Full wwPDB X-ray Structure Validation Report (i)

Jun 11, 2024 – 11:42 PM EDT

PDB ID	:	1AMO
Title	:	THREE-DIMENSIONAL STRUCTURE OF NADPH-CYTOCHROME P450 REDUCTASE: PROTOTYPE FOR FMN-AND FAD-CONTAINING ENZYMES
Authors	:	Wang, M.; Roberts, D.L.; Paschke, R.; Shea, T.M.; Masters, B.S.S.; Kim, J.J.P.
Deposited on	:	1997-06-17
Resolution	:	2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

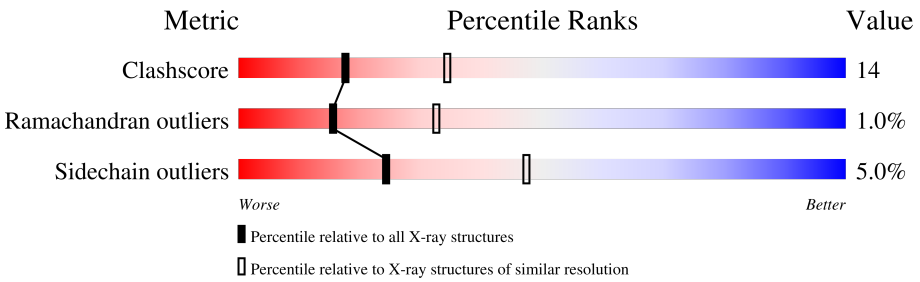
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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))
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	615	67% 27% • • •
1	B	615	67% 28% • • •

2 Entry composition (i)

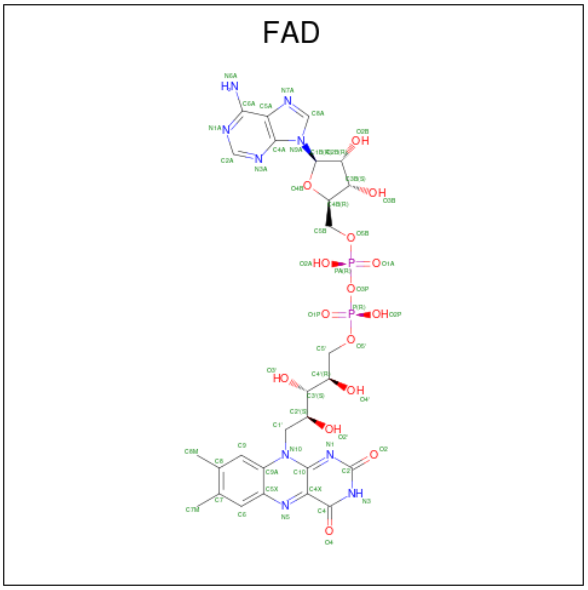
There are 5 unique types of molecules in this entry. The entry contains 10040 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

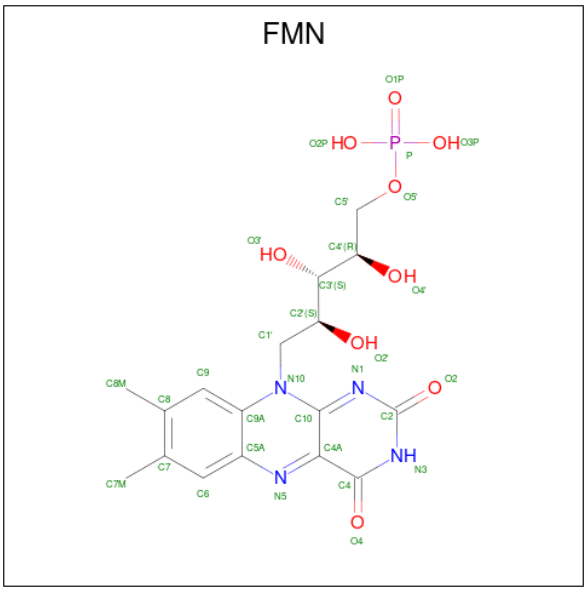
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	601	4830	3062	830	915	23	0	0	0
1	B	601	4830	3062	830	915	23	0	0	0

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



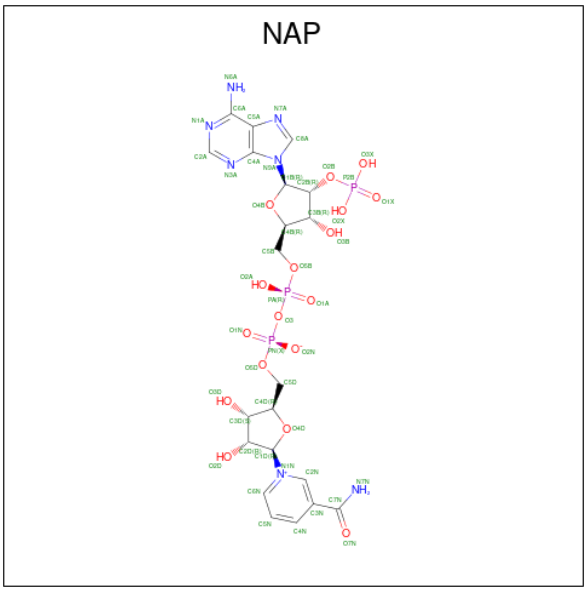
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	53	27	9	15	2	0	0
2	B	1	47	23	7	15	2	0	0

- Molecule 3 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



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

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



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

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	78	78	78	0	0
			Total	O		
5	B	78	78	78	0	0

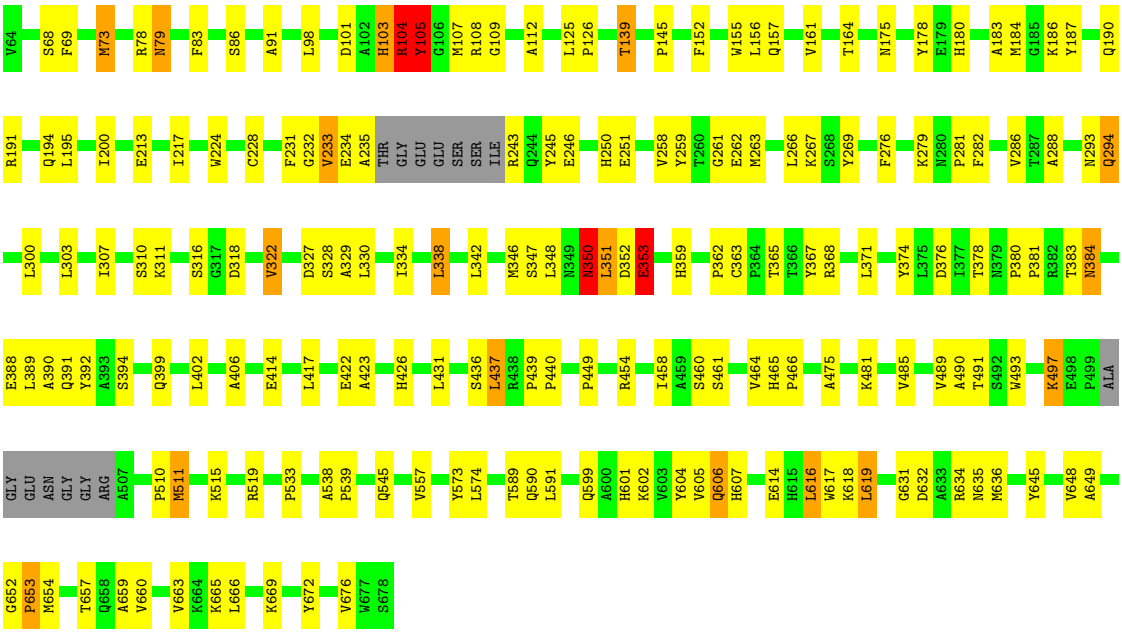
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

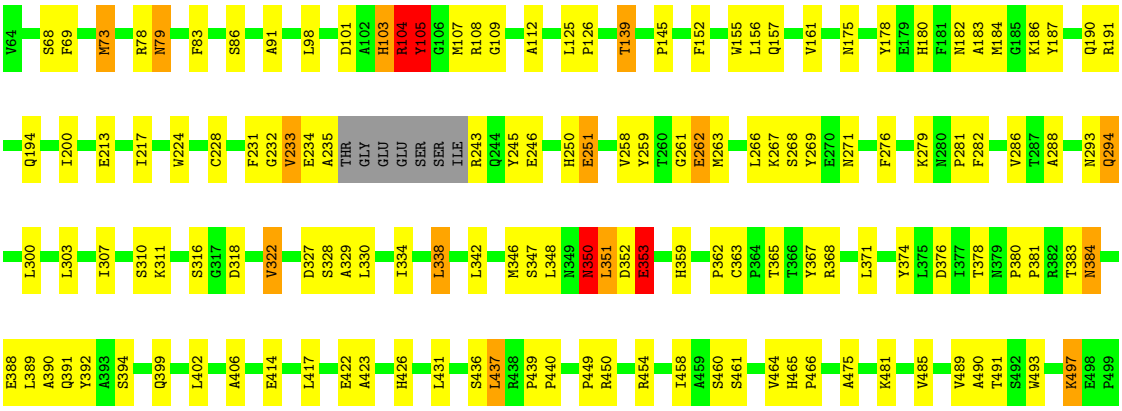
• Molecule 1: NADPH-CYTOCHROME P450 REDUCTASE

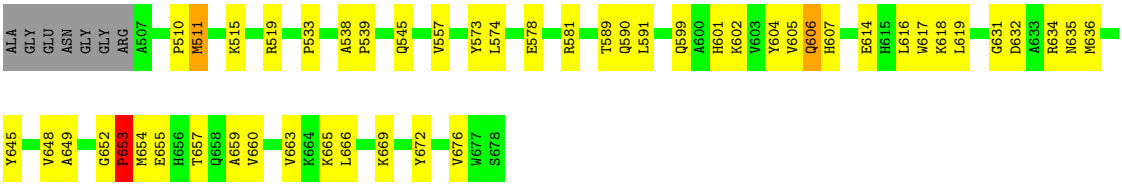
Chain A:



• Molecule 1: NADPH-CYTOCHROME P450 REDUCTASE

Chain B:




ALA
GLY
GLU
ASN
GLY
GLY
ARG
A507
P510
M511
K515
R519
P533
A538
P539
Q545
V557
Y573
L574
E578
R581
T589
Q590
L591
Q599
A600
H601
K602
V603
Y604
V605
Q606
H607
E614
H615
L616
W617
K618
L619
G631
D632
A633
R634
N635
M636
Y645
V648
A649
G652
P653
M654
E655
H656
T657
Q658
A659
V660
V663
K664
K665
L666
K669
Y672
V676
W677
S678

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	103.28Å 116.18Å 119.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.60	Depositor
% Data completeness (in resolution range)	91.5 (10.00-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.200 , 0.310	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10040	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NAP, FMN, 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	0.47	1/4945 (0.0%)	0.73	15/6691 (0.2%)
1	B	0.47	1/4945 (0.0%)	0.73	15/6691 (0.2%)
All	All	0.47	2/9890 (0.0%)	0.73	30/13382 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	350	ASN	C-N	10.97	1.59	1.34
1	B	350	ASN	C-N	10.96	1.59	1.34

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	350	ASN	C-N-CA	-9.61	97.67	121.70
1	B	350	ASN	C-N-CA	-9.60	97.69	121.70
1	A	652	GLY	N-CA-C	-8.79	91.12	113.10
1	B	652	GLY	N-CA-C	-8.79	91.12	113.10
1	A	231	PHE	C-N-CA	7.70	138.48	122.30
1	B	231	PHE	C-N-CA	7.67	138.41	122.30
1	A	243	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	B	235	ALA	CB-CA-C	-7.39	99.01	110.10
1	B	243	ARG	NE-CZ-NH2	7.38	123.99	120.30
1	A	235	ALA	CB-CA-C	-7.38	99.04	110.10
1	B	104	ARG	NE-CZ-NH2	6.88	123.74	120.30
1	A	104	ARG	NE-CZ-NH2	6.86	123.73	120.30
1	A	351	LEU	O-C-N	6.75	133.50	122.70
1	B	263	MET	CG-SD-CE	6.74	110.98	100.20
1	A	263	MET	CG-SD-CE	6.72	110.95	100.20
1	B	235	ALA	N-CA-C	6.71	129.13	111.00
1	B	351	LEU	O-C-N	6.71	133.44	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	235	ALA	N-CA-C	6.71	129.11	111.00
1	B	262	GLU	O-C-N	-6.37	112.50	122.70
1	A	652	GLY	CA-C-O	-6.35	109.17	120.60
1	A	262	GLU	O-C-N	-6.34	112.55	122.70
1	B	652	GLY	CA-C-O	-6.34	109.19	120.60
1	B	351	LEU	CA-C-N	-5.51	105.08	117.20
1	A	352	ASP	N-CA-C	-5.51	96.13	111.00
1	A	351	LEU	CA-C-N	-5.50	105.09	117.20
1	B	352	ASP	N-CA-C	-5.49	96.17	111.00
1	A	73	MET	CG-SD-CE	5.12	108.40	100.20
1	B	73	MET	CG-SD-CE	5.12	108.39	100.20
1	B	103	HIS	CA-CB-CG	-5.07	104.98	113.60
1	A	103	HIS	CA-CB-CG	-5.06	105.00	113.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4830	0	4678	129	3
1	B	4830	0	4678	130	10
2	A	53	0	31	2	0
2	B	47	0	22	9	0
3	A	31	0	19	1	0
3	B	31	0	19	1	0
4	A	31	0	11	1	0
4	B	31	0	11	1	0
5	A	78	0	0	10	0
5	B	78	0	0	11	2
All	All	10040	0	9469	266	11

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 14.

All (266) 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:A:353:GLU:OE1	1:A:353:GLU:HA	1.36	1.11
1:B:353:GLU:HA	1:B:353:GLU:OE1	1.36	1.08
2:B:750:FAD:C9	2:B:750:FAD:C7	2.39	1.00
2:B:750:FAD:C7	2:B:750:FAD:C5X	2.42	0.98
2:B:750:FAD:C7	2:B:750:FAD:C8M	2.50	0.90
2:B:750:FAD:C9	2:B:750:FAD:C8M	2.55	0.85
1:A:353:GLU:OE1	1:A:353:GLU:CA	2.25	0.84
1:A:402:LEU:HD11	1:A:437:LEU:HD21	1.63	0.81
1:B:353:GLU:OE1	1:B:353:GLU:CA	2.25	0.80
1:B:402:LEU:HD11	1:B:437:LEU:HD21	1.63	0.80
1:A:78:ARG:HG2	1:A:108:ARG:HG2	1.66	0.77
1:B:78:ARG:HG2	1:B:108:ARG:HG2	1.66	0.76
2:B:750:FAD:C4X	2:B:750:FAD:O4	2.33	0.76
1:B:461:SER:HB3	1:B:464:VAL:HG22	1.70	0.72
1:A:461:SER:HB3	1:A:464:VAL:HG22	1.70	0.71
1:B:665:LYS:HG3	1:B:669:LYS:HE3	1.77	0.66
1:A:665:LYS:HG3	1:A:669:LYS:HE3	1.77	0.65
1:A:261:GLY:HA2	1:A:374:TYR:CZ	2.32	0.64
1:B:261:GLY:HA2	1:B:374:TYR:CZ	2.32	0.64
1:A:378:THR:O	1:A:426:HIS:HB3	1.97	0.64
1:B:378:THR:O	1:B:426:HIS:HB3	1.97	0.63
1:A:86:SER:HB2	1:A:91:ALA:HB3	1.79	0.63
1:B:632:ASP:OD1	1:B:634:ARG:HG2	1.99	0.63
1:A:632:ASP:OD1	1:A:634:ARG:HG2	1.99	0.62
1:B:86:SER:HB2	1:B:91:ALA:HB3	1.79	0.62
1:A:604:TYR:H	1:A:607:HIS:CD2	2.17	0.62
1:B:156:LEU:HG	1:B:191:ARG:HG2	1.81	0.62
1:B:604:TYR:H	1:B:607:HIS:CD2	2.17	0.61
1:A:156:LEU:HG	1:A:191:ARG:HG2	1.81	0.61
1:B:213:GLU:O	1:B:217:ILE:HG13	2.00	0.61
2:B:750:FAD:C7	2:B:750:FAD:C9A	2.79	0.60
1:A:213:GLU:O	1:A:217:ILE:HG13	2.00	0.60
1:A:105:TYR:CE1	1:A:228:CYS:SG	2.95	0.60
1:B:105:TYR:CE1	1:B:228:CYS:SG	2.95	0.60
1:B:634:ARG:NH1	5:B:779:HOH:O	2.37	0.58
1:B:604:TYR:H	1:B:607:HIS:HD2	1.51	0.58
1:B:515:LYS:HD2	1:B:519:ARG:HH22	1.69	0.58
1:B:614:GLU:HG2	1:B:618:LYS:NZ	2.19	0.58
1:A:614:GLU:HG2	1:A:618:LYS:NZ	2.19	0.58
1:A:617:TRP:HZ3	1:A:648:VAL:HG22	1.69	0.57
1:B:300:LEU:HD22	1:B:574:LEU:HD21	1.86	0.57
1:A:338:LEU:HD11	1:A:439:PRO:HD2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:604:TYR:H	1:A:607:HIS:HD2	1.51	0.57
1:A:515:LYS:HD2	1:A:519:ARG:HH22	1.69	0.57
1:A:634:ARG:NH1	5:A:823:HOH:O	2.37	0.57
1:B:617:TRP:HZ3	1:B:648:VAL:HG22	1.69	0.57
1:B:105:TYR:HE1	1:B:228:CYS:SG	2.28	0.57
1:A:300:LEU:HD22	1:A:574:LEU:HD21	1.86	0.56
1:A:79:ASN:HD21	1:A:107:MET:HB3	1.70	0.56
1:B:338:LEU:HD11	1:B:439:PRO:HD2	1.87	0.56
1:A:105:TYR:HE1	1:A:228:CYS:SG	2.28	0.56
1:B:383:THR:HG23	1:B:406:ALA:HA	1.87	0.56
1:A:383:THR:HG23	1:A:406:ALA:HA	1.87	0.56
1:A:657:THR:O	1:A:660:VAL:HG22	2.06	0.56
1:B:657:THR:O	1:B:660:VAL:HG22	2.06	0.56
1:A:279:LYS:H	1:A:279:LYS:HD2	1.71	0.55
1:B:659:ALA:O	1:B:663:VAL:HG23	2.07	0.55
1:A:334:ILE:O	1:A:338:LEU:HB2	2.07	0.55
1:B:145:PRO:HB3	1:B:184:MET:SD	2.47	0.55
1:A:186:LYS:O	1:A:190:GLN:HG2	2.06	0.55
1:B:186:LYS:O	1:B:190:GLN:HG2	2.06	0.55
1:A:145:PRO:HB3	1:A:184:MET:SD	2.47	0.55
1:A:557:VAL:HG11	5:A:816:HOH:O	2.07	0.55
2:B:750:FAD:C8M	2:B:750:FAD:C7M	2.84	0.55
1:B:557:VAL:HG11	5:B:772:HOH:O	2.07	0.55
1:A:103:HIS:O	1:A:104:ARG:NE	2.30	0.54
1:A:390:ALA:HA	1:A:402:LEU:HD12	1.89	0.54
1:A:659:ALA:O	1:A:663:VAL:HG23	2.07	0.54
1:A:79:ASN:N	1:A:79:ASN:HD22	2.05	0.54
1:A:342:LEU:HB3	1:A:368:ARG:HB2	1.89	0.54
1:B:279:LYS:H	1:B:279:LYS:HD2	1.71	0.54
1:B:79:ASN:HD21	1:B:107:MET:HB3	1.70	0.54
1:B:390:ALA:HA	1:B:402:LEU:HD12	1.88	0.54
1:B:334:ILE:O	1:B:338:LEU:HB2	2.07	0.54
1:B:347:SER:OG	1:B:363:CYS:HB3	2.07	0.54
1:B:293:ASN:HB2	1:B:300:LEU:H	1.72	0.54
1:B:342:LEU:HB3	1:B:368:ARG:HB2	1.89	0.54
1:B:376:ASP:HB3	1:B:449:PRO:HG2	1.89	0.54
1:A:293:ASN:HB2	1:A:300:LEU:H	1.72	0.54
1:B:79:ASN:N	1:B:79:ASN:HD22	2.05	0.54
1:A:376:ASP:HB3	1:A:449:PRO:HG2	1.89	0.54
1:A:634:ARG:HG3	1:A:635:ASN:N	2.23	0.54
1:A:232:GLY:O	1:A:234:GLU:HB2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:GLN:O	1:B:194:GLN:HG3	2.08	0.53
1:B:232:GLY:O	1:B:234:GLU:HB2	2.08	0.53
1:B:634:ARG:HG3	1:B:635:ASN:N	2.23	0.53
1:A:190:GLN:O	1:A:194:GLN:HG3	2.08	0.53
1:A:589:THR:HG22	1:A:590:GLN:HG3	1.90	0.53
1:B:103:HIS:CE1	1:B:108:ARG:HG3	2.44	0.53
1:B:589:THR:HG22	1:B:590:GLN:HG3	1.90	0.53
1:B:101:ASP:O	1:B:105:TYR:HB2	2.09	0.53
1:B:348:LEU:O	1:B:359:HIS:HD2	1.92	0.53
1:A:347:SER:OG	1:A:363:CYS:HB3	2.07	0.53
1:B:103:HIS:O	1:B:104:ARG:NE	2.30	0.53
1:B:464:VAL:HG23	1:B:465:HIS:CD2	2.44	0.53
1:B:599:GLN:HG2	1:B:601:HIS:O	2.08	0.52
1:A:599:GLN:HG3	1:A:601:HIS:CE1	2.44	0.52
1:A:310:SER:O	1:A:311:LYS:HB2	2.09	0.52
1:A:464:VAL:HG23	1:A:465:HIS:CD2	2.44	0.52
1:A:599:GLN:HG2	1:A:601:HIS:O	2.08	0.52
1:B:599:GLN:HG3	1:B:601:HIS:CE1	2.45	0.52
1:A:103:HIS:CE1	1:A:108:ARG:HG3	2.44	0.52
1:A:368:ARG:HG3	1:A:368:ARG:HH11	1.74	0.52
1:A:101:ASP:O	1:A:105:TYR:HB2	2.09	0.52
1:B:310:SER:O	1:B:311:LYS:HB2	2.09	0.52
1:B:368:ARG:HH11	1:B:368:ARG:HG3	1.74	0.52
1:B:79:ASN:HD22	1:B:79:ASN:H	1.58	0.51
1:A:602:LYS:HD3	1:A:604:TYR:OH	2.10	0.51
1:A:604:TYR:HB3	1:A:606:GLN:OE1	2.10	0.51
1:B:604:TYR:HB3	1:B:606:GLN:OE1	2.10	0.51
1:A:68:SER:HA	5:A:775:HOH:O	2.11	0.51
1:A:348:LEU:O	1:A:359:HIS:HD2	1.92	0.51
1:B:68:SER:HA	5:B:809:HOH:O	2.11	0.51
1:A:322:VAL:HG13	1:A:493:TRP:HZ3	1.76	0.51
1:B:322:VAL:HG13	1:B:493:TRP:HZ3	1.76	0.51
1:A:557:VAL:CG1	5:A:816:HOH:O	2.60	0.50
1:B:602:LYS:HD3	1:B:604:TYR:OH	2.10	0.50
1:A:175:ASN:HB3	1:A:178:TYR:HD1	1.77	0.50
1:B:175:ASN:HB3	1:B:178:TYR:HD1	1.77	0.50
1:A:79:ASN:HD22	1:A:79:ASN:H	1.58	0.50
1:A:466:PRO:O	5:A:812:HOH:O	2.20	0.50
1:B:466:PRO:O	5:B:768:HOH:O	2.20	0.50
1:A:246:GLU:O	1:A:348:LEU:HA	2.12	0.49
1:A:125:LEU:N	1:A:126:PRO:HD2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:LEU:N	1:B:126:PRO:HD2	2.27	0.49
1:B:384:ASN:H	1:B:384:ASN:HD22	1.60	0.49
1:B:250:HIS:CE1	1:B:347:SER:HB2	2.48	0.49
1:B:246:GLU:O	1:B:348:LEU:HA	2.12	0.49
1:B:269:TYR:HE1	1:B:281:PRO:HG2	1.78	0.49
1:A:533:PRO:HG3	1:A:636:MET:HG3	1.94	0.49
1:B:475:ALA:HA	1:B:491:THR:HB	1.95	0.49
1:A:250:HIS:CE1	1:A:347:SER:HB2	2.47	0.48
1:B:318:ASP:OD1	1:B:515:LYS:HA	2.13	0.48
1:B:388:GLU:O	1:B:391:GLN:HG2	2.13	0.48
1:A:250:HIS:HE1	1:A:347:SER:HB2	1.78	0.48
1:A:388:GLU:O	1:A:391:GLN:HG2	2.13	0.48
1:A:475:ALA:HA	1:A:491:THR:HB	1.95	0.48
1:B:533:PRO:HG3	1:B:636:MET:HG3	1.94	0.48
1:A:269:TYR:HE1	1:A:281:PRO:HG2	1.78	0.48
1:B:515:LYS:HD2	1:B:519:ARG:NH2	2.29	0.48
1:B:557:VAL:CG1	5:B:772:HOH:O	2.60	0.48
1:A:318:ASP:OD1	1:A:515:LYS:HA	2.13	0.48
1:B:417:LEU:O	1:B:422:GLU:HB2	2.13	0.48
1:A:645:TYR:HD1	1:A:663:VAL:HG21	1.79	0.48
1:B:645:TYR:HD1	1:B:663:VAL:HG21	1.79	0.48
1:A:175:ASN:HB3	1:A:178:TYR:CD1	2.49	0.48
1:A:417:LEU:O	1:A:422:GLU:HB2	2.13	0.48
1:A:384:ASN:HD22	1:A:384:ASN:H	1.60	0.48
1:B:458:ILE:HG22	1:B:460:SER:H	1.79	0.48
1:A:606:GLN:OE1	4:A:752:NAP:H2A	2.14	0.48
1:B:606:GLN:OE1	4:B:752:NAP:H2A	2.14	0.48
1:A:458:ILE:HG22	1:A:460:SER:H	1.79	0.47
1:A:614:GLU:HB3	5:A:815:HOH:O	2.15	0.47
1:B:538:ALA:HB3	1:B:539:PRO:CD	2.44	0.47
1:A:316:SER:HA	5:A:787:HOH:O	2.14	0.47
1:A:515:LYS:HD2	1:A:519:ARG:NH2	2.29	0.47
1:B:175:ASN:HB3	1:B:178:TYR:CD1	2.49	0.47
1:B:250:HIS:HE1	1:B:347:SER:HB2	1.79	0.47
1:A:493:TRP:HH2	1:A:510:PRO:HD2	1.80	0.47
1:A:538:ALA:HB3	1:A:539:PRO:CD	2.44	0.47
1:A:155:TRP:HZ2	1:A:161:VAL:HG11	1.80	0.47
1:A:286:VAL:HG21	1:A:497:LYS:HB3	1.96	0.47
1:A:489:VAL:HG23	2:A:750:FAD:O1A	2.15	0.47
1:B:493:TRP:HH2	1:B:510:PRO:HD2	1.80	0.47
1:A:485:VAL:HG23	5:B:817:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:LEU:O	1:A:334:ILE:HG13	2.15	0.47
5:A:783:HOH:O	1:B:485:VAL:HG23	2.14	0.47
1:B:489:VAL:HG23	2:B:750:FAD:O1A	2.15	0.47
1:B:316:SER:HA	5:B:821:HOH:O	2.14	0.46
1:B:276:PHE:CE2	1:B:282:PHE:HD2	2.34	0.46
1:A:276:PHE:CE2	1:A:282:PHE:HD2	2.34	0.46
1:B:279:LYS:HD2	1:B:279:LYS:N	2.31	0.46
1:B:245:TYR:CE1	1:B:350:ASN:HB2	2.50	0.46
1:B:330:LEU:O	1:B:334:ILE:HG13	2.15	0.46
1:A:402:LEU:CD1	1:A:437:LEU:HD21	2.42	0.46
1:B:69:PHE:HD1	1:B:73:MET:HE2	1.80	0.46
1:B:390:ALA:O	1:B:399:GLN:HG3	2.16	0.46
1:B:98:LEU:O	1:B:224:TRP:HZ2	1.99	0.46
1:B:286:VAL:HG21	1:B:497:LYS:HB3	1.96	0.46
1:A:384:ASN:O	1:A:388:GLU:HG2	2.16	0.46
1:A:245:TYR:CE1	1:A:350:ASN:HB2	2.50	0.45
1:A:279:LYS:HD2	1:A:279:LYS:N	2.31	0.45
1:B:402:LEU:CD1	1:B:437:LEU:HD21	2.42	0.45
1:A:390:ALA:O	1:A:399:GLN:HG3	2.16	0.45
1:A:649:ALA:HA	1:A:654:MET:HB2	1.98	0.45
1:B:155:TRP:HZ2	1:B:161:VAL:HG11	1.80	0.45
1:B:649:ALA:HA	1:B:654:MET:HB2	1.98	0.45
1:A:98:LEU:O	1:A:224:TRP:HZ2	1.99	0.45
1:B:233:VAL:HG13	1:B:234:GLU:H	1.82	0.45
1:B:614:GLU:HB3	5:B:771:HOH:O	2.15	0.45
1:A:631:GLY:O	1:A:676:VAL:HA	2.17	0.45
1:B:631:GLY:O	1:B:676:VAL:HA	2.17	0.45
1:B:294:GLN:HE21	1:B:294:GLN:HB2	1.59	0.45
1:A:233:VAL:HG13	1:A:234:GLU:H	1.81	0.44
1:A:103:HIS:CD2	1:A:109:GLY:H	2.35	0.44
1:A:351:LEU:HD23	1:A:351:LEU:HA	1.90	0.44
1:A:490:ALA:O	1:A:493:TRP:HB3	2.18	0.44
1:B:327:ASP:OD2	1:B:329:ALA:HB3	2.18	0.44
1:B:103:HIS:CD2	1:B:109:GLY:H	2.35	0.44
1:B:423:ALA:HA	1:B:481:LYS:HB2	2.00	0.44
1:B:384:ASN:O	1:B:388:GLU:HG2	2.16	0.44
1:A:288:ALA:O	1:A:303:LEU:HA	2.18	0.43
1:A:104:ARG:HD3	1:A:104:ARG:HA	1.84	0.43
1:A:157:GLN:HA	1:A:187:TYR:OH	2.18	0.43
1:B:261:GLY:HA2	1:B:374:TYR:OH	2.19	0.43
1:B:338:LEU:HD13	1:B:431:LEU:HD13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:PHE:HA	1:A:112:ALA:O	2.18	0.43
1:B:288:ALA:O	1:B:303:LEU:HA	2.18	0.43
1:A:180:HIS:HB3	1:A:183:ALA:HB2	2.00	0.43
1:A:423:ALA:HA	1:A:481:LYS:HB2	2.00	0.43
1:A:511:MET:SD	1:A:511:MET:C	2.97	0.43
1:B:83:PHE:HA	1:B:112:ALA:O	2.18	0.43
1:B:511:MET:SD	1:B:511:MET:C	2.97	0.43
1:A:327:ASP:OD2	1:A:329:ALA:HB3	2.18	0.43
1:B:105:TYR:CZ	1:B:224:TRP:HB3	2.54	0.43
1:B:157:GLN:HA	1:B:187:TYR:OH	2.18	0.43
1:A:338:LEU:HD13	1:A:431:LEU:HD13	2.01	0.43
1:B:328:SER:CB	5:B:766:HOH:O	2.66	0.43
1:A:139:THR:HG23	3:A:751:FMN:C2	2.49	0.43
1:A:261:GLY:HA2	1:A:374:TYR:OH	2.19	0.43
1:B:180:HIS:HB3	1:B:183:ALA:HB2	2.00	0.43
1:B:246:GLU:HB3	1:B:351:LEU:HD12	2.01	0.43
1:B:139:THR:HG23	3:B:751:FMN:C2	2.49	0.43
1:A:380:PRO:HA	1:A:381:PRO:HD3	1.94	0.42
1:A:605:VAL:HG13	5:A:807:HOH:O	2.18	0.42
1:B:490:ALA:O	1:B:493:TRP:HB3	2.18	0.42
1:A:105:TYR:CZ	1:A:224:TRP:HB3	2.54	0.42
1:A:307:ILE:O	1:A:307:ILE:HG13	2.19	0.42
1:A:294:GLN:HE21	1:A:294:GLN:HB2	1.59	0.42
1:A:464:VAL:HG21	1:A:545:GLN:HB3	2.00	0.42
1:A:246:GLU:HB3	1:A:351:LEU:HD12	2.01	0.42
1:A:328:SER:CB	5:A:810:HOH:O	2.66	0.42
1:A:69:PHE:HD1	1:A:73:MET:HE2	1.84	0.42
1:B:303:LEU:N	1:B:303:LEU:HD12	2.35	0.42
1:A:303:LEU:HD12	1:A:303:LEU:N	2.35	0.42
1:A:392:TYR:CE1	1:A:440:PRO:HD3	2.55	0.42
1:B:307:ILE:HG13	1:B:307:ILE:O	2.19	0.42
1:B:394:SER:H	1:B:436:SER:HB3	1.85	0.42
1:B:342:LEU:HA	1:B:367:TYR:HB2	2.01	0.42
1:B:389:LEU:HA	1:B:392:TYR:HD2	1.85	0.42
1:B:491:THR:HG23	2:B:750:FAD:O2P	2.20	0.41
1:B:605:VAL:HG13	5:B:763:HOH:O	2.18	0.41
1:B:392:TYR:CE1	1:B:440:PRO:HD3	2.55	0.41
1:B:464:VAL:HG21	1:B:545:GLN:HB3	2.00	0.41
1:A:152:PHE:CZ	1:A:156:LEU:HD22	2.56	0.41
1:B:450:ARG:NH1	5:B:776:HOH:O	2.40	0.41
1:A:394:SER:H	1:A:436:SER:HB3	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:VAL:HG13	1:A:493:TRP:CZ3	2.55	0.41
1:A:619:LEU:HD12	1:A:619:LEU:HA	1.86	0.41
1:A:259:TYR:CE1	1:A:266:LEU:HG	2.56	0.41
1:A:389:LEU:HA	1:A:392:TYR:HD2	1.85	0.41
1:B:663:VAL:O	1:B:666:LEU:HB2	2.20	0.41
1:A:342:LEU:HA	1:A:367:TYR:HB2	2.01	0.41
1:A:663:VAL:O	1:A:666:LEU:HB2	2.21	0.41
1:A:491:THR:HG23	2:A:750:FAD:O2P	2.20	0.40
1:B:103:HIS:CD2	1:B:108:ARG:HA	2.57	0.40
1:B:258:VAL:HG13	1:B:365:THR:HA	2.02	0.40
1:B:262:GLU:HB2	1:B:268:SER:OG	2.21	0.40
1:B:617:TRP:CZ3	1:B:648:VAL:HA	2.56	0.40
1:A:258:VAL:HG13	1:A:365:THR:HA	2.02	0.40
1:B:104:ARG:HD3	1:B:104:ARG:HA	1.83	0.40
1:B:139:THR:HG21	1:B:182:ASN:HA	2.03	0.40
1:B:152:PHE:CZ	1:B:156:LEU:HD22	2.56	0.40
1:B:259:TYR:CE1	1:B:266:LEU:HG	2.56	0.40
1:B:380:PRO:HA	1:B:381:PRO:HD3	1.94	0.40
1:A:616:LEU:HD12	1:A:616:LEU:HA	1.95	0.40

All (11) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:ASN:OD1	1:B:655:GLU:OE2[3_655]	1.58	0.62
1:B:271:ASN:CG	1:B:655:GLU:OE2[3_655]	1.63	0.57
1:A:669:LYS:O	1:B:251:GLU:OE1[3_545]	1.64	0.56
5:B:795:HOH:O	5:B:825:HOH:O[3_645]	1.70	0.50
1:B:271:ASN:ND2	1:B:655:GLU:OE2[3_655]	1.72	0.48
1:B:266:LEU:O	5:B:796:HOH:O[3_655]	1.73	0.47
1:A:164:THR:CG2	1:B:578:GLU:OE2[2_555]	1.86	0.34
1:B:266:LEU:CD1	1:B:653:PRO:CG[3_655]	2.03	0.17
1:B:271:ASN:ND2	1:B:655:GLU:CD[3_655]	2.07	0.13
1:A:195:LEU:CA	1:B:581:ARG:NH2[2_555]	2.12	0.08
1:B:271:ASN:ND2	1:B:655:GLU:OE1[3_655]	2.17	0.03

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	595/615 (97%)	547 (92%)	42 (7%)	6 (1%)	15 32
1	B	595/615 (97%)	546 (92%)	43 (7%)	6 (1%)	15 32
All	All	1190/1230 (97%)	1093 (92%)	85 (7%)	12 (1%)	15 32

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	353	GLU
1	B	353	GLU
1	A	497	LYS
1	A	653	PRO
1	B	497	LYS
1	B	653	PRO
1	A	105	TYR
1	B	105	TYR
1	A	437	LEU
1	B	437	LEU
1	A	233	VAL
1	B	233	VAL

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	516/525 (98%)	490 (95%)	26 (5%)	24 47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	516/525 (98%)	490 (95%)	26 (5%)	24	47
All	All	1032/1050 (98%)	980 (95%)	52 (5%)	24	47

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	104	ARG
1	A	105	TYR
1	A	139	THR
1	A	200	ILE
1	A	251	GLU
1	A	267	LYS
1	A	294	GLN
1	A	322	VAL
1	A	338	LEU
1	A	346	MET
1	A	350	ASN
1	A	353	GLU
1	A	362	PRO
1	A	371	LEU
1	A	384	ASN
1	A	414	GLU
1	A	454	ARG
1	A	511	MET
1	A	573	TYR
1	A	591	LEU
1	A	606	GLN
1	A	616	LEU
1	A	619	LEU
1	A	653	PRO
1	A	672	TYR
1	B	79	ASN
1	B	104	ARG
1	B	105	TYR
1	B	139	THR
1	B	200	ILE
1	B	251	GLU
1	B	267	LYS
1	B	294	GLN
1	B	322	VAL
1	B	338	LEU

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Mol	Chain	Res	Type
1	B	346	MET
1	B	350	ASN
1	B	353	GLU
1	B	362	PRO
1	B	371	LEU
1	B	384	ASN
1	B	414	GLU
1	B	454	ARG
1	B	511	MET
1	B	573	TYR
1	B	591	LEU
1	B	606	GLN
1	B	616	LEU
1	B	619	LEU
1	B	653	PRO
1	B	672	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	180	HIS
1	A	250	HIS
1	A	293	ASN
1	A	294	GLN
1	A	302	HIS
1	A	350	ASN
1	A	384	ASN
1	A	399	GLN
1	A	465	HIS
1	A	486	ASN
1	A	607	HIS
1	B	79	ASN
1	B	180	HIS
1	B	250	HIS
1	B	293	ASN
1	B	294	GLN
1	B	302	HIS
1	B	350	ASN
1	B	384	ASN
1	B	399	GLN
1	B	465	HIS

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Mol	Chain	Res	Type
1	B	486	ASN
1	B	607	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	FAD	A	750	-	54,58,58	1.68	7 (12%)	71,89,89	1.51	14 (19%)
2	FAD	B	750	-	40,46,58	2.29	6 (15%)	41,67,89	1.32	7 (17%)
3	FMN	A	751	-	33,33,33	1.39	5 (15%)	48,50,50	1.50	8 (16%)
4	NAP	A	752	-	28,33,52	1.09	1 (3%)	40,52,80	1.70	9 (22%)
3	FMN	B	751	-	33,33,33	1.39	5 (15%)	48,50,50	1.50	9 (18%)
4	NAP	B	752	-	28,33,52	1.09	1 (3%)	40,52,80	1.69	9 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	A	750	-	-	4/30/50/50	0/6/6/6
2	FAD	B	750	-	-	4/30/50/50	0/3/4/6
3	FMN	A	751	-	-	0/18/18/18	0/3/3/3
4	NAP	A	752	-	-	7/17/37/67	0/3/3/5
3	FMN	B	751	-	-	0/18/18/18	0/3/3/3
4	NAP	B	752	-	-	7/17/37/67	0/3/3/5

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	750	FAD	C4X-N5	8.64	1.37	1.27
2	A	750	FAD	PA-O3P	-6.64	1.52	1.59
2	B	750	FAD	PA-O3P	-6.59	1.52	1.59
2	A	750	FAD	P-O3P	-6.06	1.53	1.59
2	B	750	FAD	P-O3P	-6.02	1.53	1.59
3	B	751	FMN	C9-C8	-4.33	1.33	1.39
3	A	751	FMN	C9-C8	-4.28	1.33	1.39
2	B	750	FAD	C10-C4X	-3.94	1.39	1.49
2	A	750	FAD	C4X-N5	3.27	1.37	1.30
2	B	750	FAD	C10-N10	-3.14	1.43	1.46
3	B	751	FMN	C6-C7	-3.02	1.35	1.39
3	A	751	FMN	C6-C7	-3.02	1.35	1.39
2	A	750	FAD	C10-N10	2.99	1.43	1.37
2	A	750	FAD	C6-C7	-2.46	1.36	1.39
4	A	752	NAP	P2B-O2B	-2.33	1.55	1.59
3	A	751	FMN	C4A-N5	2.31	1.35	1.30
4	B	752	NAP	P2B-O2B	-2.30	1.55	1.59
2	A	750	FAD	O4B-C1B	2.30	1.43	1.40
2	B	750	FAD	O4B-C1B	2.30	1.43	1.40
3	B	751	FMN	C4A-N5	2.30	1.35	1.30
3	A	751	FMN	C10-N10	2.21	1.42	1.37
3	B	751	FMN	C10-N10	2.21	1.42	1.37
2	A	750	FAD	C9-C9A	2.09	1.43	1.39
3	A	751	FMN	P-O3P	-2.05	1.47	1.54
3	B	751	FMN	P-O3P	-2.03	1.47	1.54

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	750	FAD	C9A-C5X-N5	4.78	127.53	122.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	751	FMN	C9A-C5A-N5	4.66	127.40	122.45
3	B	751	FMN	C9A-C5A-N5	4.62	127.36	122.45
4	A	752	NAP	C1B-N9A-C4A	-4.17	119.32	126.64
4	B	752	NAP	C1B-N9A-C4A	-4.16	119.33	126.64
3	B	751	FMN	C5A-N5-C4A	-4.11	111.43	118.09
3	A	751	FMN	C5A-N5-C4A	-4.11	111.44	118.09
2	B	750	FAD	C2B-C3B-C4B	-3.87	95.13	102.61
2	A	750	FAD	C2B-C3B-C4B	-3.85	95.18	102.61
4	A	752	NAP	N3A-C2A-N1A	-3.69	123.66	128.67
4	B	752	NAP	N3A-C2A-N1A	-3.68	123.68	128.67
2	A	750	FAD	C5X-N5-C4X	-3.32	112.71	118.09
4	A	752	NAP	C2B-C1B-N9A	3.23	119.73	112.56
4	B	752	NAP	C2B-C1B-N9A	3.23	119.73	112.56
4	A	752	NAP	O2N-PN-O1N	2.91	122.17	110.83
4	B	752	NAP	O2N-PN-O1N	2.90	122.12	110.83
3	B	751	FMN	O2-C2-N3	2.89	124.12	118.58
3	A	751	FMN	O2-C2-N3	2.89	124.12	118.58
2	A	750	FAD	O2-C2-N3	2.87	124.08	118.58
2	A	750	FAD	C10-N1-C2	2.73	122.76	116.85
3	B	751	FMN	C10-N1-C2	2.62	122.51	116.85
2	A	750	FAD	C5A-C6A-N6A	2.61	124.29	120.31
2	B	750	FAD	C5A-C6A-N6A	2.61	124.29	120.31
3	A	751	FMN	C10-N1-C2	2.61	122.49	116.85
4	A	752	NAP	O5D-PN-O3	-2.59	95.95	104.64
4	B	752	NAP	O5D-PN-O3	-2.58	95.99	104.64
2	B	750	FAD	O3'-C3'-C2'	2.46	114.53	108.93
2	A	750	FAD	C5X-C9A-N10	-2.45	115.75	117.97
2	A	750	FAD	N3-C2-N1	-2.45	114.30	119.50
2	A	750	FAD	O3'-C3'-C2'	2.45	114.49	108.93
3	A	751	FMN	O3P-P-O2P	2.29	116.40	107.80
3	B	751	FMN	O3P-P-O2P	2.29	116.39	107.80
4	B	752	NAP	C4A-C5A-N7A	2.27	111.73	109.34
4	A	752	NAP	O2A-PA-O3	2.26	113.38	107.27
4	B	752	NAP	O2A-PA-O3	2.25	113.37	107.27
4	A	752	NAP	C4A-C5A-N7A	2.25	111.72	109.34
4	B	752	NAP	C5A-C6A-N6A	2.22	123.70	120.31
2	B	750	FAD	C4B-O4B-C1B	-2.22	107.89	109.92
2	A	750	FAD	C4B-O4B-C1B	-2.21	107.90	109.92
4	A	752	NAP	C5A-C6A-N6A	2.21	123.67	120.31
3	A	751	FMN	O3'-C3'-C4'	-2.16	104.01	108.93
4	B	752	NAP	O3B-C3B-C4B	-2.16	104.87	111.08
3	B	751	FMN	O3'-C3'-C4'	-2.16	104.02	108.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	752	NAP	O3B-C3B-C4B	-2.16	104.88	111.08
2	B	750	FAD	O5'-C5'-C4'	2.13	115.06	109.36
3	B	751	FMN	N3-C2-N1	-2.13	114.98	119.50
2	A	750	FAD	O5'-C5'-C4'	2.13	115.04	109.36
3	A	751	FMN	N3-C2-N1	-2.12	114.98	119.50
2	A	750	FAD	C8M-C8-C9	2.08	123.24	119.57
2	B	750	FAD	O2A-PA-O1A	2.08	122.10	112.44
2	A	750	FAD	O2A-PA-O1A	2.07	122.06	112.44
2	A	750	FAD	O2B-C2B-C3B	-2.05	105.26	111.82
2	B	750	FAD	O2B-C2B-C3B	-2.04	105.28	111.82
3	A	751	FMN	C2'-C1'-N10	2.02	119.74	110.20
3	B	751	FMN	C2'-C1'-N10	2.02	119.73	110.20
3	B	751	FMN	C4-C4A-N5	-2.00	115.44	118.21

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	750	FAD	C1'-C2'-C3'-C4'
2	B	750	FAD	C1'-C2'-C3'-C4'
4	A	752	NAP	C5B-O5B-PA-O1A
4	A	752	NAP	C5B-O5B-PA-O2A
4	A	752	NAP	C5B-O5B-PA-O3
4	B	752	NAP	C5B-O5B-PA-O1A
4	B	752	NAP	C5B-O5B-PA-O2A
4	B	752	NAP	C5B-O5B-PA-O3
2	A	750	FAD	O2'-C2'-C3'-C4'
2	B	750	FAD	O2'-C2'-C3'-C4'
4	A	752	NAP	PN-O3-PA-O1A
4	B	752	NAP	PN-O3-PA-O1A
2	A	750	FAD	O2'-C2'-C3'-O3'
2	B	750	FAD	O2'-C2'-C3'-O3'
4	A	752	NAP	C2B-O2B-P2B-O1X
4	B	752	NAP	C2B-O2B-P2B-O1X
2	A	750	FAD	C1'-C2'-C3'-O3'
2	B	750	FAD	C1'-C2'-C3'-O3'
4	A	752	NAP	C2B-O2B-P2B-O3X
4	B	752	NAP	C2B-O2B-P2B-O3X
4	A	752	NAP	PN-O3-PA-O2A
4	B	752	NAP	PN-O3-PA-O2A

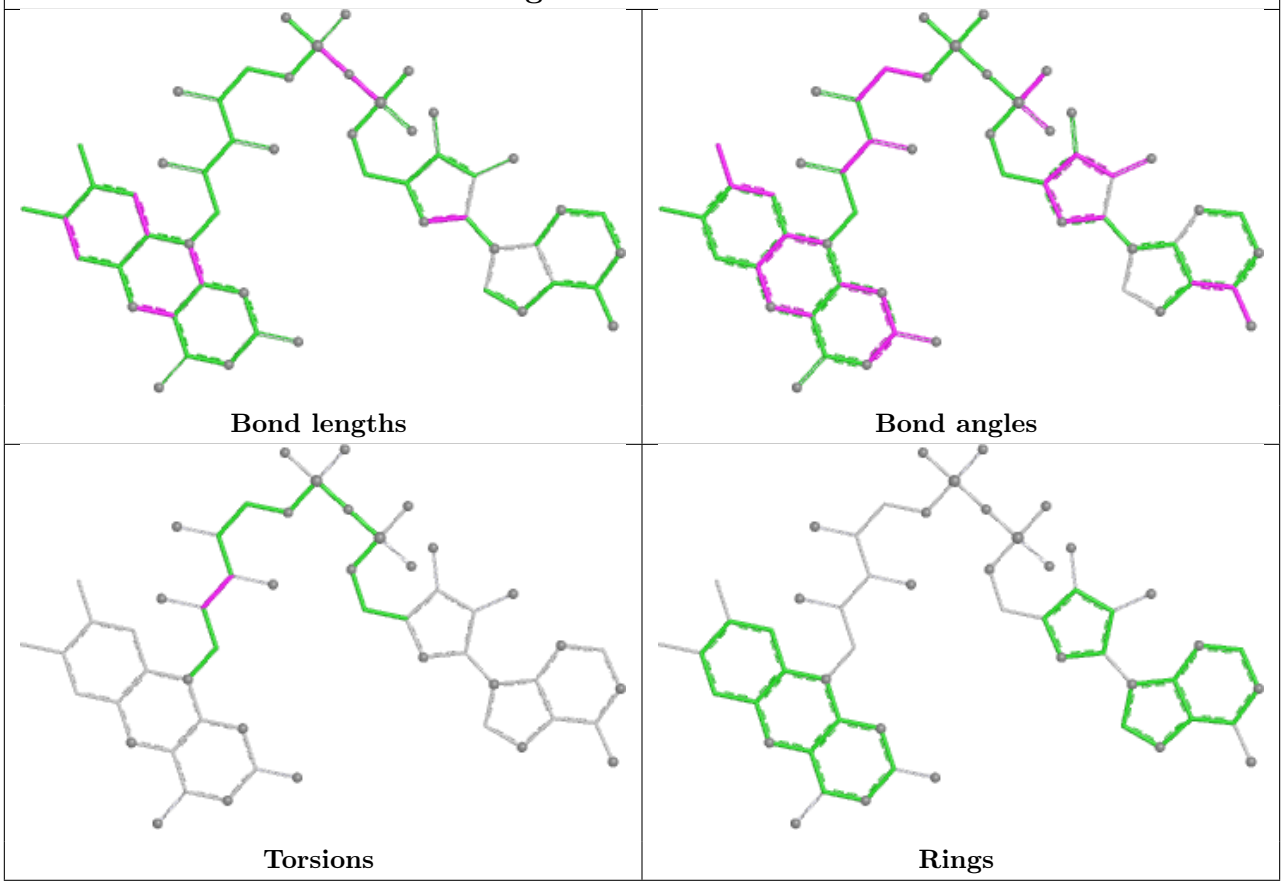
There are no ring outliers.

6 monomers are involved in 15 short contacts:

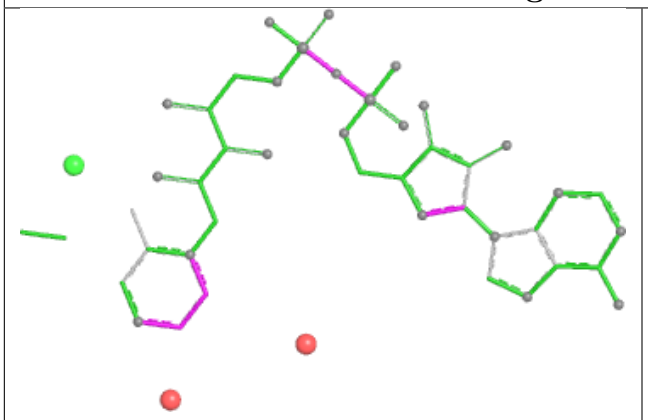
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	750	FAD	2	0
2	B	750	FAD	9	0
3	A	751	FMN	1	0
4	A	752	NAP	1	0
3	B	751	FMN	1	0
4	B	752	NAP	1	0

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

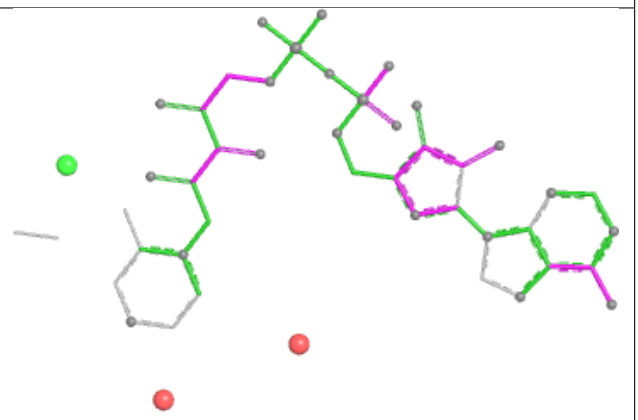
Ligand FAD A 750



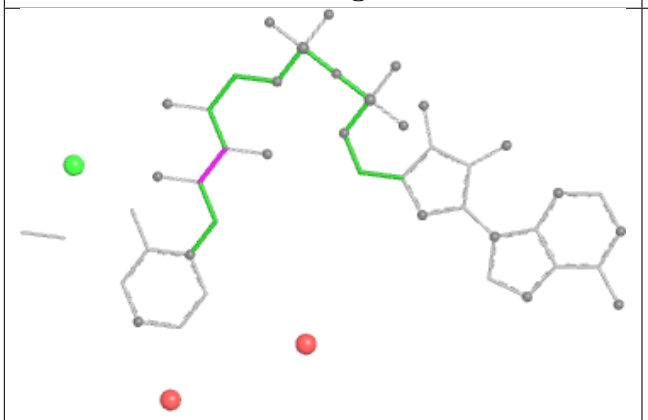
Ligand FAD B 750



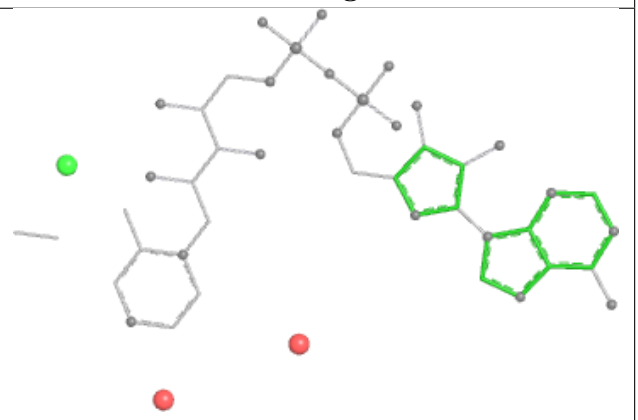
Bond lengths



Bond angles

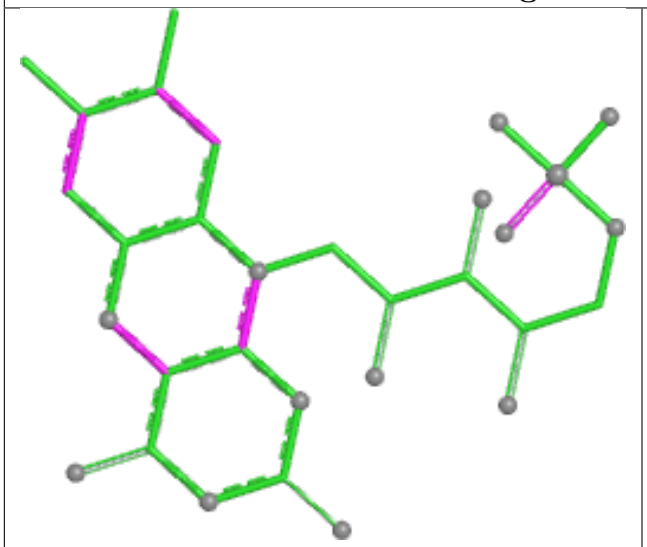


Torsions

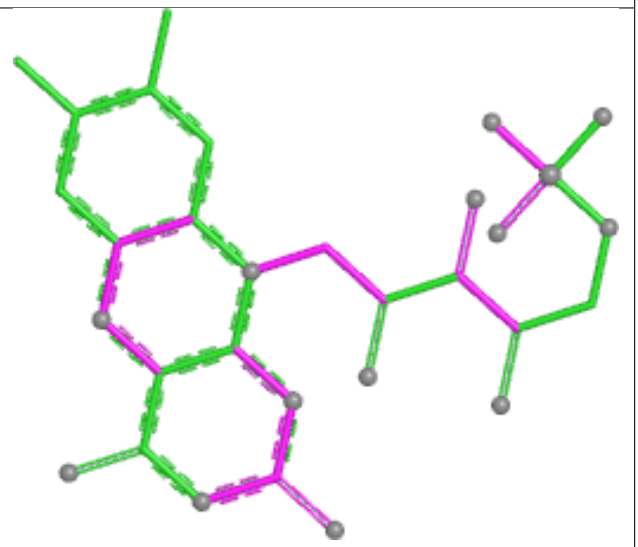


Rings

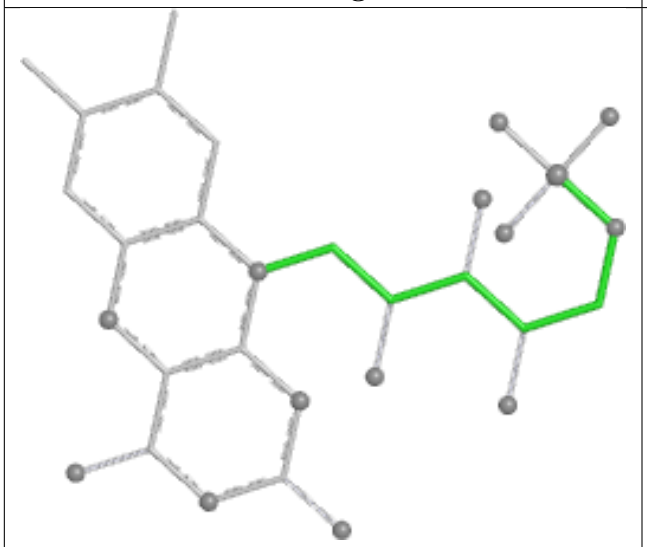
Ligand FMN A 751



Bond lengths



Bond angles



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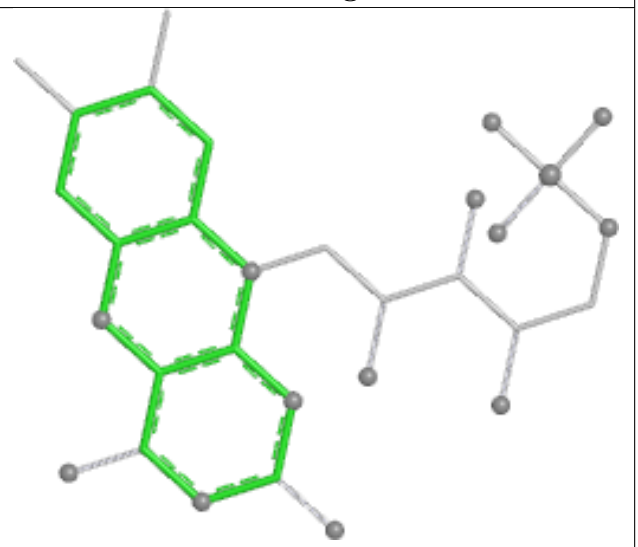
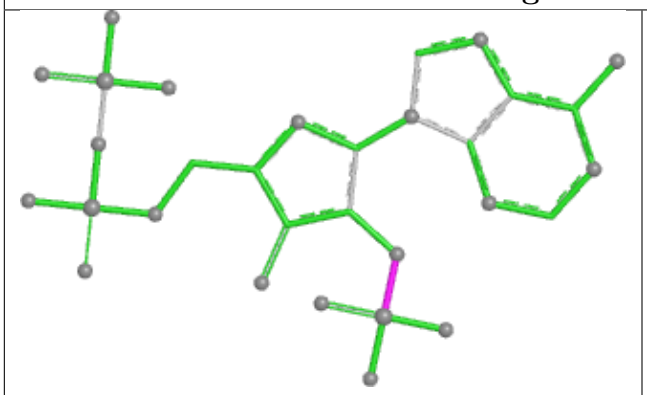
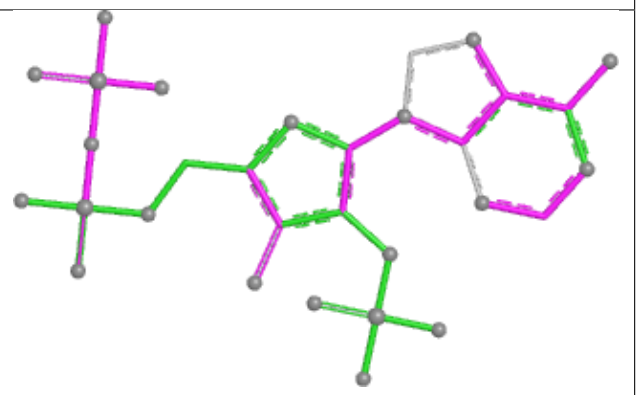


Rings

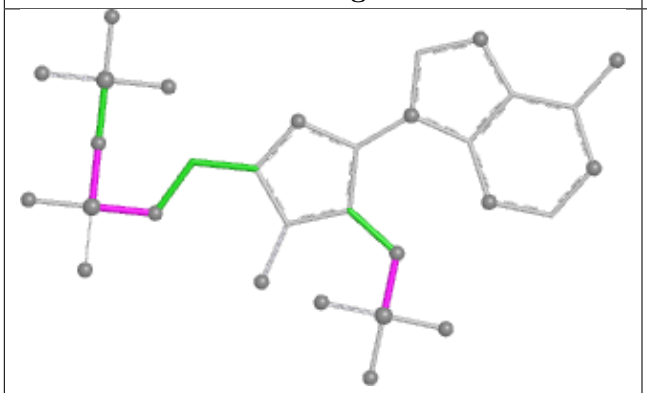
Ligand NAP A 752



Bond lengths



Bond angles



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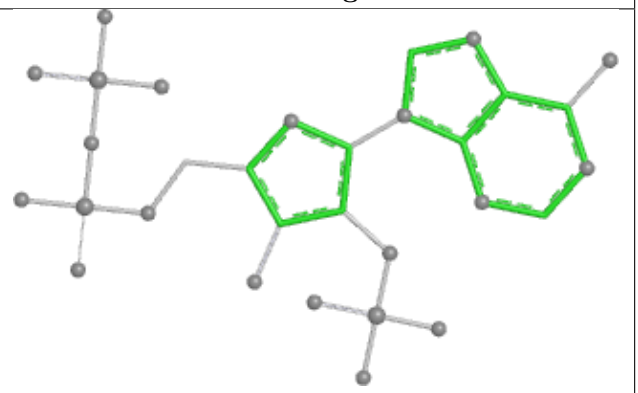
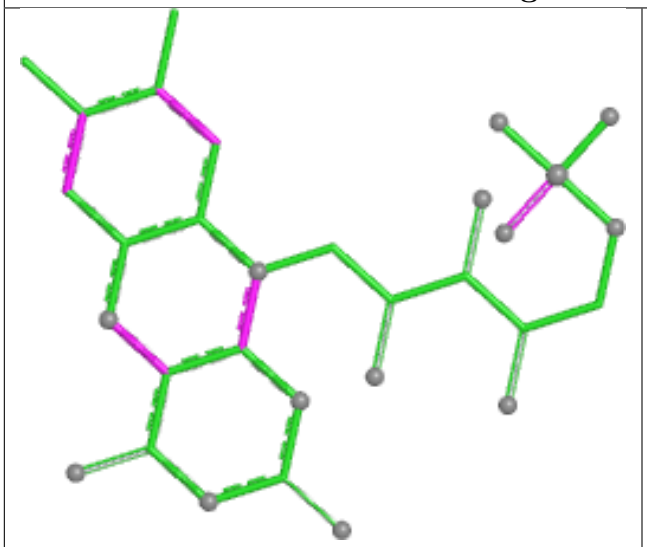
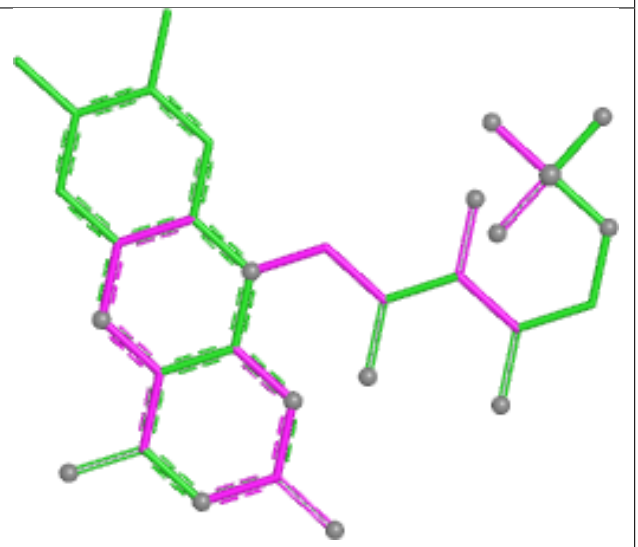


Rings

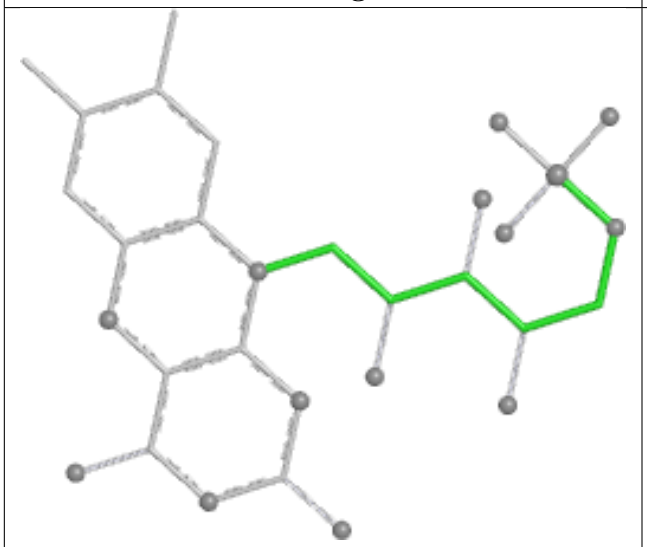
Ligand FMN B 751



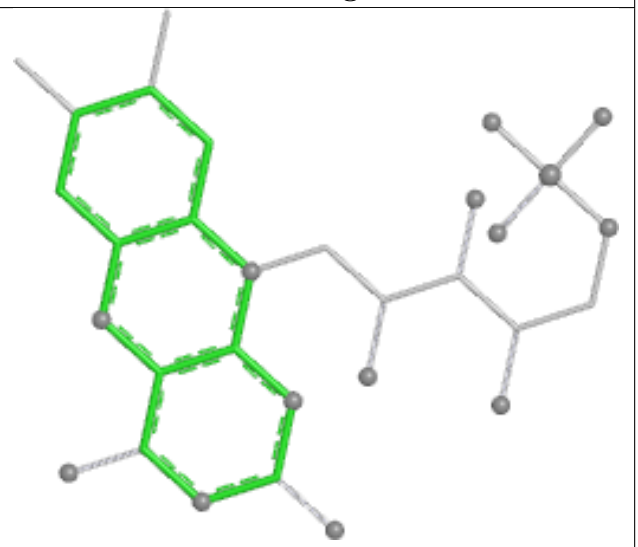
Bond lengths



Bond angles

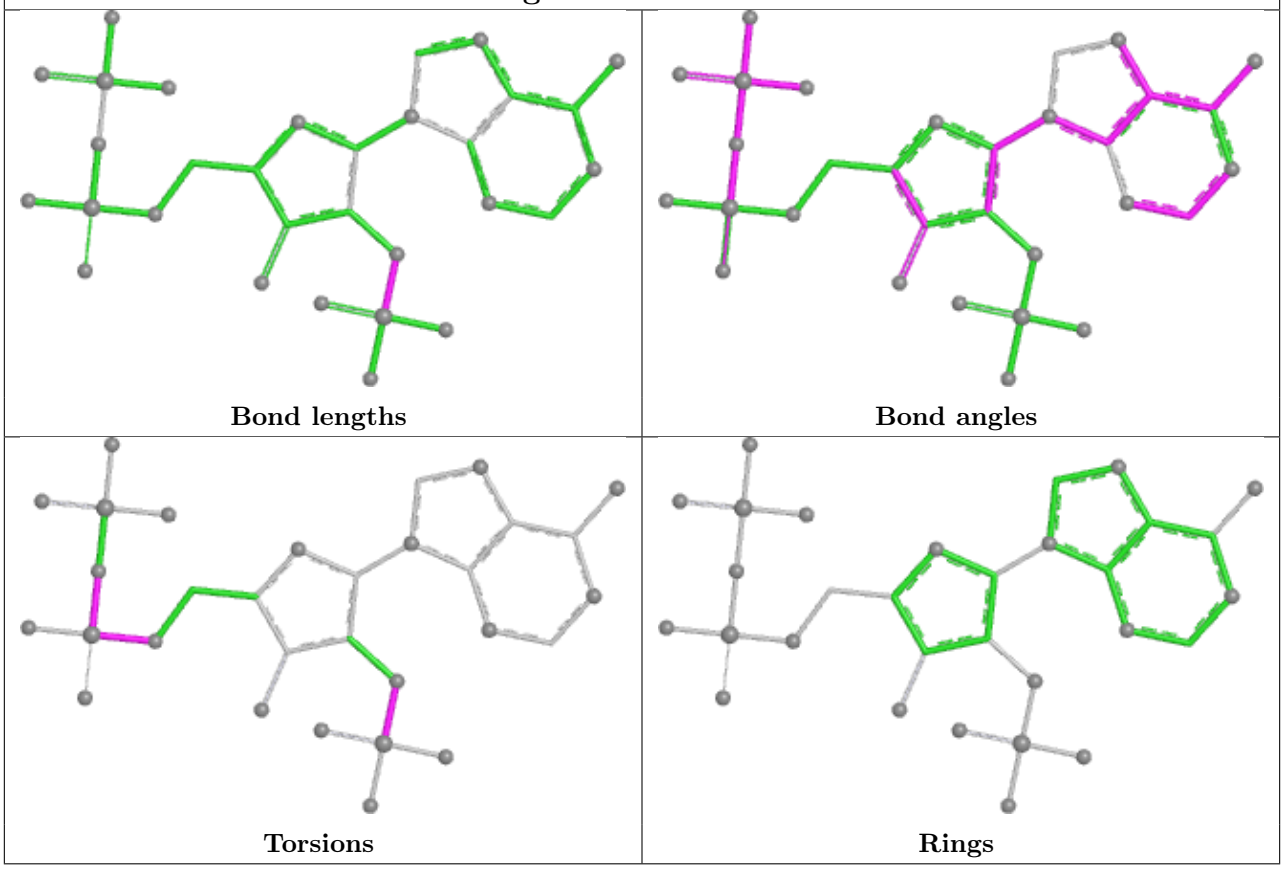


Torsions



Rings

Ligand NAP B 752



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.