:1115:HOH:O	2.19	0.41
2:B:501:FAD:H6	4:B:851:HOH:O	2.18	0.41
1:A:382:TYR:OH	1:A:386:ARG:NH2	2.53	0.41
1:A:81[B]:MSE:HB2	1:A:212:ALA:HA	2.01	0.41
1:B:55:LEU:HD21	1:B:111:ARG:NH2	2.34	0.41
1:B:319:PHE:CD2	1:B:370:TYR:HB2	2.55	0.41
1:B:101:ASP:CG	1:B:103:SER:HG	2.23	0.41
1:B:193:TYR:CE1	1:B:290:GLY:HA3	2.55	0.41
1:A:133:LEU:HD22	1:A:155:LEU:CD2	2.42	0.41
1:A:404:TYR:HE2	3:A:502:KYN:N	2.10	0.41
1:A:21:LEU:HD11	1:A:68:ALA:HB2	2.03	0.41
1:B:112[B]:ASP:OD2	1:B:116:ARG:NH2	2.52	0.41
1:B:403[A]:ARG:NH2	4:B:622:HOH:O	2.34	0.41
4:A:799:HOH:O	1:B:185:HIS:HD2	2.04	0.41
1:A:165:ALA:O	2:A:501:FAD:H52A	2.22	0.40
1:A:81[B]:MSE:SE	1:A:213:LEU:HB2	2.71	0.40
1:B:276:MSE:HE2	1:B:276:MSE:HB3	1.75	0.40
1:B:301:HIS:HE1	4:B:925:HOH:O	2.04	0.40
2:B:501:FAD:H1'1	2:B:501:FAD:H9	1.79	0.40

All (4) 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)
4:A:1128:HOH:O	4:A:1196:HOH:O[1_565]	2.09	0.11
1:B:45:GLU:OE2	1:B:82:ARG:NH2[2_655]	2.11	0.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:977:HOH:O	4:A:1106:HOH:O[2_546]	2.15	0.05
4:A:1006:HOH:O	4:A:1184:HOH:O[2_556]	2.18	0.02

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	469/473 (99%)	460 (98%)	9 (2%)	0	100	100
1	B	463/473 (98%)	448 (97%)	15 (3%)	0	100	100
All	All	932/946 (98%)	908 (97%)	24 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	383/366 (105%)	372 (97%)	11 (3%)	42	18
1	B	378/366 (103%)	363 (96%)	15 (4%)	31	10
All	All	761/732 (104%)	735 (97%)	26 (3%)	35	13

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	ARG

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Mol	Chain	Res	Type
1	A	113	ARG
1	A	155	LEU
1	A	158	ARG
1	A	256	LEU
1	A	268	ARG
1	A	291	LYS
1	A	348	LEU
1	A	383	LEU
1	A	384	LEU
1	A	437	ASP
1	B	22	VAL
1	B	49	ARG
1	B	53	ILE
1	B	77	GLU
1	B	100	ARG
1	B	101	ASP
1	B	102	ASP
1	B	151	SER
1	B	154	ARG
1	B	256	LEU
1	B	282	GLN
1	B	286	HIS
1	B	386	ARG
1	B	412	LEU
1	B	427	LEU

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

Mol	Chain	Res	Type
1	A	32	GLN
1	A	132	ASN
1	B	34	ASN

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 monosaccharides 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	$\#\|Z\| > 2$
3	KYN	B	502	-	13,15,15	1.42	3 (23%)	17,20,20	2.74	8 (47%)
3	KYN	A	502	-	13,15,15	1.49	3 (23%)	17,20,20	3.05	8 (47%)
2	FAD	B	501	-	53,58,58	1.79	10 (18%)	68,89,89	3.38	17 (25%)
2	FAD	A	501	-	53,58,58	2.21	17 (32%)	68,89,89	3.09	23 (33%)

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	KYN	B	502	-	-	6/12/12/12	0/1/1/1
3	KYN	A	502	-	-	8/12/12/12	0/1/1/1
2	FAD	B	501	-	-	2/30/50/50	0/6/6/6
2	FAD	A	501	-	-	4/30/50/50	0/6/6/6

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	FAD	O4B-C4B	7.00	1.60	1.45
2	A	501	FAD	C1'-C2'	5.03	1.59	1.52
2	B	501	FAD	O4B-C4B	4.73	1.55	1.45
2	B	501	FAD	C7M-C7	-4.68	1.41	1.51
2	B	501	FAD	C4X-N5	4.30	1.39	1.30
2	A	501	FAD	O2B-C2B	4.26	1.53	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	FAD	C4X-N5	4.14	1.38	1.30
2	B	501	FAD	O2B-C2B	3.78	1.51	1.43
2	A	501	FAD	C2A-N3A	3.74	1.38	1.32
2	B	501	FAD	C8M-C8	-3.74	1.43	1.51
2	A	501	FAD	C9A-N10	-3.61	1.34	1.41
2	A	501	FAD	O5B-C5B	3.28	1.57	1.44
2	A	501	FAD	C2A-N1A	3.16	1.39	1.33
2	A	501	FAD	C7M-C7	-3.10	1.44	1.51
3	B	502	KYN	CD2-CG	-3.01	1.37	1.41
2	B	501	FAD	C2A-N3A	2.99	1.36	1.32
2	A	501	FAD	C4A-N3A	-2.98	1.31	1.35
2	B	501	FAD	C2A-N1A	2.94	1.39	1.33
2	B	501	FAD	C10-N1	2.93	1.39	1.33
2	A	501	FAD	O3B-C3B	-2.87	1.36	1.43
2	A	501	FAD	C2'-C3'	2.81	1.58	1.53
3	A	502	KYN	CD2-CG	-2.80	1.37	1.41
2	A	501	FAD	C5'-C4'	2.73	1.55	1.51
2	B	501	FAD	C9A-N10	-2.60	1.36	1.41
2	A	501	FAD	C4'-C3'	-2.55	1.48	1.53
2	A	501	FAD	PA-O2A	-2.51	1.43	1.55
3	A	502	KYN	CB-C1	2.43	1.54	1.51
2	B	501	FAD	C5A-C4A	-2.25	1.35	1.40
3	A	502	KYN	O-C	2.25	1.29	1.22
2	A	501	FAD	C8M-C8	-2.16	1.46	1.51
3	B	502	KYN	CA-N	-2.06	1.38	1.48
2	A	501	FAD	C10-N1	2.06	1.37	1.33
3	B	502	KYN	CE2-CD2	-2.02	1.36	1.39

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	FAD	O4B-C1B-C2B	-17.93	80.72	106.93
2	A	501	FAD	O4B-C1B-C2B	-12.78	88.26	106.93
2	B	501	FAD	O4B-C4B-C5B	-8.28	82.12	109.37
2	B	501	FAD	N3A-C2A-N1A	-8.03	116.12	128.68
2	A	501	FAD	O4B-C4B-C5B	-7.98	83.13	109.37
2	A	501	FAD	O3B-C3B-C2B	7.69	136.68	111.82
2	B	501	FAD	O4B-C4B-C3B	-7.43	90.40	105.11
3	A	502	KYN	O2-C1-CD2	-7.38	107.78	120.83
2	A	501	FAD	N3A-C2A-N1A	-7.07	117.63	128.68
2	B	501	FAD	O3B-C3B-C2B	6.91	134.16	111.82
2	A	501	FAD	C3B-C2B-C1B	-6.74	90.83	100.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	502	KYN	O2-C1-CD2	-6.56	109.24	120.83
2	A	501	FAD	C5B-C4B-C3B	6.32	138.85	115.18
3	A	502	KYN	O2-C1-CB	5.78	127.46	120.76
2	B	501	FAD	O5B-C5B-C4B	-5.77	89.13	108.99
2	B	501	FAD	C5B-C4B-C3B	5.65	136.35	115.18
2	B	501	FAD	C3B-C2B-C1B	-5.37	92.89	100.98
2	A	501	FAD	O4B-C4B-C3B	-5.32	94.59	105.11
3	B	502	KYN	CA-CB-C1	-5.15	107.14	113.70
2	A	501	FAD	O5B-C5B-C4B	-4.38	93.90	108.99
3	A	502	KYN	CA-CB-C1	-4.33	108.19	113.70
3	B	502	KYN	CB-C1-CD2	4.14	125.61	119.86
2	A	501	FAD	O2B-C2B-C3B	-3.96	99.02	111.82
2	A	501	FAD	C9A-C5X-N5	-3.79	118.31	122.43
3	B	502	KYN	O2-C1-CB	3.76	125.13	120.76
3	A	502	KYN	CE2-CD2-CG	3.73	122.89	118.93
2	B	501	FAD	O2B-C2B-C3B	-3.63	100.09	111.82
2	B	501	FAD	C2A-N1A-C6A	3.61	124.93	118.75
2	A	501	FAD	C2A-N1A-C6A	3.49	124.72	118.75
3	A	502	KYN	CB-C1-CD2	3.44	124.64	119.86
2	B	501	FAD	C9A-C5X-N5	-3.13	119.03	122.43
2	A	501	FAD	C10-N1-C2	3.05	123.00	116.90
2	B	501	FAD	C1B-N9A-C4A	3.00	131.91	126.64
2	B	501	FAD	C5X-C9A-N10	2.92	120.97	117.95
2	A	501	FAD	C5X-C9A-N10	2.91	120.96	117.95
3	B	502	KYN	CD2-CG-N1	-2.90	118.84	122.67
2	A	501	FAD	C4X-C4-N3	2.89	120.54	113.19
2	A	501	FAD	C4-C4X-N5	2.79	122.20	118.23
2	A	501	FAD	C6-C5X-C9A	2.74	122.81	118.94
2	A	501	FAD	O4-C4-C4X	-2.70	119.45	126.60
2	A	501	FAD	C1B-N9A-C4A	2.63	131.26	126.64
2	A	501	FAD	O2B-C2B-C1B	-2.61	101.20	110.85
2	A	501	FAD	C6-C7-C8	-2.49	116.10	119.67
2	B	501	FAD	C6-C5X-C9A	2.48	122.44	118.94
3	B	502	KYN	OXT-C-CA	2.42	121.63	113.38
2	A	501	FAD	C9-C8-C7	2.41	123.12	119.67
2	B	501	FAD	O3'-C3'-C2'	2.30	114.37	108.81
3	A	502	KYN	CB-CA-C	-2.28	99.53	111.83
2	A	501	FAD	C9-C9A-C5X	-2.28	115.81	120.11
3	B	502	KYN	O-C-CA	-2.19	114.42	122.14
3	B	502	KYN	CD1-CG-CD2	2.16	120.19	118.10
3	A	502	KYN	CD2-CG-N1	-2.14	119.84	122.67
2	B	501	FAD	O2P-P-O1P	2.13	122.78	112.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	FAD	C10-C4X-N5	-2.13	120.33	124.86
3	A	502	KYN	CZ-CE1-CD1	2.03	123.28	120.19
2	B	501	FAD	O2-C2-N3	2.01	122.55	118.65

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	KYN	O-C-CA-N
3	B	502	KYN	O-C-CA-N
3	A	502	KYN	OXT-C-CA-N
3	B	502	KYN	OXT-C-CA-N
2	B	501	FAD	C3B-C4B-C5B-O5B
3	A	502	KYN	O2-C1-CD2-CE2
3	A	502	KYN	O2-C1-CB-CA
3	A	502	KYN	CD2-C1-CB-CA
2	A	501	FAD	O2'-C2'-C3'-C4'
3	A	502	KYN	OXT-C-CA-CB
3	B	502	KYN	O2-C1-CD2-CG
3	B	502	KYN	CD2-C1-CB-CA
2	B	501	FAD	O4B-C4B-C5B-O5B
3	B	502	KYN	O2-C1-CD2-CE2
2	A	501	FAD	O2'-C2'-C3'-O3'
3	A	502	KYN	O-C-CA-CB
2	A	501	FAD	C3B-C4B-C5B-O5B
3	A	502	KYN	O2-C1-CD2-CG
3	B	502	KYN	O2-C1-CB-CA
2	A	501	FAD	O4B-C4B-C5B-O5B

There are no ring outliers.

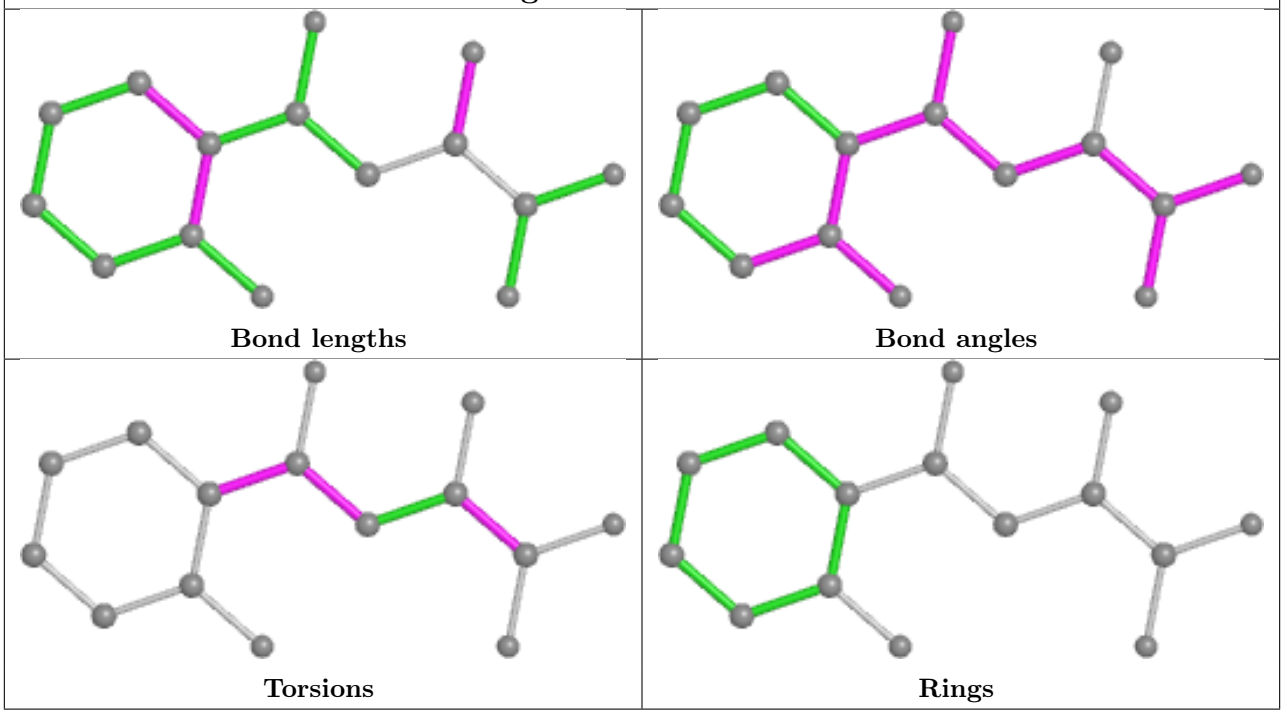
4 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	KYN	3	0
3	A	502	KYN	9	0
2	B	501	FAD	3	0
2	A	501	FAD	1	0

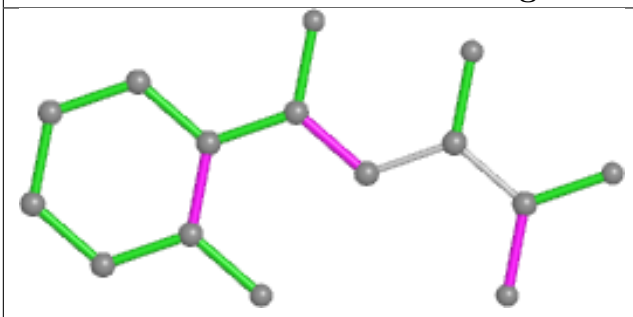
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less then 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.

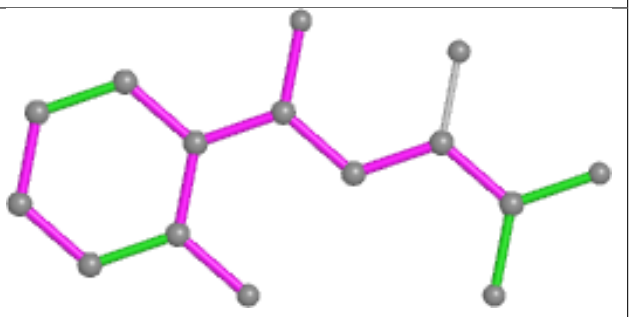
Ligand KYN B 502



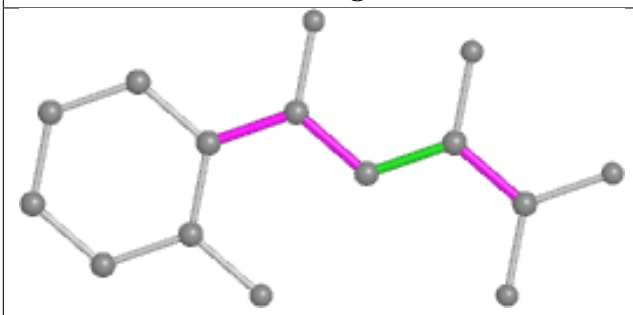
Ligand KYN A 502



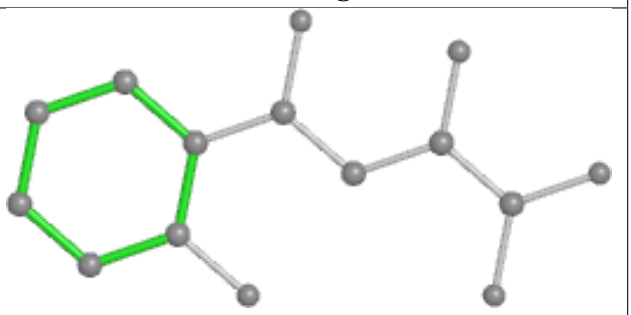
Bond lengths



Bond angles

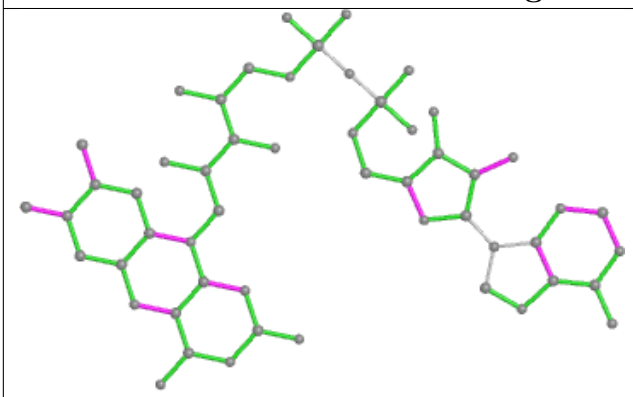


Torsions

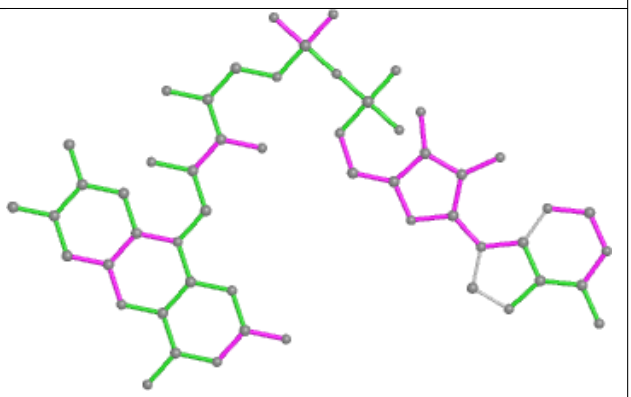


Rings

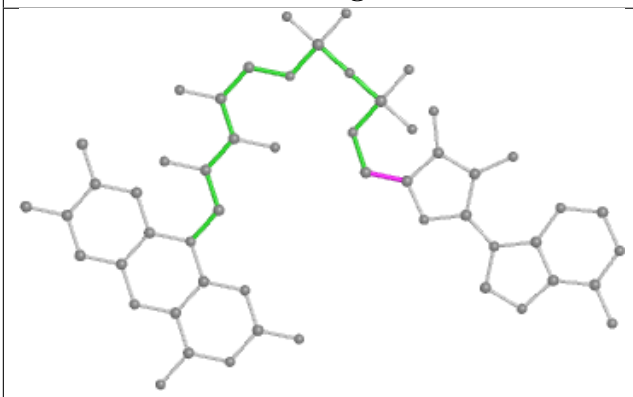
Ligand FAD B 501



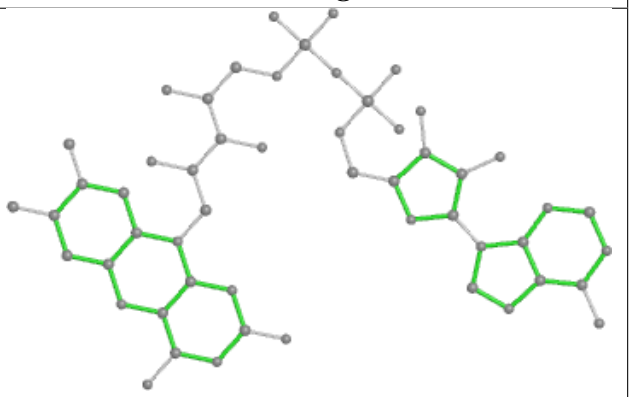
Bond lengths



Bond angles

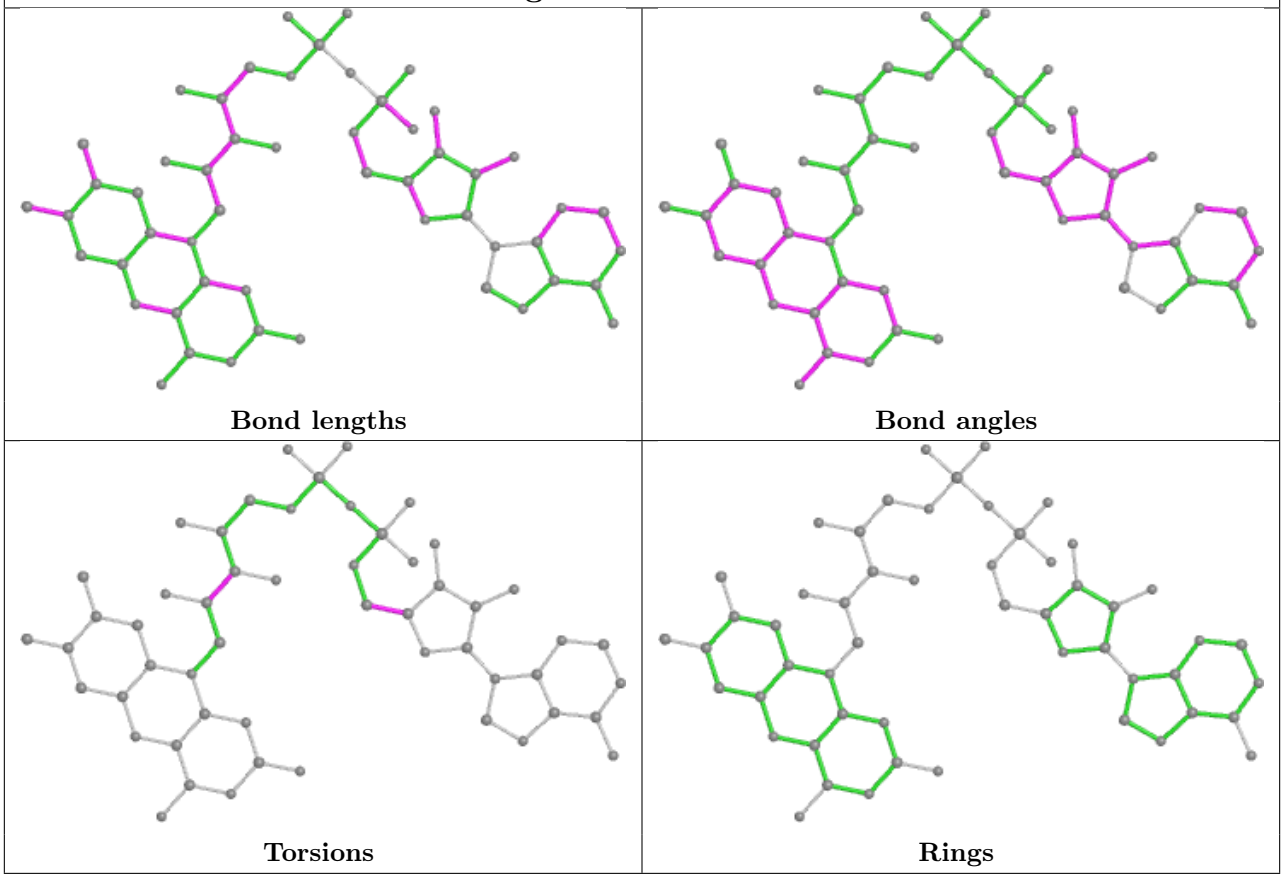


Torsions



Rings

Ligand FAD A 501



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	442/473 (93%)	-0.07	18 (4%)	37	34	11, 20, 46, 79	0
1	B	438/473 (92%)	0.22	36 (8%)	11	10	14, 27, 63, 75	0
All	All	880/946 (93%)	0.07	54 (6%)	21	19	11, 24, 57, 79	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	461	CYS	11.6
1	A	460	LEU	7.6
1	A	245	ALA	6.1
1	B	150	VAL	6.1
1	B	247	GLN	6.1
1	B	246	ALA	6.0
1	A	247	GLN	5.3
1	B	458	SER	5.3
1	B	257	VAL	5.2
1	A	246	ALA	4.9
1	B	286	HIS	4.7
1	B	245	ALA	4.5
1	A	90	GLY	4.4
1	B	260	HIS	4.3
1	B	49	ARG	4.0
1	B	100	ARG	4.0
1	B	244	PRO	3.9
1	B	438	LEU	3.9
1	B	284	PHE	3.8
1	B	249	ALA	3.8
1	B	248	PRO	3.7
1	A	91	THR	3.7
1	B	51	ARG	3.7
1	A	459	HIS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	248	PRO	3.4
1	B	262	ALA	3.3
1	A	380	PRO	3.2
1	B	258	ASP	3.1
1	A	378	ALA	3.1
1	B	259	GLY	2.9
1	B	264	ARG	2.9
1	A	439	THR	2.9
1	B	250	SER	2.8
1	B	439	THR	2.7
1	B	287	HIS	2.7
1	A	458	SER	2.7
1	B	101	ASP	2.6
1	B	154	ARG	2.6
1	B	280	LEU	2.6
1	B	282	GLN	2.6
1	A	94	ASN	2.5
1	B	152	GLY	2.5
1	B	254	ALA	2.5
1	B	283	ASP	2.4
1	A	92	PRO	2.4
1	A	379	SER	2.4
1	B	242	GLN	2.4
1	B	285	GLU	2.3
1	B	268	ARG	2.3
1	B	436	SER	2.1
1	B	437	ASP	2.1
1	A	319	PHE	2.1
1	B	55	LEU	2.1
1	A	49	ARG	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 monosaccharides in this entry.

6.4 Ligands (i)

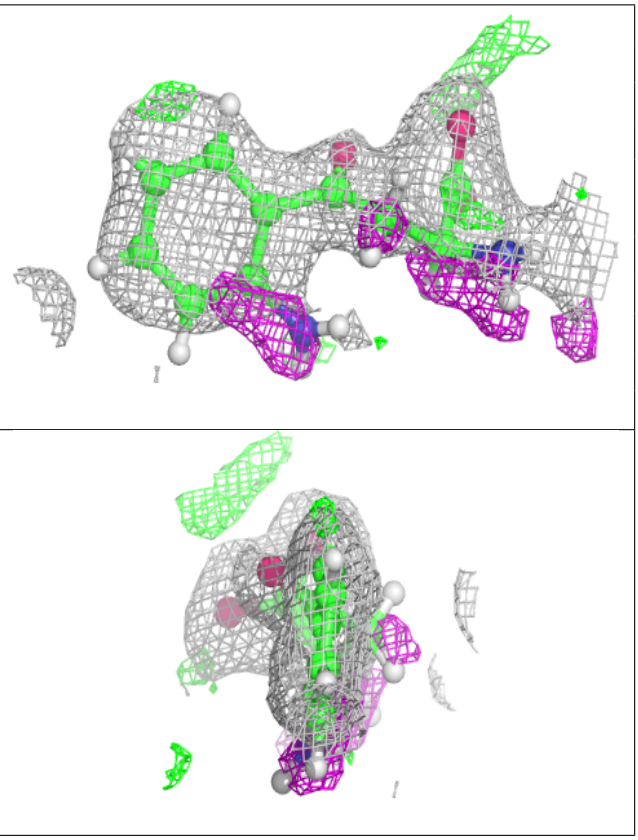
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	KYN	A	502	15/15	0.88	0.16	16,35,48,50	0
3	KYN	B	502	15/15	0.91	0.20	14,35,48,55	0
2	FAD	A	501	53/53	0.97	0.10	8,12,15,17	0
2	FAD	B	501	53/53	0.97	0.10	12,16,25,35	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.

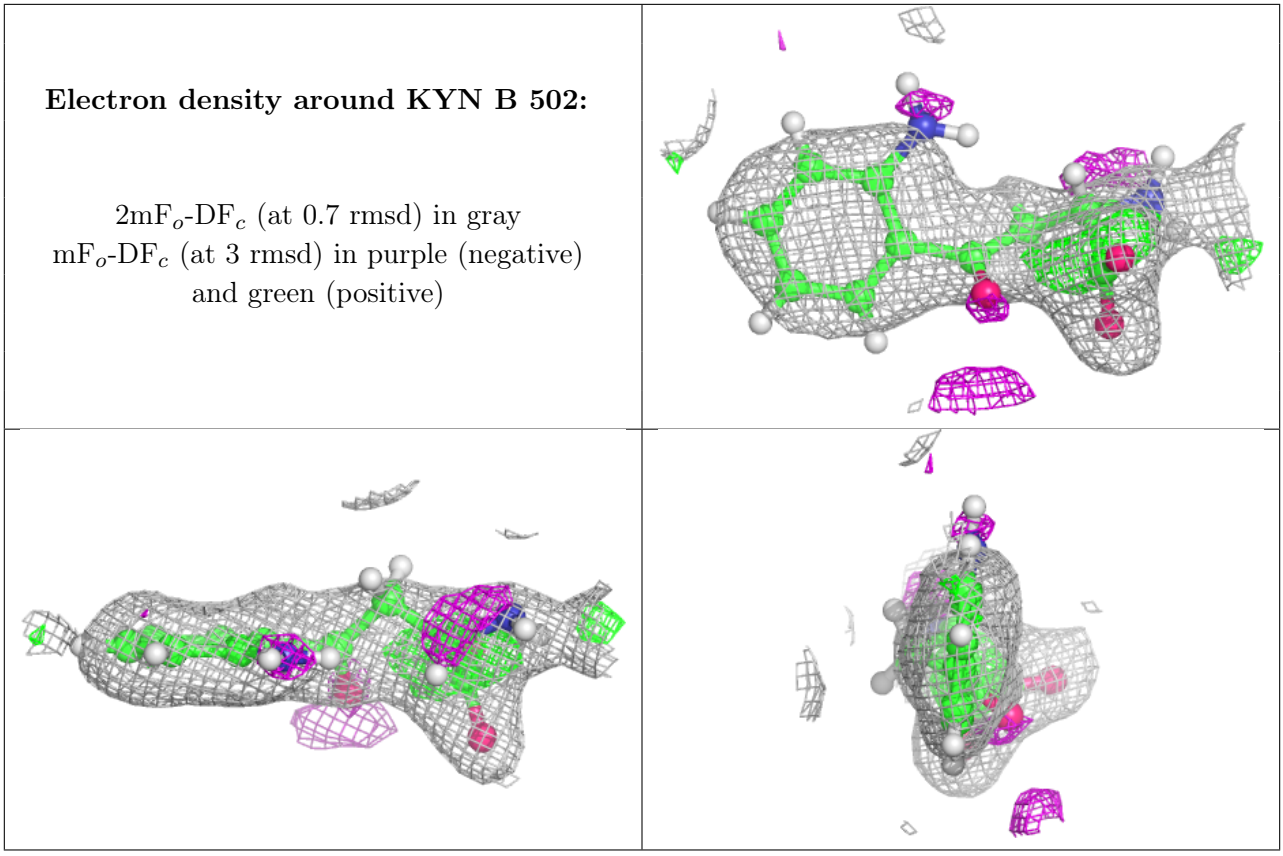
Electron density around KYN A 502:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)



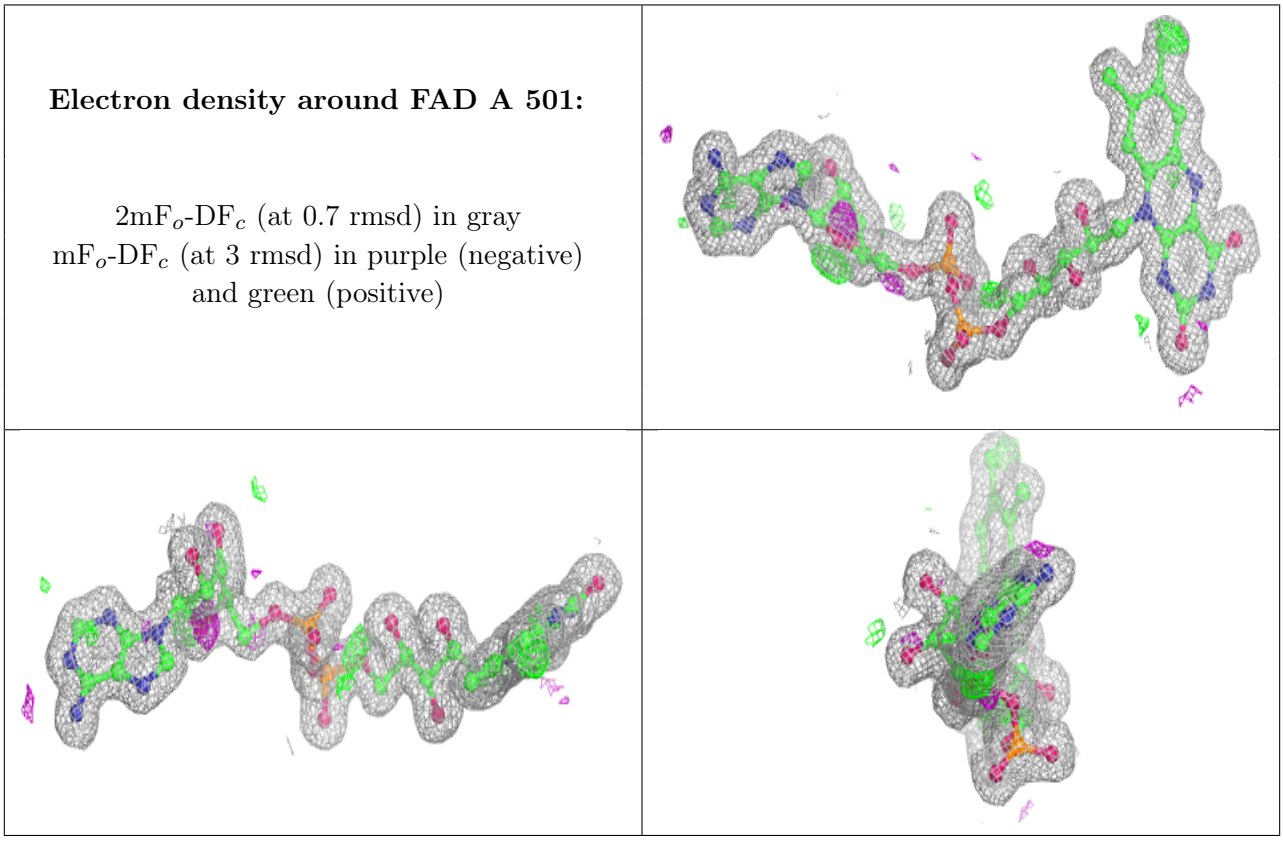
Electron density around KYN B 502:

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



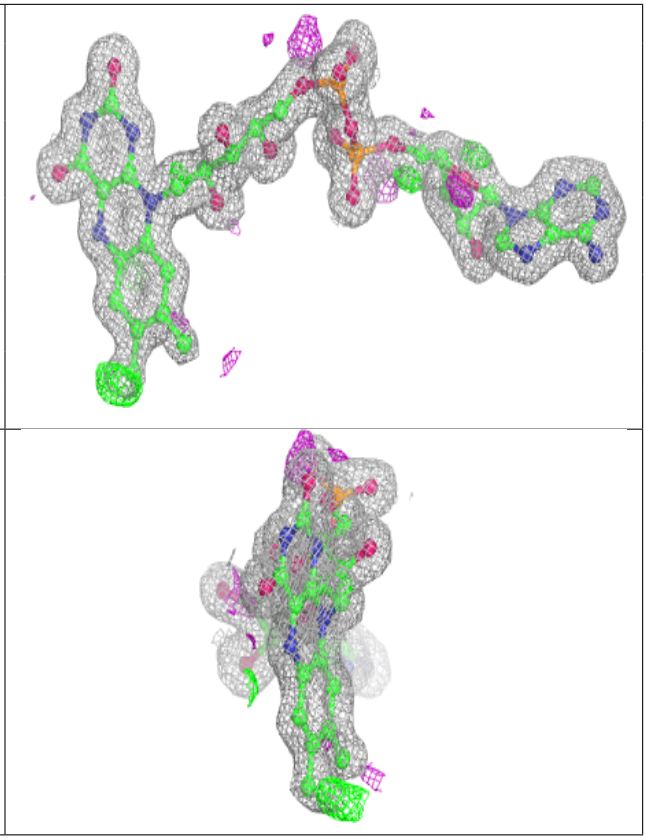
Electron density around FAD A 501:

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



Electron density around FAD B 501:

$2mF_o$-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.