



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:31 PM UTC

PDB ID : 5FN0 / pdb_00005fn0
Title : Crystal structure of *Pseudomonas fluorescens* kynurenine-3- monooxygenase (KMO) in complex with GSK180
Authors : Mole, D.J.; Webster, S.P.; Uings, I.; Zheng, X.; Binnie, M.; Wilson, K.; Hutchinson, J.P.; Mirguet, O.; Walker, A.; Beaufils, B.; Ancellin, N.; Trottet, L.; Beneton, V.; Mowat, C.G.; Wilkinson, M.; Rowland, P.; Haslam, C.; McBride, A.; Homer, N.Z.M.; Baily, J.E.; Sharp, M.G.F.; Garden, O.J.; Hughes, J.; Howie, S.E.M.; Holmes, D.; Liddle, J.; Iredale, J.P.
Deposited on : 2015-11-10
Resolution : 3.19 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)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12

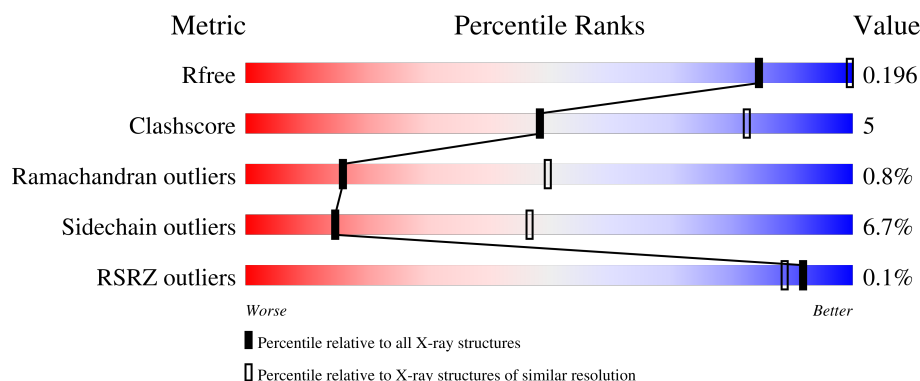
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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 $\leq 5\%$. 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	461	
1	B	461	
1	C	461	

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Ideal geometry (proteins) : Engh & Huber (2001)
 Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
1	D	461	79% 17% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14670 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.

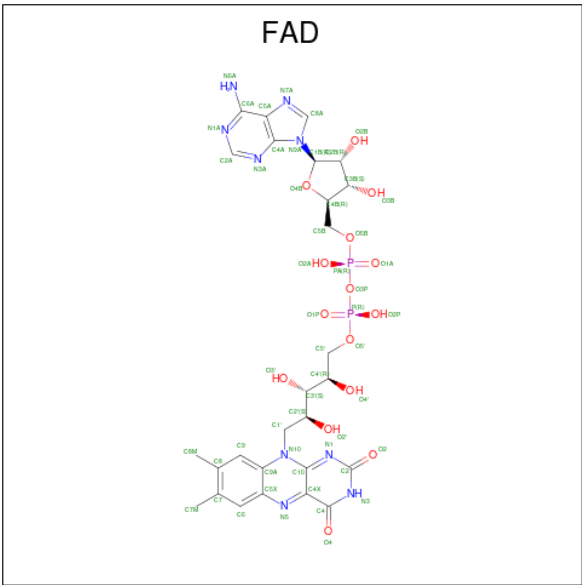
- Molecule 1 is a protein called KYNURENINE 3-MONOOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total	C	N	O	S	0	0	0
			3523	2206	654	646	17			
1	B	455	Total	C	N	O	S	0	0	0
			3523	2206	654	646	17			
1	C	455	Total	C	N	O	S	0	0	0
			3523	2206	654	646	17			
1	D	455	Total	C	N	O	S	0	0	0
			3523	2206	654	646	17			

There are 8 discrepancies between the modelled and reference sequences:

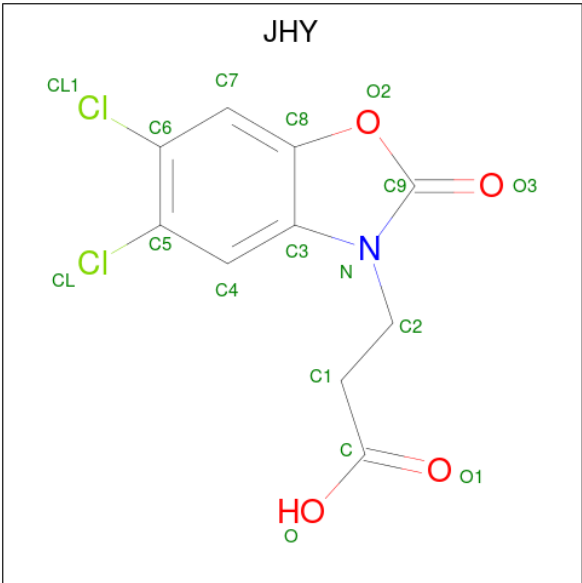
Chain	Residue	Modelled	Actual	Comment	Reference
A	461	SER	CYS	engineered mutation	UNP Q84HF5
A	252	SER	CYS	engineered mutation	UNP Q84HF5
B	461	SER	CYS	engineered mutation	UNP Q84HF5
B	252	SER	CYS	engineered mutation	UNP Q84HF5
C	461	SER	CYS	engineered mutation	UNP Q84HF5
C	252	SER	CYS	engineered mutation	UNP Q84HF5
D	461	SER	CYS	engineered mutation	UNP Q84HF5
D	252	SER	CYS	engineered mutation	UNP Q84HF5

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 3 is 3-(5,6-DICHLORO-2-OXOBENZO[D]OXAZOL-3(2H)-YL)PROPANOIC ACID (CCD ID: JHY) (formula: C₁₀H₇Cl₂NO₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			17	10	2	1	4		
3	B	1	Total	C	Cl	N	O	0	0
			17	10	2	1	4		
3	C	1	Total	C	Cl	N	O	0	0
			17	10	2	1	4		
3	D	1	Total	C	Cl	N	O	0	0
			17	10	2	1	4		

- Molecule 4 is water.

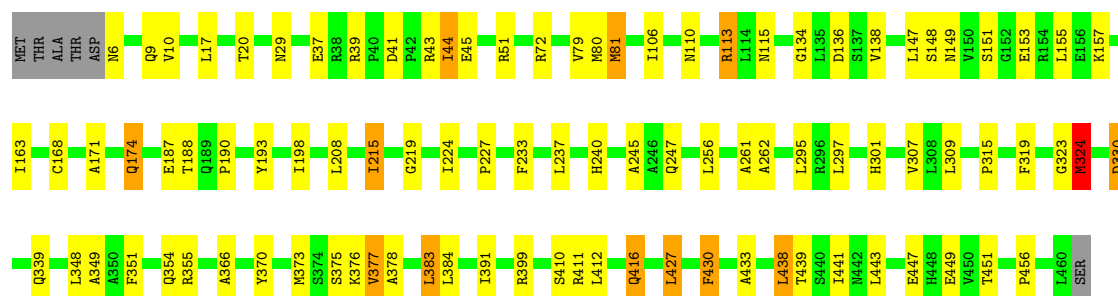
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	77	Total	O	0	0
			77	77		
4	B	73	Total	O	0	0
			73	73		
4	C	64	Total	O	0	0
			64	64		
4	D	84	Total	O	0	0
			84	84		

3 Residue-property plots

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.

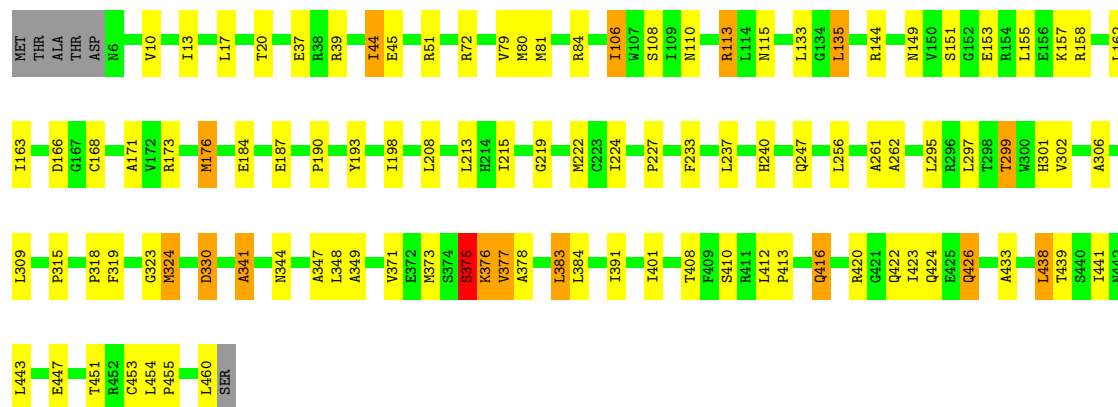
• Molecule 1: KYNURENINE 3-MONOOXYGENASE

Chain A: 



• Molecule 1: KYNURENINE 3-MONOOXYGENASE

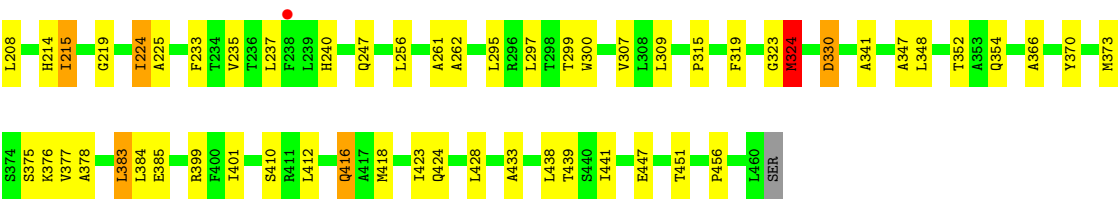
Chain B: 



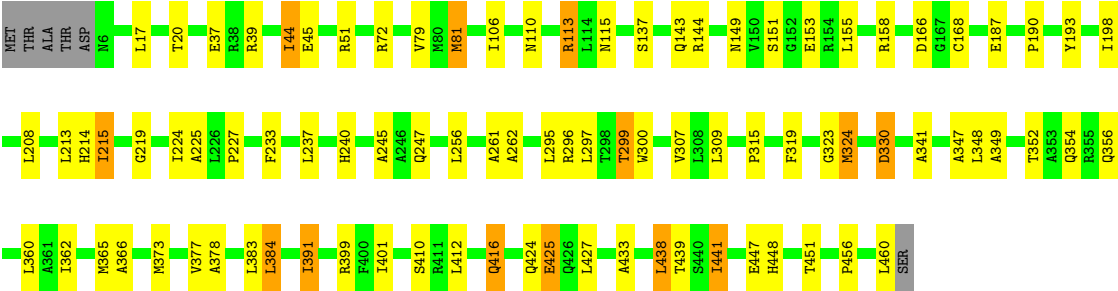
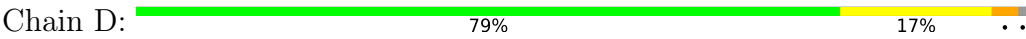
• Molecule 1: KYNURENINE 3-MONOOXYGENASE

Chain C: 





● Molecule 1: KYNURENINE 3-MONOOXYGENASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	186.85Å 105.29Å 133.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	93.43 – 3.19 93.43 – 3.19	Depositor EDS
% Data completeness (in resolution range)	99.6 (93.43-3.19) 99.9 (93.43-3.19)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.12 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.157 , 0.196 0.160 , 0.196	Depositor DCC
R_{free} test set	2244 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	89.0	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 121.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14670	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: JHY, 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	0.87	3/3600 (0.1%)	1.37	27/4891 (0.6%)
1	B	0.88	3/3600 (0.1%)	1.36	17/4891 (0.3%)
1	C	0.87	1/3600 (0.0%)	1.38	13/4891 (0.3%)
1	D	0.85	0/3600	1.36	14/4891 (0.3%)
All	All	0.87	7/14400 (0.0%)	1.37	71/19564 (0.4%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	377	VAL	CA-C	6.05	1.60	1.52
1	B	176	MET	SD-CE	-5.93	1.64	1.79
1	A	138	VAL	CA-C	5.50	1.59	1.52
1	B	299	THR	CA-C	-5.35	1.46	1.52
1	A	377	VAL	CA-C	5.28	1.59	1.52
1	C	324	MET	SD-CE	-5.25	1.66	1.79
1	A	324	MET	SD-CE	-5.03	1.67	1.79

All (71) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	330	ASP	CA-CB-CG	7.56	120.16	112.60
1	C	330	ASP	CA-CB-CG	7.41	120.01	112.60
1	A	430	PHE	CA-CB-CG	7.05	120.85	113.80
1	C	383	LEU	CA-C-N	6.94	129.46	120.44
1	C	383	LEU	C-N-CA	6.94	129.46	120.44
1	A	330	ASP	CA-CB-CG	6.70	119.30	112.60
1	A	349	ALA	CA-C-N	6.28	128.98	120.44
1	A	349	ALA	C-N-CA	6.28	128.98	120.44
1	B	330	ASP	CA-CB-CG	6.26	118.86	112.60
1	C	433	ALA	CA-C-N	6.23	133.44	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	433	ALA	C-N-CA	6.23	133.44	121.54
1	B	376	LYS	CA-C-N	6.13	133.01	121.97
1	B	376	LYS	C-N-CA	6.13	133.01	121.97
1	A	37	GLU	CA-C-N	6.06	128.73	120.54
1	A	37	GLU	C-N-CA	6.06	128.73	120.54
1	D	391	ILE	N-CA-CB	6.02	117.59	110.55
1	A	370	TYR	CA-C-N	6.00	128.34	120.60
1	A	370	TYR	C-N-CA	6.00	128.34	120.60
1	B	349	ALA	CA-C-N	5.96	128.18	120.44
1	B	349	ALA	C-N-CA	5.96	128.18	120.44
1	B	301	HIS	N-CA-C	5.94	118.10	109.07
1	A	187	GLU	CA-C-N	5.77	129.02	120.95
1	A	187	GLU	C-N-CA	5.77	129.02	120.95
1	D	425	GLU	CA-CB-CG	5.63	125.36	114.10
1	B	37	GLU	CA-C-N	5.58	127.69	120.44
1	B	37	GLU	C-N-CA	5.58	127.69	120.44
1	C	81	MET	CA-C-N	5.56	127.73	120.28
1	C	81	MET	C-N-CA	5.56	127.73	120.28
1	B	383	LEU	CA-C-N	5.55	127.71	120.28
1	B	383	LEU	C-N-CA	5.55	127.71	120.28
1	B	375	SER	CA-C-N	5.50	132.04	121.54
1	B	375	SER	C-N-CA	5.50	132.04	121.54
1	B	341	ALA	CA-C-N	5.44	128.12	120.28
1	B	341	ALA	C-N-CA	5.44	128.12	120.28
1	A	171	ALA	N-CA-C	-5.41	105.55	111.82
1	B	261	ALA	CA-C-N	5.36	127.90	120.29
1	B	261	ALA	C-N-CA	5.36	127.90	120.29
1	D	166	ASP	CA-CB-CG	5.35	117.95	112.60
1	A	261	ALA	CA-C-N	5.31	127.83	120.29
1	A	261	ALA	C-N-CA	5.31	127.83	120.29
1	D	261	ALA	CA-C-N	5.30	127.82	120.29
1	D	261	ALA	C-N-CA	5.30	127.82	120.29
1	C	423	ILE	CA-C-N	5.29	127.32	120.44
1	C	423	ILE	C-N-CA	5.29	127.32	120.44
1	A	6	ASN	CA-C-N	5.29	127.90	120.28
1	A	6	ASN	C-N-CA	5.29	127.90	120.28
1	D	349	ALA	CA-C-N	5.24	127.62	120.54
1	D	349	ALA	C-N-CA	5.24	127.62	120.54
1	A	174	GLN	N-CA-C	-5.23	105.47	111.07
1	A	81	MET	CA-C-N	5.23	127.28	120.28
1	A	81	MET	C-N-CA	5.23	127.28	120.28
1	D	81	MET	CA-C-N	5.20	127.25	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	81	MET	C-N-CA	5.20	127.25	120.28
1	A	301	HIS	N-CA-C	5.17	116.92	109.07
1	A	245	ALA	CA-C-N	5.15	127.45	120.65
1	A	245	ALA	C-N-CA	5.15	127.45	120.65
1	A	383	LEU	CA-C-N	5.15	127.44	120.44
1	A	383	LEU	C-N-CA	5.15	127.44	120.44
1	C	261	ALA	CA-C-N	5.09	127.52	120.29
1	C	261	ALA	C-N-CA	5.09	127.52	120.29
1	C	376	LYS	CA-C-N	5.02	131.01	121.97
1	C	376	LYS	C-N-CA	5.02	131.01	121.97
1	A	174	GLN	CA-C-N	5.02	127.32	120.54
1	A	174	GLN	C-N-CA	5.02	127.32	120.54
1	A	411	ARG	CA-C-N	5.01	127.47	120.65
1	A	411	ARG	C-N-CA	5.01	127.47	120.65
1	D	245	ALA	CA-C-N	5.01	127.27	120.65
1	D	245	ALA	C-N-CA	5.01	127.27	120.65
1	B	166	ASP	CA-CB-CG	5.01	117.61	112.60
1	D	37	GLU	CA-C-N	5.00	126.94	120.44
1	D	37	GLU	C-N-CA	5.00	126.94	120.44

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	3523	0	3482	37	0
1	B	3523	0	3482	46	0
1	C	3523	0	3482	36	0
1	D	3523	0	3482	37	0
2	A	53	0	31	1	0
2	B	53	0	31	2	0
2	C	53	0	31	1	0
2	D	53	0	31	1	0
3	A	17	0	6	0	0
3	B	17	0	6	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	17	0	6	1	0
3	D	17	0	6	0	0
4	A	77	0	0	0	0
4	B	73	0	0	2	0
4	C	64	0	0	0	0
4	D	84	0	0	4	0
All	All	14670	0	14076	155	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (155) 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:106:ILE:HD11	1:B:408:THR:HG21	1.64	0.78
1:C:295:LEU:HB3	1:C:315:PRO:HD2	1.73	0.70
1:A:39:ARG:HH22	1:A:115:ASN:HD21	1.44	0.66
1:D:460:LEU:C	4:D:2082:HOH:O	2.38	0.66
1:A:136:ASP:OD2	1:A:148:SER:HB3	1.96	0.66
1:D:17:LEU:HD13	1:D:324:MET:HG3	1.79	0.65
1:C:41:ASP:HB3	1:C:44:ILE:HG12	1.79	0.65
1:D:39:ARG:HH22	1:D:115:ASN:HD21	1.45	0.64
1:B:17:LEU:HD13	1:B:324:MET:HG3	1.79	0.64
1:B:295:LEU:HB3	1:B:315:PRO:HD2	1.80	0.64
1:C:17:LEU:HD13	1:C:324:MET:HG3	1.79	0.63
1:D:295:LEU:HB3	1:D:315:PRO:HD2	1.81	0.62
1:A:17:LEU:HD13	1:A:324:MET:HG3	1.81	0.62
1:C:39:ARG:HH22	1:C:115:ASN:HD21	1.47	0.62
1:B:39:ARG:HH22	1:B:115:ASN:HD21	1.47	0.61
1:A:430:PHE:HD2	1:B:426:GLN:HG2	1.66	0.61
1:B:173:ARG:HD2	1:B:297:LEU:HD11	1.81	0.61
1:C:385:GLU:HG2	1:C:428:LEU:HB3	1.82	0.61
1:D:299:THR:HA	4:D:2047:HOH:O	2.01	0.61
1:C:215:ILE:HD12	1:C:373:MET:HB2	1.83	0.61
1:B:149:ASN:HD21	1:B:153:GLU:HB2	1.66	0.60
1:B:222:MET:HE1	1:B:373:MET:HB3	1.83	0.60
1:B:378:ALA:HB1	1:B:383:LEU:HD21	1.83	0.60
1:A:29:ASN:ND2	1:A:339:GLN:HE22	2.00	0.59
1:B:81:MET:HE2	1:B:108:SER:HB2	1.84	0.59
1:A:378:ALA:HB1	1:A:383:LEU:HD21	1.86	0.58
1:D:401:ILE:H	1:D:424:GLN:HE22	1.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:LEU:HB3	1:A:315:PRO:HD2	1.85	0.57
1:A:215:ILE:HD12	1:A:373:MET:HB2	1.84	0.57
1:A:427:LEU:HD11	1:A:449:GLU:HB3	1.86	0.57
1:D:215:ILE:HD12	1:D:373:MET:HB2	1.86	0.56
1:C:378:ALA:HB1	1:C:383:LEU:HD21	1.87	0.56
1:D:378:ALA:HB1	1:D:383:LEU:HD21	1.86	0.56
1:B:110:ASN:HB3	1:B:113:ARG:HB2	1.86	0.56
1:C:224:ILE:O	1:C:235:VAL:HA	2.06	0.56
1:B:309:LEU:HD12	1:B:330:ASP:HB3	1.87	0.56
1:D:110:ASN:HB3	1:D:113:ARG:HB2	1.87	0.56
1:D:307:VAL:HB	1:D:348:LEU:HD13	1.86	0.56
1:C:309:LEU:HD12	1:C:330:ASP:HB3	1.88	0.55
1:A:51:ARG:NH2	1:A:168:CYS:HB2	2.21	0.55
1:A:110:ASN:HB3	1:A:113:ARG:HB2	1.88	0.55
1:B:44:ILE:HG13	1:B:45:GLU:H	1.72	0.55
1:D:51:ARG:NH2	1:D:168:CYS:HB2	2.22	0.55
1:B:81:MET:HE3	1:B:106:ILE:HG22	1.89	0.55
1:C:110:ASN:HB3	1:C:113:ARG:HB2	1.88	0.54
1:A:309:LEU:HD12	1:A:330:ASP:HB3	1.87	0.54
1:B:51:ARG:NH2	1:B:168:CYS:HB2	2.24	0.53
1:C:143:GLN:HG2	1:C:159:PHE:O	2.09	0.52
1:C:173:ARG:HD2	1:C:297:LEU:HD11	1.91	0.52
1:A:44:ILE:HG13	1:A:45:GLU:H	1.74	0.52
1:D:44:ILE:HG13	1:D:45:GLU:H	1.74	0.52
1:C:307:VAL:HB	1:C:348:LEU:HD13	1.91	0.52
1:A:430:PHE:HE1	1:B:453:CYS:HG	1.58	0.52
1:C:51:ARG:NH2	1:C:168:CYS:HB2	2.25	0.52
1:C:193:TYR:HA	1:C:237:LEU:O	2.10	0.51
1:B:460:LEU:C	4:B:2073:HOH:O	2.52	0.51
1:B:384:LEU:HD12	1:B:438:LEU:HD12	1.92	0.51
1:B:193:TYR:HA	1:B:237:LEU:O	2.11	0.50
1:C:214:HIS:HB2	1:C:225:ALA:HB3	1.93	0.50
1:A:29:ASN:HD22	1:A:339:GLN:HE22	1.58	0.49
1:D:256:LEU:HD22	1:D:262:ALA:HA	1.93	0.49
1:A:193:TYR:HA	1:A:237:LEU:O	2.12	0.49
1:B:135:LEU:HD23	1:B:171:ALA:HB3	1.93	0.49
1:C:256:LEU:HD22	1:C:262:ALA:HA	1.95	0.49
1:D:214:HIS:HB2	1:D:225:ALA:HB3	1.94	0.49
1:B:401:ILE:H	1:B:424:GLN:HE22	1.60	0.49
1:D:149:ASN:HD21	1:D:153:GLU:HB2	1.77	0.49
1:C:136:ASP:OD2	1:C:148:SER:HB3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:319:PHE:HA	3:C:4000:JHY:CL	2.50	0.49
1:D:193:TYR:HA	1:D:237:LEU:O	2.12	0.49
1:A:256:LEU:HD22	1:A:262:ALA:HA	1.94	0.49
1:B:256:LEU:HD22	1:B:262:ALA:HA	1.95	0.48
1:B:13:ILE:HG13	1:B:135:LEU:HD22	1.96	0.48
1:D:341:ALA:HB3	1:D:347:ALA:HB2	1.95	0.48
1:B:144:ARG:NH1	1:B:158:ARG:HE	2.12	0.48
1:A:307:VAL:HB	1:A:348:LEU:HD13	1.95	0.47
1:D:384:LEU:HG	1:D:438:LEU:HD12	1.96	0.47
1:C:224:ILE:HD11	1:C:373:MET:HE2	1.96	0.47
1:D:412:LEU:HD11	1:D:416:GLN:OE1	2.14	0.47
1:D:299:THR:HG23	4:D:2047:HOH:O	2.14	0.47
1:C:341:ALA:HB3	1:C:347:ALA:HB2	1.96	0.47
1:D:309:LEU:HD12	1:D:330:ASP:HB3	1.96	0.47
1:B:318:PRO:HA	4:B:2045:HOH:O	2.14	0.47
1:A:412:LEU:HD11	1:A:416:GLN:OE1	2.15	0.47
1:C:184:GLU:HG2	1:C:297:LEU:HD22	1.96	0.46
1:A:384:LEU:HD12	1:A:438:LEU:HD12	1.97	0.46
1:A:198:ILE:HB	1:A:233:PHE:HB2	1.98	0.46
1:D:447:GLU:O	1:D:451:THR:HG23	2.15	0.46
1:A:134:GLY:O	1:A:147:LEU:HA	2.16	0.45
1:A:190:PRO:HB2	1:A:240:HIS:CD2	2.52	0.45
1:A:29:ASN:HD22	1:A:339:GLN:NE2	2.14	0.45
1:C:319:PHE:CD2	1:C:370:TYR:HB2	2.51	0.45
1:C:447:GLU:O	1:C:451:THR:HG23	2.16	0.45
1:D:198:ILE:HB	1:D:233:PHE:HB2	1.99	0.45
1:C:198:ILE:HB	1:C:233:PHE:HB2	1.99	0.45
1:D:81:MET:HE3	1:D:106:ILE:HG13	1.99	0.45
1:D:190:PRO:HB2	1:D:240:HIS:CD2	2.52	0.45
1:B:106:ILE:HD11	1:B:408:THR:CG2	2.41	0.44
1:B:184:GLU:HG2	1:B:297:LEU:HD22	1.99	0.44
1:B:190:PRO:HB2	1:B:240:HIS:CD2	2.52	0.44
1:B:198:ILE:HB	1:B:233:PHE:HB2	2.00	0.44
1:B:412:LEU:HD11	1:B:416:GLN:OE1	2.17	0.44
1:C:190:PRO:HB2	1:C:240:HIS:CD2	2.52	0.44
1:C:401:ILE:H	1:C:424:GLN:HE22	1.64	0.44
1:B:149:ASN:ND2	1:B:153:GLU:HB2	2.29	0.44
1:B:176:MET:HE3	1:B:302:VAL:HG23	2.00	0.44
1:C:323:GLY:HA3	2:C:462:FAD:H1'2	2.00	0.44
1:A:81:MET:HB2	1:A:106:ILE:HG13	1.99	0.43
1:B:344:ASN:O	1:B:348:LEU:HG	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:LEU:HD21	1:B:157:LYS:HE3	2.01	0.43
1:D:79:VAL:HG11	1:D:227:PRO:HB2	2.01	0.43
1:D:441:ILE:H	1:D:441:ILE:HG13	1.71	0.43
1:C:44:ILE:HG13	1:C:45:GLU:N	2.34	0.43
1:C:412:LEU:HD11	1:C:416:GLN:OE1	2.19	0.43
1:C:341:ALA:CB	1:C:347:ALA:HB2	2.48	0.42
1:B:323:GLY:HA3	2:B:462:FAD:H1'2	2.01	0.42
1:A:79:VAL:HG11	1:A:227:PRO:HB2	2.00	0.42
1:B:447:GLU:O	1:B:451:THR:HG23	2.19	0.42
1:D:319:PHE:HB2	1:D:366:ALA:O	2.19	0.42
1:A:447:GLU:O	1:A:451:THR:HG23	2.20	0.42
1:B:420:ARG:HA	1:B:423:ILE:HD12	2.01	0.42
1:D:399:ARG:CZ	1:D:456:PRO:HG3	2.49	0.42
1:B:79:VAL:HG11	1:B:227:PRO:HB2	2.01	0.42
1:D:323:GLY:HA3	2:D:462:FAD:H1'2	2.00	0.42
1:A:155:LEU:HD21	1:A:157:LYS:HE3	2.02	0.42
1:A:323:GLY:HA3	2:A:462:FAD:H1'2	2.01	0.42
1:D:149:ASN:ND2	1:D:153:GLU:HB2	2.34	0.42
1:C:319:PHE:HB2	1:C:366:ALA:O	2.19	0.42
1:A:149:ASN:HD21	1:A:153:GLU:HB2	1.85	0.42
1:A:319:PHE:HB2	1:A:366:ALA:O	2.20	0.42
1:D:300:TRP:HB2	1:D:352:THR:HG23	2.01	0.42
1:A:10:VAL:HG21	1:A:163:ILE:HD12	2.01	0.42
1:C:10:VAL:HG21	1:C:163:ILE:HD12	2.02	0.42
1:B:319:PHE:HA	3:B:4000:JHY:CL	2.57	0.41
1:D:352:THR:O	1:D:356:GLN:HB2	2.19	0.41
1:A:351:PHE:CE1	1:A:355:ARG:HG3	2.55	0.41
1:D:362:ILE:HA	1:D:365:MET:HE2	2.02	0.41
1:B:162:LEU:O	1:B:306:ALA:HA	2.21	0.41
1:B:391:ILE:HG21	1:B:443:LEU:HD13	2.02	0.41
1:C:300:TRP:HB2	1:C:352:THR:HG23	2.01	0.41
1:A:399:ARG:CZ	1:A:456:PRO:HG3	2.51	0.41
1:B:133:LEU:HD23	1:B:133:LEU:HA	1.88	0.41
1:A:41:ASP:HB3	1:A:44:ILE:HG12	2.03	0.40
1:B:341:ALA:HB3	1:B:347:ALA:HB2	2.04	0.40
1:D:296:ARG:HD3	1:D:360:LEU:HD22	2.03	0.40
1:D:448:HIS:HE1	4:D:2078:HOH:O	2.04	0.40
1:A:391:ILE:HG21	1:A:443:LEU:HD13	2.03	0.40
1:A:399:ARG:HD3	1:A:451:THR:HA	2.03	0.40
1:B:454:LEU:HA	1:B:455:PRO:HD3	1.96	0.40
1:C:81:MET:HB2	1:C:106:ILE:HG13	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:10:VAL:HG21	1:B:163:ILE:HD12	2.03	0.40
2:B:462:FAD:H9	2:B:462:FAD:H1'1	1.93	0.40
1:B:413:PRO:HG2	1:B:416:GLN:HB2	2.04	0.40
1:C:399:ARG:CZ	1:C:456:PRO:HG3	2.52	0.40
1:D:144:ARG:NH1	1:D:158:ARG:HE	2.19	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	453/461 (98%)	432 (95%)	17 (4%)	4 (1%)	14	47
1	B	453/461 (98%)	431 (95%)	17 (4%)	5 (1%)	11	43
1	C	453/461 (98%)	430 (95%)	21 (5%)	2 (0%)	30	62
1	D	453/461 (98%)	431 (95%)	19 (4%)	3 (1%)	18	52
All	All	1812/1844 (98%)	1724 (95%)	74 (4%)	14 (1%)	16	50

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	433	ALA
1	B	376	LYS
1	B	433	ALA
1	D	433	ALA
1	C	377	VAL
1	D	377	VAL
1	A	219	GLY
1	A	376	LYS
1	A	377	VAL
1	B	219	GLY

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Mol	Chain	Res	Type
1	B	375	SER
1	C	219	GLY
1	D	219	GLY
1	B	377	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	367/372 (99%)	343 (94%)	24 (6%)	15	47
1	B	367/372 (99%)	341 (93%)	26 (7%)	13	44
1	C	367/372 (99%)	345 (94%)	22 (6%)	17	50
1	D	367/372 (99%)	340 (93%)	27 (7%)	13	43
All	All	1468/1488 (99%)	1369 (93%)	99 (7%)	15	47

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	20	THR
1	A	43	ARG
1	A	44	ILE
1	A	72	ARG
1	A	80	MET
1	A	113	ARG
1	A	151	SER
1	A	174	GLN
1	A	188	THR
1	A	208	LEU
1	A	215	ILE
1	A	224	ILE
1	A	247	GLN
1	A	297	LEU
1	A	324	MET
1	A	354	GLN

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Mol	Chain	Res	Type
1	A	375	SER
1	A	410	SER
1	A	416	GLN
1	A	427	LEU
1	A	438	LEU
1	A	439	THR
1	A	441	ILE
1	B	20	THR
1	B	44	ILE
1	B	72	ARG
1	B	80	MET
1	B	84	ARG
1	B	106	ILE
1	B	113	ARG
1	B	135	LEU
1	B	151	SER
1	B	187	GLU
1	B	208	LEU
1	B	213	LEU
1	B	215	ILE
1	B	224	ILE
1	B	247	GLN
1	B	299	THR
1	B	324	MET
1	B	371	VAL
1	B	375	SER
1	B	410	SER
1	B	416	GLN
1	B	422	GLN
1	B	426	GLN
1	B	438	LEU
1	B	439	THR
1	B	441	ILE
1	C	9	GLN
1	C	20	THR
1	C	45	GLU
1	C	72	ARG
1	C	80	MET
1	C	113	ARG
1	C	151	SER
1	C	208	LEU
1	C	215	ILE

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Mol	Chain	Res	Type
1	C	224	ILE
1	C	247	GLN
1	C	299	THR
1	C	324	MET
1	C	354	GLN
1	C	375	SER
1	C	384	LEU
1	C	410	SER
1	C	416	GLN
1	C	418	MET
1	C	438	LEU
1	C	439	THR
1	C	441	ILE
1	D	20	THR
1	D	44	ILE
1	D	72	ARG
1	D	113	ARG
1	D	137	SER
1	D	143	GLN
1	D	151	SER
1	D	155	LEU
1	D	187	GLU
1	D	208	LEU
1	D	213	LEU
1	D	215	ILE
1	D	224	ILE
1	D	247	GLN
1	D	297	LEU
1	D	299	THR
1	D	324	MET
1	D	354	GLN
1	D	384	LEU
1	D	391	ILE
1	D	410	SER
1	D	416	GLN
1	D	425	GLU
1	D	427	LEU
1	D	438	LEU
1	D	439	THR
1	D	441	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	32	GLN
1	A	54	ASN
1	A	94	ASN
1	A	115	ASN
1	A	143	GLN
1	A	149	ASN
1	A	197	GLN
1	A	240	HIS
1	A	305	GLN
1	A	339	GLN
1	A	390	GLN
1	A	422	GLN
1	A	459	HIS
1	B	9	GLN
1	B	54	ASN
1	B	94	ASN
1	B	115	ASN
1	B	149	ASN
1	B	197	GLN
1	B	240	HIS
1	B	339	GLN
1	B	390	GLN
1	B	424	GLN
1	B	426	GLN
1	B	448	HIS
1	B	459	HIS
1	C	9	GLN
1	C	32	GLN
1	C	54	ASN
1	C	94	ASN
1	C	115	ASN
1	C	197	GLN
1	C	240	HIS
1	C	255	GLN
1	C	305	GLN
1	C	339	GLN
1	C	424	GLN
1	D	9	GLN
1	D	32	GLN
1	D	94	ASN
1	D	115	ASN
1	D	197	GLN

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Mol	Chain	Res	Type
1	D	240	HIS
1	D	282	GLN
1	D	339	GLN
1	D	390	GLN
1	D	424	GLN
1	D	459	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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	C	462	-	58,58,58	0.34	0	85,89,89	0.61	1 (1%)
3	JHY	C	4000	-	18,18,18	0.43	0	26,26,26	0.66	0
3	JHY	A	4000	-	18,18,18	0.36	0	26,26,26	0.70	0
3	JHY	D	4000	-	18,18,18	0.37	0	26,26,26	0.73	0
3	JHY	B	4000	-	18,18,18	0.33	0	26,26,26	0.69	0
2	FAD	D	462	-	58,58,58	0.35	0	85,89,89	0.64	2 (2%)
2	FAD	B	462	-	58,58,58	0.40	0	85,89,89	0.63	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	A	462	-	58,58,58	0.37	0	85,89,89	0.64	2 (2%)

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	C	462	-	-	5/34/50/50	0/6/6/6
3	JHY	C	4000	-	-	0/5/5/5	0/2/2/2
3	JHY	A	4000	-	-	0/5/5/5	0/2/2/2
3	JHY	D	4000	-	-	0/5/5/5	0/2/2/2
3	JHY	B	4000	-	-	0/5/5/5	0/2/2/2
2	FAD	D	462	-	-	5/34/50/50	0/6/6/6
2	FAD	B	462	-	-	5/34/50/50	0/6/6/6
2	FAD	A	462	-	-	5/34/50/50	0/6/6/6

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	462	FAD	O2P-P-O5'	2.37	118.32	107.57
2	C	462	FAD	O2P-P-O5'	2.29	117.94	107.57
2	B	462	FAD	O2P-P-O5'	2.24	117.74	107.57
2	A	462	FAD	O2P-P-O5'	2.17	117.41	107.57
2	D	462	FAD	O5B-PA-O1A	2.11	117.31	108.94
2	A	462	FAD	O5B-PA-O1A	2.02	116.94	108.94

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	462	FAD	PA-O3P-P-O5'
2	B	462	FAD	PA-O3P-P-O5'
2	C	462	FAD	PA-O3P-P-O5'
2	D	462	FAD	PA-O3P-P-O5'
2	A	462	FAD	P-O3P-PA-O2A
2	B	462	FAD	P-O3P-PA-O2A
2	C	462	FAD	P-O3P-PA-O2A
2	D	462	FAD	P-O3P-PA-O2A

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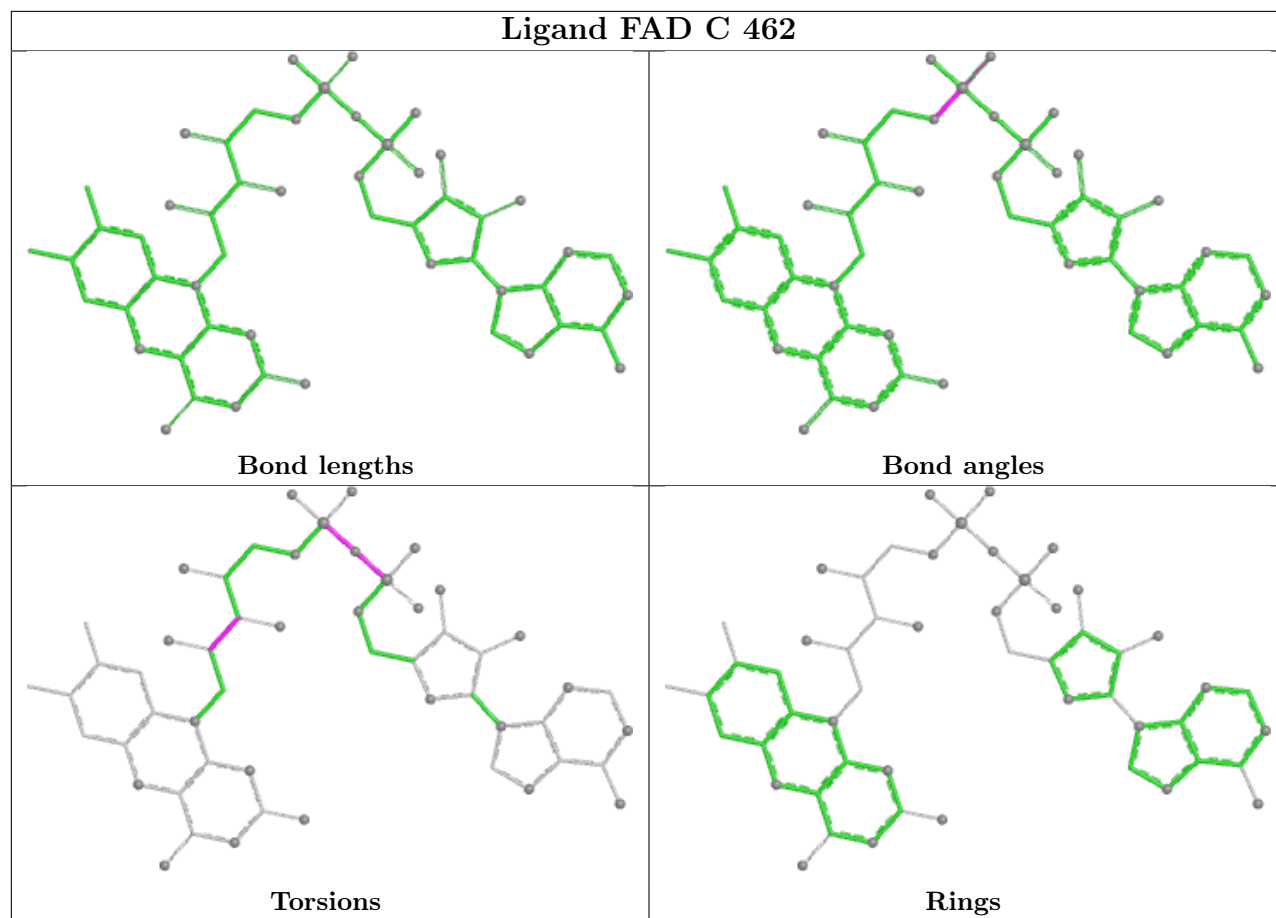
Mol	Chain	Res	Type	Atoms
2	A	462	FAD	O2'-C2'-C3'-O3'
2	B	462	FAD	O2'-C2'-C3'-O3'
2	D	462	FAD	P-O3P-PA-O1A
2	A	462	FAD	C1'-C2'-C3'-O3'
2	B	462	FAD	C1'-C2'-C3'-O3'
2	C	462	FAD	C1'-C2'-C3'-O3'
2	D	462	FAD	C1'-C2'-C3'-O3'
2	A	462	FAD	P-O3P-PA-O1A
2	C	462	FAD	O2'-C2'-C3'-O3'
2	D	462	FAD	O2'-C2'-C3'-O3'
2	B	462	FAD	P-O3P-PA-O1A
2	C	462	FAD	P-O3P-PA-O1A

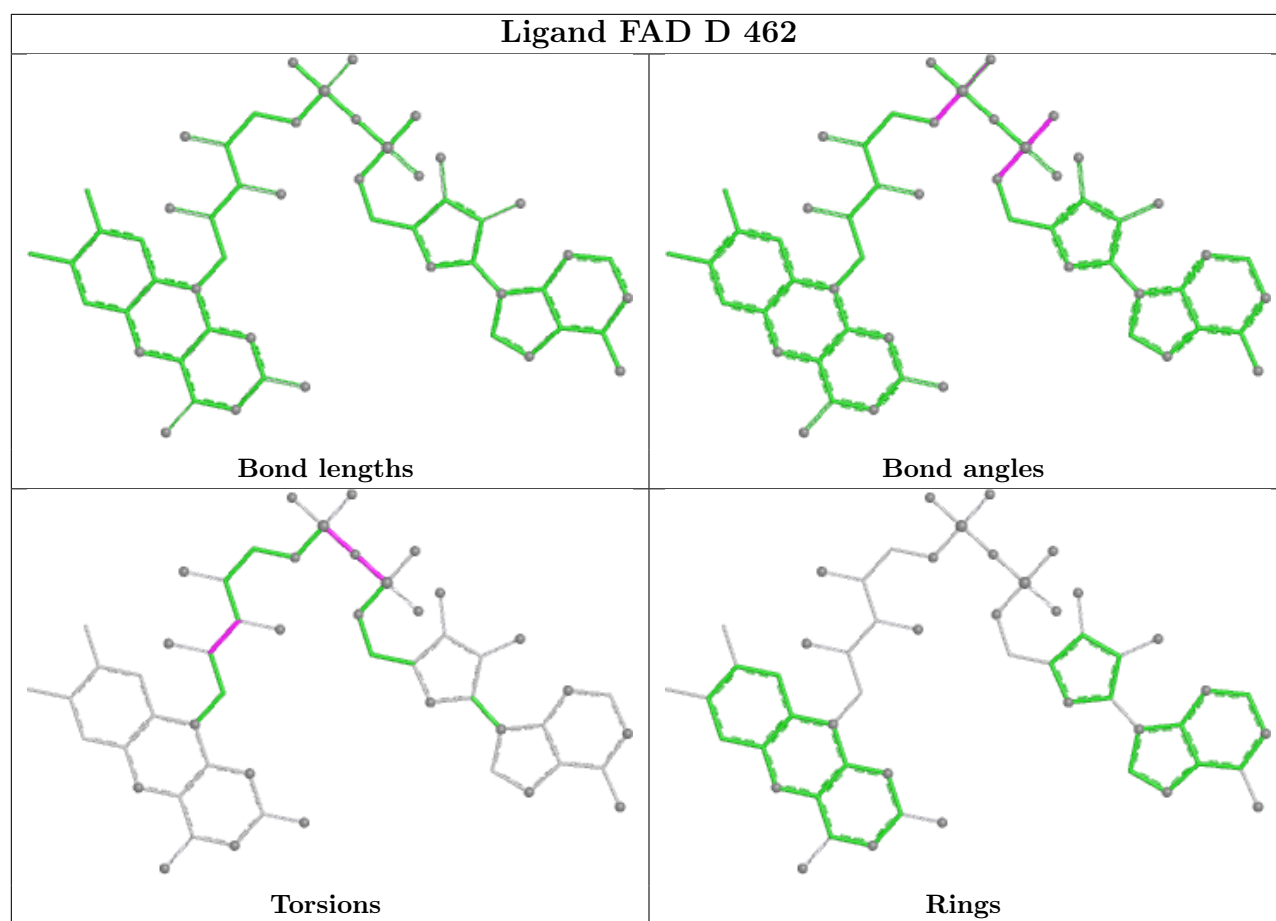
There are no ring outliers.

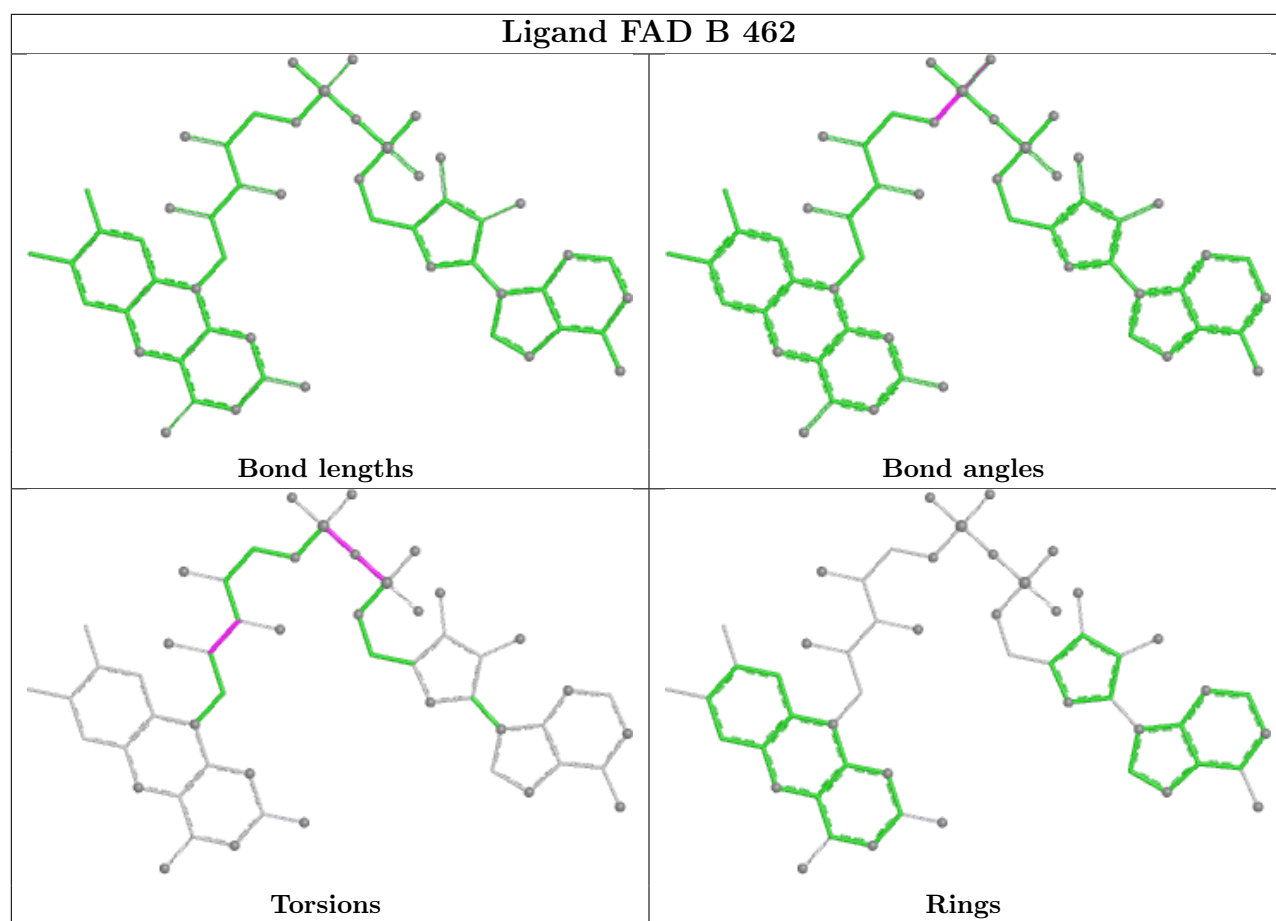
6 monomers are involved in 7 short contacts:

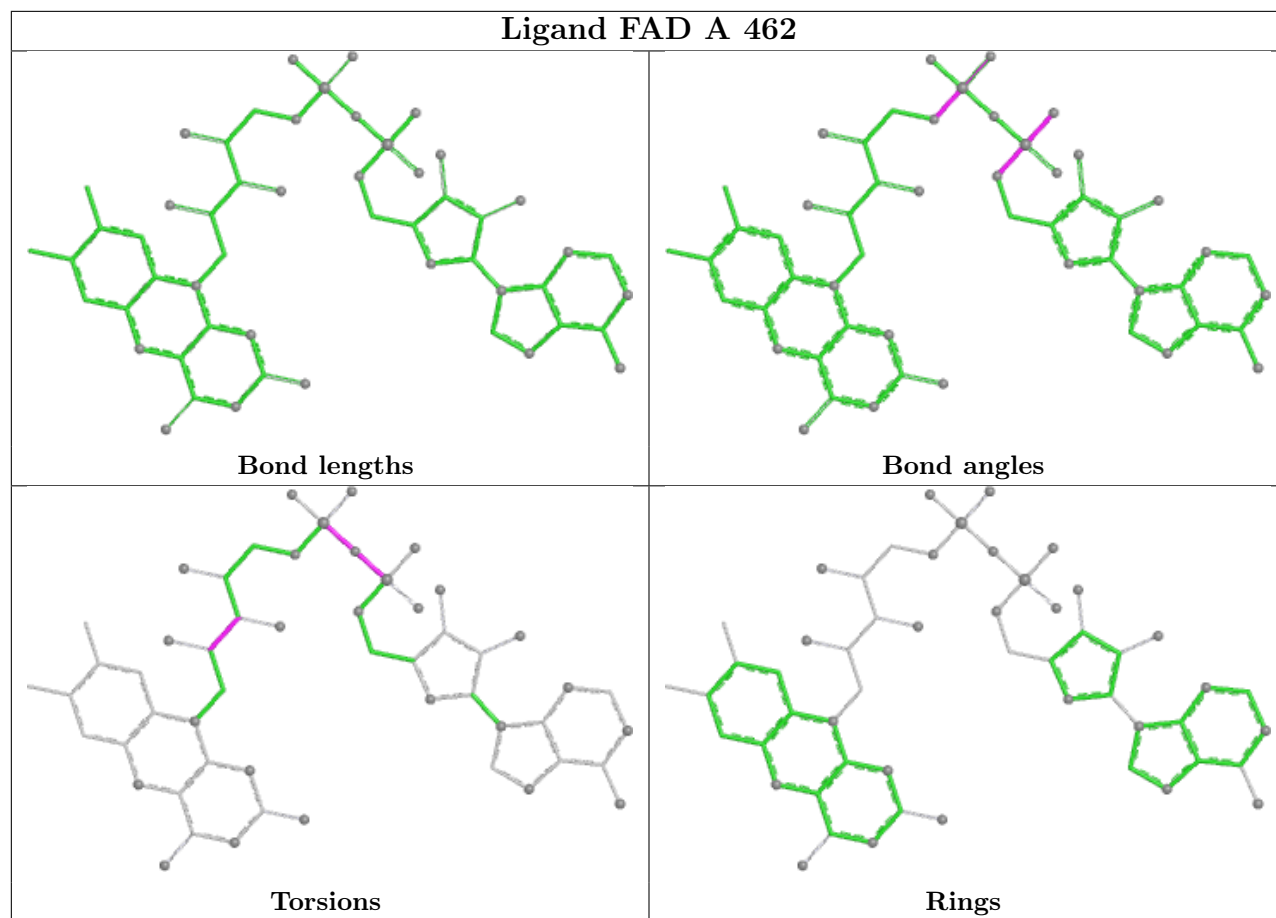
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	462	FAD	1	0
3	C	4000	JHY	1	0
3	B	4000	JHY	1	0
2	D	462	FAD	1	0
2	B	462	FAD	2	0
2	A	462	FAD	1	0

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 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 [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	455/461 (98%)	-0.69	0	100 100	49, 99, 181, 198	0
1	B	455/461 (98%)	-0.64	0	100 100	47, 103, 182, 210	0
1	C	455/461 (98%)	-0.64	1 (0%)	91 85	60, 103, 180, 194	0
1	D	455/461 (98%)	-0.67	0	100 100	56, 110, 194, 221	0
All	All	1820/1844 (98%)	-0.66	1 (0%)	92 89	47, 104, 184, 221	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	238	PHE	2.1

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 oligosaccharides 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	JHY	C	4000	17/17	0.90	0.10	105,111,115,122	0

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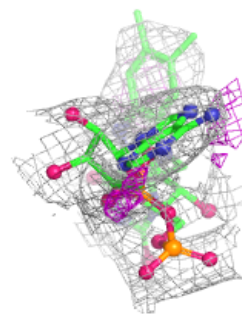
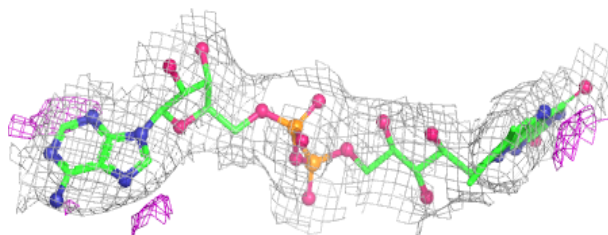
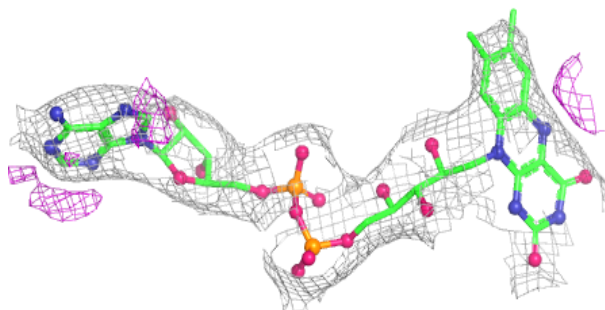
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	JHY	D	4000	17/17	0.93	0.08	103,107,121,131	0
3	JHY	B	4000	17/17	0.94	0.08	104,107,113,123	0
3	JHY	A	4000	17/17	0.95	0.07	99,105,113,122	0
2	FAD	C	462	53/53	0.97	0.06	70,85,104,106	0
2	FAD	A	462	53/53	0.98	0.05	73,86,101,112	0
2	FAD	D	462	53/53	0.98	0.04	76,87,98,102	0
2	FAD	B	462	53/53	0.98	0.05	67,80,98,105	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.

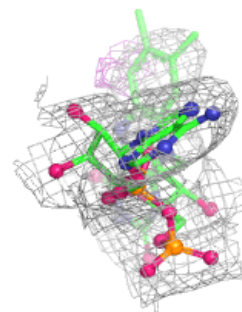
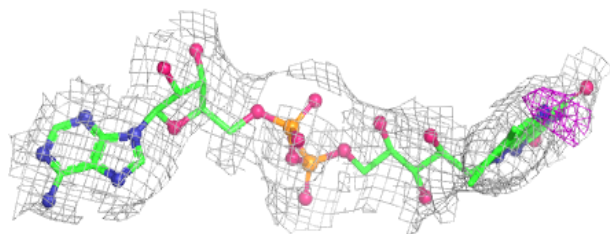
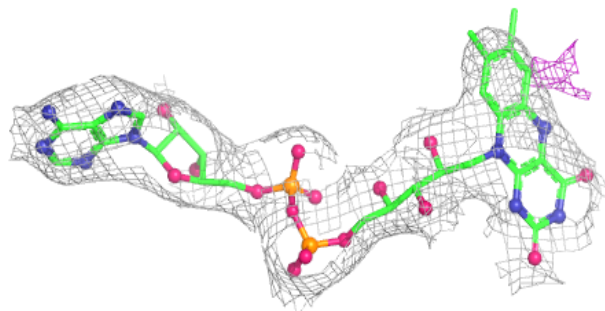
Electron density around FAD C 462:

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

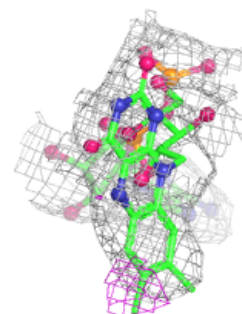
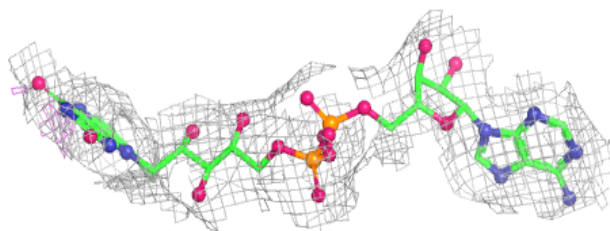
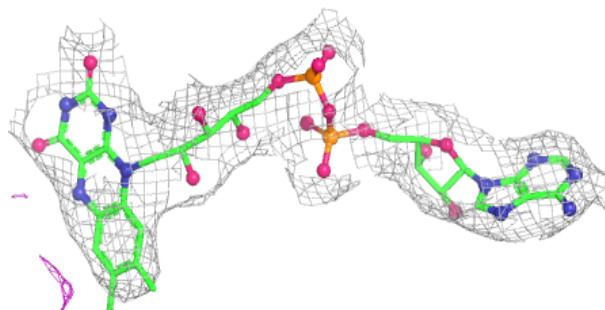


Electron density around FAD A 462:

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

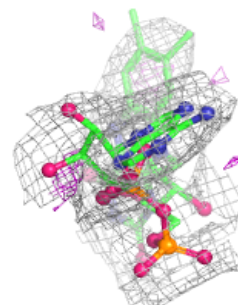
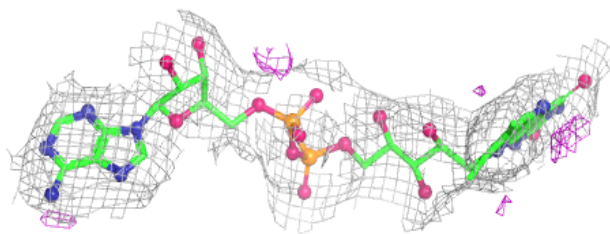
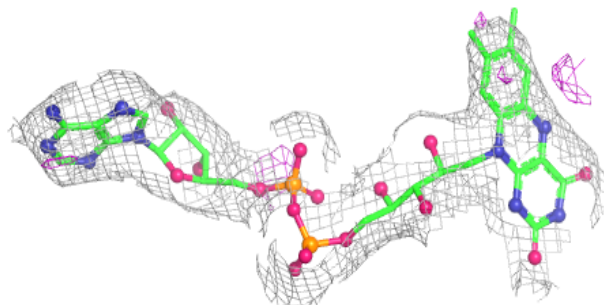
**Electron density around FAD D 462:**

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



Electron density around FAD B 462:

$2mF_o - DF_c$ (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.