



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:38 PM UTC

PDB ID : 4Y7C / pdb_00004y7c
Title : rat CYPOR mutant - G141del/E142N
Authors : Xia, C.; Kim, J.J.P.
Deposited on : 2015-02-13
Resolution : 2.20 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

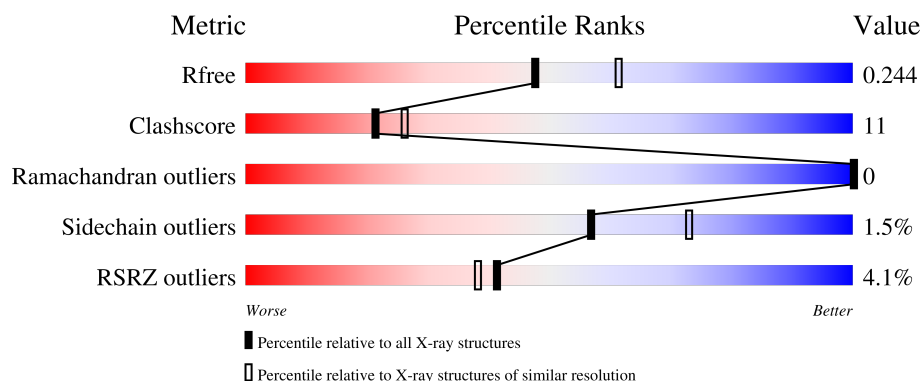
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	621	
1	B	621	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 11097 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.

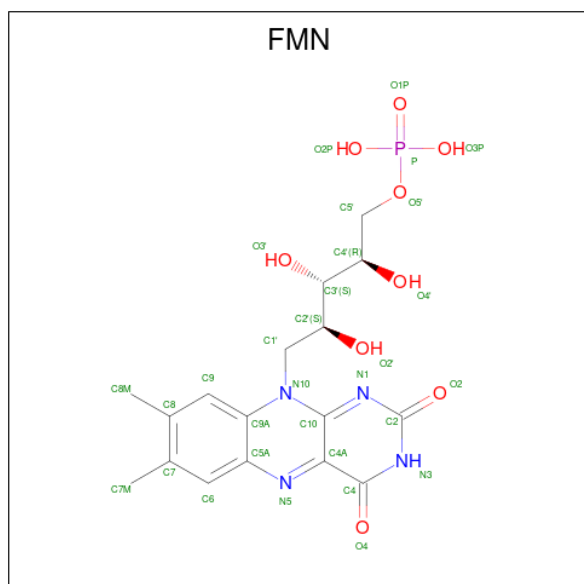
- Molecule 1 is a protein called NADPH-cytochrome P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	611	Total	C	N	O	S	0	0	0
			4895	3101	844	927	23			
1	B	605	Total	C	N	O	S	0	0	0
			4831	3059	836	913	23			

There are 4 discrepancies between the modelled and reference sequences:

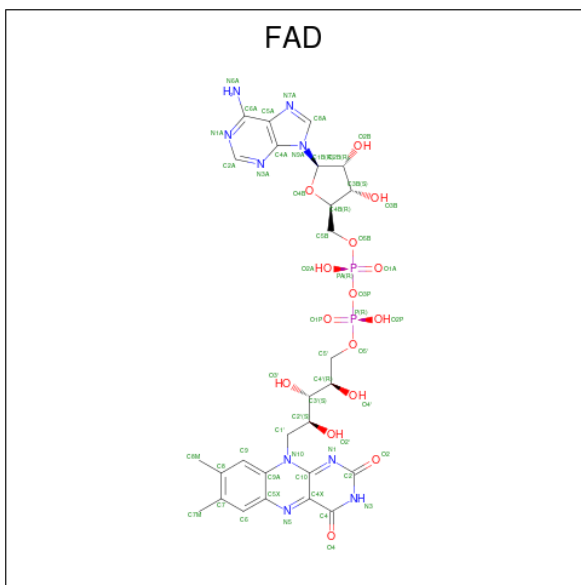
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	GLY	deletion	UNP P00388
A	142	ASN	GLU	engineered mutation	UNP P00388
B	?	-	GLY	deletion	UNP P00388
B	142	ASN	GLU	engineered mutation	UNP P00388

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



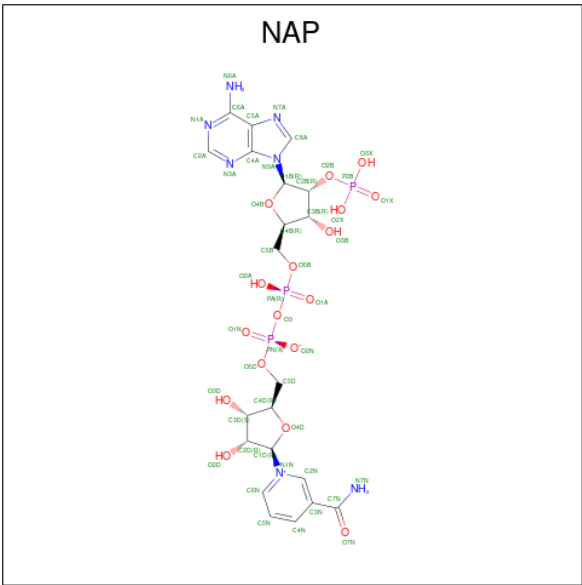
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 31	C 17	N 4	O 9	P 1	0	0
2	B	1	Total 31	C 17	N 4	O 9	P 1	0	0

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $\text{C}_{27}\text{H}_{33}\text{N}_9\text{O}_{15}\text{P}_2$).



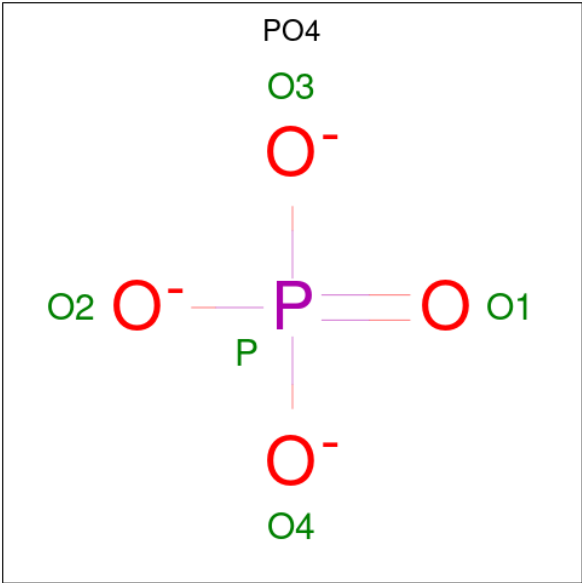
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
3	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $\text{C}_{21}\text{H}_{28}\text{N}_7\text{O}_{17}\text{P}_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
4	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	635	Total 635	O 635	0	0
6	B	501	Total 501	O 501	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.48Å 115.64Å 118.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.54 – 2.20 23.54 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.2 (23.54-2.20) 96.1 (23.54-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.19Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.192 , 0.244 0.192 , 0.244	Depositor DCC
R_{free} test set	3504 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11097	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.89% 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: PO4, FMN, NAP, FAD

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.40	0/5012	0.85	10/6782 (0.1%)
1	B	0.37	0/4946	0.86	8/6695 (0.1%)
All	All	0.38	0/9958	0.85	18/13477 (0.1%)

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	541	MET	N-CA-C	-7.20	103.51	111.36
1	B	596	SER	N-CA-C	6.97	119.73	111.71
1	B	89	GLY	N-CA-C	6.64	121.56	113.79
1	A	137	MET	N-CA-C	5.93	118.56	108.90
1	A	210	GLY	N-CA-C	-5.70	99.67	113.18
1	A	511	MET	N-CA-C	5.63	116.51	108.74
1	A	383	THR	N-CA-C	5.56	118.05	111.33
1	B	434	TYR	CA-C-N	5.51	125.86	119.47
1	B	434	TYR	C-N-CA	5.51	125.86	119.47
1	B	258	VAL	N-CA-C	5.46	115.97	107.99
1	A	596	SER	N-CA-C	5.37	118.93	112.38
1	B	263	MET	N-CA-C	5.34	117.19	111.36
1	A	346	MET	N-CA-C	5.29	115.52	108.38
1	B	297	GLU	N-CA-C	-5.27	104.95	111.33
1	B	485	VAL	N-CA-C	-5.22	99.74	107.51
1	A	482	SER	N-CA-C	-5.17	106.81	113.01
1	A	489	VAL	N-CA-C	5.12	115.27	110.30
1	A	160	ASP	N-CA-C	-5.04	105.22	113.19

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	4895	0	4739	72	0
1	B	4831	0	4642	141	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	31	0	11	0	0
4	B	31	0	11	0	0
5	B	5	0	0	0	0
6	A	635	0	0	18	0
6	B	501	0	0	16	0
All	All	11097	0	9503	213	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 11.

All (213) 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:81:ILE:HG23	1:B:133:VAL:HG23	1.42	0.97
1:B:155:TRP:O	1:B:159:THR:HG22	1.70	0.91
1:A:257:LYS:HA	1:A:266:LEU:HD21	1.54	0.90
1:B:191:ARG:HH11	1:B:195:LEU:HD11	1.35	0.89
1:B:70:VAL:HA	1:B:73:MET:HE3	1.59	0.83
1:B:86:SER:HB2	1:B:91:ALA:HB3	1.57	0.83
1:B:159:THR:OG1	1:B:161:VAL:HG13	1.81	0.81
1:B:167:LYS:HB3	1:B:200:ILE:HD11	1.62	0.81
1:B:155:TRP:O	1:B:159:THR:CG2	2.29	0.81
1:A:638:LYS:HA	1:A:638:LYS:NZ	1.99	0.77
1:A:550:LEU:HD22	1:A:555:LYS:HD2	1.67	0.77
1:A:633:ALA:HB2	1:A:676:VAL:HB	1.68	0.74
1:B:506:ARG:HB2	6:B:1074:HOH:O	1.88	0.72
1:B:214:GLU:OE1	1:B:217:ILE:HD12	1.90	0.71
1:B:81:ILE:CG2	1:B:133:VAL:HG23	2.20	0.70
1:B:222:GLN:O	1:B:225:PRO:HD2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:GLN:HB3	6:B:932:HOH:O	1.92	0.68
1:B:175:ASN:HB3	1:B:178:TYR:HD2	1.57	0.68
1:A:86:SER:HB2	1:A:91:ALA:HB3	1.73	0.68
1:A:240:SER:O	1:A:242:ILE:HG13	1.96	0.66
1:B:250:HIS:CE1	1:B:253:MET:HE3	2.30	0.65
1:B:145:PRO:HB3	1:B:184:MET:SD	2.37	0.65
1:A:506:ARG:HG2	1:A:506:ARG:HH11	1.61	0.65
1:B:107:MET:HE2	1:B:233:VAL:HG21	1.80	0.64
1:A:638:LYS:HA	1:A:638:LYS:HZ3	1.62	0.63
1:A:515:LYS:HD2	6:A:1338:HOH:O	1.99	0.63
1:B:527:PRO:HB2	1:B:625:ALA:HB2	1.81	0.63
1:A:400:GLU:OE2	1:A:404:LYS:HE3	1.99	0.63
1:A:555:LYS:HE3	6:A:1383:HOH:O	1.98	0.62
1:B:123:SER:HA	6:B:1226:HOH:O	1.98	0.62
1:A:254:ASP:O	1:A:257:LYS:HG2	1.98	0.62
1:B:222:GLN:C	1:B:225:PRO:HD2	2.25	0.62
1:B:119:LEU:HG	1:B:152:PHE:CD1	2.35	0.62
1:B:501:GLY:H	1:B:506:ARG:NH1	1.98	0.61
1:A:551:ARG:HG3	1:A:557:VAL:HG21	1.83	0.61
1:B:164:THR:HA	1:B:196:GLY:O	2.01	0.60
1:B:123:SER:O	1:B:126:PRO:HD2	2.02	0.60
1:B:188:VAL:HG12	1:B:192:LEU:HD11	1.84	0.60
1:B:506:ARG:HA	1:B:506:ARG:HE	1.65	0.60
1:B:175:ASN:HB3	1:B:178:TYR:CD2	2.37	0.60
1:A:459:ALA:HA	1:A:538:ALA:O	2.02	0.60
1:A:514:ARG:HD2	6:A:1127:HOH:O	2.02	0.58
1:A:551:ARG:HG3	1:A:557:VAL:CG2	2.33	0.58
1:B:79:ASN:HD21	1:B:107:MET:CG	2.15	0.58
1:B:660:VAL:HG12	1:B:664:LYS:HE2	1.85	0.58
1:B:175:ASN:OD1	1:B:209:ASP:HB2	2.04	0.58
1:B:626:HIS:CD2	1:B:671:ARG:HG2	2.38	0.58
1:A:156:LEU:HB3	1:A:191:ARG:HG2	1.85	0.58
1:B:233:VAL:HG12	1:B:234:GLU:N	2.19	0.58
1:B:163:LEU:HA	6:B:1226:HOH:O	2.03	0.58
1:B:279:LYS:HG3	6:B:1238:HOH:O	2.03	0.58
1:B:105:TYR:O	1:B:233:VAL:HG11	2.03	0.57
1:B:581:ARG:NH1	6:B:802:HOH:O	2.36	0.57
1:B:120:ALA:HA	1:B:155:TRP:CZ2	2.39	0.57
1:B:475:ALA:HA	1:B:491:THR:HB	1.86	0.57
1:B:506:ARG:HA	1:B:506:ARG:NE	2.19	0.57
1:B:86:SER:CB	1:B:91:ALA:HB3	2.31	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:396:PRO:HG2	6:B:863:HOH:O	2.06	0.56
1:B:266:LEU:O	1:B:267:LYS:HB2	2.06	0.56
1:A:638:LYS:HA	1:A:638:LYS:HZ2	1.71	0.55
1:A:506:ARG:N	1:A:506:ARG:HD3	2.21	0.55
1:B:353:GLU:H	1:B:353:GLU:CD	2.14	0.55
1:A:277:ASP:HB2	6:A:1196:HOH:O	2.06	0.55
1:B:217:ILE:HD11	1:B:383:THR:HG21	1.89	0.55
1:B:633:ALA:HB2	1:B:676:VAL:HB	1.88	0.55
1:B:203:LEU:HB3	6:B:871:HOH:O	2.07	0.54
1:A:300:LEU:HD22	1:A:574:LEU:HD21	1.90	0.54
1:B:223:PHE:O	1:B:227:VAL:HG23	2.07	0.54
1:B:182:ASN:O	1:B:186:LYS:HG3	2.07	0.54
1:A:78:ARG:HE	1:A:108:ARG:NE	2.06	0.54
1:B:549:TRP:O	1:B:553:GLN:NE2	2.38	0.53
1:B:548:ALA:O	1:B:552:GLU:HG2	2.08	0.53
1:B:255:VAL:HG23	6:B:1032:HOH:O	2.07	0.53
1:B:93:GLU:O	1:B:97:ARG:HG3	2.08	0.53
1:A:634:ARG:HD2	6:A:1152:HOH:O	2.09	0.53
1:B:130:LYS:HA	1:B:231:PHE:CD1	2.44	0.53
1:B:390:ALA:O	1:B:399:GLN:HG3	2.10	0.52
1:A:257:LYS:NZ	1:A:257:LYS:HB3	2.26	0.51
1:B:479:GLU:HG2	6:B:1001:HOH:O	2.10	0.51
1:B:170:VAL:O	1:B:203:LEU:HD12	2.10	0.51
1:A:581:ARG:NH1	1:A:585:ASP:OD2	2.43	0.51
1:B:385:VAL:HG13	1:B:447:LEU:HB3	1.93	0.51
1:B:73:MET:HB3	1:B:78:ARG:HB2	1.93	0.51
1:B:126:PRO:C	1:B:128:ILE:H	2.19	0.51
1:A:68:SER:HA	6:A:904:HOH:O	2.10	0.50
1:A:228:CYS:HA	1:A:233:VAL:CG2	2.41	0.50
1:B:167:LYS:HB3	1:B:200:ILE:CD1	2.37	0.50
1:B:103:HIS:C	1:B:105:TYR:H	2.20	0.50
1:B:214:GLU:OE1	1:B:214:GLU:HA	2.12	0.50
1:B:300:LEU:HD22	1:B:574:LEU:HD11	1.94	0.50
1:A:267:LYS:HA	6:A:1203:HOH:O	2.12	0.50
1:B:98:LEU:O	1:B:224:TRP:HZ2	1.95	0.50
1:B:159:THR:O	1:B:191:ARG:NH2	2.45	0.49
1:A:96:ASN:O	1:A:100:LYS:HG3	2.13	0.49
1:B:402:LEU:HD11	1:B:437:LEU:HD22	1.94	0.49
1:A:76:THR:OG1	1:A:78:ARG:HD3	2.13	0.49
1:A:321:ALA:HB2	1:A:455:TYR:CE1	2.48	0.49
1:B:195:LEU:HD12	1:B:195:LEU:N	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:214:GLU:CD	1:B:413:LYS:HE3	2.37	0.49
1:A:595:PHE:HB3	1:A:598:GLU:HG3	1.95	0.48
1:B:70:VAL:HG23	1:B:71:GLU:N	2.28	0.48
1:B:233:VAL:HG13	6:B:1126:HOH:O	2.11	0.48
1:A:84:TYR:CE2	1:A:92:GLU:HG3	2.48	0.48
1:A:119:LEU:HG	1:A:152:PHE:CD1	2.49	0.48
1:B:250:HIS:ND1	1:B:253:MET:HE3	2.29	0.48
1:B:119:LEU:O	1:B:122:LEU:HG	2.14	0.48
1:A:132:LEU:HD12	1:A:167:LYS:O	2.14	0.48
1:A:228:CYS:HA	1:A:233:VAL:HG22	1.96	0.48
1:B:84:TYR:HB3	1:B:95:ALA:HB2	1.95	0.48
1:B:135:PHE:HB2	1:B:170:VAL:HG13	1.96	0.48
1:B:242:ILE:CB	6:B:844:HOH:O	2.61	0.48
1:B:568:ARG:HB3	6:B:1124:HOH:O	2.14	0.48
1:A:262:GLU:HG2	6:A:1203:HOH:O	2.14	0.47
1:B:82:VAL:HG13	1:B:134:VAL:HB	1.96	0.47
1:A:80:ILE:O	1:A:109:GLY:HA2	2.15	0.47
1:B:139:THR:HG22	1:B:140:TYR:N	2.29	0.47
1:B:187:TYR:O	1:B:191:ARG:HB2	2.15	0.47
1:A:244:GLN:HG2	6:A:972:HOH:O	2.15	0.47
1:B:115:GLU:OE2	1:B:148:ASN:HA	2.15	0.47
1:B:660:VAL:CG1	1:B:664:LYS:HE2	2.44	0.47
1:A:584:LYS:HD3	6:A:835:HOH:O	2.14	0.46
1:B:69:PHE:HE1	1:B:121:ASP:HB2	1.80	0.46
1:B:169:ALA:HA	6:B:833:HOH:O	2.14	0.46
1:B:247:LEU:HD12	1:B:248:VAL:H	1.80	0.46
1:A:155:TRP:CZ2	1:A:161:VAL:HG11	2.50	0.46
1:B:80:ILE:HD11	1:B:107:MET:SD	2.55	0.46
1:B:246:GLU:HB2	1:B:351:LEU:HD21	1.97	0.46
1:A:629:VAL:HB	1:A:674:LEU:HD23	1.98	0.46
1:A:222:GLN:C	1:A:225:PRO:HD2	2.41	0.45
1:A:272:GLN:HA	6:A:937:HOH:O	2.16	0.45
1:B:69:PHE:CE1	1:B:121:ASP:HB2	2.52	0.45
1:B:133:VAL:O	1:B:133:VAL:CG1	2.65	0.45
1:B:656:HIS:O	1:B:660:VAL:HG23	2.15	0.45
1:A:78:ARG:NH2	1:A:352:ASP:OD2	2.48	0.45
1:A:161:VAL:HG23	6:A:1303:HOH:O	2.15	0.45
1:A:479:GLU:HB3	6:A:1222:HOH:O	2.17	0.45
1:B:547:ARG:O	1:B:551:ARG:HG3	2.17	0.45
1:A:80:ILE:HG23	1:A:109:GLY:HA3	1.99	0.45
1:A:159:THR:OG1	1:A:161:VAL:HG12	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:561:LEU:HD22	1:B:590:GLN:HB2	1.98	0.44
1:B:73:MET:CB	1:B:78:ARG:HB2	2.47	0.44
1:A:349:ASN:HB3	6:A:1071:HOH:O	2.17	0.44
1:B:660:VAL:O	1:B:664:LYS:HG3	2.17	0.44
1:A:652:GLY:N	1:A:653:PRO:HA	2.33	0.44
1:B:152:PHE:CE2	1:B:156:LEU:HD11	2.53	0.44
1:A:411:GLU:HG3	1:A:412:GLY:N	2.32	0.44
1:B:130:LYS:HA	1:B:231:PHE:HD1	1.83	0.44
1:A:184:MET:HE3	1:A:188:VAL:CG2	2.47	0.44
1:B:79:ASN:HD21	1:B:107:MET:HG2	1.83	0.44
1:B:213:GLU:O	1:B:217:ILE:HG13	2.18	0.44
1:B:152:PHE:O	1:B:156:LEU:HG	2.17	0.43
1:B:246:GLU:HB2	1:B:351:LEU:CD2	2.48	0.43
1:A:664:LYS:HE2	6:A:1223:HOH:O	2.17	0.43
1:B:101:ASP:OD2	1:B:220:ARG:NH2	2.51	0.43
1:B:126:PRO:HG3	1:B:165:GLY:C	2.44	0.43
1:B:378:THR:O	1:B:426:HIS:HB3	2.18	0.43
1:A:568:ARG:NH1	6:A:811:HOH:O	2.50	0.43
1:A:176:LYS:C	1:A:178:TYR:H	2.27	0.43
1:B:69:PHE:HA	1:B:72:LYS:HB2	2.00	0.43
1:B:133:VAL:HG12	1:B:168:PHE:HB3	2.00	0.43
1:B:84:TYR:CG	1:B:92:GLU:HA	2.53	0.43
1:B:352:ASP:C	1:B:354:GLU:H	2.25	0.43
1:A:246:GLU:HB3	1:A:351:LEU:HD21	2.01	0.42
1:B:366:THR:HG23	6:B:993:HOH:O	2.18	0.42
1:B:563:TYR:CE2	1:B:609:LEU:HD23	2.54	0.42
1:B:76:THR:HG21	1:B:354:GLU:HG2	2.02	0.42
1:B:327:ASP:OD1	1:B:329:ALA:HB3	2.19	0.42
1:B:173:LEU:HD12	1:B:173:LEU:N	2.34	0.42
1:A:633:ALA:HB2	1:A:676:VAL:CB	2.45	0.42
1:B:81:ILE:HD12	1:B:110:MET:O	2.20	0.42
1:B:349:ASN:OD1	1:B:359:HIS:CE1	2.72	0.42
1:B:188:VAL:CG1	1:B:192:LEU:HD11	2.47	0.42
1:B:224:TRP:HB2	1:B:225:PRO:HD3	2.02	0.42
1:B:233:VAL:HG12	1:B:234:GLU:H	1.82	0.42
1:A:506:ARG:HG2	1:A:506:ARG:NH1	2.29	0.42
1:B:81:ILE:HG23	1:B:81:ILE:O	2.20	0.42
1:B:395:GLU:HA	1:B:396:PRO:HD2	1.86	0.42
1:B:448:LEU:HA	1:B:449:PRO:HD3	1.92	0.42
1:A:253:MET:SD	1:A:364:PRO:HG3	2.59	0.42
1:B:186:LYS:O	1:B:190:GLN:NE2	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:VAL:O	1:B:389:LEU:HG	2.20	0.41
1:B:81:ILE:HD13	1:B:110:MET:HE2	2.02	0.41
1:B:214:GLU:OE2	1:B:413:LYS:HG3	2.20	0.41
1:B:672:TYR:HE1	1:B:674:LEU:HD21	1.85	0.41
1:A:374:TYR:HA	6:A:1028:HOH:O	2.21	0.41
1:A:519:ARG:HH11	1:A:519:ARG:HG2	1.85	0.41
1:A:559:GLU:HB3	1:A:561:LEU:CD1	2.51	0.41
1:B:529:ILE:HD13	1:B:616:LEU:HD22	2.02	0.41
1:A:617:TRP:CD1	1:A:651:PHE:HB3	2.56	0.41
1:A:651:PHE:O	1:A:652:GLY:C	2.62	0.41
1:B:376:ASP:OD2	1:B:379:ASN:ND2	2.48	0.41
1:A:564:TYR:CG	1:A:565:GLY:N	2.88	0.41
1:B:119:LEU:C	1:B:121:ASP:H	2.28	0.41
1:B:126:PRO:C	1:B:128:ILE:N	2.79	0.41
1:B:191:ARG:CD	1:B:195:LEU:HD11	2.51	0.41
1:B:84:TYR:CD2	1:B:92:GLU:HB2	2.56	0.41
1:A:409:SER:HA	1:A:413:LYS:HD3	2.03	0.41
1:B:159:THR:HG1	1:B:161:VAL:HG13	1.83	0.41
1:B:184:MET:O	1:B:188:VAL:HG23	2.20	0.41
1:B:523:LYS:NZ	6:B:828:HOH:O	2.54	0.41
1:A:155:TRP:HZ2	1:A:161:VAL:HG11	1.86	0.41
1:A:254:ASP:HB2	1:A:257:LYS:CD	2.51	0.41
1:A:68:SER:HB2	1:A:121:ASP:OD2	2.21	0.40
1:A:69:PHE:HB3	1:A:117:TYR:CD1	2.56	0.40
1:B:119:LEU:C	1:B:121:ASP:N	2.79	0.40
1:B:133:VAL:O	1:B:133:VAL:HG13	2.20	0.40
1:B:359:HIS:HB3	1:B:360:PRO:HD2	2.03	0.40
1:B:526:THR:HA	1:B:527:PRO:HD3	1.92	0.40
1:A:581:ARG:NE	6:A:835:HOH:O	2.55	0.40
1:B:133:VAL:O	1:B:168:PHE:HA	2.22	0.40
1:B:433:ASP:C	1:B:435:PRO:HD3	2.46	0.40
1:B:233:VAL:CG1	1:B:234:GLU:N	2.85	0.40
1:B:79:ASN:ND2	1:B:107:MET:HG2	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	605/621 (97%)	586 (97%)	19 (3%)	0	100	100
1	B	599/621 (96%)	566 (94%)	33 (6%)	0	100	100
All	All	1204/1242 (97%)	1152 (96%)	52 (4%)	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	522/531 (98%)	513 (98%)	9 (2%)	53	69
1	B	508/531 (96%)	502 (99%)	6 (1%)	63	78
All	All	1030/1062 (97%)	1015 (98%)	15 (2%)	57	73

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	68	SER
1	A	88	THR
1	A	179	GLU
1	A	234	GLU
1	A	257	LYS
1	A	411	GLU
1	A	484	ARG
1	A	498	GLU

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Mol	Chain	Res	Type
1	A	638	LYS
1	B	192	LEU
1	B	251	GLU
1	B	253	MET
1	B	263	MET
1	B	356	ASN
1	B	553	GLN

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	142	ASN
1	A	222	GLN
1	A	333	GLN
1	A	359	HIS
1	A	465	HIS
1	A	599	GLN
1	A	626	HIS
1	A	635	ASN
1	B	157	GLN
1	B	182	ASN
1	B	359	HIS
1	B	384	ASN
1	B	391	GLN
1	B	399	GLN
1	B	401	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

7 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	701	-	33,33,33	2.63	12 (36%)	48,50,50	1.65	12 (25%)
4	NAP	B	703	-	32,33,52	1.63	7 (21%)	50,52,80	1.91	13 (26%)
5	PO4	B	704	-	4,4,4	1.68	0	6,6,6	0.47	0
2	FMN	A	701	-	33,33,33	2.67	14 (42%)	48,50,50	1.59	12 (25%)
3	FAD	A	702	-	58,58,58	2.31	13 (22%)	85,89,89	1.95	18 (21%)
4	NAP	A	703	-	32,33,52	1.63	7 (21%)	50,52,80	1.92	13 (26%)
3	FAD	B	702	-	58,58,58	2.30	13 (22%)	85,89,89	1.95	17 (20%)

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	701	-	-	0/18/18/18	0/3/3/3
4	NAP	B	703	-	-	3/21/37/67	0/3/3/5
2	FMN	A	701	-	-	0/18/18/18	0/3/3/3
3	FAD	A	702	-	-	1/34/50/50	0/6/6/6
4	NAP	A	703	-	-	4/21/37/67	0/3/3/5
3	FAD	B	702	-	-	2/34/50/50	0/6/6/6

All (66) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	FAD	C8A-N7A	8.57	1.48	1.31
3	B	702	FAD	C8A-N7A	8.56	1.48	1.31
2	B	701	FMN	C8M-C8	-6.92	1.38	1.51
2	A	701	FMN	C8M-C8	-6.56	1.38	1.51
3	A	702	FAD	C8A-N9A	6.21	1.48	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	FAD	C8M-C8	-6.18	1.39	1.51
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	A	702	FAD	C7M-C7	-6.12	1.39	1.51
3	B	702	FAD	C7M-C7	-6.12	1.39	1.51
2	A	701	FMN	C4A-N5	5.35	1.42	1.30
2	B	701	FMN	C4A-N5	5.31	1.42	1.30
2	A	701	FMN	C10-N10	5.08	1.48	1.37
2	B	701	FMN	C10-N10	4.97	1.48	1.37
2	B	701	FMN	C9-C9A	4.87	1.47	1.39
2	A	701	FMN	C9-C9A	4.70	1.47	1.39
2	A	701	FMN	C9A-C5A	4.20	1.47	1.41
2	B	701	FMN	C9A-N10	4.13	1.48	1.41
2	B	701	FMN	C9A-C5A	4.06	1.47	1.41
2	A	701	FMN	C8-C7	3.94	1.50	1.40
2	A	701	FMN	C9A-N10	3.91	1.47	1.41
3	A	702	FAD	C9A-N10	-3.77	1.34	1.41
2	B	701	FMN	C8-C7	3.77	1.50	1.40
3	B	702	FAD	C9A-N10	-3.77	1.34	1.41
3	B	702	FAD	C2A-N3A	3.66	1.40	1.33
4	A	703	NAP	C2A-N3A	3.63	1.40	1.33
4	A	703	NAP	C2A-N1A	3.60	1.40	1.33
3	B	702	FAD	C10-N1	3.60	1.40	1.33
4	B	703	NAP	C2A-N1A	3.59	1.40	1.33
3	A	702	FAD	C2A-N3A	3.58	1.40	1.33
4	B	703	NAP	C2A-N3A	3.58	1.40	1.33
3	A	702	FAD	C10-N1	3.57	1.40	1.33
3	B	702	FAD	C2A-N1A	3.55	1.40	1.33
3	A	702	FAD	C2A-N1A	3.54	1.40	1.33
2	A	701	FMN	C1'-C2'	3.33	1.57	1.52
2	A	701	FMN	C6-C5A	3.29	1.45	1.40
2	A	701	FMN	C10-N1	3.20	1.39	1.33
2	B	701	FMN	C10-N1	3.19	1.39	1.33
4	A	703	NAP	C5A-N7A	-3.09	1.33	1.39
4	B	703	NAP	C5A-N7A	-3.06	1.33	1.39
4	A	703	NAP	C5A-C4A	-3.05	1.33	1.39
4	B	703	NAP	C5A-C4A	-3.03	1.33	1.39
3	B	702	FAD	C5A-N7A	-3.01	1.33	1.39
3	A	702	FAD	C5A-N7A	-3.01	1.33	1.39
3	A	702	FAD	C5A-C4A	-2.92	1.33	1.39
3	B	702	FAD	C5A-C4A	-2.92	1.33	1.39
2	B	701	FMN	C6-C5A	2.89	1.44	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	702	FAD	C5X-N5	-2.72	1.34	1.39
3	B	702	FAD	C5X-N5	-2.70	1.34	1.39
3	A	702	FAD	C5A-C6A	-2.67	1.33	1.41
3	B	702	FAD	C5A-C6A	-2.66	1.33	1.41
4	B	703	NAP	C5A-C6A	-2.66	1.33	1.41
4	A	703	NAP	C5A-C6A	-2.64	1.33	1.41
2	A	701	FMN	O2'-C2'	2.56	1.48	1.43
2	B	701	FMN	C4'-C3'	2.55	1.57	1.53
2	B	701	FMN	O2'-C2'	2.49	1.48	1.43
2	A	701	FMN	C4'-C3'	2.49	1.57	1.53
4	B	703	NAP	C8A-N9A	-2.39	1.33	1.37
2	A	701	FMN	C7M-C7	2.38	1.55	1.51
4	A	703	NAP	C8A-N9A	-2.36	1.33	1.37
2	A	701	FMN	C6-C7	2.36	1.42	1.39
2	B	701	FMN	C7M-C7	2.07	1.54	1.51
3	A	702	FAD	C4A-N9A	-2.05	1.33	1.37
4	A	703	NAP	C4A-N9A	-2.05	1.33	1.37
4	B	703	NAP	C4A-N9A	-2.01	1.33	1.37
3	B	702	FAD	C4A-N9A	-2.00	1.33	1.37

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	FAD	N9A-C8A-N7A	-9.46	100.52	113.94
3	B	702	FAD	N9A-C8A-N7A	-9.43	100.55	113.94
3	A	702	FAD	N3A-C2A-N1A	-7.47	117.28	128.58
3	B	702	FAD	N3A-C2A-N1A	-7.35	117.45	128.58
4	A	703	NAP	N3A-C2A-N1A	-7.11	117.82	128.58
4	B	703	NAP	N3A-C2A-N1A	-7.09	117.85	128.58
3	B	702	FAD	C5A-C4A-N3A	-4.55	120.45	126.72
4	A	703	NAP	N9A-C8A-N7A	-4.39	107.71	113.94
3	A	702	FAD	C5A-C4A-N3A	-4.36	120.71	126.72
4	B	703	NAP	N9A-C8A-N7A	-4.34	107.78	113.94
4	B	703	NAP	C5A-C4A-N3A	-4.26	120.85	126.72
4	A	703	NAP	C5A-C4A-N3A	-4.20	120.94	126.72
3	B	702	FAD	C5A-C4A-N9A	4.18	110.37	105.81
3	A	702	FAD	C5A-C4A-N9A	4.00	110.17	105.81
2	B	701	FMN	C9A-C5A-N5	3.98	126.68	122.45
2	A	701	FMN	C9A-C5A-N5	3.93	126.62	122.45
2	A	701	FMN	O4'-C4'-C3'	-3.80	100.36	109.25
2	B	701	FMN	O4'-C4'-C3'	-3.76	100.44	109.25
3	B	702	FAD	C2A-N3A-C4A	3.56	120.52	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	FAD	C2A-N3A-C4A	3.51	120.41	111.83
3	B	702	FAD	C5A-N7A-C8A	3.42	108.83	103.45
3	A	702	FAD	C5A-N7A-C8A	3.41	108.81	103.45
4	B	703	NAP	C2A-N3A-C4A	3.30	119.88	111.83
4	A	703	NAP	C2A-N3A-C4A	3.28	119.84	111.83
4	B	703	NAP	C5A-N7A-C8A	3.15	108.39	103.45
4	A	703	NAP	C5A-N7A-C8A	3.11	108.34	103.45
3	A	702	FAD	C4-N3-C2	-3.10	120.13	125.64
3	B	702	FAD	C4-N3-C2	-3.07	120.18	125.64
3	A	702	FAD	C4A-N9A-C8A	2.90	108.78	105.74
3	B	702	FAD	C5A-C6A-N6A	-2.89	116.14	123.29
3	A	702	FAD	C5A-C6A-N6A	-2.89	116.14	123.29
2	B	701	FMN	O4-C4-N3	-2.83	114.78	120.11
3	B	702	FAD	C4A-N9A-C8A	2.83	108.71	105.74
2	B	701	FMN	C5A-C9A-N10	-2.76	115.48	117.97
3	B	702	FAD	C10-C4X-N5	-2.76	119.18	124.81
2	A	701	FMN	O4-C4-N3	-2.72	115.00	120.11
3	B	702	FAD	C2B-C3B-C4B	-2.68	97.44	102.61
2	A	701	FMN	C8M-C8-C7	2.65	126.17	120.76
3	A	702	FAD	C4X-C4-N3	2.64	119.97	113.25
3	A	702	FAD	C10-C4X-N5	-2.63	119.45	124.81
4	A	703	NAP	C5A-C6A-N6A	-2.61	116.81	123.29
2	B	701	FMN	C8M-C8-C7	2.59	126.05	120.76
4	B	703	NAP	C4A-C5A-N7A	-2.59	107.62	110.58
2	B	701	FMN	C4'-C3'-C2'	-2.59	109.27	113.57
4	B	703	NAP	C5A-C6A-N6A	-2.58	116.91	123.29
2	B	701	FMN	O3P-P-O5'	-2.57	99.96	106.67
2	A	701	FMN	C5A-C9A-N10	-2.53	115.68	117.97
3	B	702	FAD	C4X-C4-N3	2.52	119.67	113.25
4	B	703	NAP	C6A-C5A-C4A	2.49	120.58	117.18
4	A	703	NAP	C4A-C5A-N7A	-2.49	107.73	110.58
2	A	701	FMN	C5A-N5-C4A	-2.49	114.06	118.09
2	B	701	FMN	C5A-N5-C4A	-2.49	114.07	118.09
2	A	701	FMN	C4'-C3'-C2'	-2.47	109.46	113.57
4	A	703	NAP	C6A-C5A-C4A	2.45	120.52	117.18
3	A	702	FAD	O4-C4-C4X	-2.37	120.28	126.53
2	A	701	FMN	C8M-C8-C9	-2.35	115.44	119.57
3	B	702	FAD	C6A-C5A-C4A	2.33	120.36	117.18
4	A	703	NAP	C3B-C2B-C1B	-2.33	98.35	102.81
2	B	701	FMN	O2P-P-O1P	2.32	119.88	110.83
2	A	701	FMN	O2P-P-O1P	2.30	119.80	110.83
2	B	701	FMN	C8M-C8-C9	-2.29	115.54	119.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	703	NAP	C4A-N9A-C8A	2.29	108.14	105.74
3	A	702	FAD	C9A-C5X-N5	-2.28	120.04	122.45
4	B	703	NAP	C3B-C2B-C1B	-2.28	98.45	102.81
2	A	701	FMN	O3P-P-O5'	-2.27	100.75	106.67
4	A	703	NAP	N3A-C4A-N9A	2.25	130.99	127.17
3	B	702	FAD	C9A-C5X-N5	-2.25	120.07	122.45
4	B	703	NAP	N3A-C4A-N9A	2.23	130.95	127.17
4	B	703	NAP	C4A-N9A-C8A	2.21	108.06	105.74
3	A	702	FAD	C6A-C5A-C4A	2.20	120.18	117.18
3	B	702	FAD	O4-C4-C4X	-2.19	120.76	126.53
3	A	702	FAD	C2B-C3B-C4B	-2.18	98.39	102.61
2	B	701	FMN	C9-C9A-N10	2.18	124.78	121.85
2	B	701	FMN	C7M-C7-C6	-2.17	115.75	119.57
4	B	703	NAP	C5A-C4A-N9A	2.14	108.14	105.81
3	A	702	FAD	C4X-C10-N10	2.12	119.52	116.48
3	B	702	FAD	C4-C4X-C10	2.12	120.56	116.93
4	B	703	NAP	O2N-PN-O1N	2.11	119.06	110.83
4	A	703	NAP	O2N-PN-O1N	2.10	119.00	110.83
2	A	701	FMN	C7M-C7-C6	-2.08	115.91	119.57
3	B	702	FAD	C4X-C10-N10	2.06	119.44	116.48
2	A	701	FMN	O4'-C4'-C5'	-2.06	105.44	109.99
3	A	702	FAD	C5X-C9A-N10	2.05	119.83	117.97
4	A	703	NAP	C5A-C4A-N9A	2.03	108.02	105.81
3	A	702	FAD	C4'-C3'-C2'	-2.02	110.21	113.57

There are no chirality outliers.

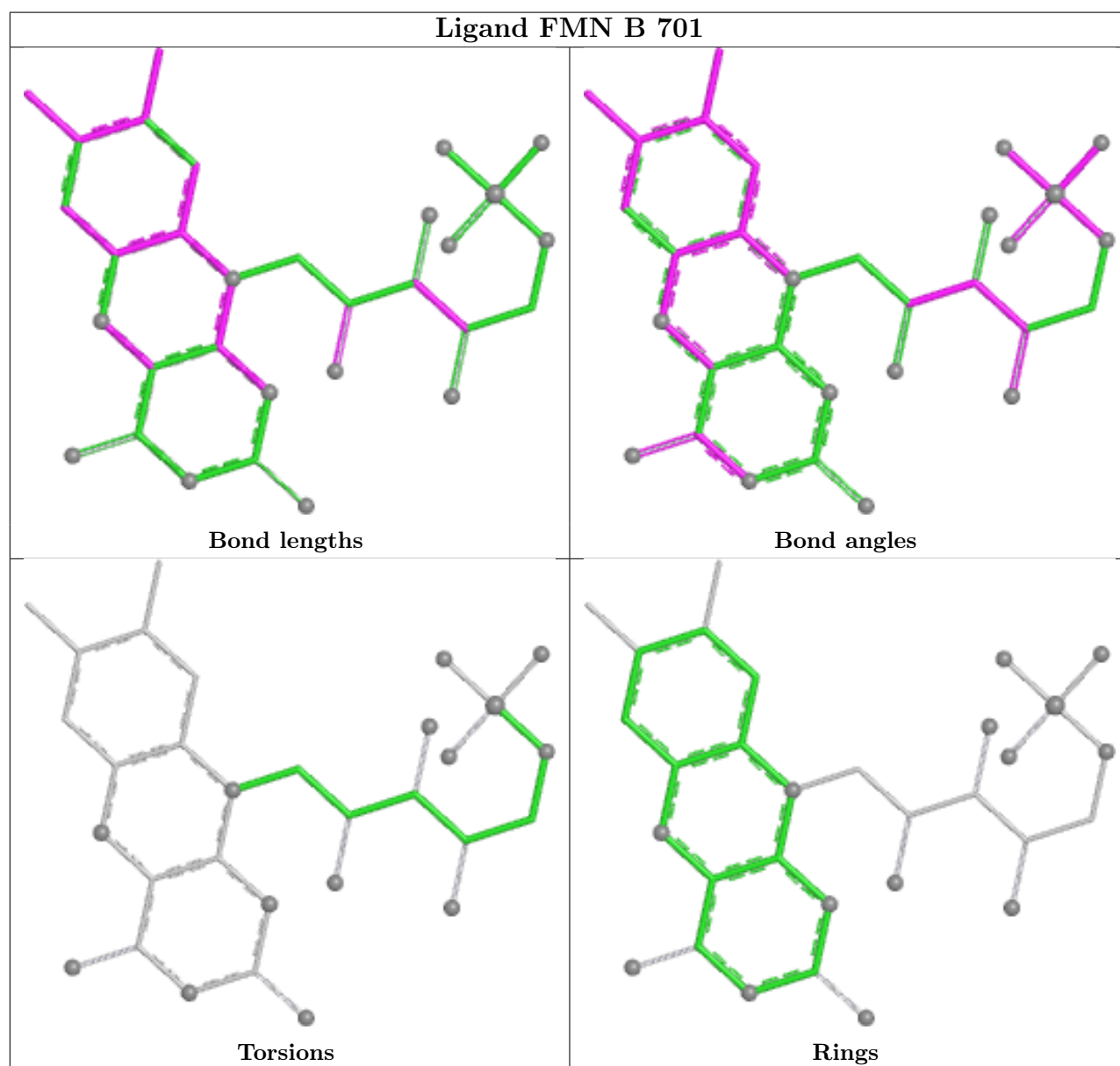
All (10) torsion outliers are listed below:

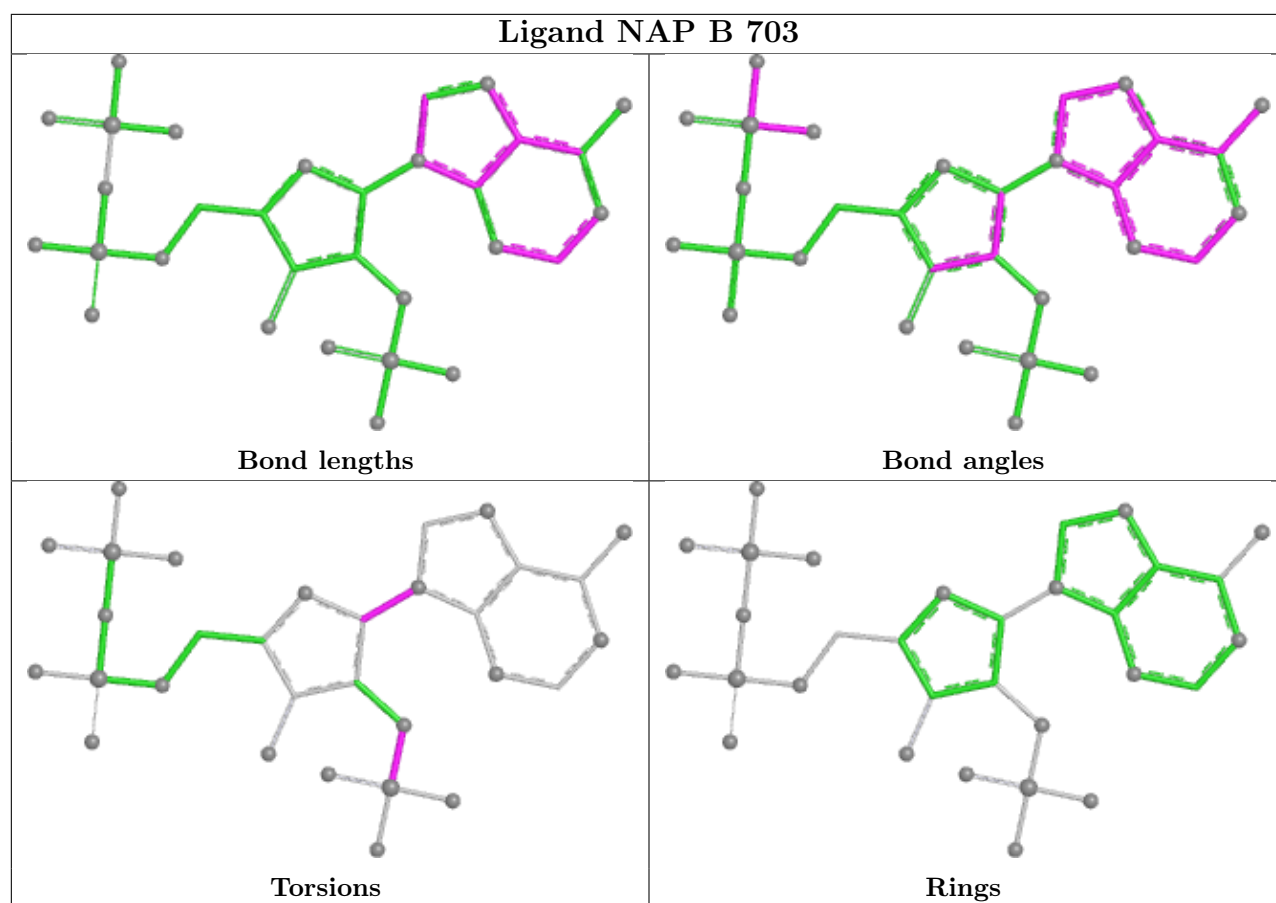
Mol	Chain	Res	Type	Atoms
4	A	703	NAP	C2B-C1B-N9A-C8A
4	B	703	NAP	C2B-C1B-N9A-C8A
4	A	703	NAP	C2B-C1B-N9A-C4A
4	B	703	NAP	C2B-C1B-N9A-C4A
4	B	703	NAP	C2B-O2B-P2B-O1X
4	A	703	NAP	C2B-O2B-P2B-O1X
3	B	702	FAD	PA-O3P-P-O1P
3	A	702	FAD	PA-O3P-P-O2P
3	B	702	FAD	PA-O3P-P-O2P
4	A	703	NAP	C2B-O2B-P2B-O3X

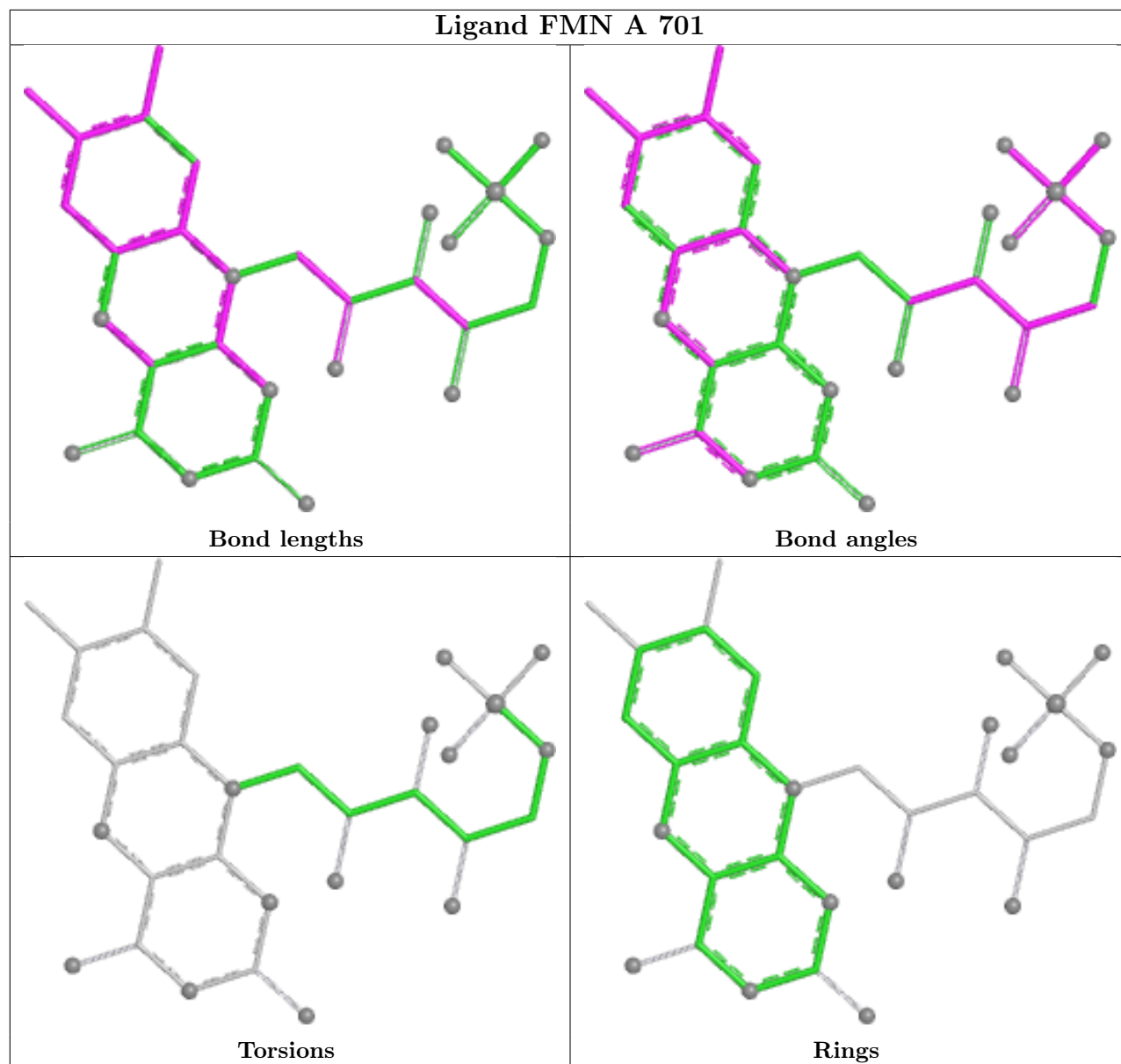
There are no ring outliers.

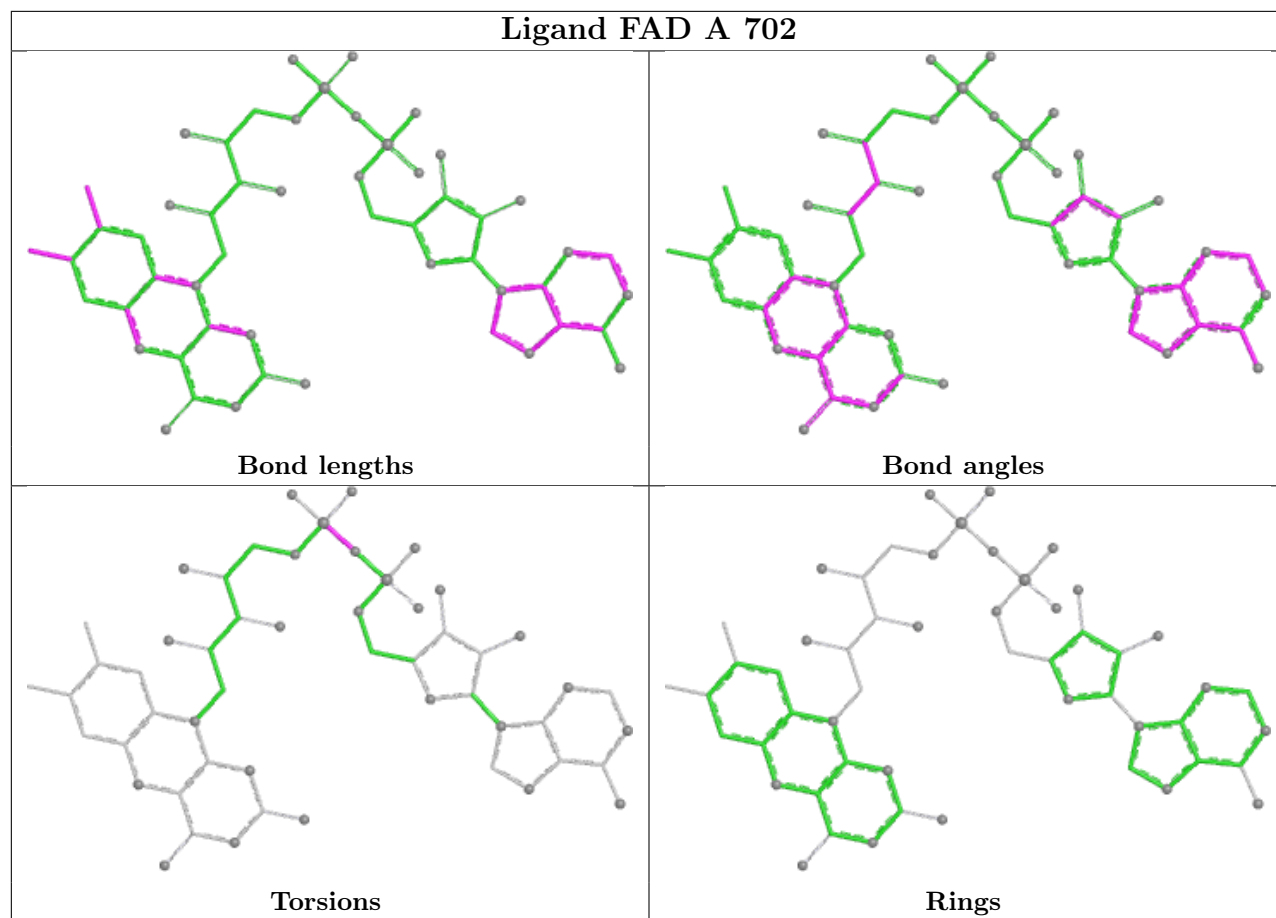
No monomer is involved in short contacts.

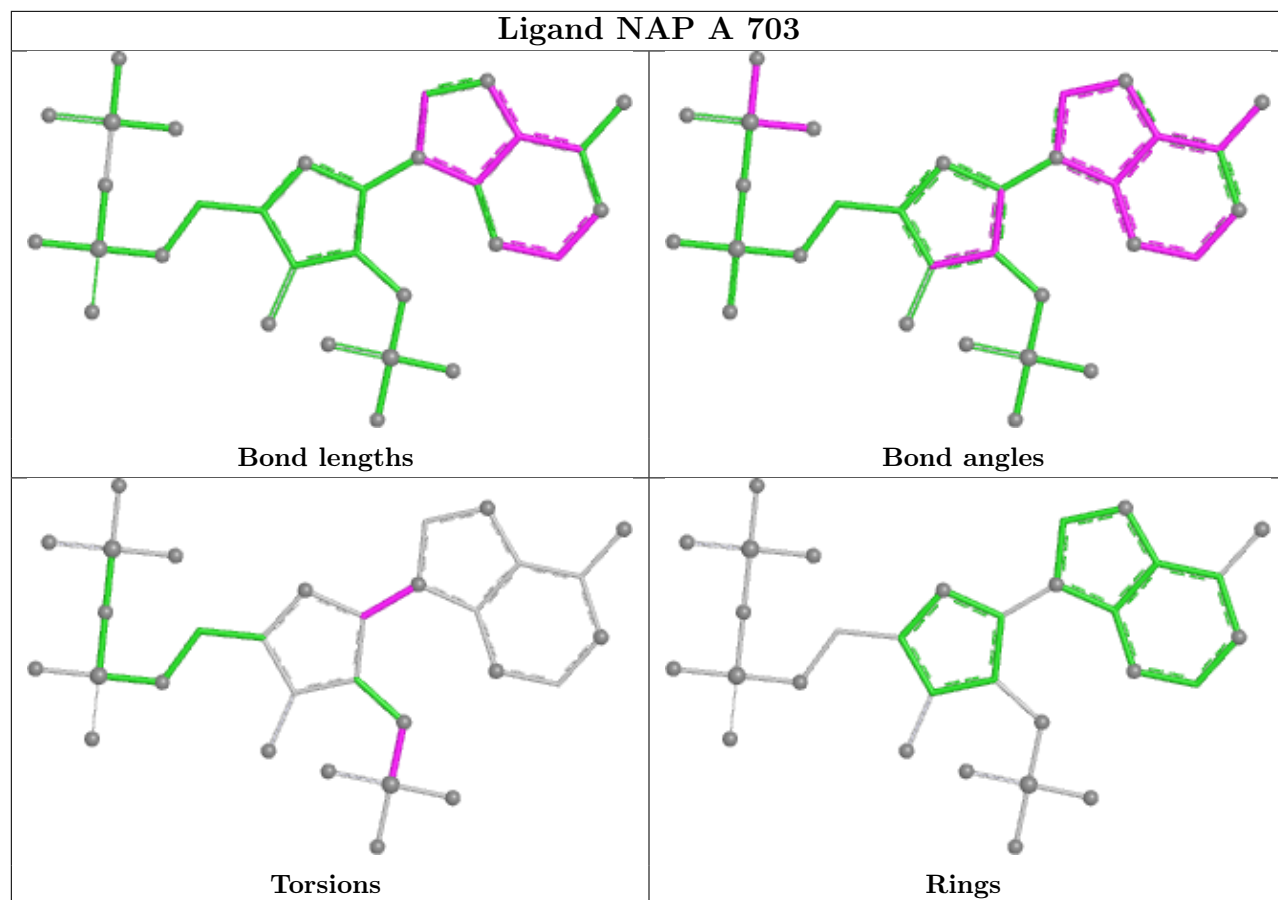
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

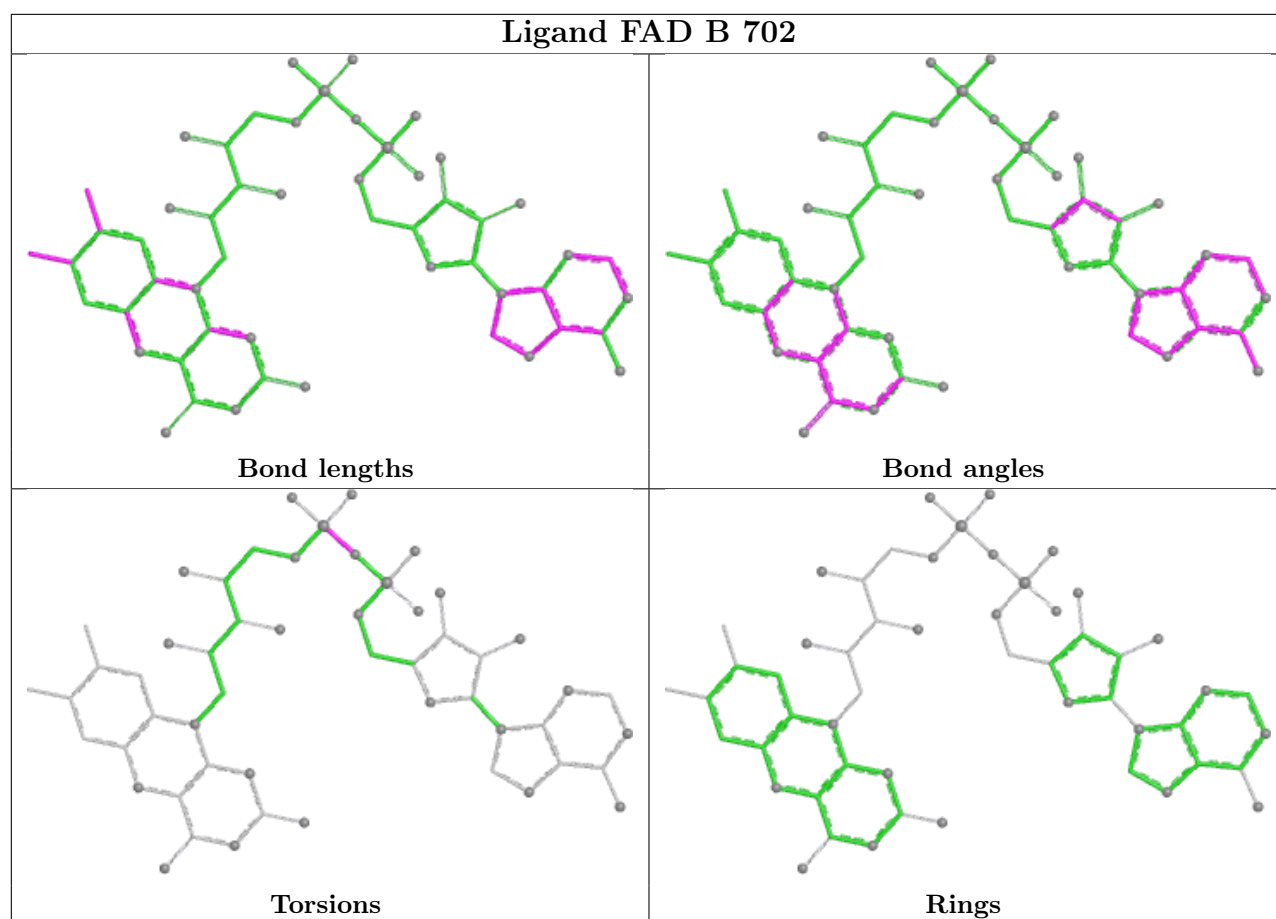












5.7 Other polymers [i](#)

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	611/621 (98%)	-0.40	1 (0%) 91 90	15, 28, 46, 86	0
1	B	605/621 (97%)	0.25	49 (8%) 18 15	16, 35, 86, 99	0
All	All	1216/1242 (97%)	-0.08	50 (4%) 41 38	15, 30, 82, 99	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	64	VAL	4.4
1	B	192	LEU	3.9
1	B	236	THR	3.7
1	B	161	VAL	3.7
1	B	181	PHE	3.5
1	B	138	ALA	3.4
1	B	235	ALA	3.4
1	B	173	LEU	3.3
1	B	67	SER	3.2
1	B	85	GLY	3.1
1	B	163	LEU	3.1
1	B	197	ALA	2.9
1	B	66	GLU	2.8
1	B	233	VAL	2.8
1	B	500	ALA	2.8
1	B	255	VAL	2.8
1	B	125	LEU	2.7
1	B	152	PHE	2.6
1	B	69	PHE	2.6
1	B	70	VAL	2.6
1	B	206	GLY	2.6
1	B	120	ALA	2.6
1	B	254	ASP	2.6
1	B	82	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	355	SER	2.5
1	B	164	THR	2.5
1	B	188	VAL	2.5
1	B	149	ALA	2.4
1	B	136	CYS	2.4
1	B	71	GLU	2.4
1	B	80	ILE	2.3
1	B	139	THR	2.3
1	B	106	GLY	2.3
1	B	122	LEU	2.2
1	A	506	ARG	2.2
1	B	171	PHE	2.2
1	B	182	ASN	2.2
1	B	68	SER	2.2
1	B	102	ALA	2.2
1	B	128	ILE	2.1
1	B	252	ASP	2.1
1	B	166	VAL	2.1
1	B	232	GLY	2.1
1	B	177	THR	2.1
1	B	112	ALA	2.1
1	B	103	HIS	2.1
1	B	403	HIS	2.1
1	B	185	GLY	2.1
1	B	140	TYR	2.0
1	B	156	LEU	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](#)

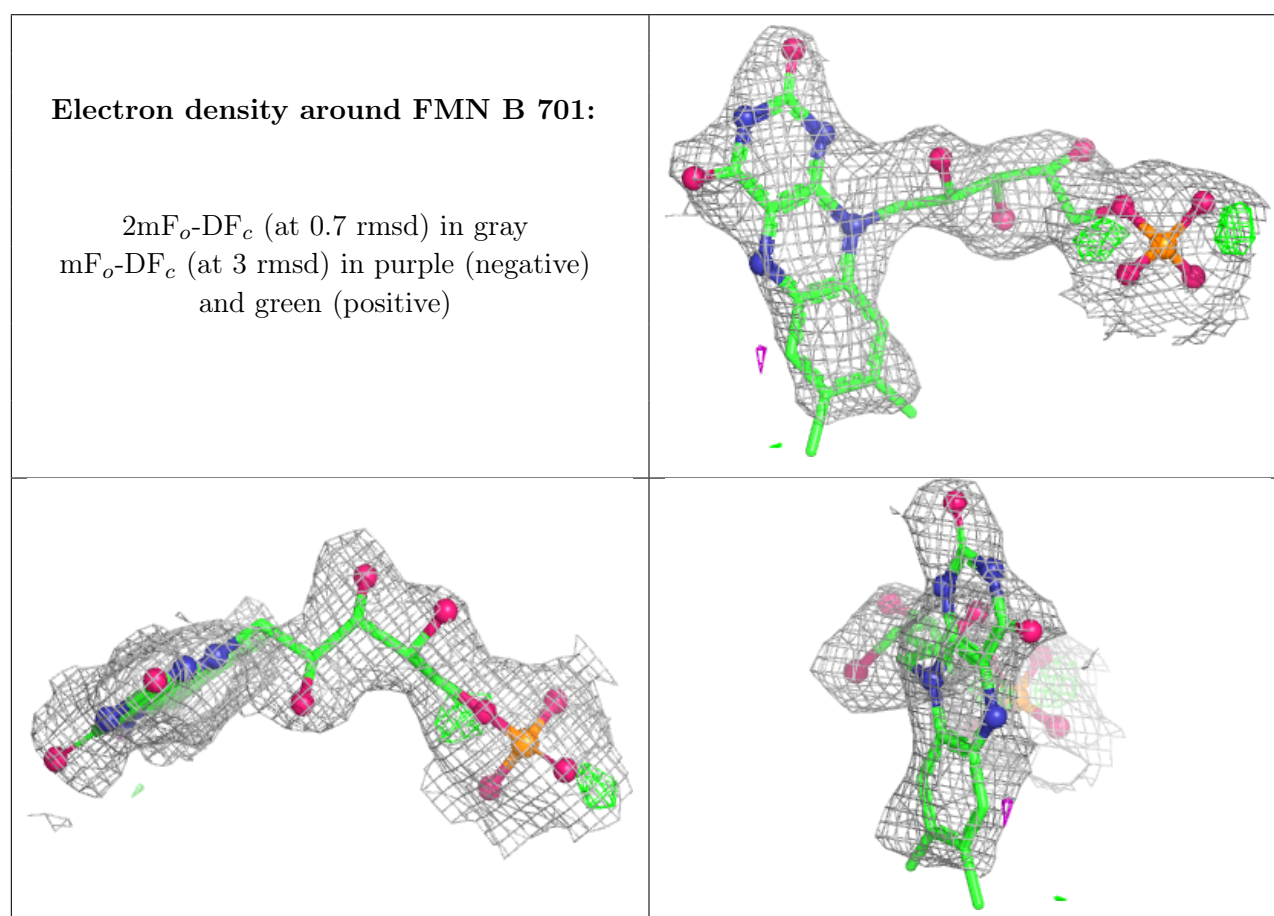
There are no oligosaccharides in this entry.

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.

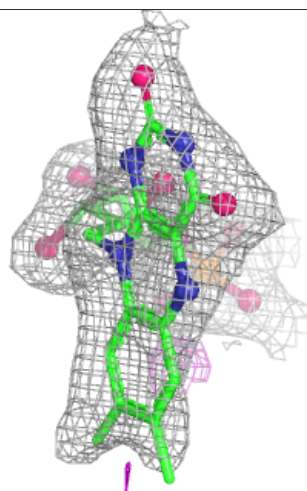
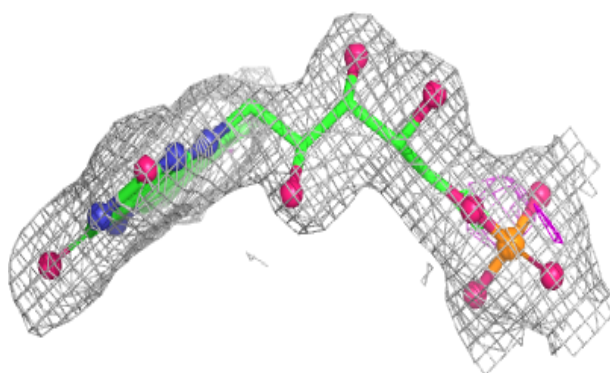
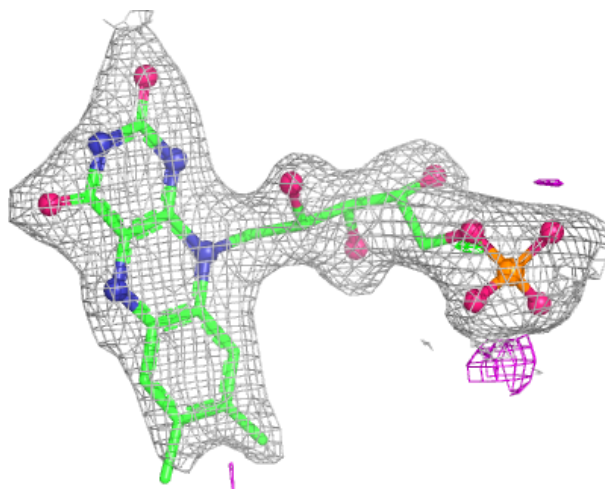
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	PO4	B	704	5/5	0.66	0.16	98,99,99,99	0
2	FMN	B	701	31/31	0.88	0.13	57,73,74,74	0
2	FMN	A	701	31/31	0.95	0.09	29,35,39,40	0
4	NAP	B	703	31/48	0.96	0.06	17,21,43,48	0
4	NAP	A	703	31/48	0.97	0.06	17,20,44,50	0
3	FAD	B	702	53/53	0.98	0.05	12,19,25,29	0
3	FAD	A	702	53/53	0.98	0.05	16,20,28,30	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.



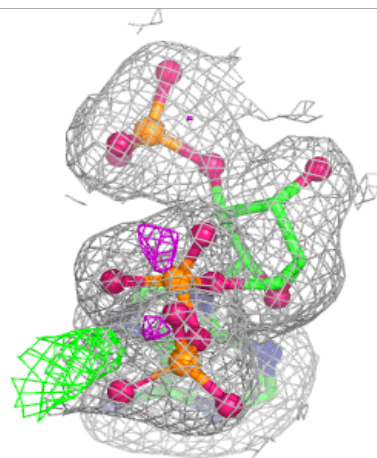
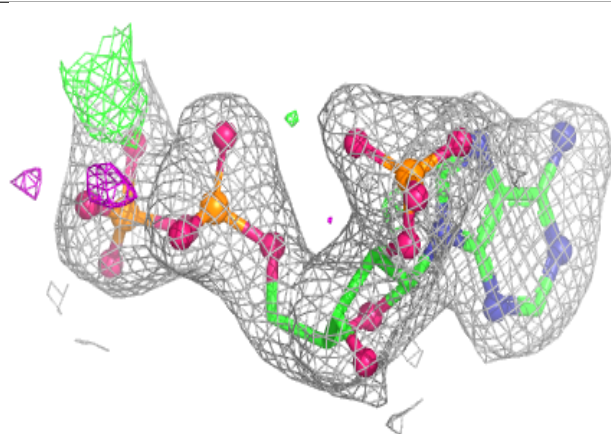
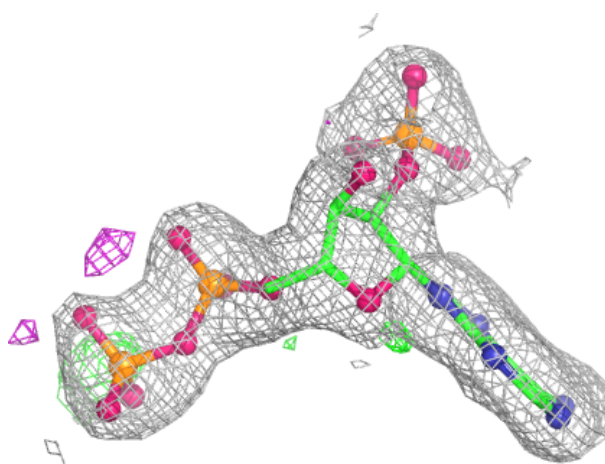
Electron density around FMN A 701:

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



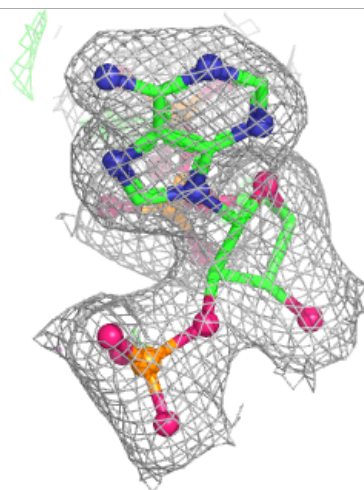
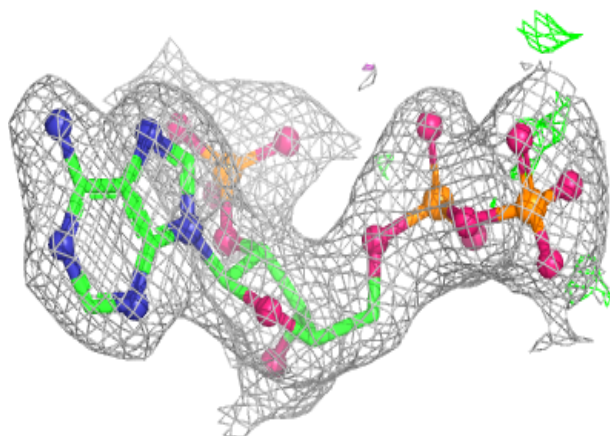
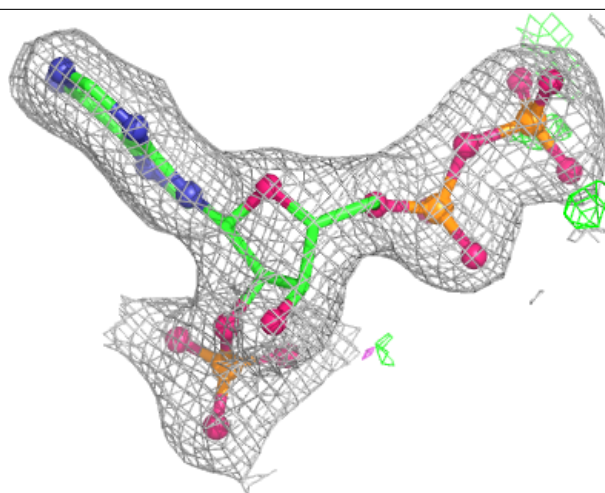
Electron density around NAP B 703:

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



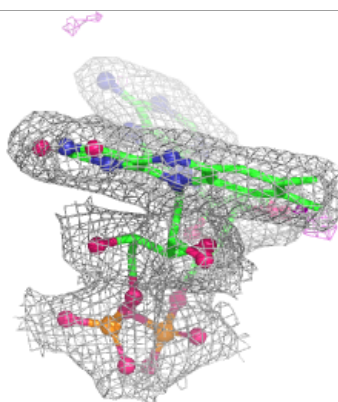
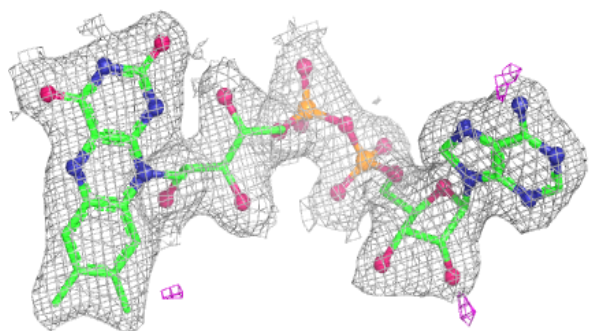
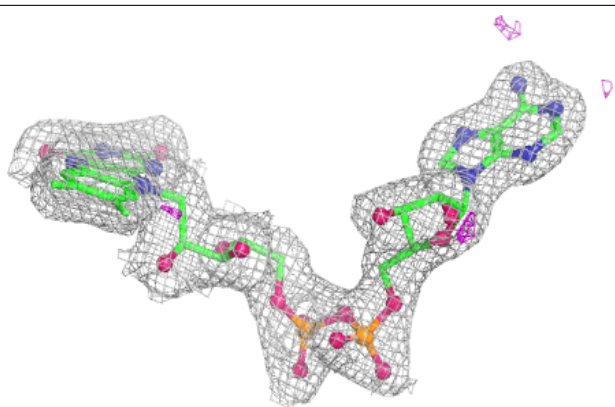
Electron density around NAP A 703:

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

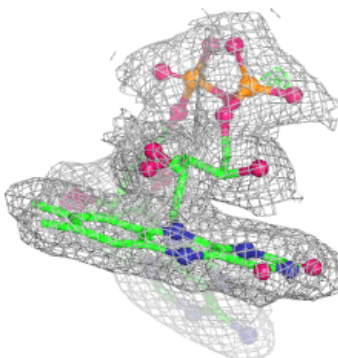
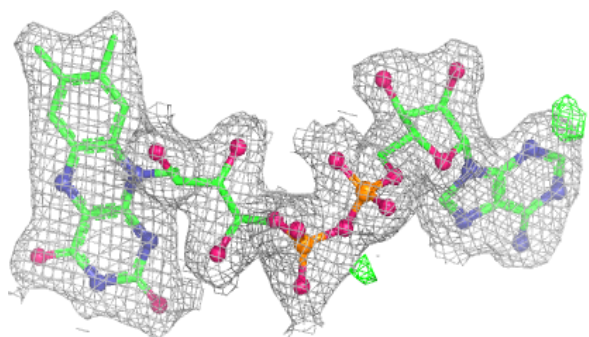
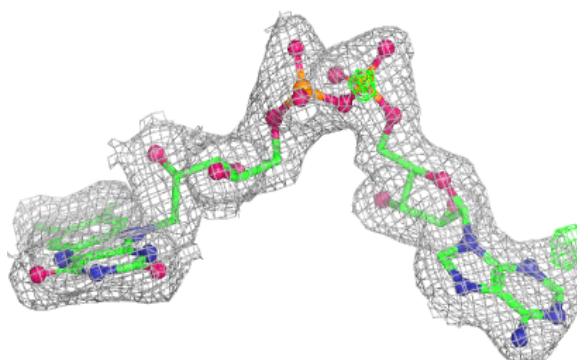


Electron density around FAD B 702:

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

**Electron density around FAD A 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.