



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:00 AM UTC

PDB ID : 6PH3 / pdb_00006ph3
Title : LOV-PAS construct from the LOV-HK sensory protein from Brucella abortus (dark-adapted, construct 15-273)
Authors : Rinaldi, J.; Otero, L.H.; Fernandez, I.; Goldbaum, F.A.; Shin, H.; Yang, X.; Klinke, S.
Deposited on : 2019-06-25
Resolution : 2.74 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
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

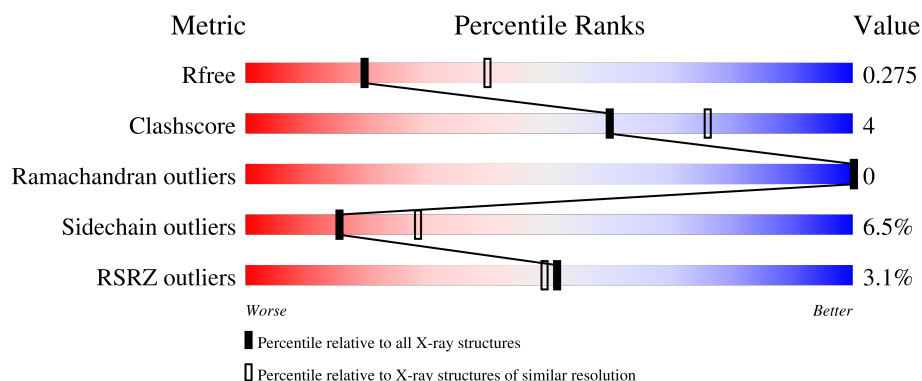
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	266	0% 80% 15% 5%
1	B	266	2% 81% 15% 4%
1	C	266	8% 70% 18% 4% 9%
1	D	266	2% 79% 15% 4%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8249 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 Blue-light-activated histidine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			2037	1293	357	379	8			
1	B	257	Total	C	N	O	S	0	0	0
			2052	1301	360	383	8			
1	C	241	Total	C	N	O	S	0	0	0
			1927	1225	336	358	8			
1	D	256	Total	C	N	O	S	0	0	0
			2047	1299	360	380	8			

There are 28 discrepancies between the modelled and reference sequences:

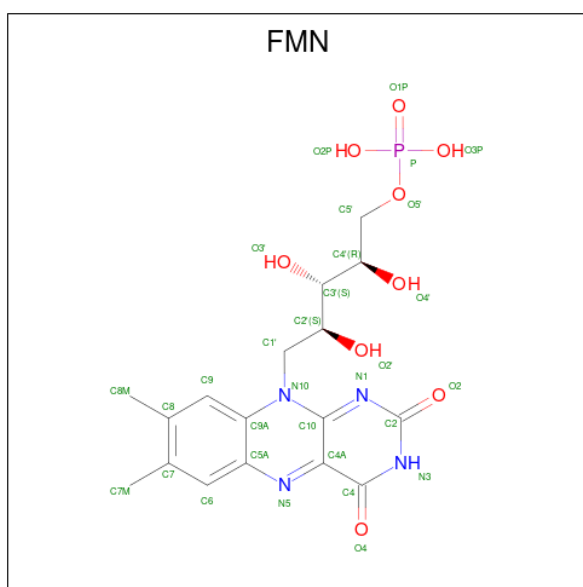
Chain	Residue	Modelled	Actual	Comment	Reference
A	14	MET	-	initiating methionine	UNP Q8YC53
A	274	HIS	-	expression tag	UNP Q8YC53
A	275	HIS	-	expression tag	UNP Q8YC53
A	276	HIS	-	expression tag	UNP Q8YC53
A	277	HIS	-	expression tag	UNP Q8YC53
A	278	HIS	-	expression tag	UNP Q8YC53
A	279	HIS	-	expression tag	UNP Q8YC53
B	14	MET	-	initiating methionine	UNP Q8YC53
B	274	HIS	-	expression tag	UNP Q8YC53
B	275	HIS	-	expression tag	UNP Q8YC53
B	276	HIS	-	expression tag	UNP Q8YC53
B	277	HIS	-	expression tag	UNP Q8YC53
B	278	HIS	-	expression tag	UNP Q8YC53
B	279	HIS	-	expression tag	UNP Q8YC53
C	14	MET	-	initiating methionine	UNP Q8YC53
C	274	HIS	-	expression tag	UNP Q8YC53
C	275	HIS	-	expression tag	UNP Q8YC53
C	276	HIS	-	expression tag	UNP Q8YC53
C	277	HIS	-	expression tag	UNP Q8YC53
C	278	HIS	-	expression tag	UNP Q8YC53
C	279	HIS	-	expression tag	UNP Q8YC53

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Chain	Residue	Modelled	Actual	Comment	Reference
D	14	MET	-	initiating methionine	UNP Q8YC53
D	274	HIS	-	expression tag	UNP Q8YC53
D	275	HIS	-	expression tag	UNP Q8YC53
D	276	HIS	-	expression tag	UNP Q8YC53
D	277	HIS	-	expression tag	UNP Q8YC53
D	278	HIS	-	expression tag	UNP Q8YC53
D	279	HIS	-	expression tag	UNP Q8YC53

- Molecule 2 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: $C_{17}H_{21}N_4O_9P$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	20	Total	O	0	0
			20	20		

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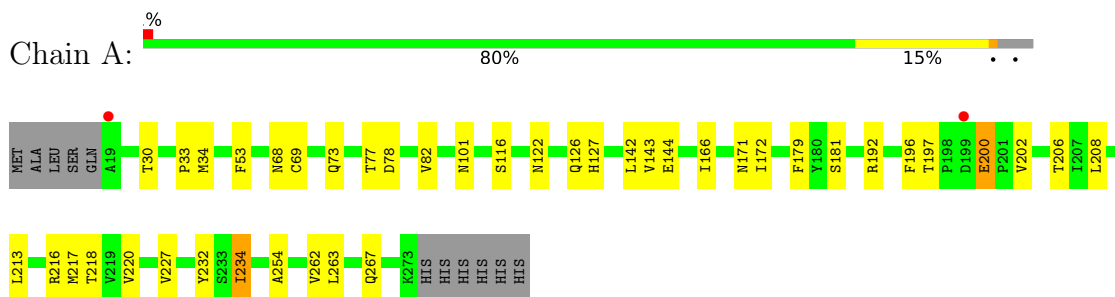
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	16	Total	O	0	0
			16	16		
3	C	4	Total	O	0	0
			4	4		
3	D	22	Total	O	0	0
			22	22		

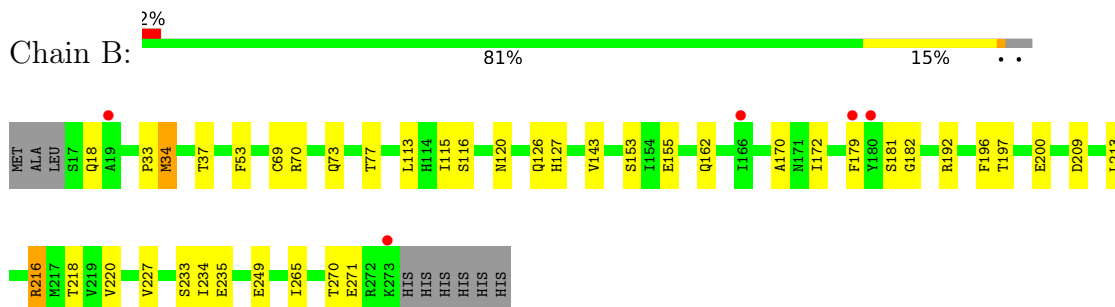
3 Residue-property plots

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.

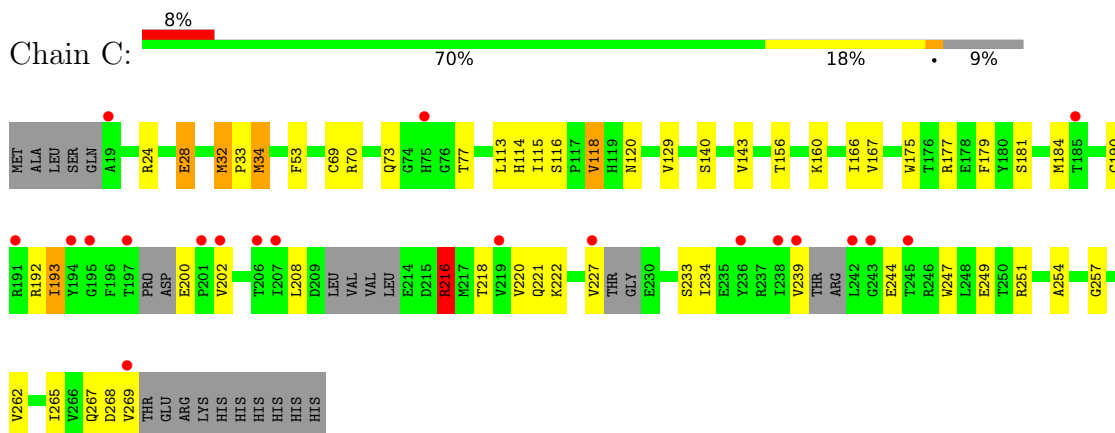
- Molecule 1: Blue-light-activated histidine kinase



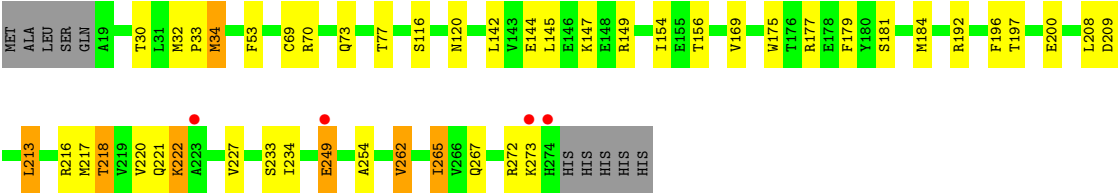
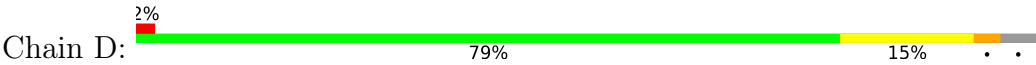
- Molecule 1: Blue-light-activated histidine kinase



- Molecule 1: Blue-light-activated histidine kinase



- Molecule 1: Blue-light-activated histidine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	108.42Å 56.93Å 114.60Å 90.00° 103.36° 90.00°	Depositor
Resolution (Å)	47.70 – 2.74 47.70 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.70-2.74) 99.2 (47.70-2.74)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 2.73Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.224 , 0.269 0.235 , 0.275	Depositor DCC
R_{free} test set	1801 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.663	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8249	wwPDB-VP
Average B, all atoms (Å ²)	74.0	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 23.45 % 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 4.6494e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: 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	0.85	0/2081	1.40	10/2825 (0.4%)
1	B	0.91	2/2096 (0.1%)	1.42	15/2845 (0.5%)
1	C	0.88	2/1966 (0.1%)	1.43	12/2662 (0.5%)
1	D	0.89	1/2092 (0.0%)	1.40	11/2840 (0.4%)
All	All	0.88	5/8235 (0.1%)	1.41	48/11172 (0.4%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	34	MET	SD-CE	-7.74	1.60	1.79
1	D	34	MET	SD-CE	-7.50	1.60	1.79
1	B	34	MET	SD-CE	-6.13	1.64	1.79
1	C	32	MET	SD-CE	-5.39	1.66	1.79
1	B	37	THR	CA-C	5.19	1.59	1.52

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	270	THR	CA-C-N	7.04	131.38	120.82
1	B	270	THR	C-N-CA	7.04	131.38	120.82
1	B	170	ALA	CA-C-N	-7.04	109.93	122.32
1	B	170	ALA	C-N-CA	-7.04	109.93	122.32
1	C	69	CYS	CA-C-N	6.80	130.31	120.38
1	C	69	CYS	C-N-CA	6.80	130.31	120.38
1	B	196	PHE	CA-C-N	6.68	129.98	120.49
1	B	196	PHE	C-N-CA	6.68	129.98	120.49
1	A	122	ASN	CA-CB-CG	6.66	119.26	112.60
1	B	69	CYS	CA-C-N	6.56	129.96	120.38
1	B	69	CYS	C-N-CA	6.56	129.96	120.38
1	C	216	ARG	CA-C-N	6.50	129.23	120.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	216	ARG	C-N-CA	6.50	129.23	120.65
1	B	120	ASN	CA-CB-CG	6.09	118.69	112.60
1	C	193	ILE	N-CA-C	-6.07	104.82	110.53
1	A	101	ASN	CA-CB-CG	6.04	118.64	112.60
1	D	192	ARG	CA-C-N	5.98	128.09	120.56
1	D	192	ARG	C-N-CA	5.98	128.09	120.56
1	B	170	ALA	CA-C-O	-5.88	113.99	120.69
1	A	82	VAL	CA-C-N	5.82	128.07	120.28
1	A	82	VAL	C-N-CA	5.82	128.07	120.28
1	A	196	PHE	CA-C-N	5.77	128.69	120.49
1	A	196	PHE	C-N-CA	5.77	128.69	120.49
1	A	68	ASN	N-CA-C	-5.75	102.40	110.50
1	C	268	ASP	CA-C-N	5.72	132.00	121.70
1	C	268	ASP	C-N-CA	5.72	132.00	121.70
1	C	120	ASN	CA-CB-CG	5.70	118.30	112.60
1	D	196	PHE	CA-C-N	5.68	128.56	120.49
1	D	196	PHE	C-N-CA	5.68	128.56	120.49
1	C	118	VAL	N-CA-CB	5.66	118.99	111.64
1	C	184	MET	N-CA-C	5.59	117.11	109.18
1	D	69	CYS	CA-C-N	5.58	128.53	120.38
1	D	69	CYS	C-N-CA	5.58	128.53	120.38
1	D	272	ARG	N-CA-C	-5.43	100.76	109.39
1	D	120	ASN	CA-CB-CG	5.23	117.83	112.60
1	A	69	CYS	CA-C-N	5.23	128.01	120.38
1	A	69	CYS	C-N-CA	5.23	128.01	120.38
1	B	271	GLU	CA-C-N	5.19	127.19	120.44
1	B	271	GLU	C-N-CA	5.19	127.19	120.44
1	C	257	GLY	CA-C-N	5.14	127.68	120.28
1	C	257	GLY	C-N-CA	5.14	127.68	120.28
1	D	265	ILE	N-CA-CB	5.14	118.32	111.64
1	B	170	ALA	N-CA-C	-5.12	101.61	109.76
1	D	154	ILE	CA-C-N	5.09	127.06	120.44
1	D	154	ILE	C-N-CA	5.09	127.06	120.44
1	A	78	ASP	CA-CB-CG	5.05	117.65	112.60
1	B	182	GLY	CA-C-N	5.05	128.49	121.02
1	B	182	GLY	C-N-CA	5.05	128.49	121.02

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	2037	0	2033	19	0
1	B	2052	0	2046	15	0
1	C	1927	0	1909	24	0
1	D	2047	0	2040	23	0
2	A	31	0	19	0	0
2	B	31	0	19	1	0
2	C	31	0	19	1	0
2	D	31	0	19	1	0
3	A	20	0	0	0	0
3	B	16	0	0	0	0
3	C	4	0	0	0	0
3	D	22	0	0	0	0
All	All	8249	0	8104	70	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 4.

All (70) 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:C:156:THR:HG23	1:D:156:THR:HG23	1.64	0.78
1:B:197:THR:HB	1:B:200:GLU:HB3	1.67	0.76
1:C:34:MET:HE1	1:D:34:MET:HE1	1.71	0.72
1:A:232:TYR:HE2	1:A:234:ILE:HD11	1.53	0.72
1:D:197:THR:HB	1:D:200:GLU:HB2	1.70	0.71
1:C:70:ARG:HD3	2:C:300:FMN:O3P	1.94	0.67
1:A:34:MET:HE1	1:B:34:MET:HE1	1.78	0.66
1:C:24:ARG:HH12	1:C:28:GLU:HG2	1.60	0.64
1:C:156:THR:CG2	1:D:156:THR:HG23	2.29	0.63
1:D:254:ALA:HA	1:D:262:VAL:HA	1.81	0.63
1:B:70:ARG:HD3	2:B:300:FMN:O3P	1.97	0.63
1:A:202:VAL:HG13	1:A:206:THR:OG1	2.00	0.62
1:B:113:LEU:HD21	1:B:115:ILE:HD11	1.83	0.60
1:C:113:LEU:HD21	1:C:115:ILE:HD11	1.83	0.60
1:C:177:ARG:HB3	1:C:262:VAL:CG2	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:216:ARG:O	1:D:220:VAL:HG23	2.03	0.58
1:B:235:GLU:HG3	1:B:249:GLU:HG3	1.89	0.55
1:C:32:MET:HB2	1:D:32:MET:HE2	1.89	0.55
1:A:232:TYR:CE2	1:A:234:ILE:HD11	2.39	0.54
1:A:216:ARG:O	1:A:220:VAL:HG23	2.07	0.54
1:D:177:ARG:HB3	1:D:262:VAL:HG22	1.89	0.54
1:C:239:VAL:HA	1:C:244:GLU:O	2.08	0.53
1:A:254:ALA:HA	1:A:262:VAL:HA	1.90	0.53
1:C:34:MET:HE2	1:D:30:THR:OG1	2.08	0.53
1:A:197:THR:HB	1:A:200:GLU:HB3	1.91	0.53
1:B:216:ARG:O	1:B:220:VAL:HG23	2.09	0.53
1:C:177:ARG:HB3	1:C:262:VAL:HG23	1.90	0.53
1:B:213:LEU:HA	1:B:216:ARG:HG3	1.90	0.52
1:A:208:LEU:HD22	1:A:216:ARG:HD3	1.92	0.52
1:C:216:ARG:O	1:C:220:VAL:HG23	2.10	0.51
1:D:213:LEU:HA	1:D:216:ARG:HG3	1.92	0.51
1:C:73:GLN:HG2	1:C:77:THR:HG21	1.93	0.51
1:A:33:PRO:HB3	1:A:53:PHE:HA	1.93	0.51
1:B:126:GLN:HE21	1:B:127:HIS:HD2	1.57	0.51
1:C:254:ALA:HB2	1:C:262:VAL:HG12	1.92	0.51
1:C:221:GLN:HE21	1:C:222:LYS:HD2	1.77	0.50
1:A:126:GLN:HE21	1:A:127:HIS:HD2	1.59	0.49
1:B:33:PRO:HB3	1:B:53:PHE:HA	1.93	0.49
1:B:73:GLN:HG2	1:B:77:THR:HG21	1.94	0.49
1:A:213:LEU:HA	1:A:216:ARG:HG3	1.94	0.49
1:C:166:ILE:HG13	1:C:167:VAL:N	2.28	0.49
1:D:73:GLN:HG2	1:D:77:THR:HG21	1.95	0.48
1:D:208:LEU:HD11	1:D:220:VAL:HG22	1.94	0.48
1:A:73:GLN:HG2	1:A:77:THR:HG21	1.96	0.48
1:C:33:PRO:HB3	1:C:53:PHE:HA	1.95	0.48
1:C:251:ARG:HD2	1:D:169:VAL:HG13	1.95	0.47
1:C:208:LEU:HD11	1:C:220:VAL:HG22	1.97	0.47
1:C:114:HIS:NE2	1:C:116:SER:HB2	2.30	0.47
1:A:172:ILE:HD11	1:B:172:ILE:HD11	1.97	0.47
1:C:175:TRP:HZ3	1:C:262:VAL:HG23	1.80	0.46
1:D:175:TRP:HZ3	1:D:262:VAL:HG23	1.80	0.46
1:D:249:GLU:HG3	1:D:267:GLN:HB2	1.98	0.46
1:A:197:THR:HB	1:A:200:GLU:CB	2.45	0.46
1:D:33:PRO:HB3	1:D:53:PHE:HA	1.96	0.45
1:D:208:LEU:HD22	1:D:216:ARG:HD3	1.98	0.45
1:A:217:MET:HE3	1:A:217:MET:HB2	1.87	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:ARG:HD3	2:D:300:FMN:O3P	2.18	0.44
1:C:247:TRP:HB2	1:C:269:VAL:HG22	1.99	0.44
1:D:177:ARG:HB2	1:D:184:MET:HG2	2.00	0.44
1:C:190:CYS:HA	1:C:193:ILE:HD12	2.00	0.44
1:C:179:PHE:CZ	1:C:227:VAL:HB	2.54	0.43
1:D:175:TRP:CZ3	1:D:262:VAL:HG23	2.54	0.43
1:A:30:THR:OG1	1:B:34:MET:HE2	2.19	0.42
1:B:179:PHE:CZ	1:B:227:VAL:HB	2.54	0.42
1:A:263:LEU:HD22	1:B:162:GLN:HG2	2.02	0.42
1:A:166:ILE:HA	1:B:265:ILE:HD11	2.01	0.42
1:D:218:THR:O	1:D:222:LYS:HB2	2.18	0.42
1:D:179:PHE:CZ	1:D:227:VAL:HB	2.55	0.42
1:A:179:PHE:CZ	1:A:227:VAL:HB	2.55	0.42
1:D:217:MET:O	1:D:221:GLN:HG2	2.19	0.41

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	253/266 (95%)	247 (98%)	6 (2%)	0	100	100
1	B	255/266 (96%)	249 (98%)	6 (2%)	0	100	100
1	C	231/266 (87%)	221 (96%)	10 (4%)	0	100	100
1	D	254/266 (96%)	246 (97%)	8 (3%)	0	100	100
All	All	993/1064 (93%)	963 (97%)	30 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	222/232 (96%)	211 (95%)	11 (5%)	22	40
1	B	224/232 (97%)	212 (95%)	12 (5%)	20	36
1	C	209/232 (90%)	192 (92%)	17 (8%)	11	20
1	D	223/232 (96%)	206 (92%)	17 (8%)	12	22
All	All	878/928 (95%)	821 (94%)	57 (6%)	15	28

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	116	SER
1	A	142	LEU
1	A	143	VAL
1	A	144	GLU
1	A	171	ASN
1	A	181	SER
1	A	192	ARG
1	A	200	GLU
1	A	218	THR
1	A	234	ILE
1	A	267	GLN
1	B	18	GLN
1	B	116	SER
1	B	143	VAL
1	B	153	SER
1	B	155	GLU
1	B	181	SER
1	B	192	ARG
1	B	209	ASP
1	B	216	ARG
1	B	218	THR
1	B	233	SER
1	B	234	ILE
1	C	28	GLU

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Mol	Chain	Res	Type
1	C	118	VAL
1	C	129	VAL
1	C	140	SER
1	C	143	VAL
1	C	160	LYS
1	C	181	SER
1	C	192	ARG
1	C	200	GLU
1	C	202	VAL
1	C	216	ARG
1	C	218	THR
1	C	233	SER
1	C	234	ILE
1	C	249	GLU
1	C	265	ILE
1	C	267	GLN
1	D	116	SER
1	D	142	LEU
1	D	144	GLU
1	D	145	LEU
1	D	147	LYS
1	D	149	ARG
1	D	181	SER
1	D	209	ASP
1	D	213	LEU
1	D	218	THR
1	D	222	LYS
1	D	233	SER
1	D	234	ILE
1	D	249	GLU
1	D	262	VAL
1	D	265	ILE
1	D	273	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	HIS
1	A	127	HIS
1	A	171	ASN
1	B	38	ASN
1	B	119	HIS

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Mol	Chain	Res	Type
1	B	127	HIS
1	B	162	GLN
1	C	119	HIS
1	C	162	GLN
1	C	221	GLN
1	C	267	GLN
1	D	259	ASN
1	D	267	GLN

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 oligosaccharides 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	C	300	-	33,33,33	0.46	0	48,50,50	0.59	1 (2%)
2	FMN	B	300	-	33,33,33	0.40	0	48,50,50	0.58	1 (2%)
2	FMN	D	300	-	33,33,33	0.40	0	48,50,50	0.63	1 (2%)
2	FMN	A	300	-	33,33,33	0.41	0	48,50,50	0.61	1 (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	FMN	C	300	-	-	1/18/18/18	0/3/3/3
2	FMN	B	300	-	-	1/18/18/18	0/3/3/3
2	FMN	D	300	-	-	1/18/18/18	0/3/3/3
2	FMN	A	300	-	-	1/18/18/18	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	300	FMN	O5'-P-O1P	2.58	113.41	106.44
2	A	300	FMN	O5'-P-O1P	2.56	113.37	106.44
2	C	300	FMN	O5'-P-O1P	2.26	112.56	106.44
2	B	300	FMN	O5'-P-O1P	2.21	112.42	106.44

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	300	FMN	C4'-C5'-O5'-P
2	B	300	FMN	C4'-C5'-O5'-P
2	C	300	FMN	C4'-C5'-O5'-P
2	D	300	FMN	C4'-C5'-O5'-P

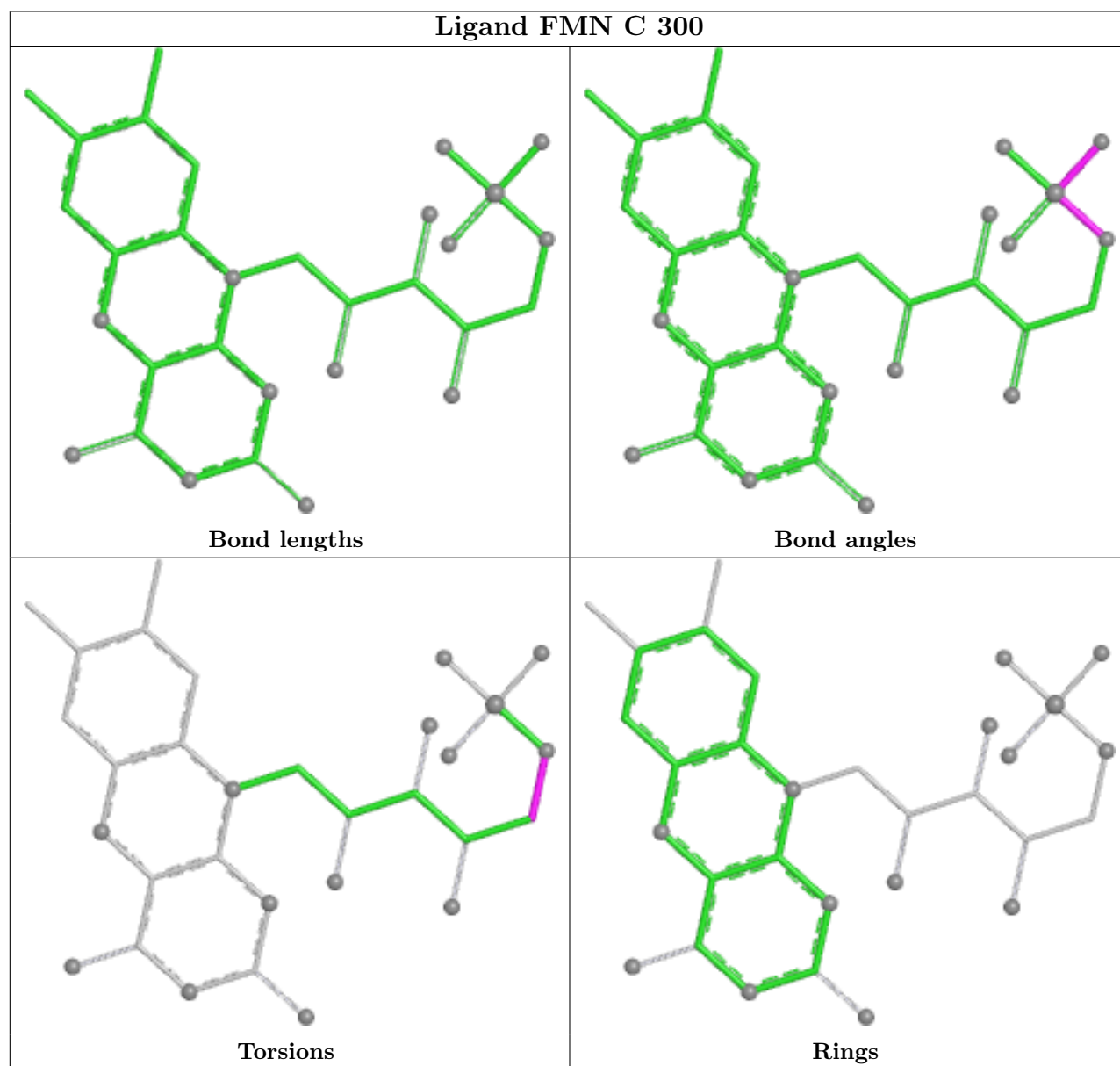
There are no ring outliers.

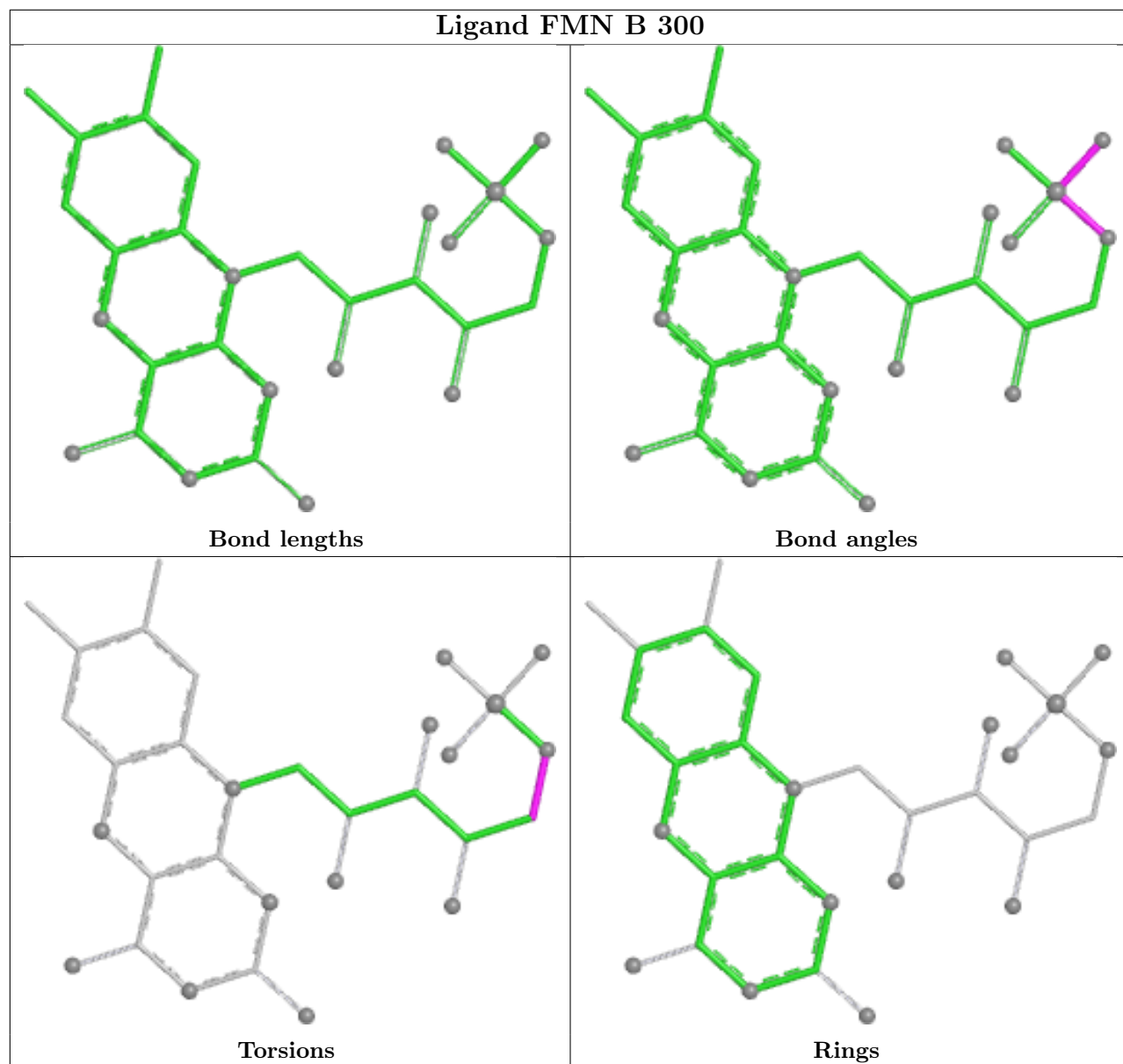
3 monomers are involved in 3 short contacts:

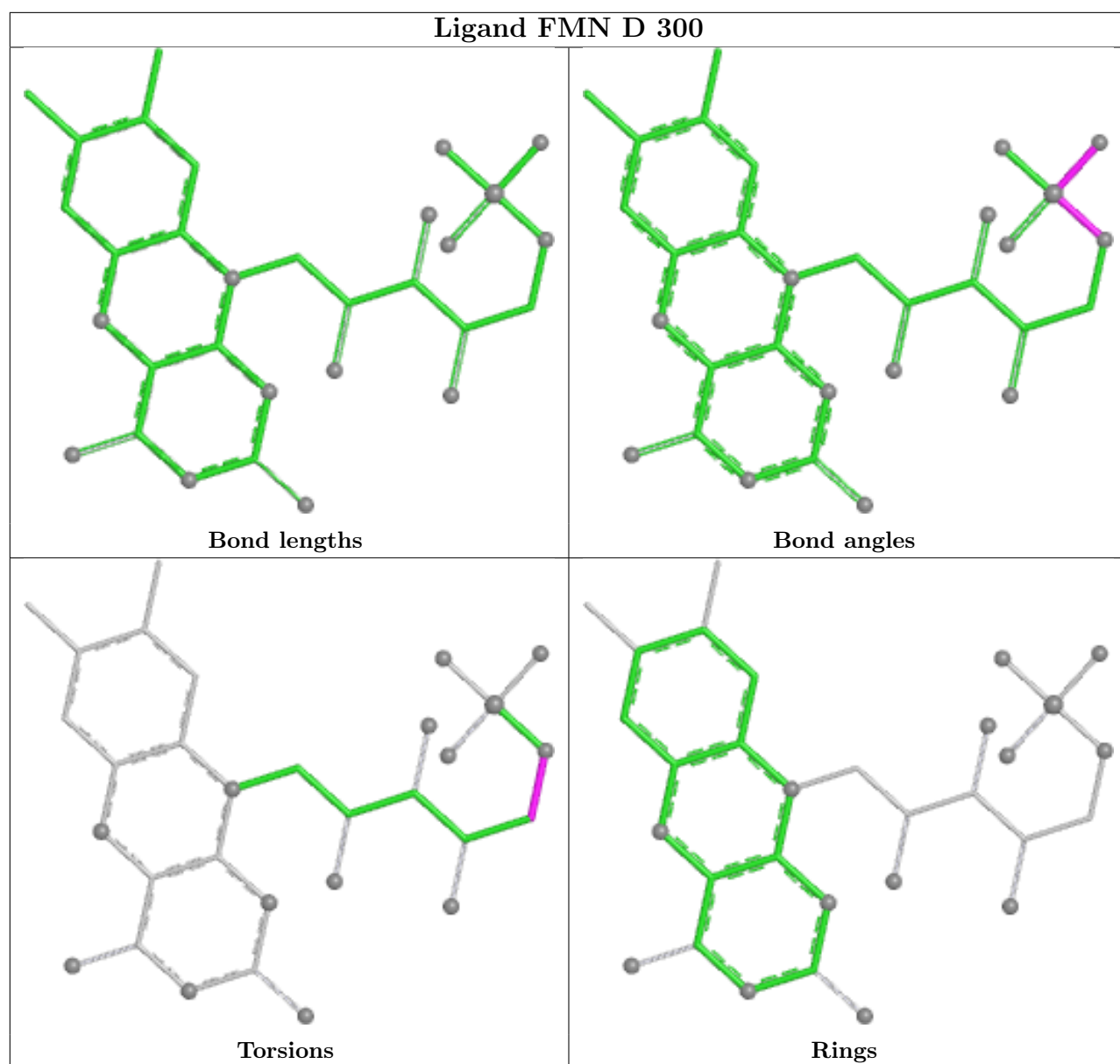
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	300	FMN	1	0
2	B	300	FMN	1	0
2	D	300	FMN	1	0

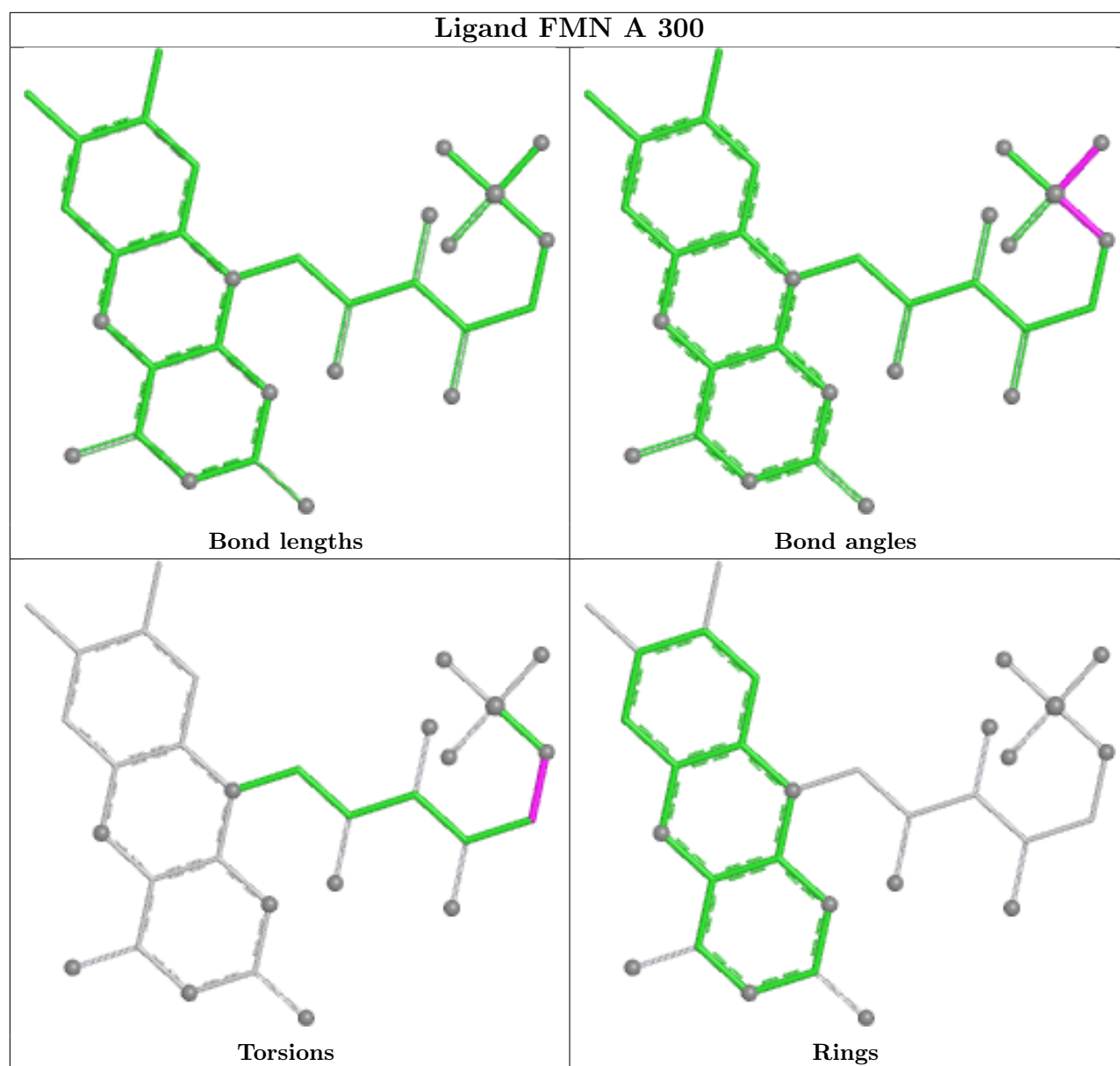
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	255/266 (95%)	0.25	2 (0%) 82 82	45, 65, 94, 114	0
1	B	257/266 (96%)	0.32	5 (1%) 66 63	44, 68, 95, 130	0
1	C	241/266 (90%)	0.72	20 (8%) 17 17	51, 81, 136, 147	0
1	D	256/266 (96%)	0.36	4 (1%) 70 68	48, 68, 95, 121	0
All	All	1009/1064 (94%)	0.41	31 (3%) 51 49	44, 69, 117, 147	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	274	HIS	4.5
1	C	242	LEU	4.4
1	C	236	TYR	3.6
1	C	219	VAL	3.5
1	C	227	VAL	3.4
1	C	197	THR	3.2
1	B	180	TYR	3.2
1	C	191	ARG	3.1
1	C	207	ILE	3.1
1	A	199	ASP	3.0
1	C	238	ILE	2.9
1	C	245	THR	2.9
1	C	201	PRO	2.9
1	C	202	VAL	2.9
1	B	166	ILE	2.8
1	B	273	LYS	2.8
1	B	19	ALA	2.8
1	D	273	LYS	2.7
1	D	223	ALA	2.6
1	C	194	TYR	2.5
1	C	206	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	239	VAL	2.5
1	C	243	GLY	2.4
1	C	269	VAL	2.4
1	C	195	GLY	2.4
1	B	179	PHE	2.3
1	D	249	GLU	2.2
1	C	185	THR	2.2
1	A	19	ALA	2.2
1	C	19	ALA	2.1
1	C	75	HIS	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

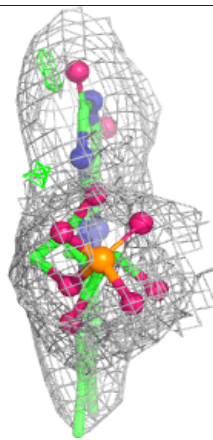
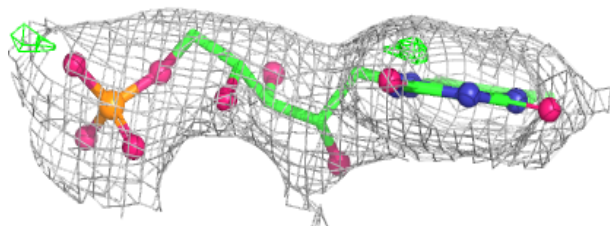
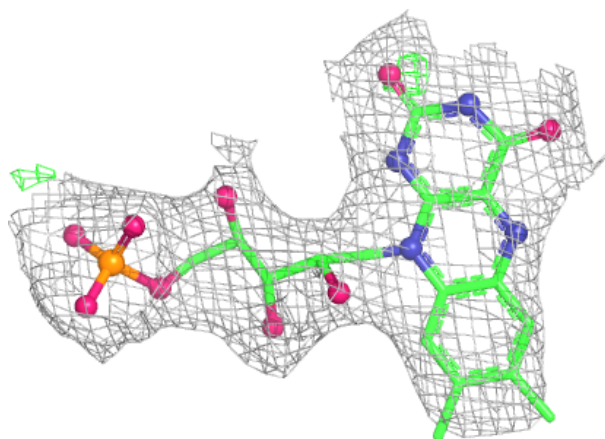
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	FMN	C	300	31/31	0.93	0.11	67,71,76,78	0
2	FMN	D	300	31/31	0.93	0.09	43,59,73,77	0
2	FMN	A	300	31/31	0.94	0.09	48,61,69,74	0
2	FMN	B	300	31/31	0.94	0.09	53,59,79,84	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.

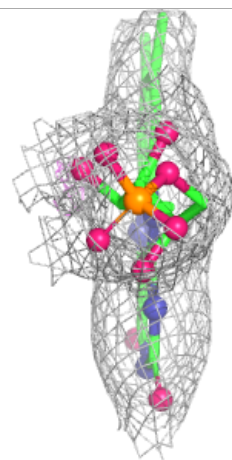
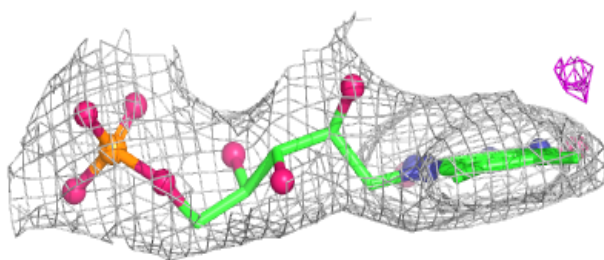
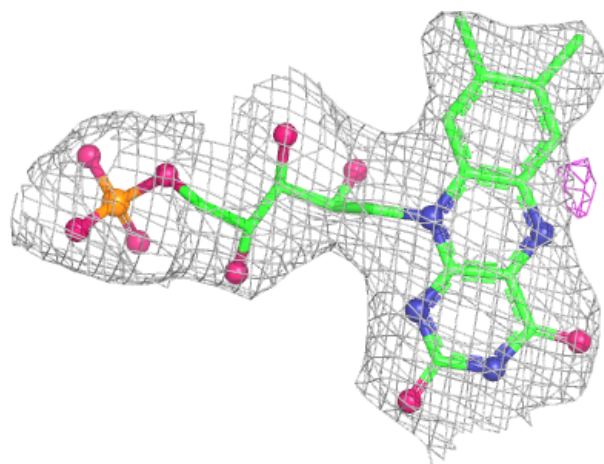
Electron density around FMN C 300:

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



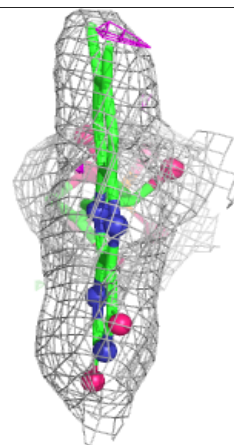
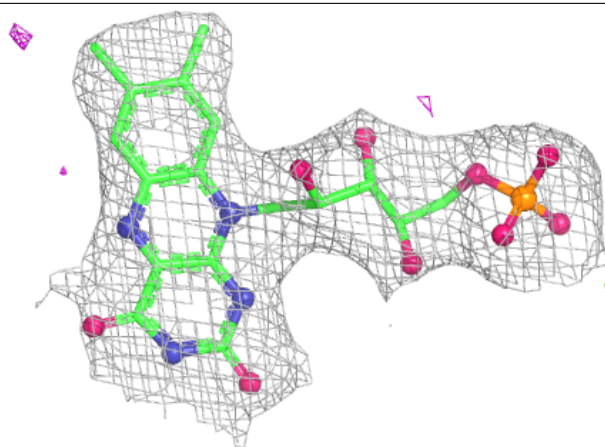
Electron density around FMN D 300:

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



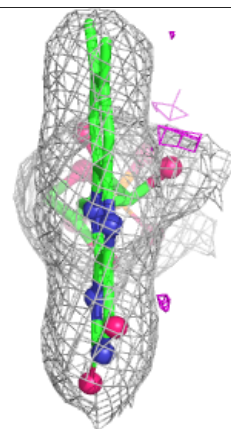
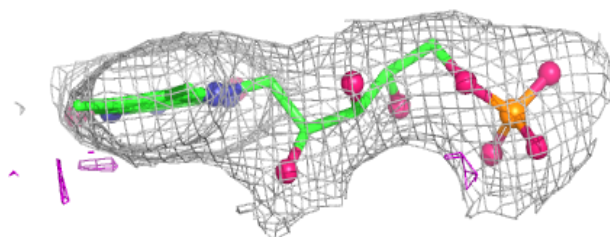
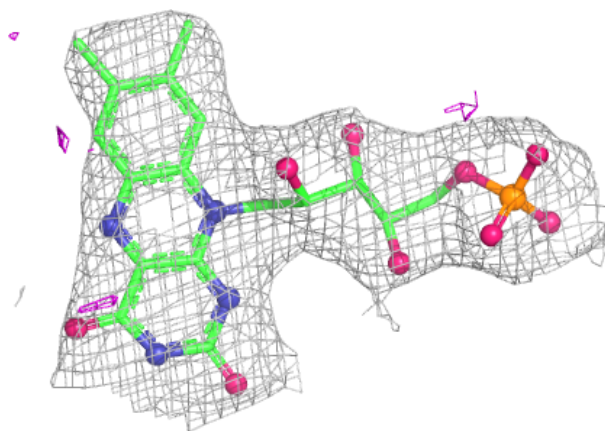
Electron density around FMN A 300:

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



Electron density around FMN B 300:

$2mF_o-DF_c$ (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.