Full wwPDB X-ray Structure Validation Report (i)

Apr 28, 2024 – 05:05 pm BST

PDB ID	:	2JBS
Title	:	Structure of the monooxygenase component of p-hydroxyphenylacetate hydroxylase from Acinetobacter baumannii
Authors	:	Alfieri, A.; Mattevi, A.
Deposited on	:	2006-12-11
Resolution	:	2.80 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:

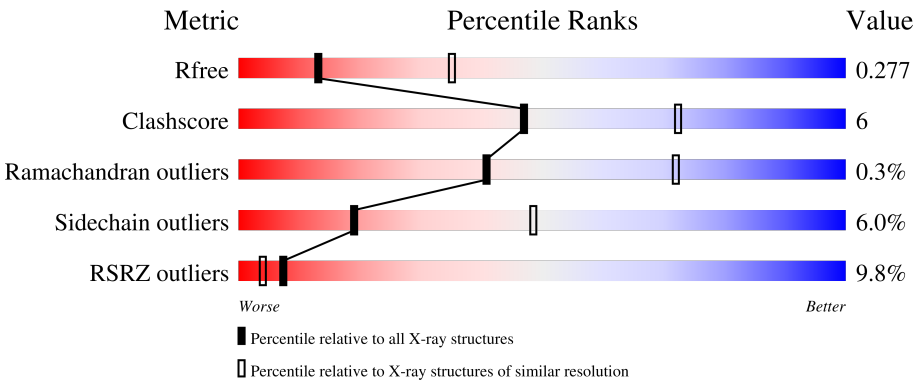
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	422	5% 77% 17% • 5%
1	B	422	23% 80% 12% • 5%
1	C	422	6% 79% 14% • 5%
1	D	422	3% 78% 14% • 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
2	FMN	B	1423	-	-	-	X

2 Entry composition (i)

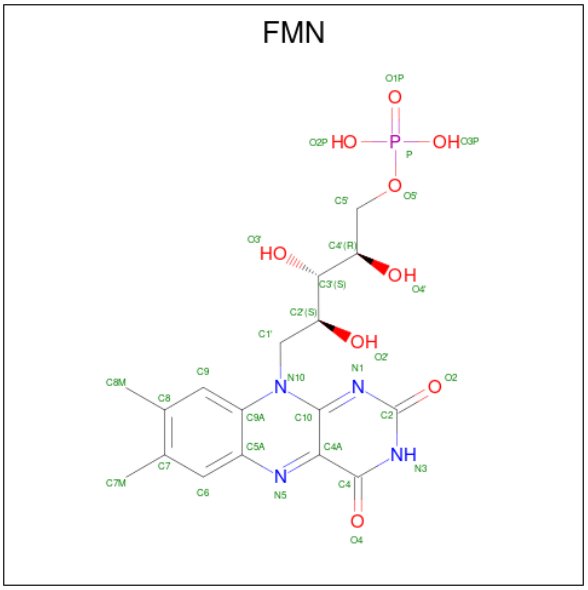
There are 2 unique types of molecules in this entry. The entry contains 12624 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 P-HYDROXYPHENYLACETATE HYDROXYLASE C2\:O XYGENASE COMPONENT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	399	3121	1980	538	582	21	0	0	0
1	B	399	3121	1980	538	582	21	0	0	0
1	C	400	3129	1986	539	583	21	0	0	0
1	D	400	3129	1986	539	583	21	0	0	0

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	31	17	4	9	1	0	0
2	B	1	31	17	4	9	1	0	0

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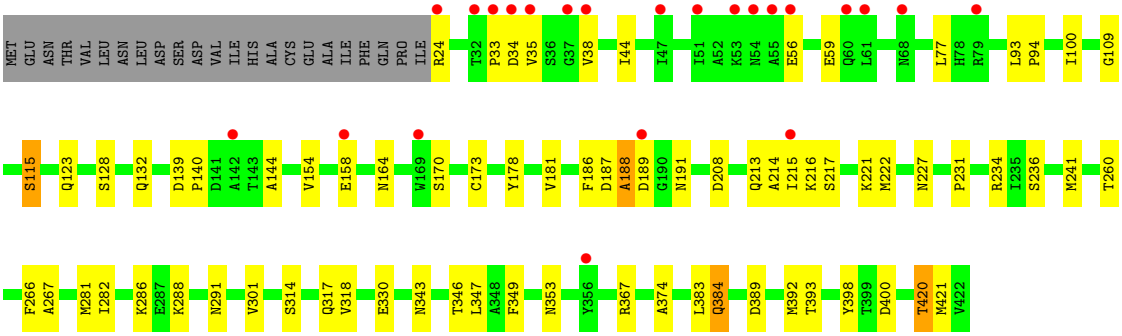
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	C	1	31	17	4	9	1	0	0
			Total	C	N	O	P		
2	D	1	31	17	4	9	1	0	0

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.

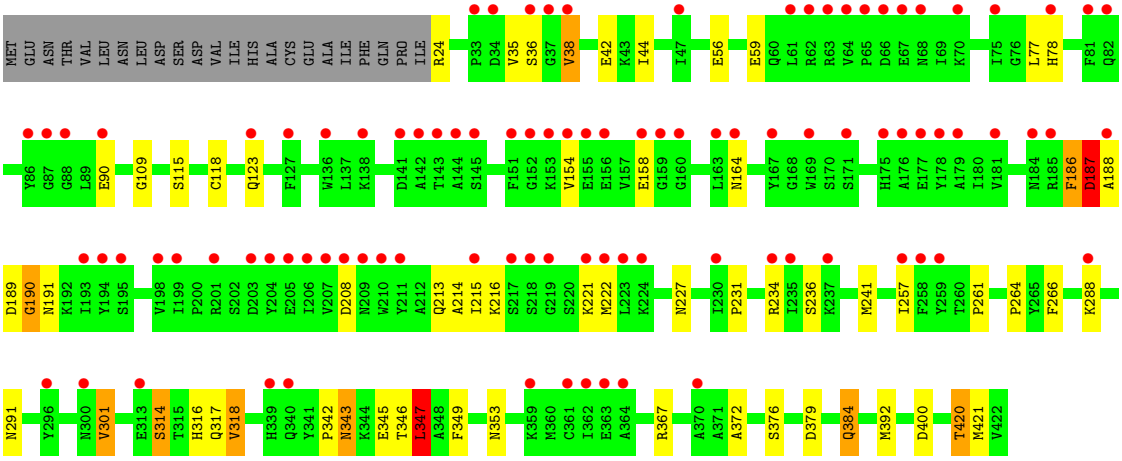
• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\:OXYGENASE COMPONENT

Chain A:



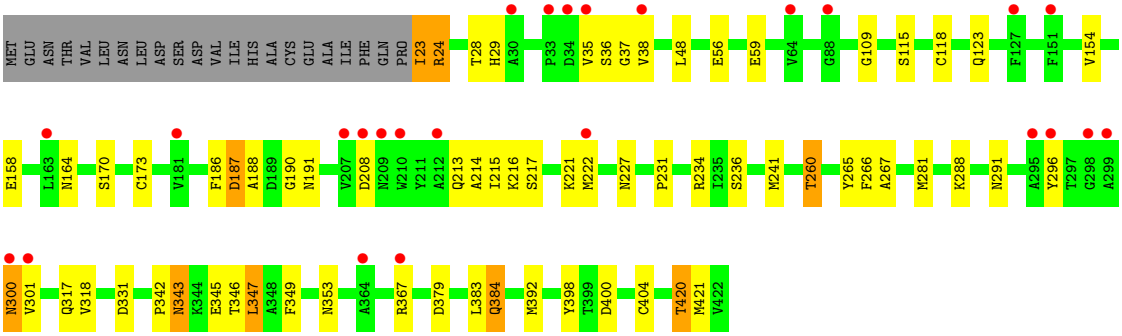
• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\:OXYGENASE COMPONENT

Chain B:



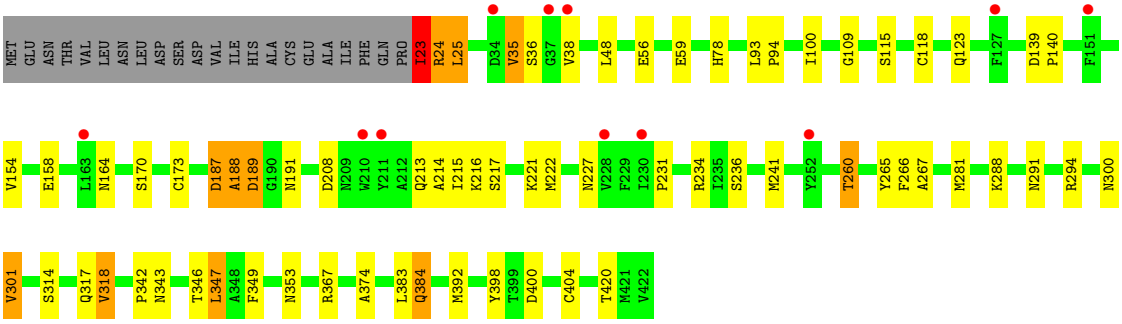
• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\:OXYGENASE COMPONENT

Chain C:



• Molecule 1: P-HYDROXYPHENYLACETATE HYDROXYLASE C2\:OXYGENASE COMPONENT

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	94.32Å 183.43Å 284.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.23 – 2.80 54.03 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (54.23-2.80) 99.4 (54.03-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.91 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.251 , 0.287 0.240 , 0.277	Depositor DCC
R_{free} test set	634 reflections (1.04%)	wwPDB-VP
Wilson B-factor (Å^2)	78.4	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 96.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	12624	wwPDB-VP
Average B, all atoms (Å^2)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: FMN

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	0.74	1/3193 (0.0%)	0.69	1/4318 (0.0%)
1	B	0.71	6/3193 (0.2%)	0.68	2/4318 (0.0%)
1	C	0.77	9/3201 (0.3%)	0.70	3/4329 (0.1%)
1	D	0.90	6/3201 (0.2%)	0.72	3/4329 (0.1%)
All	All	0.79	22/12788 (0.2%)	0.70	9/17294 (0.1%)

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	C	0	1
1	D	0	2
All	All	0	4

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	23	ILE	C-N	15.14	1.68	1.34
1	D	189	ASP	C-O	14.57	1.51	1.23
1	A	188	ALA	CA-CB	14.26	1.82	1.52
1	D	25	LEU	N-CA	12.66	1.71	1.46
1	D	188	ALA	CA-CB	12.44	1.78	1.52
1	C	188	ALA	CA-CB	8.19	1.69	1.52
1	C	37	GLY	C-O	8.15	1.36	1.23
1	C	404	CYS	CB-SG	-7.92	1.68	1.82
1	C	296	TYR	CE2-CZ	7.24	1.48	1.38
1	D	24	ARG	C-N	-7.05	1.17	1.34
1	C	296	TYR	CG-CD1	6.90	1.48	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	189	ASP	C-O	6.87	1.36	1.23
1	C	296	TYR	CG-CD2	6.54	1.47	1.39
1	B	188	ALA	CA-CB	6.52	1.66	1.52
1	C	118	CYS	CB-SG	-6.39	1.71	1.82
1	B	189	ASP	C-N	6.38	1.44	1.33
1	C	296	TYR	CE1-CZ	6.08	1.46	1.38
1	D	404	CYS	CB-SG	-5.85	1.72	1.81
1	B	186	PHE	CG-CD2	5.83	1.47	1.38
1	B	187	ASP	CG-OD2	5.80	1.38	1.25
1	B	186	PHE	CE1-CZ	5.55	1.47	1.37
1	C	300	ASN	CG-OD1	5.43	1.35	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	25	LEU	O-C-N	6.71	133.44	122.70
1	D	23	ILE	O-C-N	-6.29	112.63	122.70
1	A	188	ALA	N-CA-CB	-5.78	102.01	110.10
1	C	24	ARG	N-CA-C	5.76	126.57	111.00
1	D	24	ARG	CB-CA-C	5.61	121.61	110.40
1	C	24	ARG	CB-CA-C	-5.41	99.59	110.40
1	B	187	ASP	CB-CG-OD2	5.22	123.00	118.30
1	C	187	ASP	CB-CA-C	-5.20	100.01	110.40
1	B	347	LEU	CB-CG-CD2	-5.04	102.44	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	35	VAL	Peptide
1	C	36	SER	Peptide
1	D	187	ASP	Peptide
1	D	35	VAL	Peptide

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	3121	0	3060	41	0
1	B	3121	0	3060	40	0
1	C	3129	0	3071	37	0
1	D	3129	0	3071	48	0
2	A	31	0	19	1	0
2	B	31	0	19	0	0
2	C	31	0	19	0	0
2	D	31	0	19	1	0
All	All	12624	0	12338	154	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 6.

All (154) 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:D:188:ALA:CA	1:D:188:ALA:CB	1.78	1.56
1:A:188:ALA:CA	1:A:188:ALA:CB	1.82	1.52
1:D:25:LEU:CA	1:D:25:LEU:N	1.71	1.47
1:D:23:ILE:C	1:D:24:ARG:N	1.68	1.46
1:D:23:ILE:O	1:D:23:ILE:HG13	1.62	1.00
1:C:24:ARG:O	1:C:24:ARG:HG3	1.63	0.96
1:D:23:ILE:O	1:D:23:ILE:CG1	2.17	0.92
1:D:24:ARG:C	1:D:25:LEU:CA	2.45	0.83
1:A:187:ASP:OD2	1:A:191:ASN:HB2	1.84	0.77
1:C:23:ILE:O	1:C:23:ILE:HG13	1.89	0.72
1:C:186:PHE:HB3	1:C:190:GLY:HA2	1.73	0.71
1:B:59:GLU:HG3	1:B:216:LYS:HD2	1.73	0.70
1:D:59:GLU:HG3	1:D:216:LYS:HD2	1.74	0.69
1:A:164:ASN:ND2	1:A:227:ASN:H	1.90	0.69
1:D:164:ASN:ND2	1:D:227:ASN:H	1.94	0.66
1:C:123:GLN:HG2	1:C:241:MET:HE1	1.78	0.66
1:B:213:GLN:HG3	1:C:384:GLN:CG	2.26	0.65
1:A:59:GLU:HG3	1:A:216:LYS:HD2	1.77	0.65
1:A:231:PRO:HG2	1:A:234:ARG:HG3	1.78	0.65
1:C:59:GLU:HG3	1:C:216:LYS:HD2	1.78	0.65
1:C:164:ASN:ND2	1:C:227:ASN:H	1.97	0.62
1:B:123:GLN:HG2	1:B:241:MET:HE1	1.81	0.62
1:A:188:ALA:CB	1:A:188:ALA:N	2.63	0.62
1:B:317:GLN:OE1	1:B:367:ARG:NH2	2.31	0.61
1:B:384:GLN:CG	1:C:213:GLN:HG3	2.31	0.61
1:A:188:ALA:CB	1:A:188:ALA:C	2.68	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:ASN:ND2	1:B:227:ASN:H	1.98	0.61
1:C:317:GLN:OE1	1:C:367:ARG:NH2	2.28	0.61
1:A:164:ASN:HD21	1:A:227:ASN:H	1.48	0.60
2:A:1423:FMN:O1P	1:D:374:ALA:N	2.34	0.60
1:A:123:GLN:HG2	1:A:241:MET:HE1	1.82	0.60
1:A:215:ILE:HG22	1:A:215:ILE:O	2.02	0.60
1:B:213:GLN:HG3	1:C:384:GLN:CD	2.22	0.60
1:D:23:ILE:O	1:D:23:ILE:CD1	2.49	0.60
1:D:188:ALA:CB	1:D:188:ALA:C	2.68	0.60
1:B:213:GLN:HG3	1:C:384:GLN:HG3	1.84	0.59
1:D:188:ALA:CB	1:D:188:ALA:N	2.64	0.59
1:B:186:PHE:HB3	1:B:190:GLY:HA2	1.85	0.59
1:D:123:GLN:HG2	1:D:241:MET:HE1	1.85	0.59
1:D:215:ILE:O	1:D:215:ILE:HG22	2.01	0.58
1:A:400:ASP:OD1	1:A:400:ASP:C	2.42	0.58
1:A:189:ASP:OD2	1:A:191:ASN:ND2	2.37	0.58
1:C:215:ILE:HG22	1:C:215:ILE:O	2.04	0.57
1:D:35:VAL:HG12	1:D:36:SER:N	2.19	0.57
1:B:215:ILE:HG22	1:B:215:ILE:O	2.04	0.57
1:B:384:GLN:HG3	1:C:213:GLN:HG3	1.86	0.57
1:B:384:GLN:CD	1:C:213:GLN:HG3	2.25	0.56
1:A:420:THR:HG22	1:A:421:MET:HE3	1.87	0.56
1:B:231:PRO:HG2	1:B:234:ARG:HG3	1.86	0.56
1:A:317:GLN:OE1	1:A:367:ARG:NH2	2.38	0.56
1:B:109:GLY:HA2	1:B:215:ILE:HG23	1.88	0.56
1:A:24:ARG:O	1:A:24:ARG:HG3	2.05	0.55
1:A:214:ALA:HB3	1:A:392:MET:HB2	1.89	0.55
1:B:214:ALA:HB3	1:B:392:MET:HB2	1.89	0.55
1:B:400:ASP:C	1:B:400:ASP:OD1	2.46	0.55
1:D:187:ASP:OD2	1:D:191:ASN:HB2	2.07	0.55
1:D:231:PRO:HG2	1:D:234:ARG:HG3	1.88	0.55
1:D:25:LEU:N	1:D:25:LEU:CB	2.64	0.54
1:D:24:ARG:O	1:D:25:LEU:CA	2.55	0.53
1:D:400:ASP:OD1	1:D:400:ASP:C	2.47	0.53
1:D:25:LEU:N	1:D:25:LEU:C	2.57	0.53
1:B:314:SER:O	1:B:318:VAL:HG13	2.08	0.52
1:D:317:GLN:OE1	1:D:367:ARG:NH2	2.38	0.52
1:C:400:ASP:OD1	1:C:400:ASP:C	2.46	0.52
1:D:164:ASN:HD21	1:D:227:ASN:H	1.58	0.52
1:C:231:PRO:HG2	1:C:234:ARG:HG3	1.92	0.52
1:C:208:ASP:HA	1:C:221:LYS:HG2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:23:ILE:O	1:D:23:ILE:HD12	2.09	0.51
1:A:109:GLY:HA2	1:A:215:ILE:HG23	1.91	0.51
1:B:123:GLN:HG2	1:B:241:MET:CE	2.41	0.51
1:B:420:THR:HG22	1:B:421:MET:HE3	1.93	0.51
1:C:342:PRO:HG2	1:C:347:LEU:HD22	1.92	0.50
1:A:374:ALA:N	2:D:1423:FMN:O1P	2.44	0.50
1:C:214:ALA:CB	1:C:392:MET:HB2	2.41	0.50
1:B:90:GLU:HG2	1:B:257:ILE:HD11	1.94	0.50
1:C:123:GLN:HG2	1:C:241:MET:CE	2.42	0.50
1:A:330:GLU:HA	1:A:330:GLU:OE1	2.12	0.49
1:B:349:PHE:HD2	1:B:353:ASN:HD21	1.58	0.49
1:B:379:ASP:OD1	1:C:213:GLN:HG2	2.13	0.49
1:B:164:ASN:HD21	1:B:227:ASN:H	1.61	0.48
1:B:261:PRO:O	1:B:264:PRO:HD2	2.13	0.48
1:A:214:ALA:CB	1:A:392:MET:HB2	2.43	0.48
1:C:349:PHE:HD2	1:C:353:ASN:HD21	1.60	0.48
1:A:144:ALA:HA	1:A:178:TYR:O	2.14	0.48
1:B:213:GLN:HG2	1:C:379:ASP:OD1	2.14	0.47
1:D:78:HIS:HA	1:D:118:CYS:SG	2.53	0.47
1:B:187:ASP:N	1:B:187:ASP:OD1	2.47	0.47
1:C:260:THR:HG21	1:C:265:TYR:CE1	2.49	0.47
1:D:35:VAL:HG12	1:D:36:SER:O	2.13	0.47
1:C:109:GLY:HA2	1:C:215:ILE:HG23	1.95	0.47
1:B:44:ILE:HD11	1:B:77:LEU:HD22	1.96	0.46
1:D:342:PRO:HG2	1:D:347:LEU:HD22	1.97	0.46
1:D:109:GLY:HA2	1:D:215:ILE:HG23	1.96	0.46
1:D:214:ALA:HB3	1:D:392:MET:HB2	1.96	0.46
1:B:187:ASP:OD2	1:B:191:ASN:HB2	2.16	0.46
1:C:420:THR:HG22	1:C:421:MET:HE3	1.97	0.46
1:B:208:ASP:HA	1:B:221:LYS:HG2	1.97	0.46
1:B:214:ALA:CB	1:B:392:MET:HB2	2.46	0.46
1:C:214:ALA:HB3	1:C:392:MET:HB2	1.96	0.46
1:A:123:GLN:HG2	1:A:241:MET:CE	2.46	0.46
1:A:353:ASN:HD22	1:B:316:HIS:CD2	2.34	0.46
1:B:24:ARG:O	1:B:24:ARG:HG3	2.15	0.46
1:D:214:ALA:CB	1:D:392:MET:HB2	2.45	0.46
1:D:260:THR:HG21	1:D:265:TYR:CE1	2.51	0.46
1:D:314:SER:O	1:D:318:VAL:HG13	2.16	0.45
1:C:48:LEU:HD12	1:C:48:LEU:HA	1.82	0.45
1:C:164:ASN:HD21	1:C:227:ASN:H	1.64	0.45
1:D:349:PHE:HD2	1:D:353:ASN:HD21	1.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:267:ALA:HB1	1:D:398:TYR:CD2	2.51	0.45
1:C:28:THR:O	1:C:29:HIS:HB3	2.17	0.45
1:B:78:HIS:HA	1:B:118:CYS:SG	2.57	0.45
1:A:128:SER:O	1:A:132:GLN:HG3	2.17	0.44
1:B:301:VAL:O	1:B:301:VAL:HG13	2.17	0.44
1:D:139:ASP:HA	1:D:140:PRO:HD3	1.84	0.44
1:A:389:ASP:O	1:A:393:THR:HG23	2.18	0.44
1:A:281:MET:CE	1:A:383:LEU:HD22	2.47	0.44
1:D:208:ASP:HA	1:D:221:LYS:HG2	2.00	0.43
1:C:281:MET:CE	1:C:383:LEU:HD22	2.48	0.43
1:D:123:GLN:HG2	1:D:241:MET:CE	2.48	0.43
1:C:186:PHE:HA	1:C:191:ASN:O	2.18	0.43
1:A:349:PHE:HD2	1:A:353:ASN:HD21	1.66	0.43
1:B:343:ASN:HD22	1:B:345:GLU:H	1.66	0.43
1:C:109:GLY:HA3	1:C:217:SER:OG	2.18	0.43
1:A:44:ILE:HD11	1:A:77:LEU:HD22	2.00	0.43
1:A:267:ALA:HB1	1:A:398:TYR:CE2	2.54	0.43
1:C:170:SER:HB3	1:C:173:CYS:HB3	2.00	0.43
1:D:48:LEU:HD12	1:D:48:LEU:HA	1.91	0.43
1:B:342:PRO:HG2	1:B:347:LEU:HD22	2.00	0.43
1:A:267:ALA:HB1	1:A:398:TYR:CD2	2.54	0.42
1:B:35:VAL:HG12	1:B:36:SER:O	2.18	0.42
1:A:93:LEU:O	1:A:94:PRO:C	2.57	0.42
1:D:301:VAL:HG13	1:D:301:VAL:O	2.20	0.42
1:D:100:ILE:HG13	1:D:115:SER:HB2	2.01	0.42
1:A:139:ASP:HA	1:A:140:PRO:HD3	1.80	0.42
1:A:186:PHE:HA	1:A:191:ASN:O	2.20	0.42
1:A:100:ILE:HG13	1:A:115:SER:HB2	2.02	0.41
1:D:109:GLY:HA3	1:D:217:SER:OG	2.20	0.41
1:D:170:SER:HB3	1:D:173:CYS:HB3	2.02	0.41
1:D:187:ASP:HB2	1:D:189:ASP:N	2.33	0.41
1:A:384:GLN:CG	1:D:213:GLN:HG3	2.50	0.41
1:C:343:ASN:HD22	1:C:345:GLU:H	1.68	0.41
1:A:208:ASP:HA	1:A:221:LYS:HG2	2.03	0.41
1:C:267:ALA:HB1	1:C:398:TYR:CD2	2.55	0.41
1:B:38:VAL:HG12	1:B:42:GLU:OE2	2.21	0.41
1:D:93:LEU:HB3	1:D:94:PRO:HD3	2.02	0.41
1:A:109:GLY:HA3	1:A:217:SER:OG	2.21	0.41
1:B:186:PHE:HA	1:B:191:ASN:O	2.21	0.41
1:D:281:MET:CE	1:D:383:LEU:HD22	2.51	0.41
1:A:282:ILE:O	1:A:286:LYS:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:372:ALA:HB1	1:B:376:SER:OG	2.21	0.40
1:A:170:SER:HB3	1:A:173:CYS:HB3	2.03	0.40
1:A:213:GLN:HG3	1:D:384:GLN:CG	2.52	0.40
1:A:314:SER:O	1:A:318:VAL:HG13	2.22	0.40
1:C:331:ASP:HB2	1:D:25:LEU:HD12	2.03	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	397/422 (94%)	376 (95%)	19 (5%)	2 (0%)	29	61
1	B	397/422 (94%)	383 (96%)	13 (3%)	1 (0%)	41	72
1	C	398/422 (94%)	373 (94%)	24 (6%)	1 (0%)	41	72
1	D	398/422 (94%)	387 (97%)	11 (3%)	0	100	100
All	All	1590/1688 (94%)	1519 (96%)	67 (4%)	4 (0%)	41	72

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	190	GLY
1	A	34	ASP
1	C	35	VAL
1	A	33	PRO

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	324/345 (94%)	306 (94%)	18 (6%)	21	51
1	B	324/345 (94%)	305 (94%)	19 (6%)	19	49
1	C	325/345 (94%)	304 (94%)	21 (6%)	17	44
1	D	325/345 (94%)	305 (94%)	20 (6%)	18	47
All	All	1298/1380 (94%)	1220 (94%)	78 (6%)	19	48

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	VAL
1	A	56	GLU
1	A	115	SER
1	A	154	VAL
1	A	158	GLU
1	A	181	VAL
1	A	222	MET
1	A	236	SER
1	A	260	THR
1	A	266	PHE
1	A	288	LYS
1	A	291	ASN
1	A	301	VAL
1	A	343	ASN
1	A	346	THR
1	A	347	LEU
1	A	384	GLN
1	A	420	THR
1	B	38	VAL
1	B	56	GLU
1	B	115	SER
1	B	154	VAL
1	B	158	GLU
1	B	187	ASP
1	B	222	MET
1	B	236	SER
1	B	266	PHE
1	B	288	LYS
1	B	291	ASN
1	B	301	VAL

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Mol	Chain	Res	Type
1	B	314	SER
1	B	318	VAL
1	B	343	ASN
1	B	346	THR
1	B	347	LEU
1	B	384	GLN
1	B	420	THR
1	C	23	ILE
1	C	38	VAL
1	C	56	GLU
1	C	115	SER
1	C	154	VAL
1	C	158	GLU
1	C	187	ASP
1	C	222	MET
1	C	236	SER
1	C	260	THR
1	C	266	PHE
1	C	288	LYS
1	C	291	ASN
1	C	300	ASN
1	C	301	VAL
1	C	318	VAL
1	C	343	ASN
1	C	346	THR
1	C	347	LEU
1	C	384	GLN
1	C	420	THR
1	D	23	ILE
1	D	38	VAL
1	D	56	GLU
1	D	154	VAL
1	D	158	GLU
1	D	222	MET
1	D	236	SER
1	D	260	THR
1	D	266	PHE
1	D	288	LYS
1	D	291	ASN
1	D	294	ARG
1	D	300	ASN
1	D	301	VAL

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Mol	Chain	Res	Type
1	D	318	VAL
1	D	343	ASN
1	D	346	THR
1	D	347	LEU
1	D	384	GLN
1	D	420	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	164	ASN
1	A	184	ASN
1	A	191	ASN
1	A	227	ASN
1	A	343	ASN
1	A	353	ASN
1	B	46	GLN
1	B	164	ASN
1	B	184	ASN
1	B	227	ASN
1	B	343	ASN
1	B	353	ASN
1	C	46	GLN
1	C	164	ASN
1	C	184	ASN
1	C	191	ASN
1	C	227	ASN
1	C	353	ASN
1	D	46	GLN
1	D	164	ASN
1	D	184	ASN
1	D	227	ASN
1	D	343	ASN
1	D	353	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	$\#\|Z\| > 2$
2	FMN	B	1423	-	33,33,33	1.10	2 (6%)	48,50,50	1.55	12 (25%)
2	FMN	C	1423	-	33,33,33	1.19	2 (6%)	48,50,50	1.57	12 (25%)
2	FMN	D	1423	-	33,33,33	1.24	2 (6%)	48,50,50	1.67	9 (18%)
2	FMN	A	1423	-	33,33,33	1.26	3 (9%)	48,50,50	1.64	12 (25%)

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	FMN	B	1423	-	-	5/18/18/18	0/3/3/3
2	FMN	C	1423	-	-	5/18/18/18	0/3/3/3
2	FMN	D	1423	-	-	5/18/18/18	0/3/3/3
2	FMN	A	1423	-	-	5/18/18/18	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1423	FMN	C4A-N5	4.69	1.39	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1423	FMN	C4A-N5	4.66	1.39	1.30
2	A	1423	FMN	C4A-N5	4.62	1.39	1.30
2	B	1423	FMN	C4A-N5	4.17	1.38	1.30
2	D	1423	FMN	C10-N1	3.22	1.39	1.33
2	A	1423	FMN	C10-N1	3.05	1.39	1.33
2	C	1423	FMN	C10-N1	2.96	1.39	1.33
2	B	1423	FMN	C10-N1	2.91	1.39	1.33
2	A	1423	FMN	C9-C8	2.16	1.42	1.39

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	1423	FMN	C8M-C8-C9	-4.21	111.71	119.49
2	C	1423	FMN	C8M-C8-C9	-3.75	112.56	119.49
2	A	1423	FMN	P-O5'-C5'	-3.50	108.66	118.30
2	D	1423	FMN	C4A-C10-N10	3.43	121.50	116.48
2	B	1423	FMN	C8M-C8-C9	-3.37	113.26	119.49
2	D	1423	FMN	C4-N3-C2	-3.19	119.75	125.64
2	C	1423	FMN	C4A-C10-N10	3.14	121.07	116.48
2	B	1423	FMN	C4'-C3'-C2'	-3.07	106.99	113.36
2	A	1423	FMN	C8M-C8-C9	-2.94	114.06	119.49
2	B	1423	FMN	C4A-C10-N10	2.93	120.77	116.48
2	A	1423	FMN	C4-N3-C2	-2.90	120.28	125.64
2	D	1423	FMN	C7M-C7-C6	-2.88	114.17	119.49
2	A	1423	FMN	C4A-C10-N10	2.87	120.68	116.48
2	D	1423	FMN	C10-C4A-N5	-2.82	118.86	124.86
2	C	1423	FMN	C10-C4A-N5	-2.81	118.90	124.86
2	C	1423	FMN	C4-N3-C2	-2.79	120.50	125.64
2	A	1423	FMN	C10-C4A-N5	-2.76	119.00	124.86
2	B	1423	FMN	C4-N3-C2	-2.70	120.66	125.64
2	D	1423	FMN	C4A-C4-N3	2.68	119.99	113.19
2	D	1423	FMN	C4-C4A-N5	2.66	122.02	118.23
2	A	1423	FMN	C4A-C4-N3	2.65	119.93	113.19
2	B	1423	FMN	C10-C4A-N5	-2.62	119.31	124.86
2	A	1423	FMN	C4A-C10-N1	-2.61	118.67	124.73
2	A	1423	FMN	C7M-C7-C6	-2.47	114.93	119.49
2	A	1423	FMN	C4'-C3'-C2'	-2.43	108.30	113.36
2	C	1423	FMN	C4A-C4-N3	2.41	119.30	113.19
2	C	1423	FMN	C4-C4A-N5	2.39	121.63	118.23
2	B	1423	FMN	C4A-C4-N3	2.37	119.22	113.19
2	D	1423	FMN	O2-C2-N1	-2.36	117.92	121.83
2	D	1423	FMN	C4A-C10-N1	-2.35	119.28	124.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1423	FMN	C4A-C10-N1	-2.32	119.34	124.73
2	B	1423	FMN	O4-C4-C4A	-2.32	120.45	126.60
2	C	1423	FMN	C5'-C4'-C3'	-2.31	107.74	112.20
2	A	1423	FMN	O4'-C4'-C5'	-2.20	104.97	109.92
2	C	1423	FMN	C7M-C7-C6	-2.19	115.44	119.49
2	B	1423	FMN	C7M-C7-C6	-2.19	115.44	119.49
2	A	1423	FMN	C5A-C9A-N10	2.19	120.21	117.95
2	B	1423	FMN	O2-C2-N1	-2.13	118.31	121.83
2	B	1423	FMN	C4A-C10-N1	-2.09	119.87	124.73
2	C	1423	FMN	C10-N1-C2	2.09	121.09	116.90
2	C	1423	FMN	O4-C4-C4A	-2.06	121.14	126.60
2	B	1423	FMN	C5'-C4'-C3'	-2.03	108.28	112.20
2	C	1423	FMN	O2-C2-N1	-2.03	118.47	121.83
2	B	1423	FMN	O3'-C3'-C2'	2.02	113.69	108.81
2	A	1423	FMN	C8M-C8-C7	-2.01	116.62	120.74

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1423	FMN	O4'-C4'-C5'-O5'
2	A	1423	FMN	C5'-O5'-P-O1P
2	A	1423	FMN	C5'-O5'-P-O2P
2	A	1423	FMN	C5'-O5'-P-O3P
2	B	1423	FMN	O4'-C4'-C5'-O5'
2	B	1423	FMN	C5'-O5'-P-O2P
2	B	1423	FMN	C5'-O5'-P-O3P
2	C	1423	FMN	C3'-C4'-C5'-O5'
2	C	1423	FMN	O4'-C4'-C5'-O5'
2	C	1423	FMN	C5'-O5'-P-O1P
2	C	1423	FMN	C5'-O5'-P-O2P
2	C	1423	FMN	C5'-O5'-P-O3P
2	D	1423	FMN	O4'-C4'-C5'-O5'
2	D	1423	FMN	C5'-O5'-P-O1P
2	D	1423	FMN	C5'-O5'-P-O2P
2	D	1423	FMN	C5'-O5'-P-O3P
2	A	1423	FMN	C3'-C4'-C5'-O5'
2	D	1423	FMN	C3'-C4'-C5'-O5'
2	B	1423	FMN	C5'-O5'-P-O1P
2	B	1423	FMN	C3'-C4'-C5'-O5'

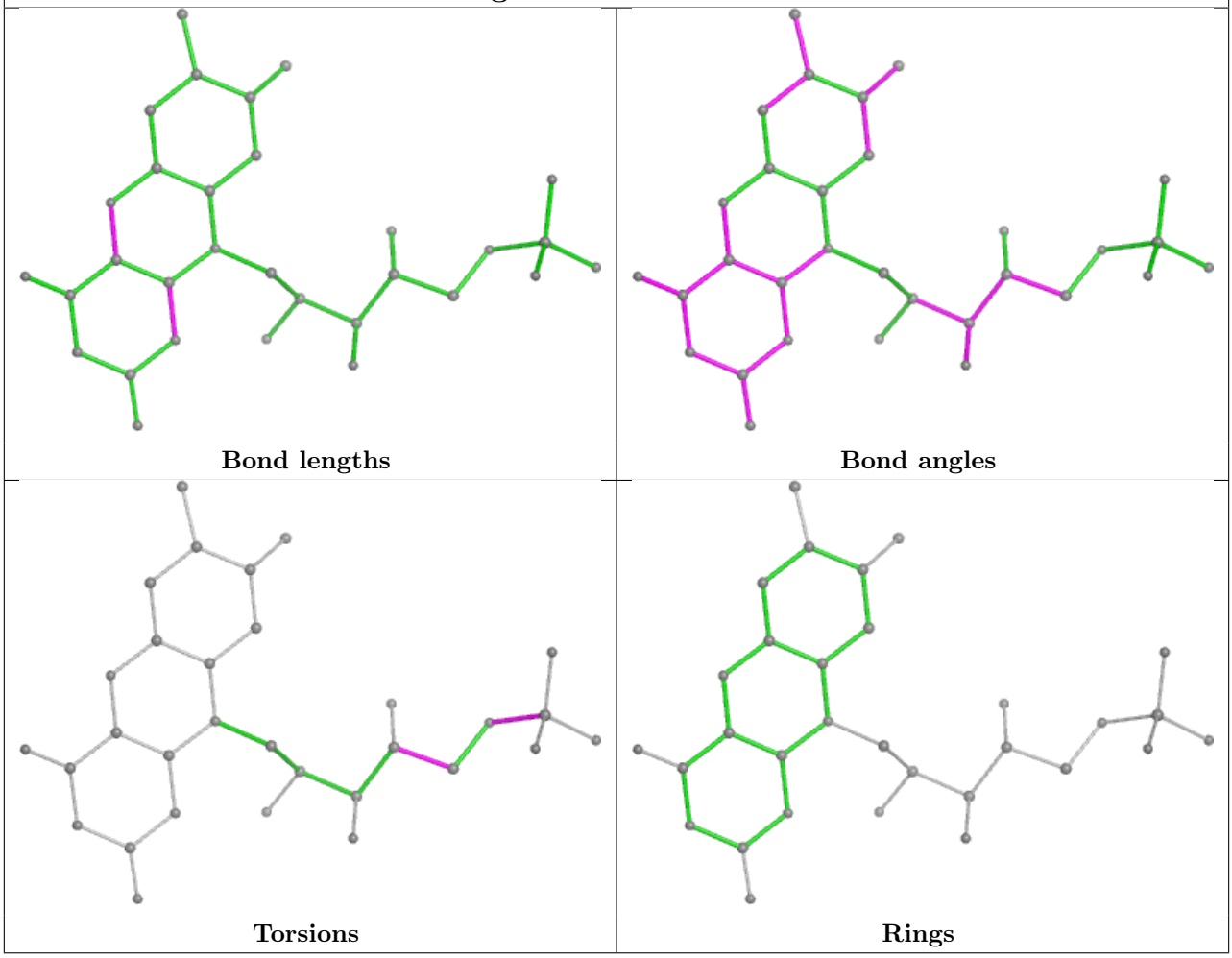
There are no ring outliers.

2 monomers are involved in 2 short contacts:

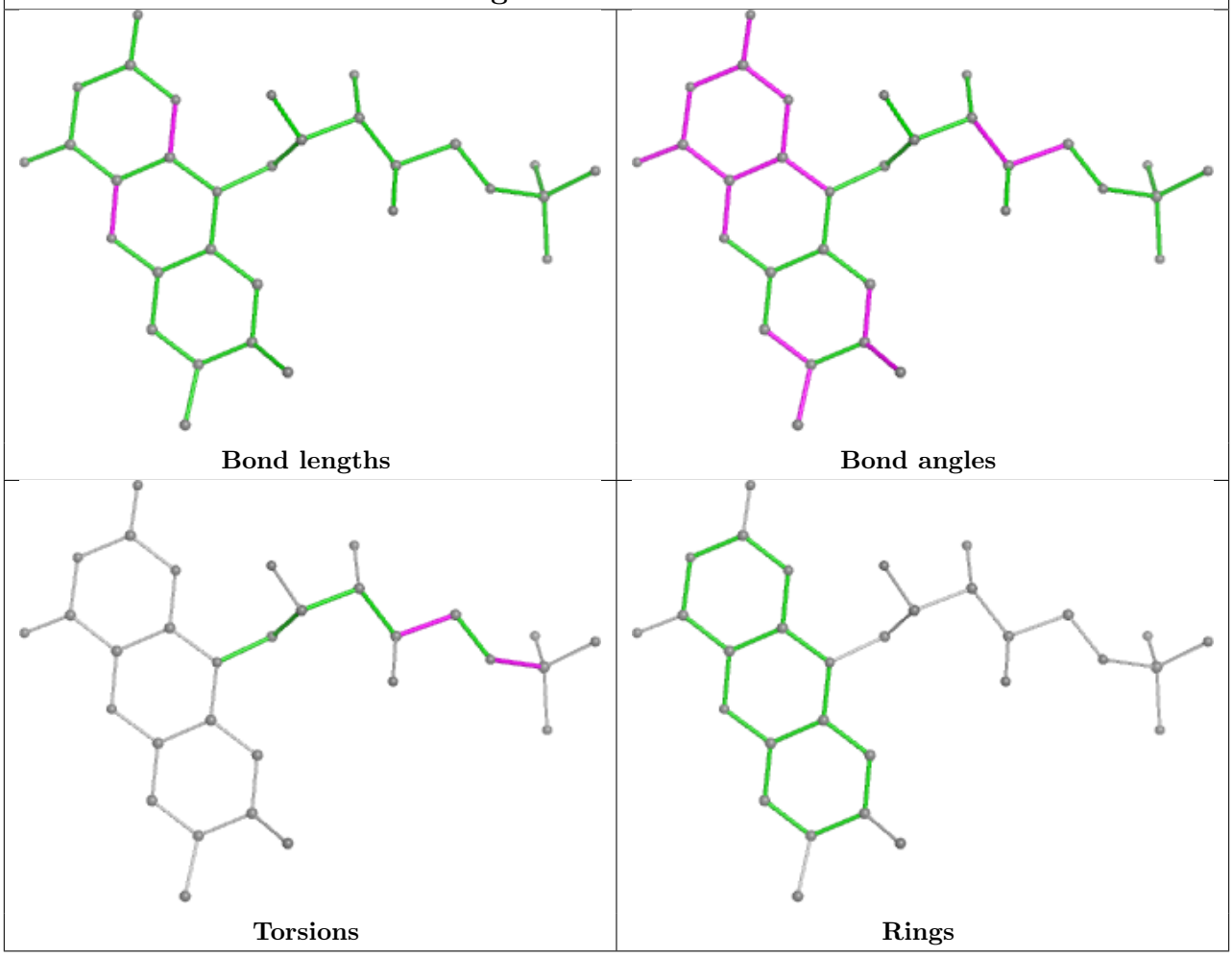
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1423	FMN	1	0
2	A	1423	FMN	1	0

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.

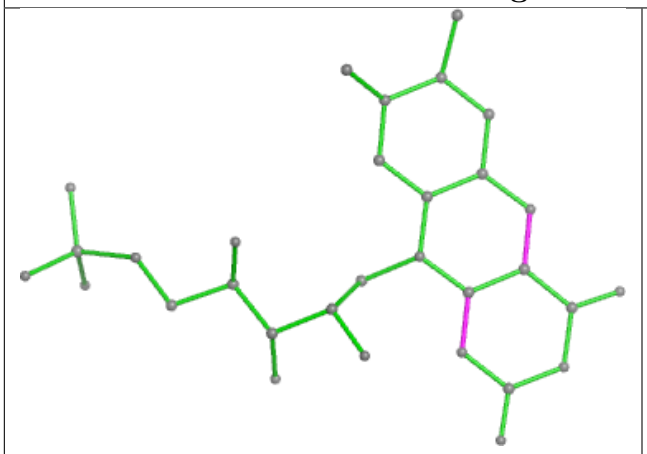
Ligand FMN B 1423



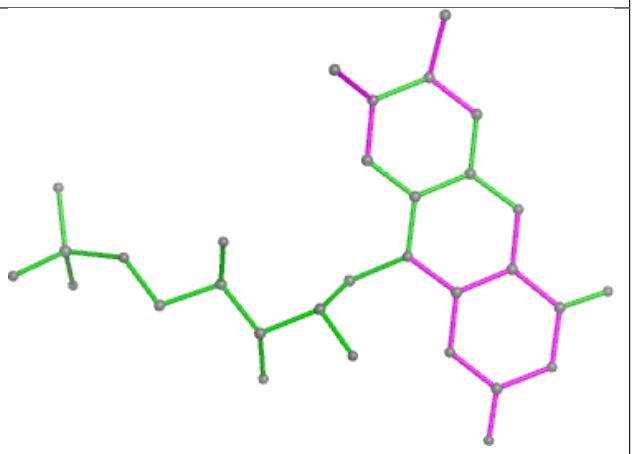
Ligand FMN C 1423



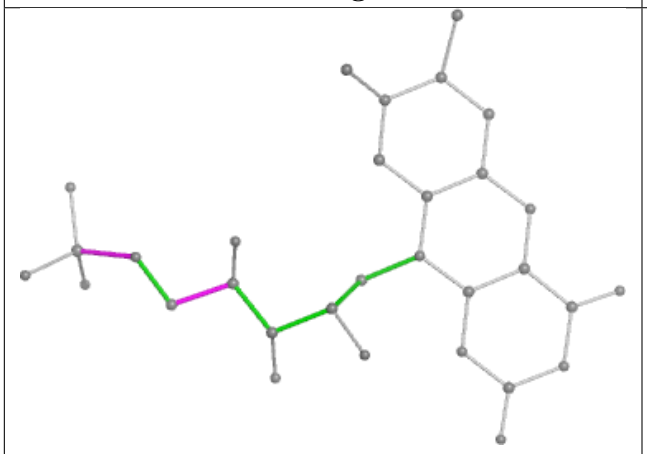
Ligand FMN D 1423



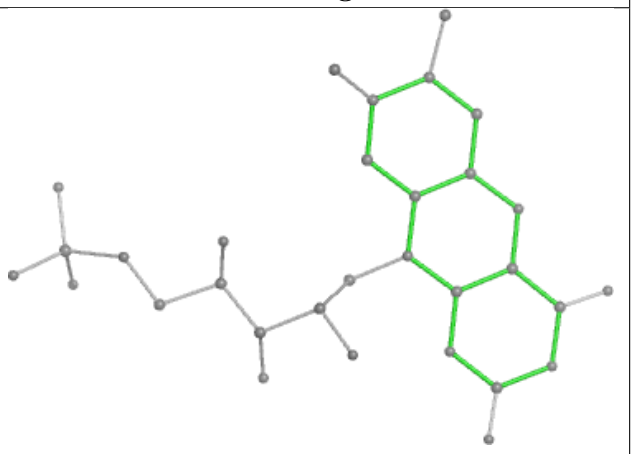
Bond lengths



Bond angles

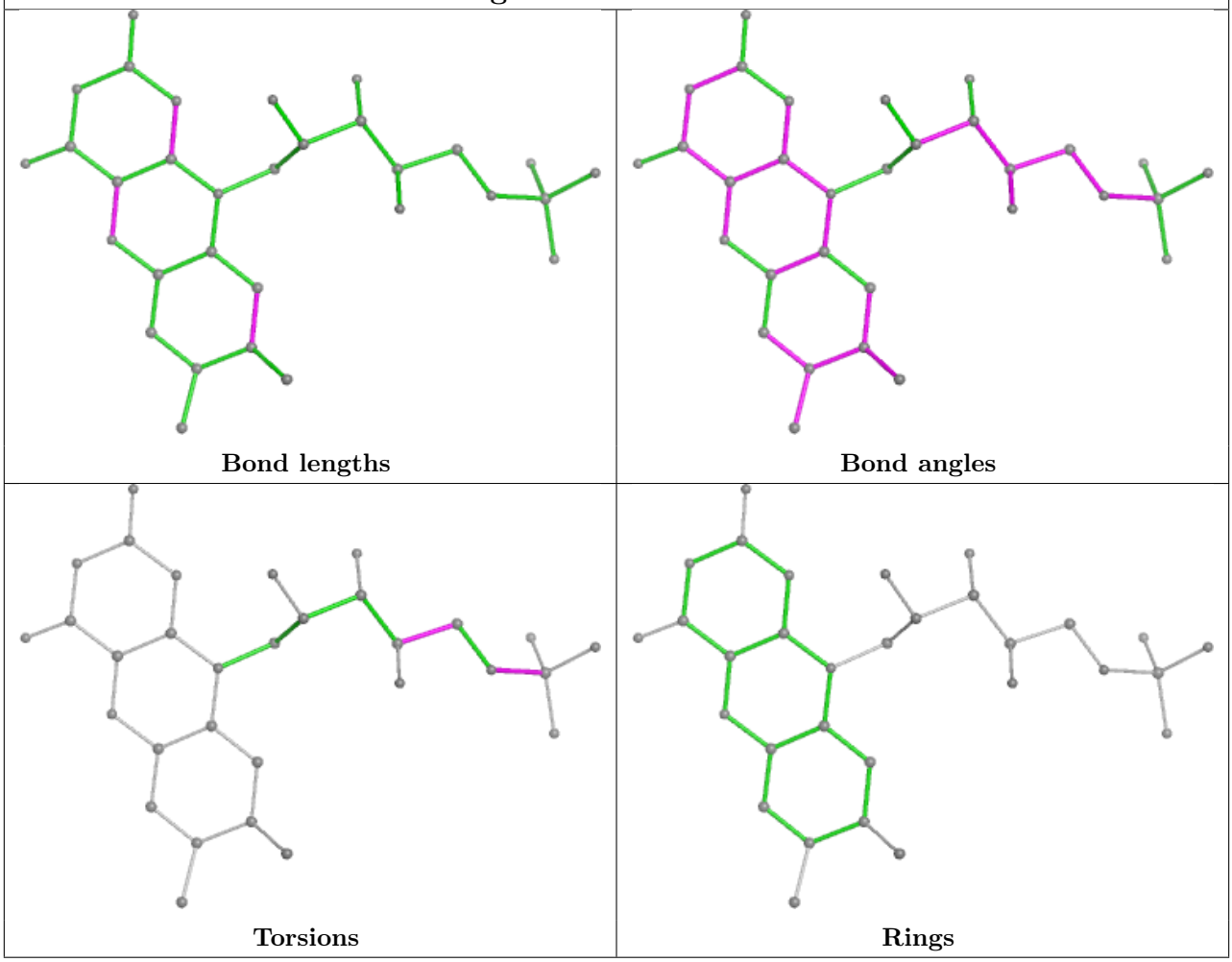


Torsions



Rings

Ligand FMN A 1423



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	23:ILE	C	24:ARG	N	1.68
1	D	24:ARG	C	25:LEU	N	1.17

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	399/422 (94%)	0.37	23 (5%)	23	15	79, 82, 87, 100	0
1	B	399/422 (94%)	1.23	97 (24%)	0	0	79, 82, 87, 98	0
1	C	400/422 (94%)	0.60	25 (6%)	20	12	79, 82, 87, 105	0
1	D	400/422 (94%)	0.28	11 (2%)	53	43	79, 82, 86, 99	0
All	All	1598/1688 (94%)	0.62	156 (9%)	7	4	79, 82, 87, 105	0

All (156) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	223	LEU	9.2
1	B	152	GLY	8.1
1	B	151	PHE	6.9
1	C	295	ALA	6.8
1	C	299	ALA	5.8
1	A	38	VAL	5.7
1	B	206	ILE	5.5
1	B	154	VAL	5.4
1	B	36	SER	5.3
1	B	37	GLY	5.3
1	B	204	TYR	5.2
1	C	301	VAL	5.1
1	B	259	TYR	5.0
1	B	219	GLY	5.0
1	B	195	SER	4.9
1	B	224	LYS	4.7
1	B	88	GLY	4.7
1	B	179	ALA	4.7
1	B	339	HIS	4.6
1	B	163	LEU	4.6
1	B	136	TRP	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	65	PRO	4.4
1	B	33	PRO	4.3
1	B	127	PHE	4.3
1	C	34	ASP	4.2
1	C	298	GLY	4.2
1	B	209	ASN	4.1
1	B	211	TYR	4.1
1	B	143	THR	4.1
1	B	218	SER	4.1
1	C	35	VAL	4.0
1	C	38	VAL	4.0
1	B	175	HIS	4.0
1	B	142	ALA	3.8
1	D	163	LEU	3.8
1	B	82	GLN	3.7
1	B	188	ALA	3.6
1	C	33	PRO	3.6
1	B	208	ASP	3.6
1	B	167	TYR	3.6
1	B	222	MET	3.6
1	B	230	ILE	3.6
1	B	138	LYS	3.6
1	B	199	ILE	3.6
1	B	68	ASN	3.6
1	B	155	GLU	3.6
1	A	33	PRO	3.5
1	B	207	VAL	3.5
1	B	66	ASP	3.5
1	B	178	TYR	3.4
1	B	194	TYR	3.4
1	B	61	LEU	3.3
1	D	228	VAL	3.3
1	B	221	LYS	3.3
1	A	47	ILE	3.3
1	B	123	GLN	3.3
1	B	64	VAL	3.2
1	B	340	GLN	3.2
1	C	207	VAL	3.1
1	C	209	ASN	3.1
1	B	201	ARG	3.1
1	B	70	LYS	3.1
1	D	230	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	38	VAL	3.1
1	D	151	PHE	3.0
1	C	208	ASP	3.0
1	B	176	ALA	3.0
1	B	81	PHE	3.0
1	B	141	ASP	3.0
1	B	235	ILE	3.0
1	D	252	TYR	3.0
1	C	212	ALA	2.9
1	B	363	GLU	2.9
1	C	64	VAL	2.9
1	B	86	TYR	2.9
1	A	68	ASN	2.9
1	B	217	SER	2.9
1	B	205	GLU	2.9
1	B	87	GLY	2.9
1	A	37	GLY	2.9
1	A	60	GLN	2.9
1	C	210	TRP	2.9
1	B	169	TRP	2.8
1	D	210	TRP	2.8
1	B	234	ARG	2.8
1	D	37	GLY	2.8
1	D	34	ASP	2.8
1	B	156	GLU	2.8
1	B	203	ASP	2.7
1	C	88	GLY	2.7
1	B	215	ILE	2.7
1	B	164	ASN	2.7
1	B	63	ARG	2.7
1	B	153	LYS	2.6
1	A	34	ASP	2.6
1	A	35	VAL	2.6
1	B	144	ALA	2.6
1	C	364	ALA	2.6
1	B	75	ILE	2.6
1	A	189	ASP	2.6
1	C	296	TYR	2.6
1	C	127	PHE	2.6
1	B	362	ILE	2.5
1	B	160	GLY	2.5
1	C	163	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	30	ALA	2.5
1	C	181	VAL	2.5
1	A	215	ILE	2.5
1	B	359	LYS	2.5
1	B	171	SER	2.4
1	A	32	THR	2.4
1	D	127	PHE	2.4
1	B	300	ASN	2.4
1	A	54	ASN	2.4
1	B	257	ILE	2.4
1	A	142	ALA	2.4
1	D	211	TYR	2.3
1	A	56	GLU	2.3
1	A	61	LEU	2.3
1	B	62	ARG	2.3
1	A	51	ILE	2.3
1	B	181	VAL	2.3
1	B	159	GLY	2.3
1	B	361	CYS	2.3
1	B	184	ASN	2.3
1	B	237	LYS	2.3
1	B	258	PHE	2.2
1	C	151	PHE	2.2
1	B	177	GLU	2.2
1	B	193	ILE	2.2
1	B	185	ARG	2.2
1	A	158	GLU	2.2
1	B	47	ILE	2.2
1	B	296	TYR	2.2
1	A	55	ALA	2.2
1	A	53	LYS	2.2
1	B	158	GLU	2.1
1	B	210	TRP	2.1
1	C	222	MET	2.1
1	B	34	ASP	2.1
1	B	67	GLU	2.1
1	B	78	HIS	2.1
1	C	300	ASN	2.1
1	B	90	GLU	2.1
1	B	313	GLU	2.1
1	B	288	LYS	2.1
1	A	79	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	38	VAL	2.1
1	A	24	ARG	2.0
1	C	367	ARG	2.0
1	B	364	ALA	2.0
1	B	370	ALA	2.0
1	B	198	VAL	2.0
1	B	145	SER	2.0
1	A	169	TRP	2.0
1	A	356	TYR	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)

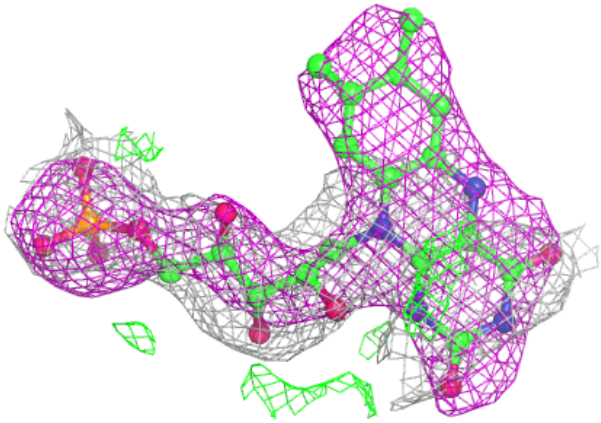
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

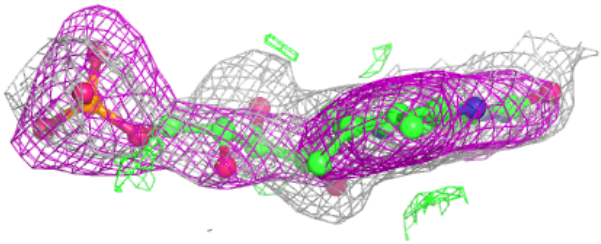
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FMN	B	1423	31/31	0.78	0.57	84,86,91,95	0
2	FMN	D	1423	31/31	0.91	0.29	84,86,92,95	0
2	FMN	C	1423	31/31	0.92	0.26	84,86,91,95	0
2	FMN	A	1423	31/31	0.93	0.17	84,86,91,95	0

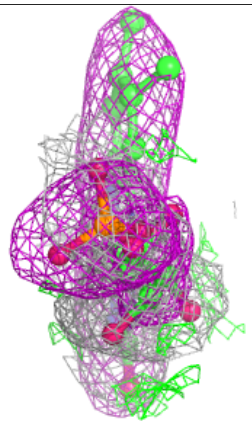
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around FMN B 1423:

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

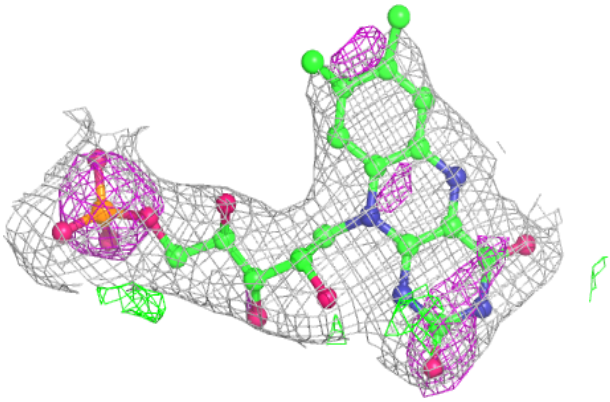


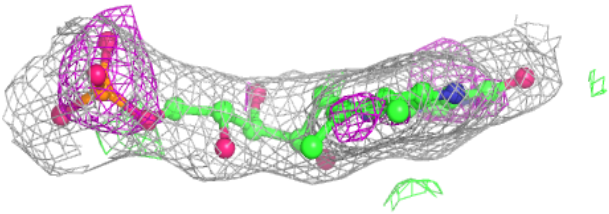


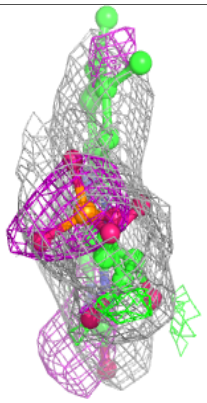


Electron density around FMN D 1423:

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

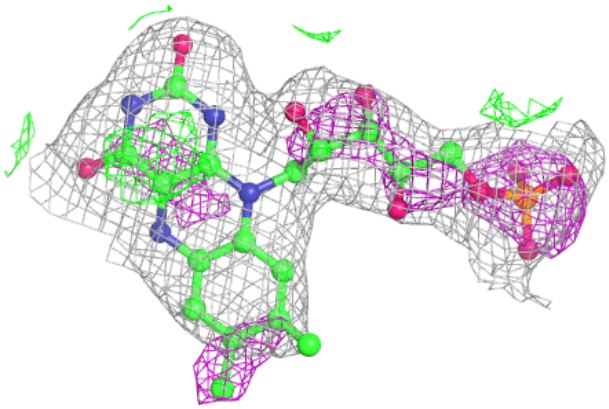


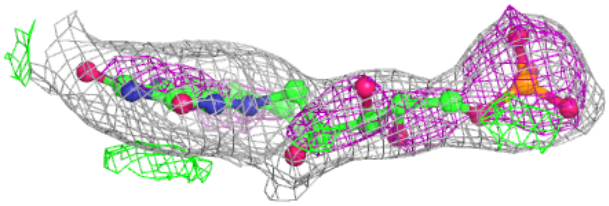


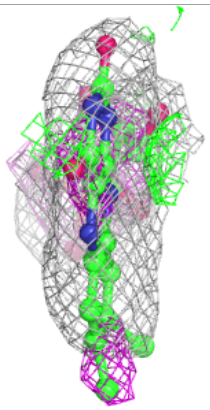


Electron density around FMN C 1423:

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

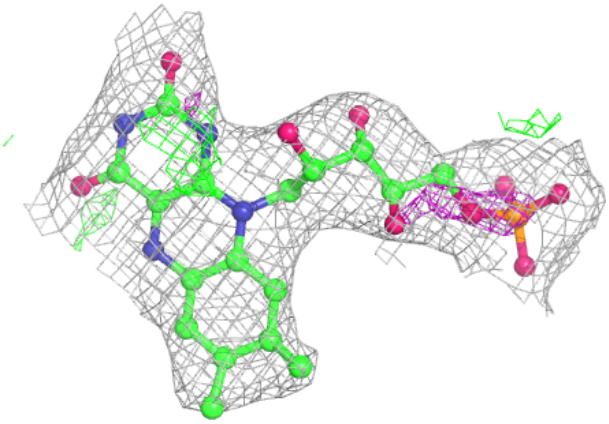


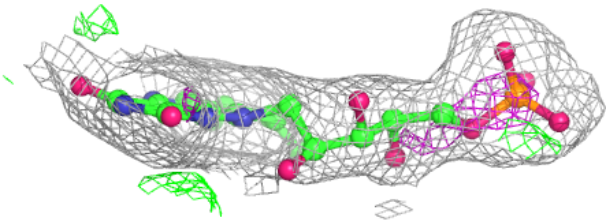


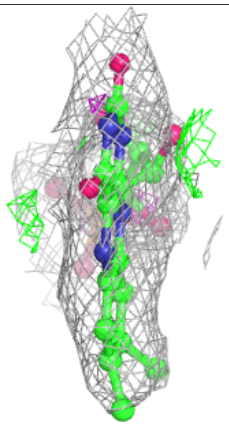


Electron density around FMN A 1423:

$2mF_o$-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.