



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:58 PM UTC

PDB ID : 4YAL / pdb_00004yal
Title : Reduced CYPOR with 2'-AMP
Authors : Xia, C.; Kim, J.J.P.
Deposited on : 2015-02-17
Resolution : 1.88 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

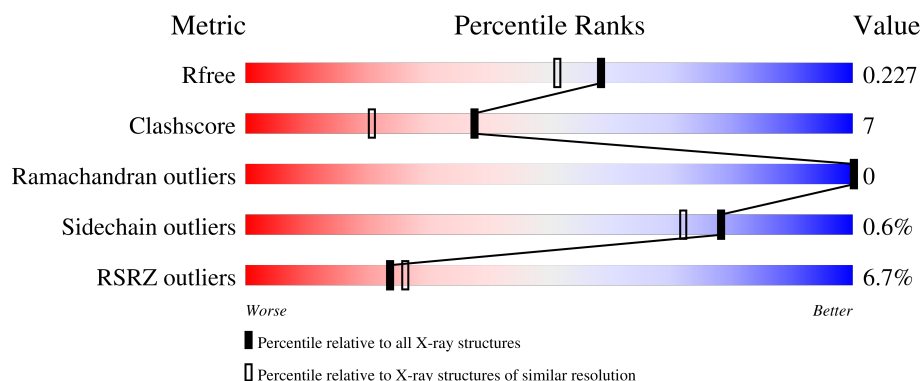
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	622	3% 84% 13% .
1	B	622	10% 79% 18% .

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 11182 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 NADPH–cytochrome P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	608	Total	C	N	O	S	0	0	0
			4868	3087	837	921	23			
1	B	608	Total	C	N	O	S	0	0	0
			4857	3080	837	917	23			

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



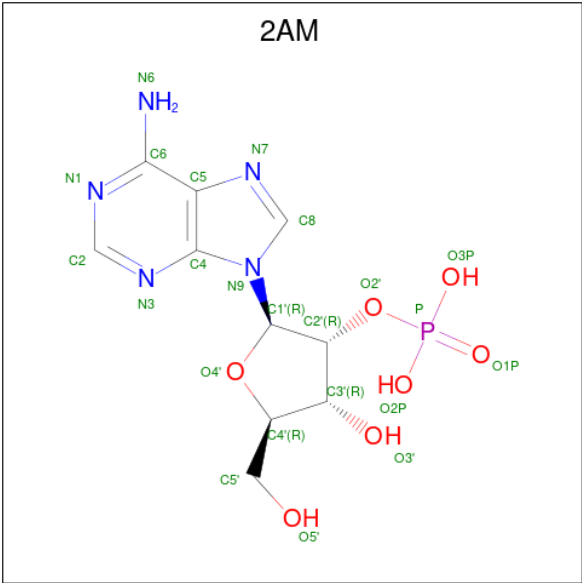
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 4 is ADENOSINE-2'-MONOPHOSPHATE (CCD ID: 2AM) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
4	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		
5	A	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		

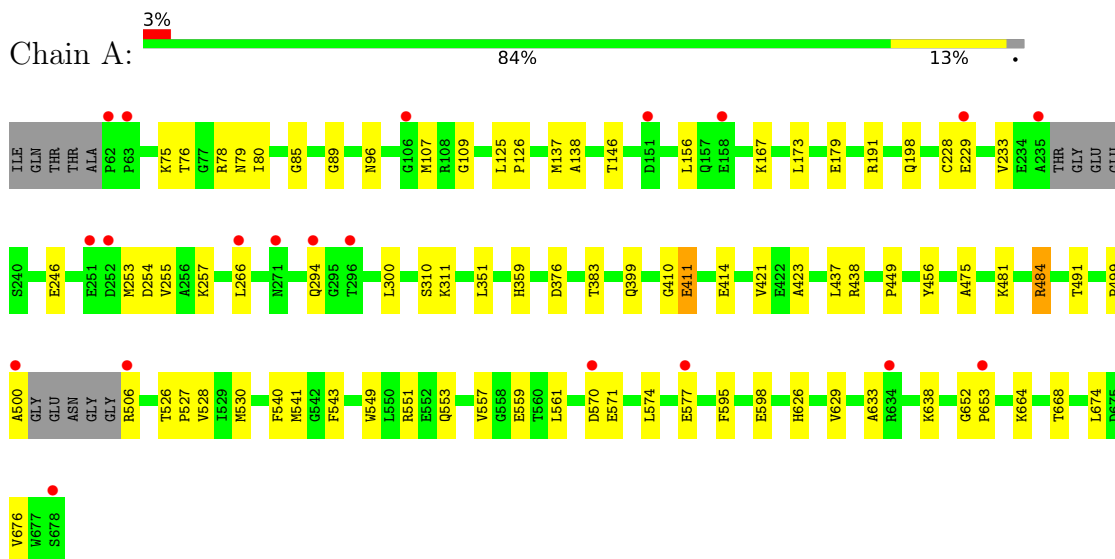
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	674	Total	O	0	0
			674	674		
6	B	549	Total	O	0	0
			549	549		

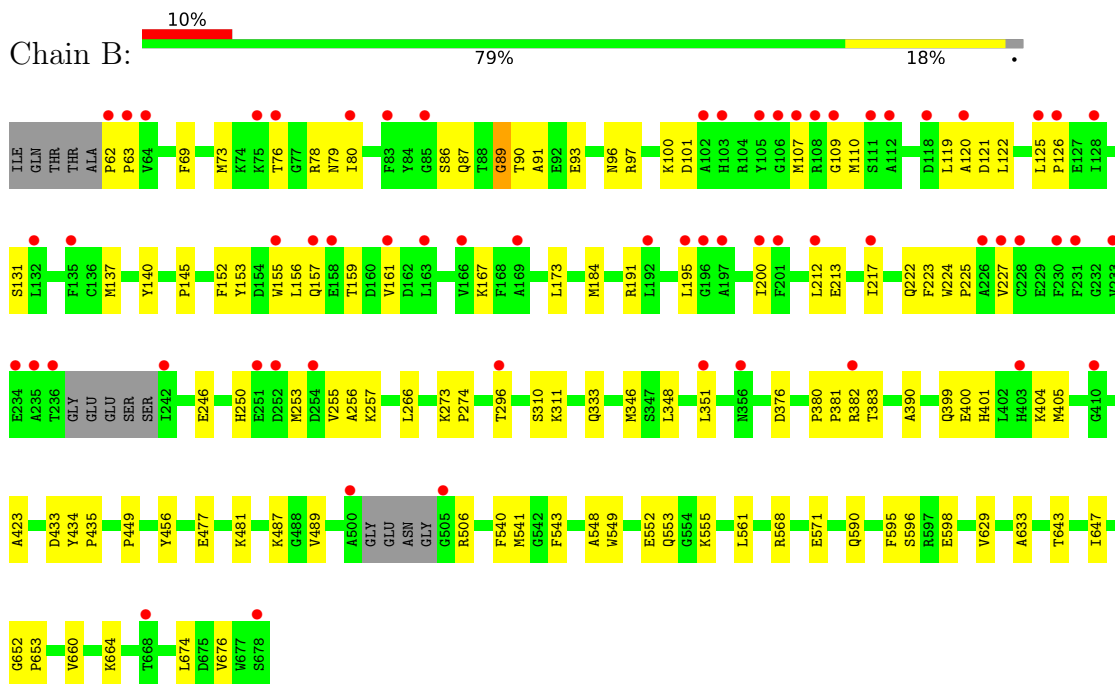
3 Residue-property plots [i](#)

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: NADPH-cytochrome P450 reductase



- Molecule 1: NADPH-cytochrome P450 reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.91Å 115.42Å 118.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.61 – 1.88 46.61 – 1.88	Depositor EDS
% Data completeness (in resolution range)	99.6 (46.61-1.88) 99.8 (46.61-1.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.23 (at 1.88Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.196 , 0.228 0.196 , 0.227	Depositor DCC
R_{free} test set	5717 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.007 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11182	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.34% 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: 2AM, FMN, FAD, PO4

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	0.38	1/4985 (0.0%)	0.84	7/6746 (0.1%)
1	B	0.34	0/4974	0.84	7/6731 (0.1%)
All	All	0.36	1/9959 (0.0%)	0.84	14/13477 (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	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	359	HIS	C-O	-8.11	1.15	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	THR	N-CA-C	6.48	119.16	111.71
1	A	541	MET	N-CA-C	-6.21	104.59	111.36
1	A	421	VAL	N-CA-C	6.05	116.11	110.42
1	B	89	GLY	N-CA-C	6.00	123.32	115.40
1	A	89	GLY	N-CA-C	5.92	123.23	115.36
1	B	434	TYR	CA-C-N	5.47	125.81	119.47
1	B	434	TYR	C-N-CA	5.47	125.81	119.47
1	A	137	MET	N-CA-C	5.45	117.40	108.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	541	MET	N-CA-C	-5.33	105.47	111.28
1	B	137	MET	N-CA-C	5.33	117.20	108.52
1	A	146	THR	N-CA-C	-5.25	103.60	110.53
1	B	489	VAL	N-CA-C	5.17	115.32	110.30
1	B	596	SER	N-CA-C	5.16	118.67	112.38
1	A	85	GLY	N-CA-C	-5.05	101.22	113.18

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	456	TYR	Sidechain
1	B	456	TYR	Sidechain

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	4868	0	4716	54	0
1	B	4857	0	4693	83	0
2	A	31	0	19	0	0
2	B	31	0	19	0	0
3	A	53	0	31	0	0
3	B	53	0	31	0	0
4	A	23	0	12	0	0
4	B	23	0	12	0	0
5	A	10	0	0	0	0
5	B	10	0	0	0	0
6	A	674	0	0	8	0
6	B	549	0	0	2	0
All	All	11182	0	9533	137	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 7.

All (137) 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:346:MET:HE1	1:B:348:LEU:HG	1.43	1.00
1:B:167:LYS:HG3	1:B:200:ILE:HD11	1.44	0.99
1:B:595:PHE:HB3	1:B:598:GLU:HG3	1.60	0.83
1:A:577:GLU:H	1:A:577:GLU:CD	1.89	0.79
1:A:254:ASP:HB3	1:A:257:LYS:HD2	1.65	0.76
1:B:86:SER:HB2	1:B:91:ALA:HB3	1.71	0.72
1:B:79:ASN:ND2	1:B:107:MET:HE2	2.03	0.72
1:B:213:GLU:O	1:B:217:ILE:HG13	1.93	0.69
1:B:561:LEU:HD22	1:B:590:GLN:HB2	1.74	0.68
1:B:660:VAL:O	1:B:664:LYS:HG2	1.94	0.68
1:B:477:GLU:HG3	1:B:487:LYS:HD2	1.74	0.67
1:B:549:TRP:O	1:B:553:GLN:HG2	1.95	0.67
1:B:76:THR:HG22	1:B:78:ARG:HE	1.62	0.65
1:A:257:LYS:HA	1:A:266:LEU:HD21	1.78	0.65
1:B:400:GLU:CD	1:B:404:LYS:HE2	2.22	0.65
1:A:399:GLN:HG2	6:A:1326:HOH:O	1.97	0.64
1:B:376:ASP:HB3	1:B:449:PRO:HG2	1.79	0.64
1:A:551:ARG:HG3	1:A:557:VAL:HG21	1.80	0.64
1:A:79:ASN:ND2	1:A:107:MET:HE2	2.13	0.64
1:B:125:LEU:N	1:B:126:PRO:HD2	2.15	0.61
1:A:376:ASP:HB3	1:A:449:PRO:HG2	1.82	0.61
1:A:626:HIS:HE1	6:A:1363:HOH:O	1.82	0.61
1:B:246:GLU:HB2	1:B:351:LEU:HD21	1.82	0.61
1:A:79:ASN:HD21	1:A:107:MET:HE2	1.65	0.60
1:B:79:ASN:HD22	1:B:107:MET:HE2	1.67	0.60
1:A:633:ALA:HB2	1:A:676:VAL:HB	1.83	0.59
1:B:76:THR:HG21	1:B:78:ARG:HH21	1.68	0.58
1:B:62:PRO:HG2	1:B:63:PRO:HD3	1.85	0.58
1:B:629:VAL:HB	1:B:674:LEU:HD23	1.86	0.58
1:A:76:THR:OG1	1:A:78:ARG:HD3	2.03	0.58
1:B:119:LEU:HG	1:B:152:PHE:CD1	2.39	0.57
1:B:382:ARG:HG2	1:B:382:ARG:HH11	1.72	0.55
1:B:79:ASN:HD21	1:B:107:MET:HB3	1.72	0.55
1:B:145:PRO:HB3	1:B:184:MET:SD	2.46	0.55
1:A:506:ARG:NH1	1:A:506:ARG:HB3	2.22	0.55
1:B:87:GLN:HG2	1:B:140:TYR:CE1	2.42	0.55
1:B:76:THR:HG21	1:B:78:ARG:NH2	2.23	0.54
1:A:437:LEU:O	1:A:438:ARG:HD3	2.07	0.54
1:A:561:LEU:HD12	1:A:561:LEU:N	2.23	0.53
1:B:253:MET:HE2	1:B:257:LYS:HB2	1.91	0.53
1:A:570:ASP:OD1	1:A:571:GLU:HG3	2.08	0.53
1:A:228:CYS:HA	1:A:233:VAL:CG2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:TRP:O	1:B:159:THR:HG22	2.10	0.52
1:B:79:ASN:O	1:B:131:SER:HA	2.09	0.52
1:B:90:THR:HG23	1:B:382:ARG:HH21	1.74	0.52
1:B:568:ARG:HD2	1:B:571:GLU:CD	2.34	0.52
1:B:222:GLN:C	1:B:225:PRO:HD2	2.35	0.51
1:B:595:PHE:CB	1:B:598:GLU:HG3	2.37	0.51
1:B:253:MET:CE	1:B:257:LYS:HD2	2.40	0.51
1:B:555:LYS:N	1:B:555:LYS:HD2	2.26	0.51
1:B:548:ALA:O	1:B:552:GLU:HG2	2.11	0.50
1:B:120:ALA:HA	1:B:155:TRP:CZ2	2.46	0.50
1:B:159:THR:OG1	1:B:161:VAL:HG23	2.11	0.50
1:A:179:GLU:OE2	1:A:664:LYS:HE2	2.13	0.49
1:A:78:ARG:NH1	6:A:805:HOH:O	2.44	0.49
1:A:78:ARG:HH11	1:A:78:ARG:HG2	1.76	0.49
1:B:90:THR:O	1:B:93:GLU:HB2	2.11	0.49
1:B:253:MET:HE3	6:B:1227:HOH:O	2.12	0.49
1:B:433:ASP:C	1:B:435:PRO:HD3	2.38	0.49
1:A:499:PRO:O	1:A:500:ALA:HB3	2.13	0.49
1:B:62:PRO:N	1:B:63:PRO:HD2	2.28	0.49
1:B:310:SER:O	1:B:311:LYS:HB2	2.13	0.49
1:A:528:VAL:HG12	1:A:530:MET:HG3	1.95	0.48
1:B:223:PHE:O	1:B:227:VAL:HG23	2.12	0.48
1:A:254:ASP:O	1:A:257:LYS:HG2	2.12	0.48
1:A:638:LYS:HG2	6:A:1021:HOH:O	2.12	0.48
1:A:506:ARG:CB	1:A:506:ARG:HH11	2.26	0.48
1:B:96:ASN:O	1:B:100:LYS:HG3	2.14	0.48
1:B:78:ARG:HH11	1:B:110:MET:HB3	1.79	0.48
1:B:80:ILE:O	1:B:109:GLY:HA2	2.13	0.47
1:B:153:TYR:CE2	1:B:157:GLN:OE1	2.67	0.47
1:B:191:ARG:O	1:B:195:LEU:HD13	2.14	0.47
1:B:69:PHE:CE1	1:B:121:ASP:HB2	2.49	0.47
1:B:222:GLN:O	1:B:225:PRO:HD2	2.14	0.47
1:B:257:LYS:HA	1:B:266:LEU:HD21	1.96	0.47
1:B:633:ALA:HB2	1:B:676:VAL:HB	1.96	0.47
1:B:477:GLU:HG3	1:B:487:LYS:CD	2.45	0.47
1:B:119:LEU:O	1:B:122:LEU:HG	2.16	0.46
1:A:652:GLY:N	1:A:653:PRO:HA	2.30	0.46
1:B:346:MET:SD	1:B:346:MET:C	2.99	0.46
1:B:346:MET:CE	1:B:348:LEU:HG	2.30	0.46
1:B:390:ALA:O	1:B:399:GLN:HG3	2.16	0.46
1:A:75:LYS:HD2	6:A:970:HOH:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:HIS:O	1:B:405:MET:HG2	2.16	0.45
1:B:255:VAL:HG23	1:B:256:ALA:N	2.31	0.45
1:A:506:ARG:NH1	1:A:506:ARG:CB	2.80	0.45
1:B:62:PRO:N	1:B:63:PRO:CD	2.80	0.45
1:B:76:THR:HG22	1:B:76:THR:O	2.18	0.44
1:B:296:THR:O	1:B:296:THR:HG22	2.16	0.44
1:A:253:MET:HG3	1:A:257:LYS:HD3	2.00	0.44
1:A:540:PHE:HA	1:A:543:PHE:HB2	1.99	0.44
1:A:246:GLU:HB3	1:A:351:LEU:HD21	1.99	0.44
1:B:152:PHE:CZ	1:B:156:LEU:HD11	2.51	0.44
1:B:89:GLY:O	1:B:93:GLU:HG2	2.16	0.44
1:B:101:ASP:HB2	1:B:224:TRP:CZ2	2.52	0.44
1:B:246:GLU:HB2	1:B:351:LEU:CD2	2.47	0.44
1:A:228:CYS:HA	1:A:233:VAL:HG22	1.99	0.44
1:A:310:SER:O	1:A:311:LYS:HB2	2.18	0.44
1:B:69:PHE:O	1:B:73:MET:HG3	2.18	0.44
1:A:125:LEU:N	1:A:126:PRO:CD	2.81	0.44
1:A:96:ASN:ND2	6:A:806:HOH:O	2.51	0.43
1:A:411:GLU:HA	1:A:414:GLU:OE1	2.18	0.43
1:A:475:ALA:HA	1:A:491:THR:HB	1.99	0.43
1:A:595:PHE:HB3	1:A:598:GLU:HG3	1.98	0.43
1:A:423:ALA:HA	1:A:481:LYS:HB2	2.00	0.43
1:A:629:VAL:HB	1:A:674:LEU:HD23	1.99	0.43
1:B:253:MET:HE2	1:B:257:LYS:HD2	2.00	0.43
1:B:125:LEU:N	1:B:126:PRO:CD	2.81	0.43
1:B:250:HIS:HB3	1:B:253:MET:HB2	2.00	0.43
1:A:300:LEU:HD13	1:A:574:LEU:HD21	1.99	0.43
1:B:173:LEU:HD12	1:B:173:LEU:N	2.34	0.43
1:A:138:ALA:HA	1:A:173:LEU:HB2	2.01	0.42
1:B:423:ALA:HA	1:B:481:LYS:HB2	2.00	0.42
1:A:257:LYS:HB2	6:A:873:HOH:O	2.18	0.42
1:B:78:ARG:NH1	1:B:110:MET:HB3	2.34	0.42
1:B:382:ARG:HH11	1:B:382:ARG:CG	2.32	0.42
1:B:643:THR:O	1:B:647:ILE:HG13	2.19	0.42
1:B:540:PHE:HA	1:B:543:PHE:HB2	2.01	0.42
1:A:410:GLY:O	1:A:414:GLU:HG3	2.20	0.42
1:A:577:GLU:CD	1:A:577:GLU:N	2.66	0.42
1:A:664:LYS:O	1:A:668:THR:HG23	2.20	0.42
1:A:167:LYS:HE3	1:A:198:GLN:OE1	2.19	0.41
1:A:549:TRP:O	1:A:553:GLN:HG2	2.20	0.41
1:A:559:GLU:HB3	1:A:561:LEU:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:VAL:HG23	6:A:1415:HOH:O	2.20	0.41
1:B:93:GLU:O	1:B:97:ARG:HG3	2.20	0.41
1:B:87:GLN:HG2	1:B:140:TYR:CZ	2.56	0.41
1:B:333:GLN:NE2	6:B:816:HOH:O	2.53	0.41
1:B:273:LYS:HA	1:B:274:PRO:HD3	1.95	0.41
1:B:652:GLY:N	1:B:653:PRO:HA	2.36	0.41
1:A:484:ARG:N	1:A:484:ARG:HD2	2.36	0.41
1:A:526:THR:HA	1:A:527:PRO:HD3	1.94	0.40
1:B:217:ILE:HD11	1:B:383:THR:HG21	2.04	0.40
1:A:78:ARG:HD2	1:A:78:ARG:N	2.37	0.40
1:A:156:LEU:HB3	1:A:191:ARG:HG2	2.04	0.40
1:A:80:ILE:O	1:A:109:GLY:HA2	2.21	0.40
1:B:380:PRO:HA	1:B:381:PRO:HD3	1.98	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	602/622 (97%)	589 (98%)	13 (2%)	0	100	100
1	B	602/622 (97%)	578 (96%)	24 (4%)	0	100	100
All	All	1204/1244 (97%)	1167 (97%)	37 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	519/531 (98%)	515 (99%)	4 (1%)	73	68
1	B	514/531 (97%)	512 (100%)	2 (0%)	84	79
All	All	1033/1062 (97%)	1027 (99%)	6 (1%)	78	72

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	229	GLU
1	A	294	GLN
1	A	411	GLU
1	A	484	ARG
1	B	212	LEU
1	B	506	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	HIS
1	A	359	HIS
1	A	467	ASN
1	A	592	ASN
1	B	180	HIS
1	B	190	GLN
1	B	194	GLN
1	B	244	GLN
1	B	391	GLN
1	B	403	HIS
1	B	467	ASN
1	B	590	GLN
1	B	592	ASN
1	B	599	GLN
1	B	635	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

10 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	$\# Z > 2$
5	PO4	B	705	-	4,4,4	1.70	0	6,6,6	0.45	0
4	2AM	A	703	-	25,25,25	1.71	7 (28%)	38,38,38	2.14	12 (31%)
5	PO4	A	705	-	4,4,4	1.70	0	6,6,6	0.45	0
2	FMN	B	701	-	33,33,33	2.59	12 (36%)	48,50,50	1.63	12 (25%)
5	PO4	B	704	-	4,4,4	1.59	1 (25%)	6,6,6	0.46	0
5	PO4	A	704	-	4,4,4	1.60	2 (50%)	6,6,6	0.47	0
3	FAD	B	702	-	58,58,58	2.31	13 (22%)	85,89,89	1.96	16 (18%)
2	FMN	A	701	-	33,33,33	2.59	12 (36%)	48,50,50	1.61	13 (27%)
4	2AM	B	703	-	25,25,25	1.71	7 (28%)	38,38,38	2.15	13 (34%)
3	FAD	A	702	-	58,58,58	2.30	13 (22%)	85,89,89	1.96	16 (18%)

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
4	2AM	A	703	-	-	4/11/27/27	0/3/3/3
2	FMN	B	701	-	-	0/18/18/18	0/3/3/3
3	FAD	B	702	-	-	2/34/50/50	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FMN	A	701	-	-	0/18/18/18	0/3/3/3
4	2AM	B	703	-	-	4/11/27/27	0/3/3/3
3	FAD	A	702	-	-	3/34/50/50	0/6/6/6

All (67) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	702	FAD	C8A-N7A	8.56	1.48	1.31
3	A	702	FAD	C8A-N7A	8.54	1.48	1.31
2	A	701	FMN	C8M-C8	-6.77	1.38	1.51
2	B	701	FMN	C8M-C8	-6.74	1.38	1.51
3	A	702	FAD	C8A-N9A	6.18	1.48	1.37
3	B	702	FAD	C8A-N9A	6.17	1.48	1.37
3	B	702	FAD	C8M-C8	-6.13	1.39	1.51
3	B	702	FAD	C7M-C7	-6.13	1.39	1.51
3	A	702	FAD	C8M-C8	-6.13	1.39	1.51
3	A	702	FAD	C7M-C7	-6.11	1.39	1.51
2	B	701	FMN	C4A-N5	5.24	1.42	1.30
2	A	701	FMN	C4A-N5	5.11	1.41	1.30
2	A	701	FMN	C10-N10	4.97	1.48	1.37
2	B	701	FMN	C10-N10	4.87	1.47	1.37
2	B	701	FMN	C9-C9A	4.76	1.47	1.39
2	A	701	FMN	C9-C9A	4.75	1.47	1.39
2	A	701	FMN	C9A-C5A	4.21	1.48	1.41
2	B	701	FMN	C9A-N10	4.17	1.48	1.41
2	A	701	FMN	C9A-N10	4.08	1.48	1.41
2	B	701	FMN	C9A-C5A	4.07	1.47	1.41
2	A	701	FMN	C8-C7	3.96	1.50	1.40
2	B	701	FMN	C8-C7	3.92	1.50	1.40
3	A	702	FAD	C9A-N10	-3.82	1.34	1.41
3	B	702	FAD	C9A-N10	-3.78	1.34	1.41
4	A	703	2AM	C2-N3	3.67	1.40	1.33
4	B	703	2AM	C2-N1	3.63	1.40	1.33
3	A	702	FAD	C2A-N3A	3.61	1.40	1.33
3	B	702	FAD	C2A-N1A	3.60	1.40	1.33
3	B	702	FAD	C2A-N3A	3.59	1.40	1.33
3	A	702	FAD	C2A-N1A	3.59	1.40	1.33
4	B	703	2AM	C2-N3	3.59	1.40	1.33
4	A	703	2AM	C2-N1	3.58	1.40	1.33
3	B	702	FAD	C10-N1	3.58	1.40	1.33
3	A	702	FAD	C10-N1	3.49	1.40	1.33
2	B	701	FMN	C10-N1	3.24	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	703	2AM	C5-C4	-3.14	1.33	1.39
2	A	701	FMN	C10-N1	3.09	1.39	1.33
4	A	703	2AM	C5-N7	-3.05	1.33	1.39
4	B	703	2AM	C5-N7	-3.05	1.33	1.39
3	A	702	FAD	C5A-N7A	-3.05	1.33	1.39
3	B	702	FAD	C5A-N7A	-3.05	1.33	1.39
4	A	703	2AM	C5-C4	-3.01	1.33	1.39
2	A	701	FMN	C6-C5A	2.94	1.44	1.40
3	A	702	FAD	C5A-C4A	-2.90	1.33	1.39
3	B	702	FAD	C5A-C4A	-2.89	1.33	1.39
2	B	701	FMN	C6-C5A	2.79	1.44	1.40
3	A	702	FAD	C5A-C6A	-2.69	1.33	1.41
3	B	702	FAD	C5X-N5	-2.68	1.34	1.39
3	B	702	FAD	C5A-C6A	-2.67	1.33	1.41
4	B	703	2AM	C5-C6	-2.65	1.33	1.41
4	A	703	2AM	C5-C6	-2.63	1.33	1.41
3	A	702	FAD	C5X-N5	-2.63	1.34	1.39
2	A	701	FMN	C1'-C2'	2.55	1.56	1.52
2	A	701	FMN	O2'-C2'	2.55	1.48	1.43
2	B	701	FMN	O2'-C2'	2.46	1.48	1.43
4	A	703	2AM	C8-N9	-2.43	1.33	1.37
2	B	701	FMN	C4'-C3'	2.40	1.57	1.53
4	B	703	2AM	C8-N9	-2.40	1.33	1.37
2	A	701	FMN	C7M-C7	2.12	1.55	1.51
2	B	701	FMN	C7M-C7	2.08	1.54	1.51
5	A	704	PO4	P-O3	-2.06	1.48	1.54
4	A	703	2AM	C4-N9	-2.05	1.33	1.37
3	A	702	FAD	C4A-N9A	-2.03	1.33	1.37
4	B	703	2AM	C4-N9	-2.02	1.33	1.37
5	B	704	PO4	P-O4	-2.02	1.48	1.54
3	B	702	FAD	C4A-N9A	-2.02	1.33	1.37
5	A	704	PO4	P-O2	-2.00	1.48	1.54

All (82) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	702	FAD	N9A-C8A-N7A	-9.41	100.59	113.94
3	A	702	FAD	N9A-C8A-N7A	-9.34	100.69	113.94
3	A	702	FAD	N3A-C2A-N1A	-7.34	117.47	128.58
3	B	702	FAD	N3A-C2A-N1A	-7.34	117.47	128.58
4	B	703	2AM	N3-C2-N1	-7.20	117.69	128.58
4	A	703	2AM	N3-C2-N1	-7.12	117.80	128.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	702	FAD	C5A-C4A-N3A	-4.54	120.47	126.72
3	A	702	FAD	C5A-C4A-N3A	-4.48	120.54	126.72
4	B	703	2AM	N9-C8-N7	-4.42	107.67	113.94
4	A	703	2AM	N9-C8-N7	-4.32	107.81	113.94
3	A	702	FAD	C5A-C4A-N9A	4.27	110.46	105.81
4	A	703	2AM	C5-C4-N3	-4.18	120.96	126.72
4	B	703	2AM	C5-C4-N3	-4.17	120.98	126.72
3	B	702	FAD	C5A-C4A-N9A	4.14	110.33	105.81
2	A	701	FMN	C9A-C5A-N5	4.04	126.74	122.45
2	B	701	FMN	C9A-C5A-N5	3.85	126.54	122.45
2	A	701	FMN	O4'-C4'-C3'	-3.58	100.88	109.25
3	B	702	FAD	C2A-N3A-C4A	3.55	120.50	111.83
3	A	702	FAD	C2A-N3A-C4A	3.53	120.45	111.83
2	B	701	FMN	O4'-C4'-C3'	-3.44	101.20	109.25
3	A	702	FAD	C5A-N7A-C8A	3.43	108.83	103.45
3	B	702	FAD	C5A-N7A-C8A	3.39	108.77	103.45
2	B	701	FMN	C4'-C3'-C2'	-3.38	107.95	113.57
4	B	703	2AM	C2-N3-C4	3.30	119.90	111.83
4	A	703	2AM	C2-N3-C4	3.27	119.81	111.83
2	A	701	FMN	C4'-C3'-C2'	-3.22	108.21	113.57
4	B	703	2AM	C5-N7-C8	3.12	108.36	103.45
3	A	702	FAD	C4A-N9A-C8A	3.12	109.01	105.74
3	B	702	FAD	C4-N3-C2	-3.06	120.20	125.64
4	A	703	2AM	C5-N7-C8	3.05	108.24	103.45
3	A	702	FAD	C5A-C6A-N6A	-3.03	115.79	123.29
3	A	702	FAD	C4-N3-C2	-2.94	120.42	125.64
3	B	702	FAD	C5A-C6A-N6A	-2.91	116.09	123.29
3	B	702	FAD	C4A-N9A-C8A	2.90	108.78	105.74
3	A	702	FAD	C10-C4X-N5	-2.87	118.95	124.81
3	B	702	FAD	C10-C4X-N5	-2.82	119.06	124.81
4	A	703	2AM	C5-C6-N6	-2.67	116.68	123.29
2	B	701	FMN	C5A-C9A-N10	-2.67	115.56	117.97
3	B	702	FAD	C4X-C4-N3	2.66	120.02	113.25
4	B	703	2AM	C5-C6-N6	-2.63	116.78	123.29
2	B	701	FMN	C8M-C8-C7	2.61	126.09	120.76
2	A	701	FMN	C5A-C9A-N10	-2.60	115.61	117.97
2	B	701	FMN	O4-C4-N3	-2.59	115.25	120.11
3	A	702	FAD	C4X-C4-N3	2.57	119.80	113.25
2	A	701	FMN	C5A-N5-C4A	-2.55	113.97	118.09
2	A	701	FMN	C8M-C8-C7	2.54	125.95	120.76
3	B	702	FAD	C2B-C3B-C4B	-2.53	97.71	102.61
2	A	701	FMN	O4-C4-N3	-2.53	115.36	120.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	FMN	C5A-N5-C4A	-2.50	114.05	118.09
4	B	703	2AM	C4-C5-N7	-2.48	107.75	110.58
4	A	703	2AM	C6-C5-C4	2.46	120.54	117.18
4	B	703	2AM	C6-C5-C4	2.45	120.52	117.18
4	A	703	2AM	C4-C5-N7	-2.44	107.79	110.58
4	A	703	2AM	O4'-C1'-C2'	-2.42	102.43	106.59
2	B	701	FMN	O2P-P-O1P	2.39	120.15	110.83
3	B	702	FAD	C9A-C5X-N5	-2.34	119.97	122.45
4	B	703	2AM	C4-N9-C8	2.34	108.19	105.74
4	A	703	2AM	C4-N9-C8	2.34	108.19	105.74
2	B	701	FMN	O3P-P-O5'	-2.34	100.58	106.67
2	B	701	FMN	C8M-C8-C9	-2.33	115.48	119.57
2	A	701	FMN	C8M-C8-C9	-2.32	115.48	119.57
3	B	702	FAD	C6A-C5A-C4A	2.32	120.34	117.18
3	A	702	FAD	C2B-C3B-C4B	-2.31	98.14	102.61
3	A	702	FAD	C6A-C5A-C4A	2.29	120.30	117.18
2	A	701	FMN	O2P-P-O1P	2.29	119.74	110.83
4	B	703	2AM	C3'-C2'-C1'	-2.28	98.44	102.81
4	A	703	2AM	N3-C4-N9	2.28	131.04	127.17
3	B	702	FAD	O4-C4-C4X	-2.25	120.58	126.53
4	B	703	2AM	N3-C4-N9	2.24	130.97	127.17
2	A	701	FMN	O3P-P-O5'	-2.23	100.84	106.67
3	A	702	FAD	C4X-C10-N10	2.23	119.68	116.48
3	A	702	FAD	O4-C4-C4X	-2.19	120.74	126.53
2	B	701	FMN	C7M-C7-C6	-2.19	115.72	119.57
3	B	702	FAD	C4X-C10-N10	2.18	119.60	116.48
3	A	702	FAD	C9A-C5X-N5	-2.17	120.15	122.45
4	A	703	2AM	C5'-C4'-C3'	-2.17	109.98	115.10
2	A	701	FMN	C7M-C7-C6	-2.16	115.77	119.57
2	B	701	FMN	C9-C9A-N10	2.14	124.73	121.85
4	B	703	2AM	O4'-C1'-C2'	-2.04	103.07	106.59
4	B	703	2AM	C5-C4-N9	2.02	108.01	105.81
2	A	701	FMN	C9-C9A-N10	2.02	124.57	121.85
2	A	701	FMN	C6-C5A-N5	-2.01	115.11	118.44

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	703	2AM	C2'-C1'-N9-C4
4	B	703	2AM	C2'-C1'-N9-C4
4	A	703	2AM	C2'-C1'-N9-C8

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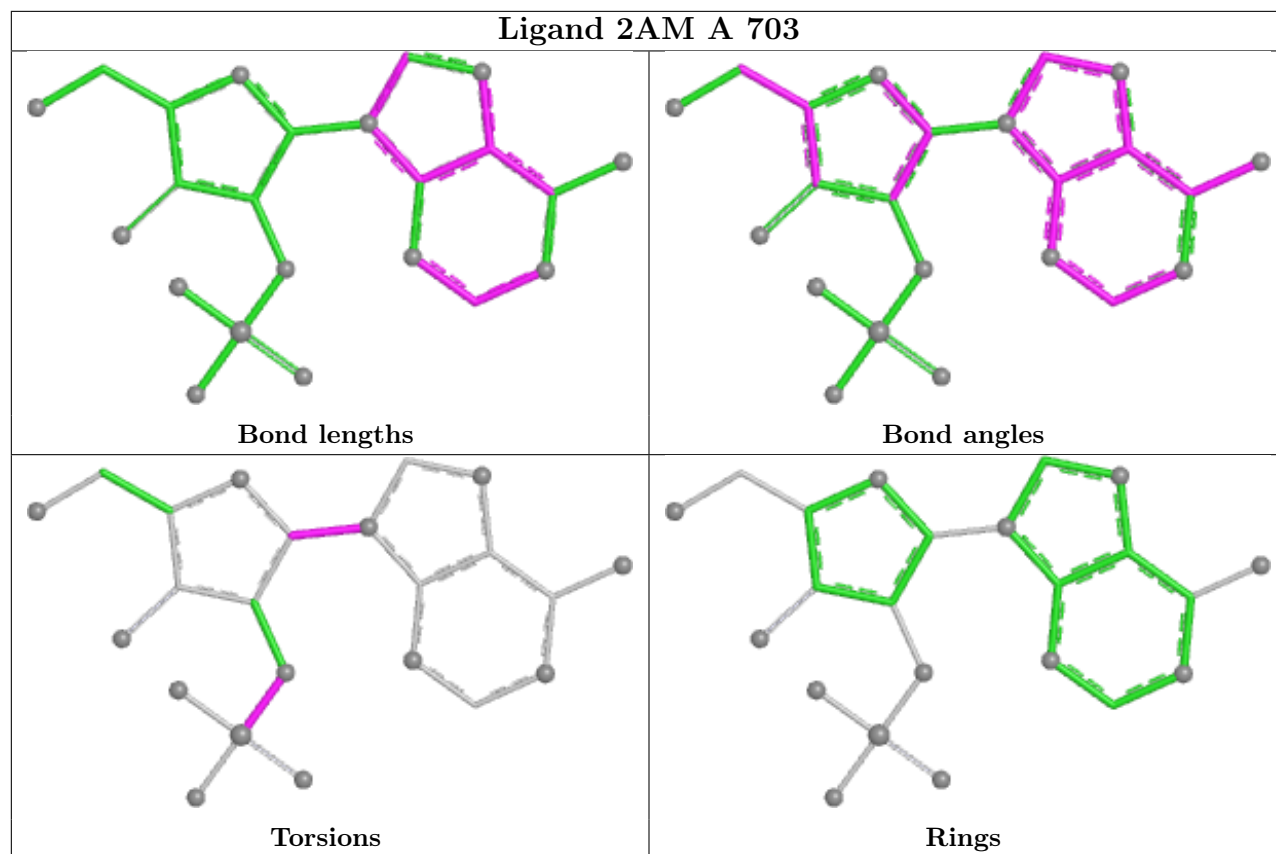
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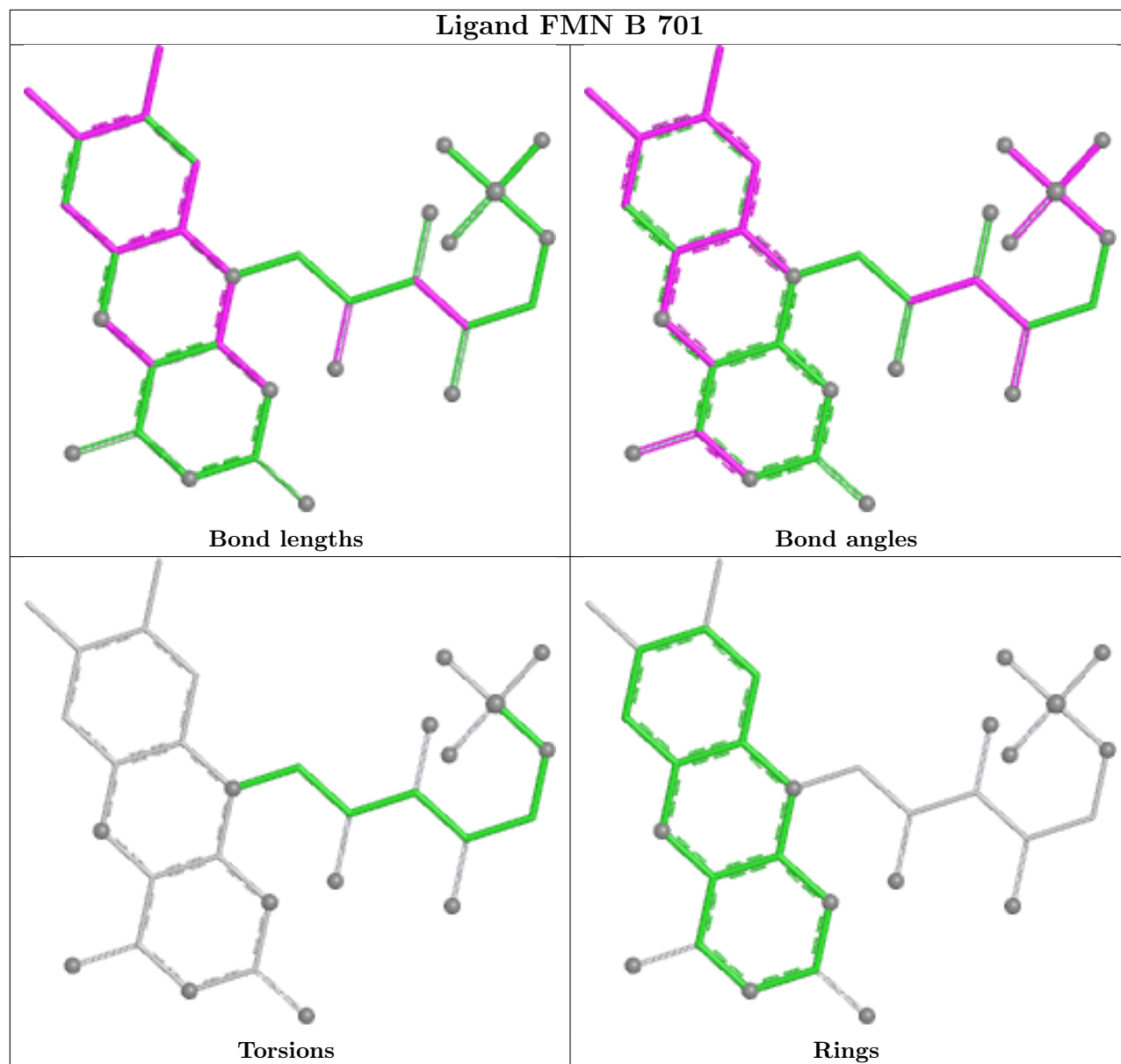
Mol	Chain	Res	Type	Atoms
4	B	703	2AM	C2'-C1'-N9-C8
4	A	703	2AM	C2'-O2'-P-O1P
4	B	703	2AM	C2'-O2'-P-O1P
3	A	702	FAD	C5'-O5'-P-O1P
3	A	702	FAD	PA-O3P-P-O1P
3	B	702	FAD	PA-O3P-P-O1P
4	A	703	2AM	O4'-C1'-N9-C8
3	B	702	FAD	PA-O3P-P-O2P
4	B	703	2AM	O4'-C1'-N9-C8
3	A	702	FAD	PA-O3P-P-O2P

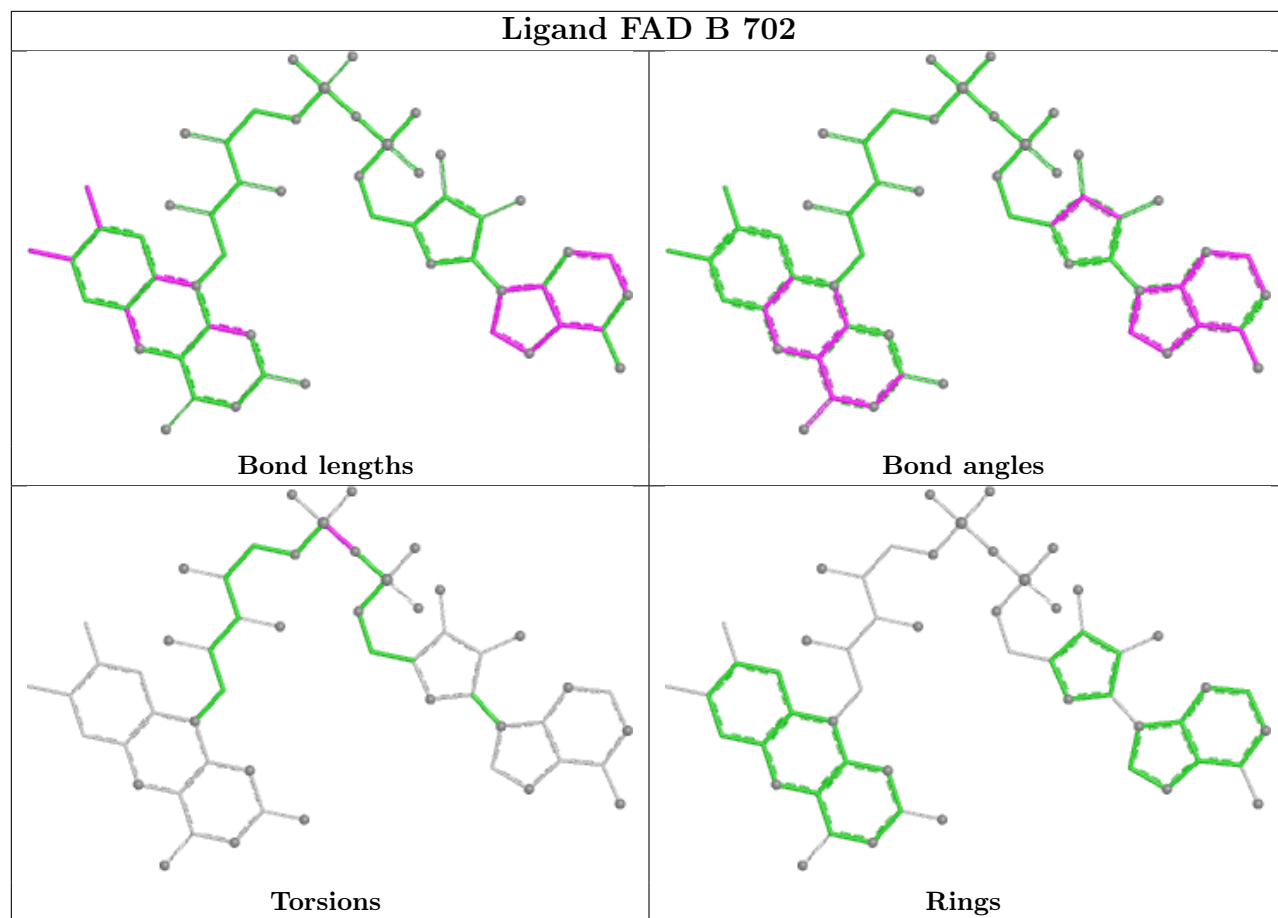
There are no ring outliers.

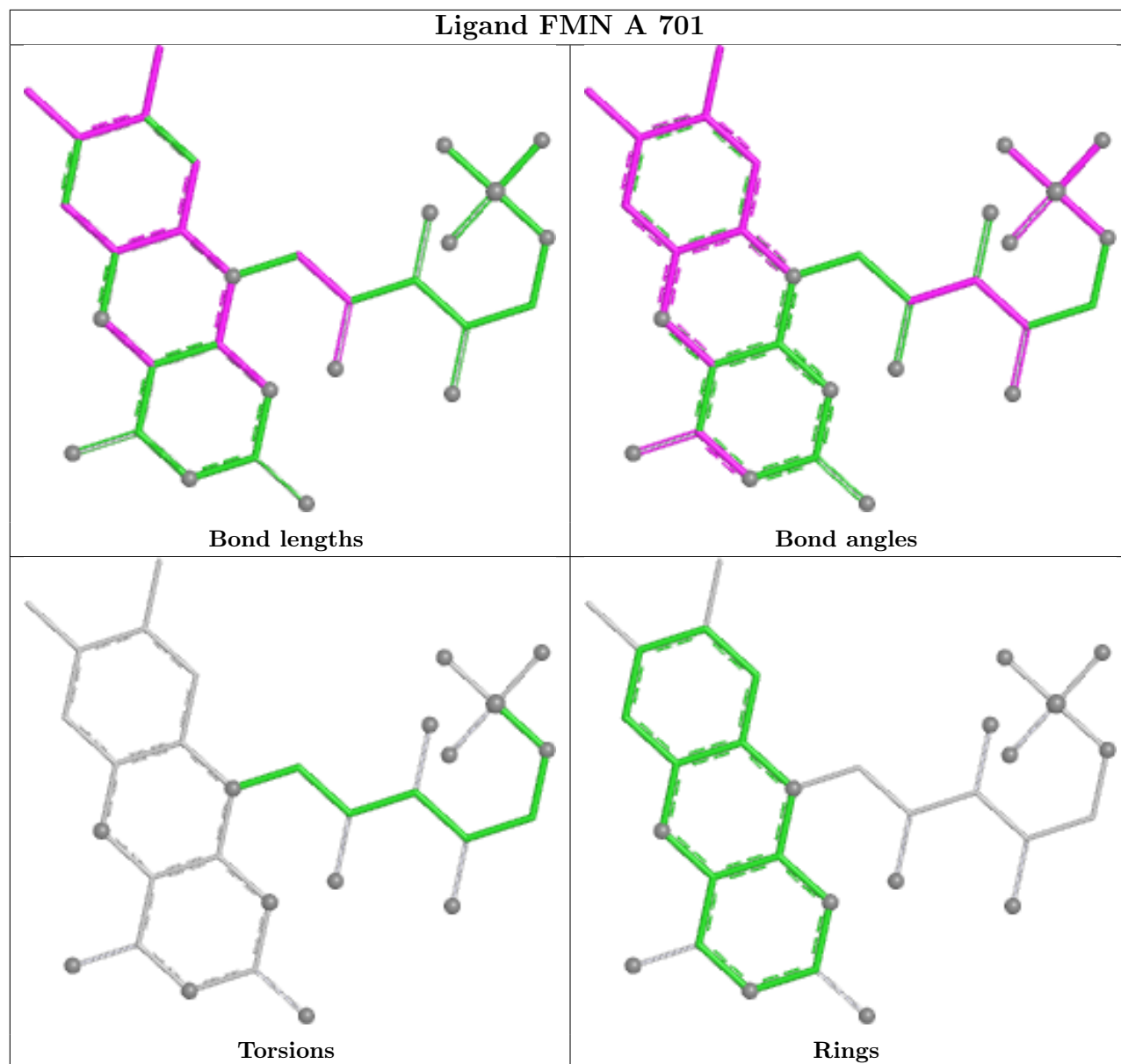
No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

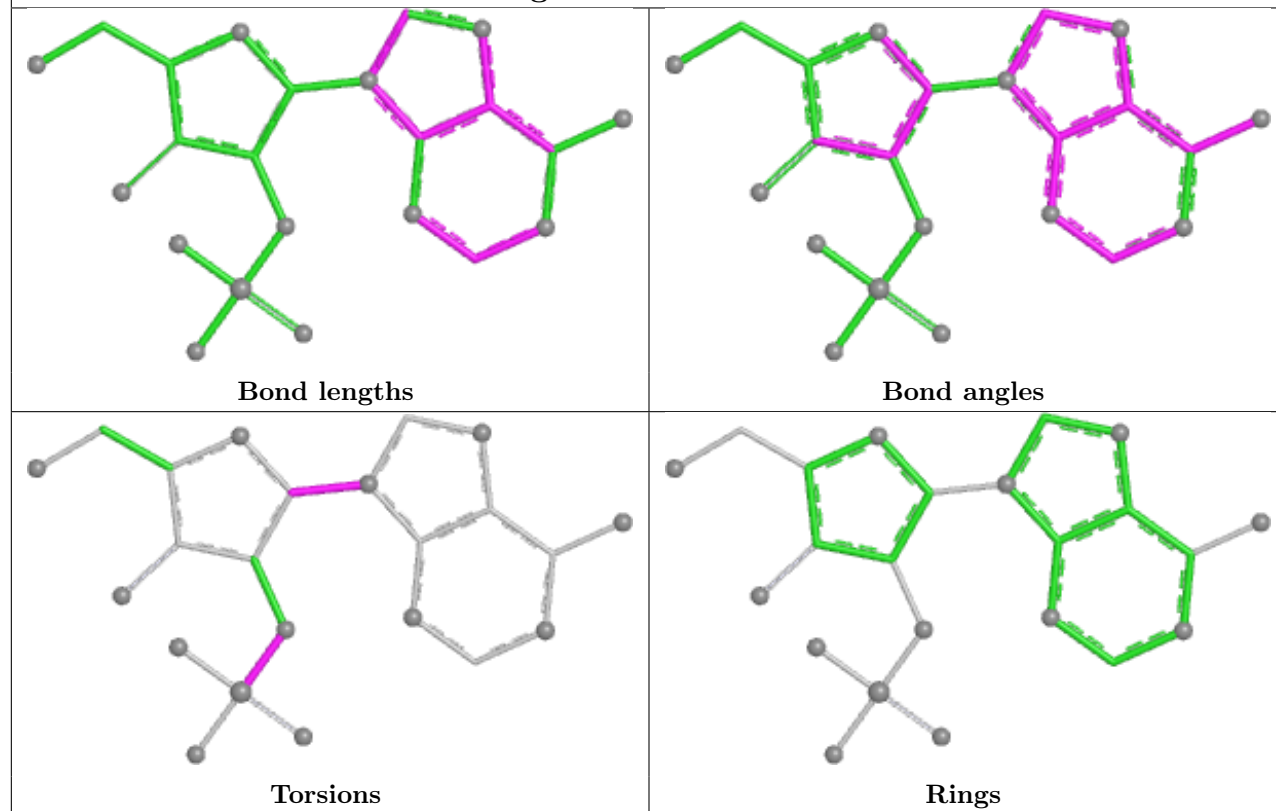




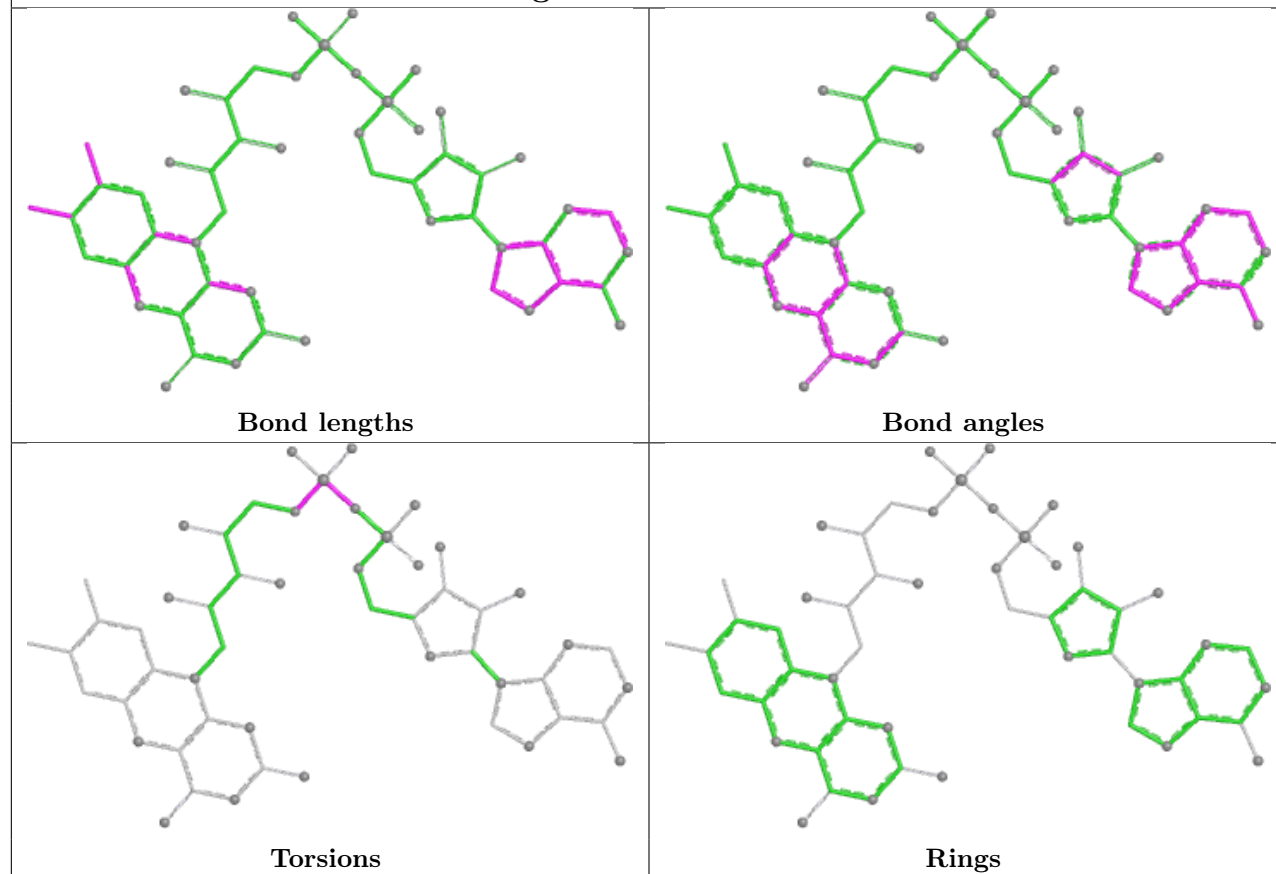




Ligand 2AM B 703



Ligand FAD A 702



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	608/622 (97%)	0.19	20 (3%) 49 54	16, 27, 41, 59	0
1	B	608/622 (97%)	0.67	62 (10%) 12 14	15, 32, 58, 72	0
All	All	1216/1244 (97%)	0.43	82 (6%) 24 26	15, 28, 54, 72	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	THR	5.6
1	B	233	VAL	4.9
1	B	118	ASP	4.6
1	B	200	ILE	4.6
1	A	106	GLY	4.4
1	B	227	VAL	4.3
1	B	252	ASP	4.0
1	A	678	SER	3.6
1	B	158	GLU	3.5
1	B	192	LEU	3.5
1	B	62	PRO	3.4
1	B	500	ALA	3.4
1	B	63	PRO	3.4
1	A	500	ALA	3.3
1	B	235	ALA	3.3
1	B	230	PHE	3.3
1	B	76	THR	3.3
1	B	64	VAL	3.3
1	B	195	LEU	3.3
1	B	102	ALA	3.2
1	A	252	ASP	3.2
1	B	242	ILE	3.0
1	B	111	SER	3.0
1	B	103	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	201	PHE	2.9
1	A	62	PRO	2.9
1	B	217	ILE	2.9
1	B	231	PHE	2.8
1	B	128	ILE	2.8
1	A	229	GLU	2.8
1	A	634	ARG	2.8
1	B	356	ASN	2.7
1	B	75	LYS	2.7
1	B	85	GLY	2.7
1	A	296	THR	2.7
1	A	235	ALA	2.7
1	B	505	GLY	2.7
1	B	125	LEU	2.7
1	B	197	ALA	2.6
1	A	294	GLN	2.6
1	B	668	THR	2.6
1	A	63	PRO	2.6
1	B	107	MET	2.6
1	B	163	LEU	2.6
1	B	106	GLY	2.6
1	B	169	ALA	2.6
1	B	226	ALA	2.5
1	B	126	PRO	2.5
1	A	158	GLU	2.5
1	A	251	GLU	2.5
1	B	234	GLU	2.5
1	B	678	SER	2.4
1	B	254	ASP	2.4
1	B	120	ALA	2.4
1	A	577	GLU	2.4
1	B	109	GLY	2.4
1	B	132	LEU	2.3
1	B	351	LEU	2.3
1	B	410	GLY	2.3
1	A	506	ARG	2.3
1	B	157	GLN	2.3
1	B	296	THR	2.3
1	B	382	ARG	2.3
1	B	161	VAL	2.3
1	B	251	GLU	2.3
1	A	266	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	105	TYR	2.3
1	B	108	ARG	2.3
1	B	228	CYS	2.3
1	B	212	LEU	2.3
1	A	653	PRO	2.2
1	B	196	GLY	2.2
1	B	112	ALA	2.2
1	A	271	ASN	2.1
1	B	135	PHE	2.1
1	B	80	ILE	2.1
1	B	155	TRP	2.1
1	B	403	HIS	2.1
1	B	83	PHE	2.1
1	A	570	ASP	2.1
1	A	151	ASP	2.0
1	B	166	VAL	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($\text{\AA}^2$)	Q<0.9
5	PO4	A	705	5/5	0.79	0.16	73,74,74,75	0
5	PO4	B	705	5/5	0.83	0.12	57,59,60,60	0
5	PO4	A	704	5/5	0.85	0.13	27,32,34,34	0
5	PO4	B	704	5/5	0.92	0.10	30,32,33,34	0
2	FMN	A	701	31/31	0.93	0.10	23,28,32,34	0
2	FMN	B	701	31/31	0.93	0.10	32,36,38,40	0
4	2AM	B	703	23/23	0.95	0.07	24,27,29,35	0

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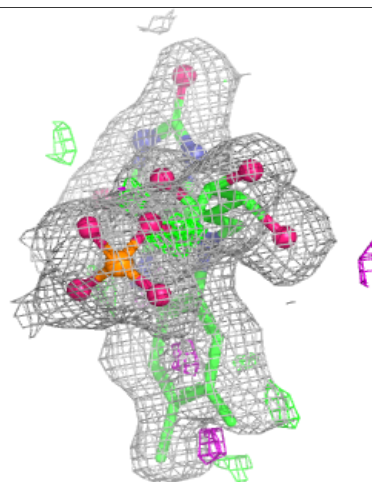
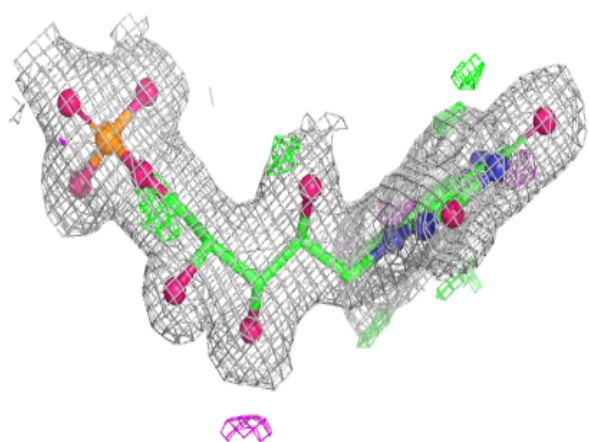
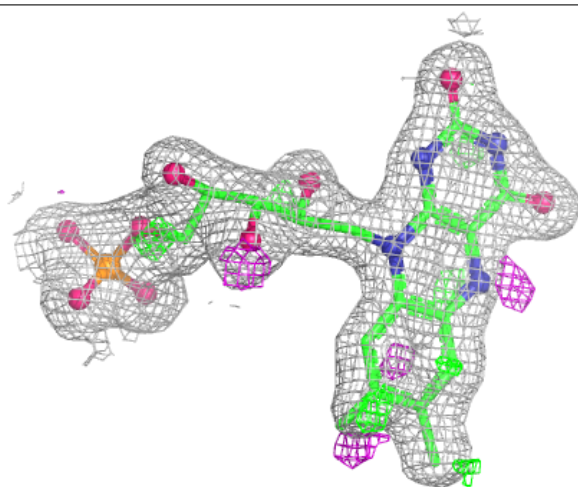
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	2AM	A	703	23/23	0.96	0.07	22,26,28,34	0
3	FAD	A	702	53/53	0.97	0.06	14,18,28,29	0
3	FAD	B	702	53/53	0.98	0.05	13,18,30,31	0

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.

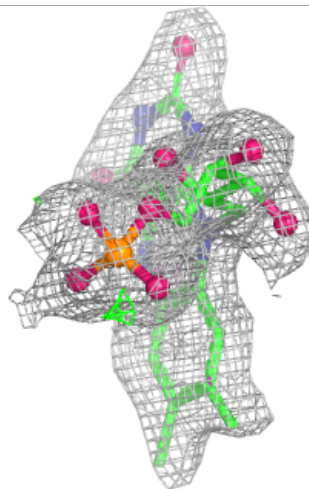
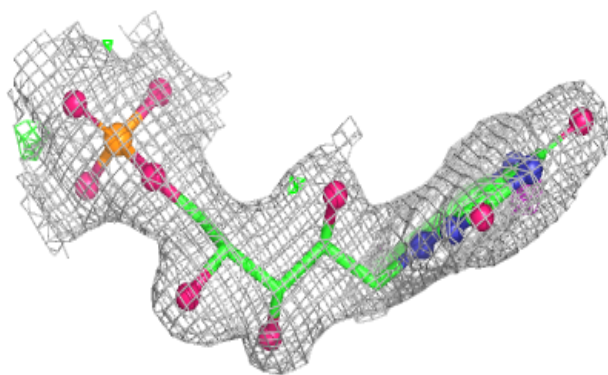
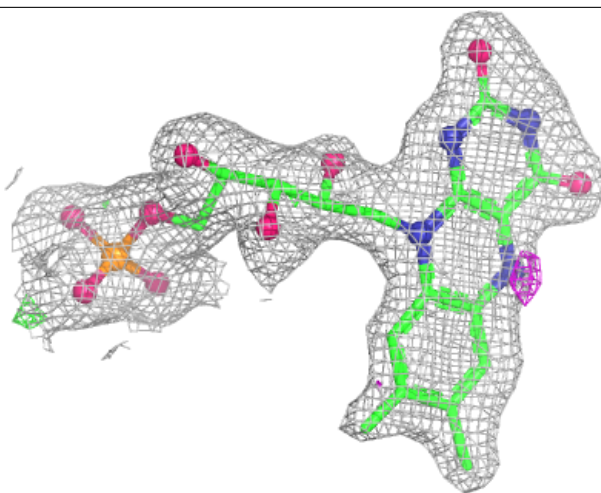
Electron density around FMN A 701:

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



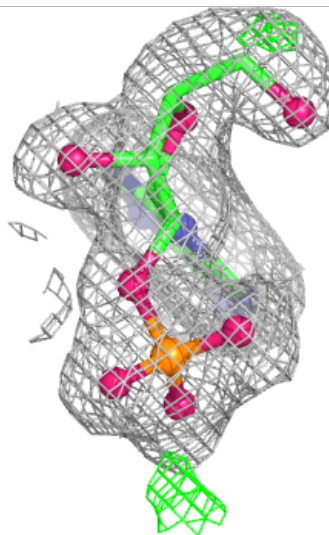
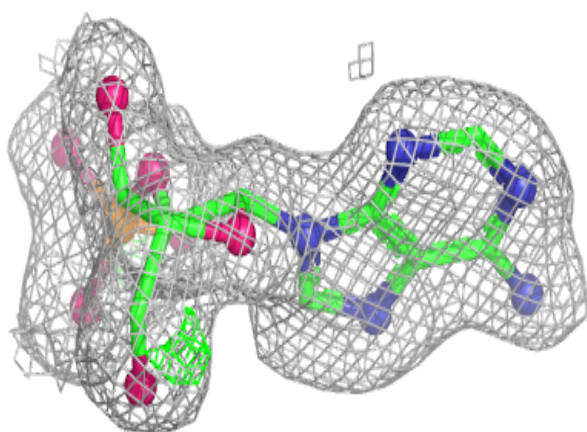
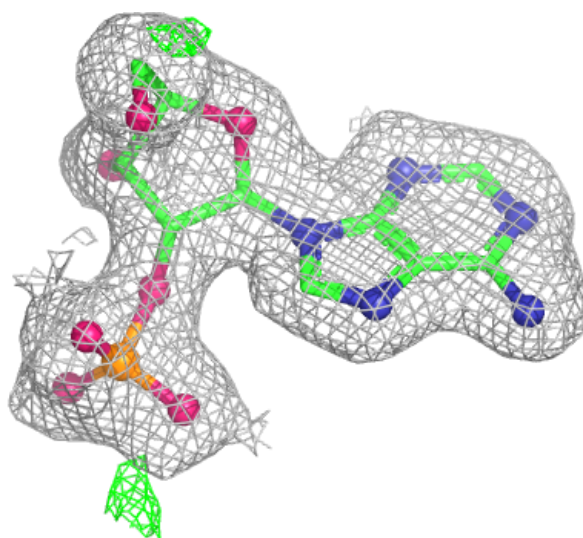
Electron density around FMN B 701:

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



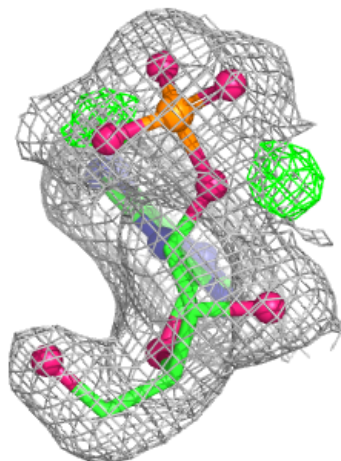
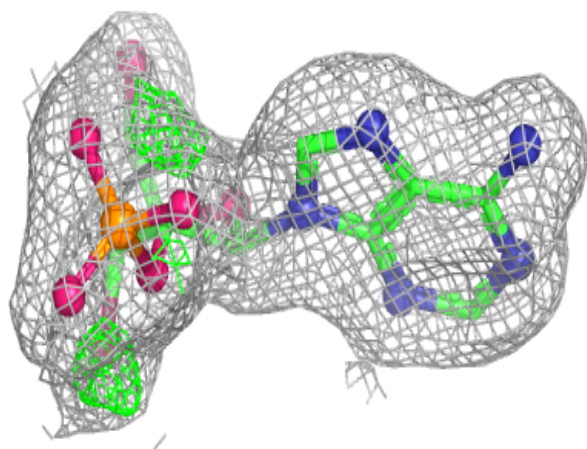
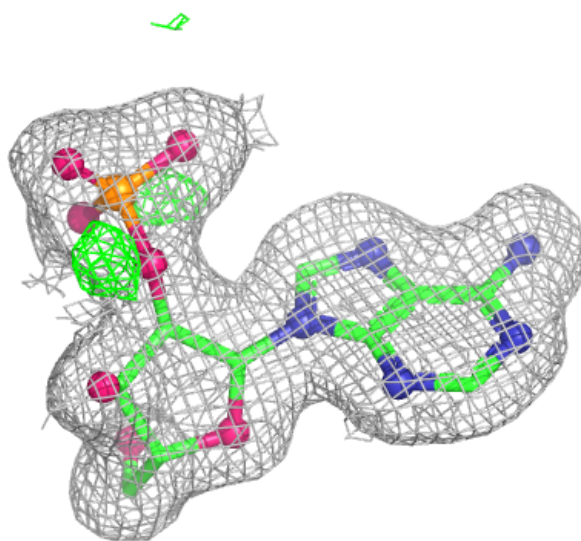
Electron density around 2AM B 703:

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



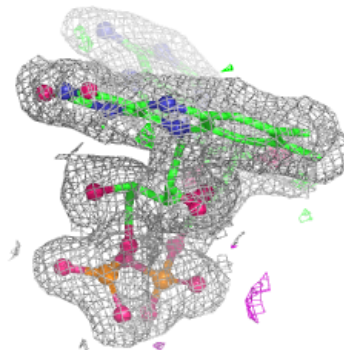
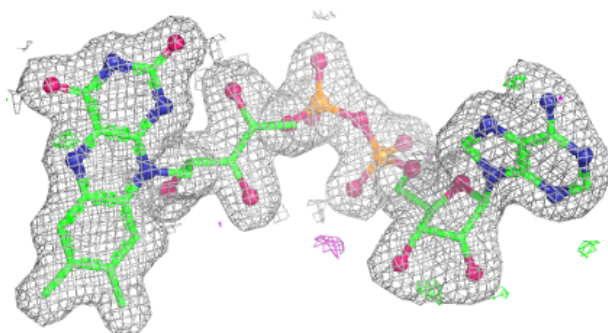
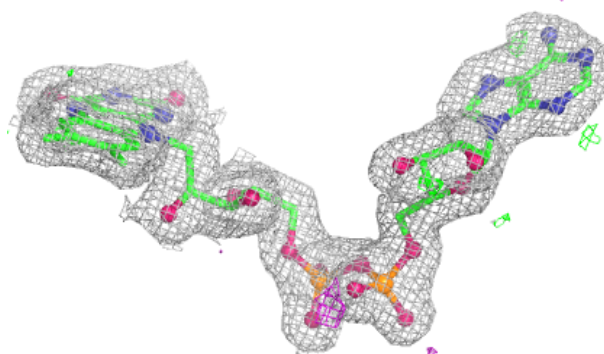
Electron density around 2AM A 703:

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

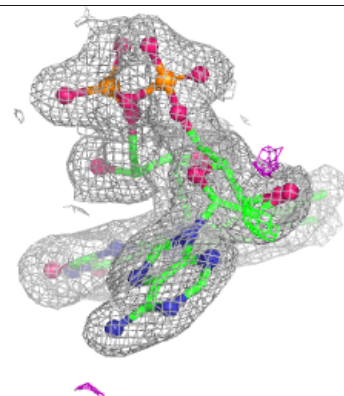
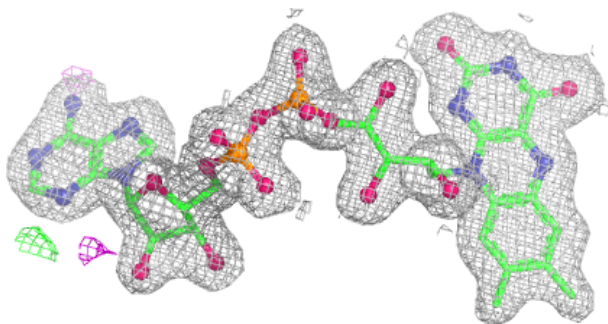
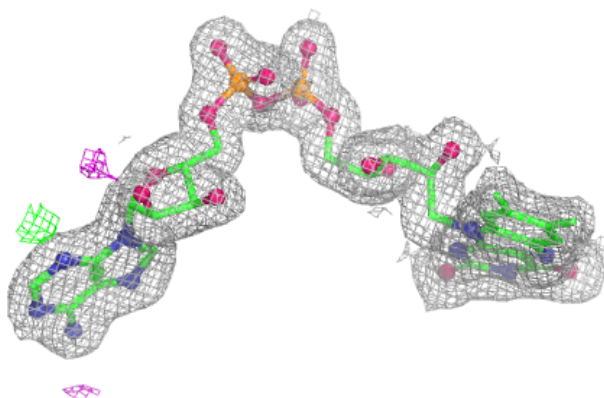


Electron density around FAD A 702:

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

**Electron density around FAD B 702:**

$2mF_o-DF_c$ (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.