



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:07 AM UTC

PDB ID : 3QFC / pdb_00003qfc
Title : Crystal Structure of Human NADPH-Cytochrome P450 (V492E mutant)
Authors : Xia, C.; Marohnic, C.; Panda, S.P.; Masters, B.S.; Kim, J.-J.P.
Deposited on : 2011-01-21
Resolution : 1.80 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ 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

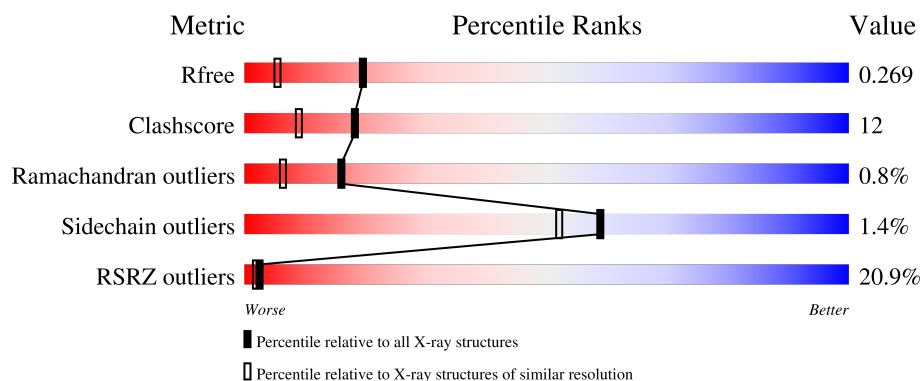
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	618	5% 80% 18% ..
1	B	618	36% 65% 29% ..

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10573 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	607	Total	C	N	O	S	0	0	0
			4844	3064	837	920	23			
1	B	595	Total	C	N	O	S	0	0	0
			4761	3014	822	902	23			

There are 14 discrepancies between the modelled and reference sequences:

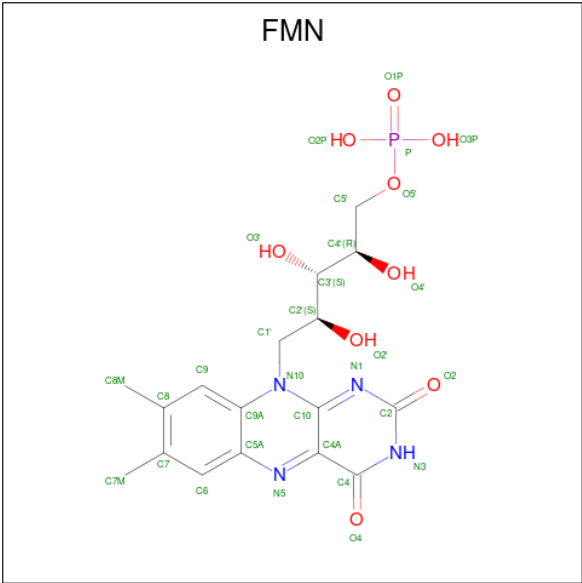
Chain	Residue	Modelled	Actual	Comment	Reference
A	63	GLY	-	expression tag	UNP P16435
A	64	SER	-	expression tag	UNP P16435
A	65	HIS	-	expression tag	UNP P16435
A	66	MET	-	expression tag	UNP P16435
A	228	LEU	PRO	SEE REMARK 999	UNP P16435
A	492	GLU	VAL	engineered mutation	UNP P16435
A	503	VAL	ALA	SEE REMARK 999	UNP P16435
B	63	GLY	-	expression tag	UNP P16435
B	64	SER	-	expression tag	UNP P16435
B	65	HIS	-	expression tag	UNP P16435
B	66	MET	-	expression tag	UNP P16435
B	228	LEU	PRO	SEE REMARK 999	UNP P16435
B	492	GLU	VAL	engineered mutation	UNP P16435
B	503	VAL	ALA	SEE REMARK 999	UNP P16435

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



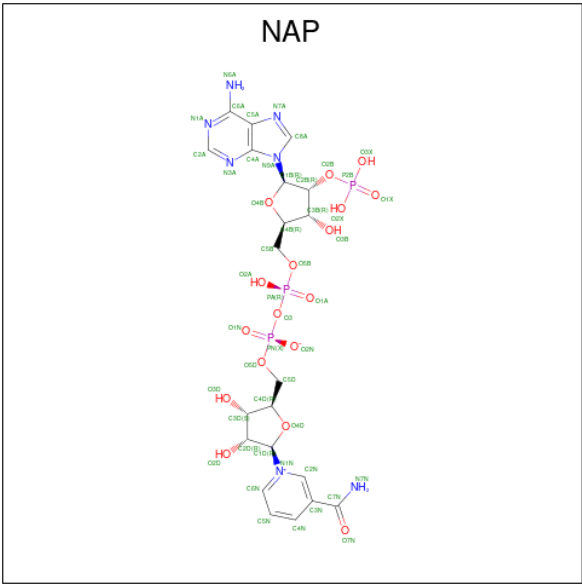
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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



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

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



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
			48	21	7	17	3		

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		

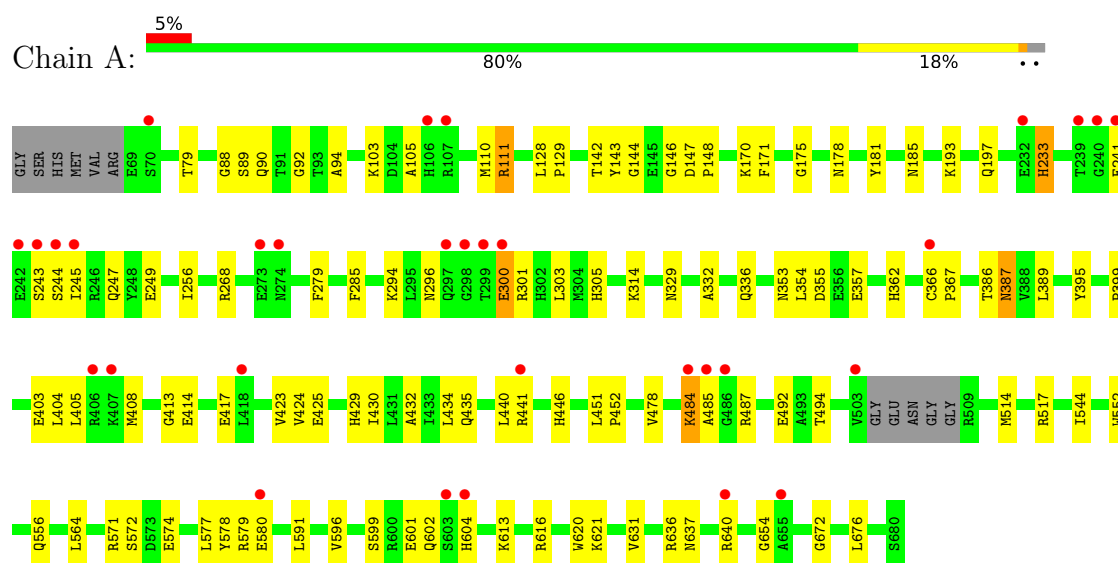
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	567	Total	O	0	0
			567	567		
6	B	153	Total	O	0	0
			153	153		

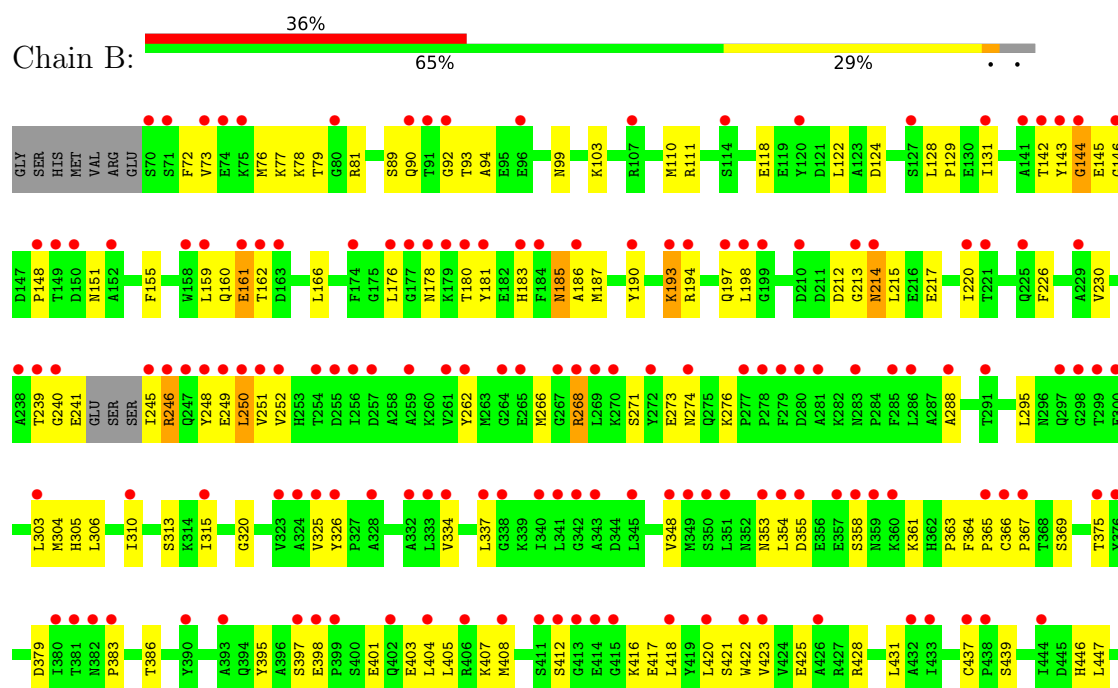
3 Residue-property plots [i](#)

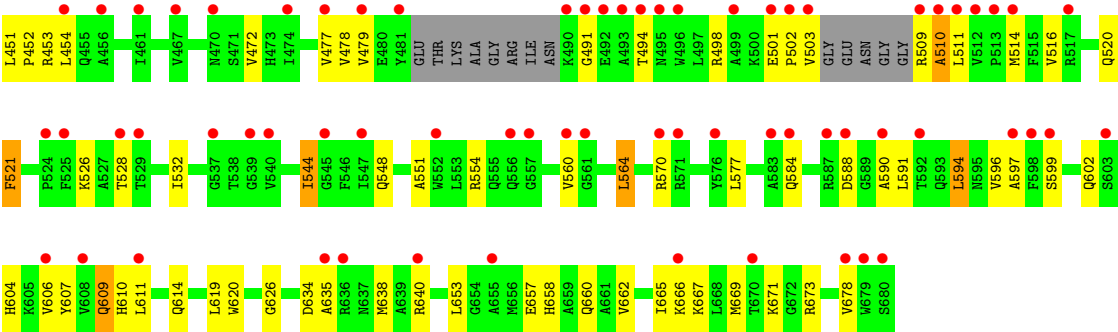
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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	70.72Å 118.51Å 155.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.42 – 1.80 45.42 – 1.80	Depositor EDS
% Data completeness (in resolution range)	93.1 (45.42-1.80) 93.2 (45.42-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 1.79Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.232 , 0.265 0.238 , 0.269	Depositor DCC
R_{free} test set	5702 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10573	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.87% 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: FAD, FMN, CA, NAP

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	0.39	0/4953	0.87	14/6704 (0.2%)
1	B	0.31	0/4868	0.79	3/6587 (0.0%)
All	All	0.35	0/9821	0.83	17/13291 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	424	VAL	N-CA-C	7.71	117.67	110.42
1	B	92	GLY	N-CA-C	7.39	124.81	115.42
1	A	599	SER	N-CA-C	6.83	119.74	111.82
1	A	386	THR	N-CA-C	6.82	119.55	111.71
1	A	92	GLY	N-CA-C	6.60	124.11	115.40
1	A	544	ILE	N-CA-C	-6.39	104.10	110.62
1	A	484	LYS	N-CA-C	-5.67	106.21	113.01
1	B	544	ILE	N-CA-C	-5.61	105.06	110.72
1	A	430	ILE	N-CA-C	5.55	116.28	110.62
1	A	88	GLY	N-CA-C	-5.45	100.27	113.18
1	B	599	SER	N-CA-C	5.44	119.18	112.54
1	A	596	VAL	N-CA-C	5.34	115.58	108.11
1	A	636	ARG	N-CA-C	5.30	116.74	111.07
1	A	514	MET	N-CA-C	5.28	116.02	108.74
1	A	171	PHE	N-CA-C	5.24	116.95	109.14
1	A	233	HIS	N-CA-C	5.22	117.05	111.36
1	A	256	ILE	N-CA-C	5.12	116.61	109.80

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	4844	0	4699	85	0
1	B	4761	0	4621	153	0
2	A	53	0	31	0	0
2	B	53	0	31	0	0
3	A	31	0	19	0	0
3	B	31	0	19	1	0
4	A	31	0	11	0	0
4	B	48	0	25	1	0
5	A	1	0	0	0	0
6	A	567	0	0	8	0
6	B	153	0	0	3	0
All	All	10573	0	9456	233	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 12.

All (233) 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:412:SER:HA	1:B:416:LYS:HD2	1.33	1.04
1:B:479:VAL:H	1:B:494:THR:HG21	1.33	0.93
1:B:180:THR:HA	1:B:640:ARG:HH21	1.34	0.92
1:B:315:ILE:HD11	1:B:516:VAL:HG23	1.55	0.88
1:A:300:GLU:HG3	1:A:301:ARG:H	1.38	0.88
1:B:609:GLN:HE22	4:B:753:NAP:H2A	1.38	0.87
1:A:405:LEU:HA	1:A:408:MET:HE3	1.57	0.84
1:A:435:GLN:HE21	1:A:487:ARG:HH11	1.27	0.82
1:B:607:TYR:H	1:B:610:HIS:HD2	1.22	0.82
1:A:89:SER:HB2	1:A:94:ALA:HB3	1.66	0.77
1:A:637:ASN:ND2	1:A:640:ARG:HH22	1.82	0.77
1:A:425:GLU:HG3	1:A:484:LYS:HZ1	1.50	0.76
1:B:412:SER:HA	1:B:416:LYS:CD	2.13	0.76
1:B:502:PRO:O	1:B:503:VAL:HG23	1.84	0.76
1:A:247:GLN:HE21	1:A:353:ASN:HD21	1.32	0.76
1:B:408:MET:HE1	1:B:418:LEU:HB3	1.66	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:VAL:HG12	1:B:514:MET:HB3	1.68	0.76
1:B:89:SER:HB2	1:B:94:ALA:HB3	1.66	0.75
1:A:637:ASN:HD22	1:A:640:ARG:HH22	1.34	0.75
1:B:250:LEU:HD13	1:B:251:VAL:N	2.03	0.73
1:B:609:GLN:H	1:B:609:GLN:HE21	1.36	0.72
1:A:404:LEU:HG	1:A:408:MET:HE2	1.69	0.72
1:A:404:LEU:HD22	1:B:274:ASN:HD21	1.55	0.71
1:A:621:LYS:HD2	6:A:1061:HOH:O	1.91	0.71
1:B:397:SER:HB2	1:B:439:SER:HB2	1.72	0.70
1:B:361:LYS:HE2	1:B:361:LYS:HA	1.73	0.70
1:B:353:ASN:HD22	1:B:354:LEU:N	1.89	0.70
1:B:310:ILE:HD13	1:B:472:VAL:HG23	1.74	0.69
1:A:435:GLN:NE2	1:A:487:ARG:HH11	1.90	0.68
1:B:240:GLY:O	1:B:241:GLU:HB2	1.92	0.68
1:B:194:ARG:HH11	1:B:197:GLN:NE2	1.92	0.68
1:A:429:HIS:HD2	1:A:432:ALA:H	1.42	0.67
1:B:180:THR:HG22	1:B:640:ARG:NH2	2.09	0.67
1:B:509:ARG:O	1:B:511:LEU:HG	1.94	0.66
1:B:81:ARG:HD3	1:B:111:ARG:HB3	1.77	0.66
1:B:268:ARG:HB3	1:B:271:SER:HB3	1.75	0.66
1:B:544:ILE:O	1:B:548:GLN:HG3	1.95	0.66
1:B:379:ASP:HB3	1:B:452:PRO:HG2	1.78	0.65
1:A:414:GLU:CD	1:B:276:LYS:HE2	2.22	0.64
1:A:440:LEU:O	1:A:441:ARG:HD3	1.97	0.64
1:B:313:SER:OG	1:B:315:ILE:HG22	1.97	0.64
1:B:671:LYS:NZ	1:B:673:ARG:HH11	1.96	0.63
1:B:607:TYR:H	1:B:610:HIS:CD2	2.10	0.63
1:A:362:HIS:CD2	1:A:366:CYS:SG	2.92	0.63
1:B:478:VAL:HA	1:B:494:THR:OG1	1.99	0.62
1:B:245:ILE:HG13	1:B:245:ILE:O	1.99	0.62
1:B:657:GLU:H	1:B:660:GLN:HE21	1.44	0.62
1:A:245:ILE:N	1:A:245:ILE:HD12	2.13	0.62
1:A:429:HIS:CD2	1:A:432:ALA:H	2.18	0.62
1:A:616:ARG:HD2	6:A:57:HOH:O	1.99	0.62
1:B:183:HIS:HB3	1:B:186:ALA:HB2	1.82	0.61
1:B:78:LYS:HD3	1:B:78:LYS:O	2.00	0.61
1:A:305:HIS:HD2	1:A:578:TYR:OH	1.84	0.61
1:B:609:GLN:H	1:B:609:GLN:NE2	1.97	0.61
1:B:178:ASN:OD1	1:B:212:ASP:HB2	2.01	0.61
1:A:399:PRO:O	1:A:403:GLU:HG3	2.01	0.61
1:B:90:GLN:HB2	1:B:143:TYR:CZ	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ALA:HB3	6:A:992:HOH:O	2.01	0.59
1:A:353:ASN:ND2	1:A:355:ASP:H	2.00	0.59
1:B:564:LEU:HD12	1:B:564:LEU:N	2.16	0.59
1:A:613:LYS:HE2	6:A:1192:HOH:O	2.02	0.59
1:A:425:GLU:HG3	1:A:484:LYS:NZ	2.16	0.59
1:B:93:THR:HG21	1:B:215:LEU:HD23	1.84	0.59
1:B:353:ASN:ND2	1:B:355:ASP:H	2.00	0.59
1:B:355:ASP:OD2	1:B:358:SER:HB2	2.03	0.59
1:B:401:GLU:OE2	1:B:437:CYS:HA	2.04	0.58
1:A:193:LYS:O	1:A:197:GLN:HG3	2.04	0.57
1:B:160:GLN:HG3	1:B:161:GLU:OE2	2.04	0.57
1:B:239:THR:HG23	1:B:240:GLY:H	1.70	0.57
1:A:571:ARG:HH12	1:A:601:GLU:HA	1.70	0.57
1:A:314:LYS:NZ	1:A:314:LYS:HB3	2.19	0.57
1:A:485:ALA:HB3	6:A:814:HOH:O	2.04	0.57
1:B:250:LEU:HD13	1:B:251:VAL:H	1.68	0.57
1:B:597:ALA:HB1	1:B:606:VAL:HG12	1.87	0.57
1:A:247:GLN:HE21	1:A:353:ASN:ND2	2.02	0.57
1:B:662:VAL:O	1:B:666:LYS:HG2	2.05	0.57
1:B:310:ILE:HD12	1:B:315:ILE:HG23	1.86	0.57
1:B:403:GLU:O	1:B:407:LYS:HG3	2.05	0.56
1:B:610:HIS:O	1:B:614:GLN:HG3	2.05	0.56
1:B:159:LEU:HB3	1:B:194:ARG:HG2	1.88	0.56
1:B:160:GLN:HB3	1:B:190:TYR:OH	2.06	0.56
1:A:353:ASN:HD22	1:A:354:LEU:N	2.03	0.56
1:B:609:GLN:HE21	1:B:609:GLN:N	2.04	0.56
1:B:303:LEU:HD22	1:B:577:LEU:HD21	1.88	0.55
1:B:288:ALA:HA	1:B:510:ALA:O	2.07	0.55
1:A:294:LYS:HE3	1:A:296:ASN:O	2.06	0.55
1:B:239:THR:HG23	1:B:240:GLY:N	2.21	0.54
1:B:303:LEU:N	1:B:303:LEU:HD12	2.22	0.54
1:A:413:GLY:O	1:A:417:GLU:HG2	2.08	0.54
1:A:425:GLU:HG3	1:A:484:LYS:CE	2.37	0.54
1:B:110:MET:HE1	1:B:230:VAL:HG11	1.90	0.54
1:B:142:THR:CG2	1:B:146:GLY:HA2	2.38	0.54
1:A:389:LEU:HD11	1:A:423:VAL:HG21	1.89	0.54
1:B:262:TYR:CD2	1:B:365:PRO:HB3	2.42	0.54
1:A:362:HIS:CG	1:A:366:CYS:SG	3.01	0.54
1:A:395:TYR:OH	1:A:446:HIS:HD2	1.91	0.54
1:B:671:LYS:HZ1	1:B:673:ARG:HH11	1.54	0.54
1:B:491:GLY:HA3	1:B:494:THR:HG22	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:304:MET:HG3	1:B:306:LEU:HD11	1.89	0.54
1:B:249:GLU:HB3	1:B:354:LEU:HD21	1.89	0.53
1:A:366:CYS:HB2	1:A:367:PRO:HA	1.88	0.53
1:B:303:LEU:CD2	1:B:577:LEU:HD21	2.38	0.53
1:B:401:GLU:O	1:B:405:LEU:HG	2.09	0.53
1:B:295:LEU:HD21	1:B:305:HIS:HB2	1.91	0.52
1:A:353:ASN:HD22	1:A:355:ASP:H	1.57	0.52
1:A:572:SER:O	1:A:579:ARG:NH1	2.43	0.52
1:B:273:GLU:HG2	6:B:807:HOH:O	2.09	0.52
1:A:170:LYS:NZ	1:A:233:HIS:HE1	2.08	0.51
1:B:220:ILE:HD11	1:B:386:THR:HG21	1.92	0.51
1:A:332:ALA:O	1:A:336:GLN:HG3	2.11	0.51
1:B:145:GLU:OE1	1:B:520:GLN:HG2	2.10	0.51
1:B:78:LYS:HD3	1:B:78:LYS:C	2.36	0.51
1:B:348:VAL:HA	1:B:369:SER:HA	1.93	0.51
1:B:315:ILE:HD11	1:B:516:VAL:CG2	2.36	0.50
1:A:404:LEU:CD2	1:B:274:ASN:HD21	2.23	0.50
1:B:268:ARG:HE	1:B:268:ARG:HA	1.76	0.50
1:A:147:ASP:HB3	1:A:148:PRO:HD2	1.94	0.50
1:B:144:GLY:H	3:B:751:FMN:H6	1.77	0.50
1:B:404:LEU:C	1:B:404:LEU:HD23	2.37	0.50
1:A:303:LEU:HD22	1:A:577:LEU:HD21	1.94	0.49
1:B:176:LEU:C	1:B:185:ASN:HD21	2.21	0.49
1:B:358:SER:O	1:B:361:LYS:HE3	2.12	0.49
1:B:353:ASN:HD22	1:B:353:ASN:C	2.18	0.49
1:A:564:LEU:N	1:A:564:LEU:HD12	2.28	0.48
1:A:631:VAL:HB	1:A:676:LEU:HD23	1.94	0.48
1:B:99:ASN:O	1:B:103:LYS:HG3	2.13	0.48
1:A:387:ASN:H	1:A:387:ASN:HD22	1.62	0.48
1:B:122:LEU:HG	1:B:155:PHE:CD1	2.48	0.48
1:B:607:TYR:N	1:B:610:HIS:HD2	2.01	0.48
1:B:128:LEU:N	1:B:129:PRO:CD	2.77	0.48
1:A:175:GLY:C	1:A:185:ASN:HD21	2.22	0.47
1:A:672:GLY:HA2	6:A:1032:HOH:O	2.15	0.47
1:B:477:VAL:O	1:B:477:VAL:HG13	2.14	0.47
1:B:526:LYS:N	1:B:526:LYS:HD2	2.30	0.47
1:B:337:LEU:HD23	1:B:431:LEU:HD13	1.96	0.47
1:B:602:GLN:HG2	1:B:604:HIS:O	2.15	0.47
1:A:268:ARG:HB2	1:A:268:ARG:NH1	2.29	0.47
1:B:532:ILE:HD13	1:B:619:LEU:HD22	1.96	0.47
1:A:249:GLU:HB3	1:A:354:LEU:HD21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:ASN:HB2	1:A:574:GLU:O	2.14	0.47
1:B:118:GLU:HB2	1:B:151:ASN:O	2.14	0.47
1:B:214:ASN:ND2	1:B:420:LEU:HD22	2.30	0.46
1:A:79:THR:HG22	1:A:357:GLU:CD	2.40	0.46
1:A:425:GLU:CG	1:A:484:LYS:HZ1	2.23	0.46
1:A:435:GLN:HE21	1:A:487:ARG:NH1	2.02	0.46
1:A:366:CYS:HA	1:A:367:PRO:C	2.41	0.46
1:A:640:ARG:NH1	1:A:640:ARG:HB3	2.31	0.46
1:B:337:LEU:HD21	1:B:431:LEU:HB2	1.98	0.46
1:B:570:ARG:HG3	1:B:570:ARG:HH11	1.79	0.46
1:A:103:LYS:HE2	1:A:247:GLN:OE1	2.16	0.46
1:A:571:ARG:NH1	1:A:601:GLU:HA	2.29	0.46
1:A:602:GLN:HG2	1:A:604:HIS:O	2.16	0.45
1:B:635:ALA:HB2	1:B:678:VAL:HB	1.97	0.45
1:A:414:GLU:OE1	1:B:276:LYS:HE2	2.14	0.45
1:A:110:MET:O	1:A:111:ARG:HD2	2.16	0.45
1:B:526:LYS:HB3	1:B:528:THR:HG22	1.98	0.45
1:B:594:LEU:HD22	1:B:596:VAL:CG2	2.47	0.45
1:B:383:PRO:HB3	1:B:423:VAL:HG12	1.97	0.45
1:A:241:GLU:C	1:A:243:SER:H	2.24	0.45
1:B:334:VAL:HG21	1:B:375:THR:OG1	2.17	0.45
1:B:422:TRP:CE2	1:B:428:ARG:NH1	2.85	0.45
1:B:304:MET:HG3	1:B:306:LEU:CD1	2.47	0.45
1:B:364:PHE:HB2	1:B:365:PRO:CD	2.47	0.45
1:B:420:LEU:O	1:B:425:GLU:HB2	2.17	0.45
1:B:479:VAL:O	1:B:491:GLY:N	2.44	0.45
1:B:193:LYS:NZ	1:B:193:LYS:HB3	2.32	0.45
1:B:159:LEU:O	1:B:162:THR:HG22	2.17	0.45
1:B:246:ARG:NH1	1:B:246:ARG:HG2	2.32	0.45
1:B:408:MET:HE1	1:B:418:LEU:HD23	1.98	0.45
1:A:128:LEU:N	1:A:129:PRO:CD	2.80	0.44
1:A:451:LEU:HA	1:A:452:PRO:HD3	1.89	0.44
1:A:552:TRP:O	1:A:556:GLN:HG2	2.17	0.44
1:B:77:LYS:HG2	1:B:131:ILE:HG12	1.98	0.44
1:A:434:LEU:HD23	1:A:440:LEU:HD23	1.99	0.44
1:A:300:GLU:HG3	1:A:301:ARG:N	2.17	0.44
1:A:329:ASN:ND2	1:A:492:GLU:OE2	2.51	0.44
1:B:250:LEU:HD12	1:B:252:VAL:HG23	1.99	0.44
1:B:395:TYR:OH	1:B:446:HIS:HD2	2.00	0.44
1:B:554:ARG:HG2	1:B:560:VAL:HG21	1.99	0.44
1:B:611:LEU:N	1:B:611:LEU:HD12	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:GLY:HA3	1:B:521:PHE:O	2.18	0.44
1:B:551:ALA:HB2	1:B:590:ALA:HB1	1.98	0.44
1:B:143:TYR:O	1:B:144:GLY:C	2.60	0.44
1:A:404:LEU:HG	1:A:408:MET:CE	2.44	0.44
1:A:90:GLN:HB3	1:A:143:TYR:CZ	2.52	0.43
1:B:73:VAL:O	1:B:77:LYS:HG3	2.18	0.43
1:A:429:HIS:HE1	6:A:855:HOH:O	2.00	0.43
1:B:626:GLY:HA2	1:B:673:ARG:NH2	2.33	0.43
1:B:266:MET:HE3	1:B:326:TYR:CD2	2.53	0.43
1:B:180:THR:CA	1:B:640:ARG:HH21	2.17	0.43
1:A:620:TRP:CZ2	1:A:654:GLY:HA2	2.53	0.43
1:B:366:CYS:HA	1:B:367:PRO:C	2.43	0.43
1:B:620:TRP:CD1	1:B:653:LEU:HB3	2.54	0.43
1:A:353:ASN:HD22	1:A:353:ASN:C	2.23	0.43
1:B:246:ARG:HG2	1:B:246:ARG:HH11	1.83	0.43
1:A:279:PHE:CG	1:A:285:PHE:HB2	2.54	0.42
1:A:404:LEU:HD13	1:B:274:ASN:HD22	1.84	0.42
1:B:584:GLN:NE2	1:B:588:ASP:OD1	2.52	0.42
1:B:364:PHE:HB2	1:B:365:PRO:HD2	2.00	0.42
1:B:217:GLU:OE1	1:B:416:LYS:HE3	2.18	0.42
1:B:72:PHE:O	1:B:76:MET:HG3	2.19	0.42
1:B:166:LEU:HG	1:B:198:LEU:HB3	2.02	0.42
1:B:181:TYR:HB3	6:B:705:HOH:O	2.18	0.42
1:B:421:SER:HA	1:B:425:GLU:HB3	2.02	0.42
1:B:226:PHE:O	1:B:230:VAL:HG23	2.20	0.42
1:B:665:ILE:HG22	1:B:669:MET:HE3	2.02	0.42
1:A:329:ASN:HD22	1:A:492:GLU:CD	2.28	0.41
1:A:517:ARG:HG2	6:A:850:HOH:O	2.20	0.41
1:B:315:ILE:HG23	1:B:315:ILE:O	2.19	0.41
1:A:425:GLU:HG3	1:A:484:LYS:HE3	2.02	0.41
1:B:634:ASP:HB3	1:B:638:MET:HB3	2.02	0.41
1:A:178:ASN:HB3	1:A:181:TYR:HD2	1.85	0.41
1:B:148:PRO:HB3	1:B:187:MET:SD	2.60	0.41
1:B:194:ARG:HH11	1:B:197:GLN:HE22	1.67	0.41
1:B:478:VAL:HG21	1:B:498:ARG:HD2	2.03	0.41
1:B:398:GLU:H	1:B:439:SER:CB	2.34	0.41
1:A:142:THR:CG2	1:A:146:GLY:HA2	2.50	0.41
1:B:79:THR:OG1	1:B:81:ARG:HG2	2.21	0.41
1:A:314:LYS:HB3	1:A:314:LYS:HZ3	1.85	0.41
1:B:363:PRO:O	1:B:453:ARG:HD3	2.21	0.41
1:B:671:LYS:HE2	1:B:673:ARG:HD3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:ARG:HH11	1:A:111:ARG:HG3	1.86	0.40
1:B:501:GLU:HA	1:B:502:PRO:HD3	1.95	0.40
1:B:594:LEU:HD22	1:B:596:VAL:HG23	2.02	0.40
1:B:667:LYS:HG3	6:B:818:HOH:O	2.20	0.40
1:B:408:MET:CE	1:B:418:LEU:HD23	2.50	0.40
1:B:72:PHE:CE1	1:B:124:ASP:HB2	2.57	0.40
1:A:478:VAL:HA	1:A:494:THR:HB	2.03	0.40
1:B:248:TYR:CE2	1:B:363:PRO:HD3	2.56	0.40
1:B:416:LYS:O	1:B:417:GLU:C	2.65	0.40
1:B:416:LYS:HG3	1:B:417:GLU:N	2.35	0.40
1:B:447:LEU:O	1:B:451:LEU:HG	2.22	0.40
1:B:658:HIS:O	1:B:662:VAL:HG23	2.21	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	603/618 (98%)	584 (97%)	16 (3%)	3 (0%)	24	14
1	B	587/618 (95%)	548 (93%)	33 (6%)	6 (1%)	12	4
All	All	1190/1236 (96%)	1132 (95%)	49 (4%)	9 (1%)	16	6

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	144	GLY
1	A	144	GLY
1	B	510	ALA
1	A	244	SER
1	A	300	GLU
1	B	214	ASN

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Mol	Chain	Res	Type
1	B	246	ARG
1	B	521	PHE
1	B	213	GLY

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	514/526 (98%)	510 (99%)	4 (1%)	73	70
1	B	507/526 (96%)	497 (98%)	10 (2%)	48	38
All	All	1021/1052 (97%)	1007 (99%)	14 (1%)	59	52

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	ARG
1	A	387	ASN
1	A	580	GLU
1	A	591	LEU
1	B	161	GLU
1	B	185	ASN
1	B	193	LYS
1	B	250	LEU
1	B	268	ARG
1	B	454	LEU
1	B	564	LEU
1	B	591	LEU
1	B	594	LEU
1	B	609	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	183	HIS

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Mol	Chain	Res	Type
1	A	185	ASN
1	A	233	HIS
1	A	253	HIS
1	A	274	ASN
1	A	283	ASN
1	A	305	HIS
1	A	329	ASN
1	A	353	ASN
1	A	362	HIS
1	A	387	ASN
1	A	402	GLN
1	A	429	HIS
1	A	435	GLN
1	A	446	HIS
1	A	455	GLN
1	A	489	ASN
1	A	555	GLN
1	A	556	GLN
1	A	584	GLN
1	A	637	ASN
1	A	643	GLN
1	B	90	GLN
1	B	133	ASN
1	B	160	GLN
1	B	183	HIS
1	B	185	ASN
1	B	197	GLN
1	B	247	GLN
1	B	274	ASN
1	B	329	ASN
1	B	336	GLN
1	B	353	ASN
1	B	387	ASN
1	B	394	GLN
1	B	435	GLN
1	B	446	HIS
1	B	455	GLN
1	B	470	ASN
1	B	555	GLN
1	B	556	GLN
1	B	595	ASN
1	B	609	GLN

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Mol	Chain	Res	Type
1	B	610	HIS
1	B	628	HIS
1	B	660	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
4	NAP	B	753	-	50,52,52	2.63	20 (40%)	71,80,80	1.85	16 (22%)
2	FAD	A	752	-	58,58,58	2.63	20 (34%)	85,89,89	1.50	15 (17%)
2	FAD	B	752	-	58,58,58	2.99	21 (36%)	85,89,89	1.51	14 (16%)
4	NAP	A	753	-	32,33,52	2.62	9 (28%)	50,52,80	1.97	12 (24%)
3	FMN	A	751	-	33,33,33	2.43	11 (33%)	48,50,50	1.47	5 (10%)
3	FMN	B	751	-	33,33,33	2.34	12 (36%)	48,50,50	1.52	10 (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
4	NAP	B	753	-	-	12/35/67/67	0/5/5/5
2	FAD	A	752	-	-	2/34/50/50	0/6/6/6
2	FAD	B	752	-	-	4/34/50/50	0/6/6/6
4	NAP	A	753	-	-	5/21/37/67	0/3/3/5
3	FMN	A	751	-	-	0/18/18/18	0/3/3/3
3	FMN	B	751	-	-	0/18/18/18	0/3/3/3

All (93) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	752	FAD	PA-O3P	-9.97	1.48	1.59
4	B	753	NAP	PA-O3	-9.02	1.49	1.59
4	B	753	NAP	C4A-N3A	8.65	1.50	1.34
4	A	753	NAP	C4A-N3A	8.18	1.49	1.34
4	A	753	NAP	PA-O3	-8.15	1.50	1.59
2	A	752	FAD	PA-O3P	-7.84	1.51	1.59
2	A	752	FAD	C1'-C2'	7.22	1.62	1.52
2	B	752	FAD	C1'-C2'	6.51	1.61	1.52
2	B	752	FAD	C4A-N3A	6.26	1.46	1.34
2	B	752	FAD	C5'-C4'	-6.04	1.43	1.51
2	A	752	FAD	C4A-N3A	5.83	1.45	1.34
3	A	751	FMN	C10-N10	5.66	1.49	1.37
3	B	751	FMN	C10-N10	5.53	1.49	1.37
2	B	752	FAD	C9A-C5X	5.40	1.49	1.41
2	A	752	FAD	C9A-C5X	5.29	1.49	1.41
2	B	752	FAD	C9A-N10	5.25	1.50	1.41
3	A	751	FMN	C4A-N5	4.90	1.41	1.30
3	A	751	FMN	C9A-N10	4.89	1.49	1.41
4	B	753	NAP	C5A-C4A	4.85	1.47	1.39
3	A	751	FMN	C9-C9A	4.84	1.47	1.39
3	B	751	FMN	C4A-N5	4.81	1.41	1.30
2	B	752	FAD	C6-C5X	4.76	1.47	1.40
2	B	752	FAD	C4X-N5	4.71	1.40	1.30
3	B	751	FMN	C9A-N10	4.69	1.49	1.41
2	A	752	FAD	C2'-C3'	4.52	1.61	1.53
2	B	752	FAD	C2'-C3'	4.51	1.61	1.53
2	A	752	FAD	C5'-C4'	-4.46	1.45	1.51
3	B	751	FMN	C9-C9A	4.44	1.46	1.39
2	A	752	FAD	C9A-N10	4.42	1.48	1.41
4	A	753	NAP	C5A-C4A	4.36	1.46	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	751	FMN	C9A-C5A	4.25	1.48	1.41
2	B	752	FAD	C8A-N7A	4.19	1.39	1.31
3	A	751	FMN	C9A-C5A	4.06	1.47	1.41
4	B	753	NAP	C6N-N1N	4.01	1.44	1.35
2	A	752	FAD	C4X-N5	3.99	1.39	1.30
3	A	751	FMN	C6-C5A	3.98	1.46	1.40
2	A	752	FAD	C6-C5X	3.93	1.46	1.40
2	B	752	FAD	C5A-C4A	3.91	1.46	1.39
2	B	752	FAD	P-O3P	-3.86	1.55	1.59
2	B	752	FAD	C10-N1	3.83	1.41	1.33
4	B	753	NAP	C4N-C3N	3.80	1.45	1.39
4	A	753	NAP	P2B-O1X	3.68	1.61	1.50
2	B	752	FAD	C8-C7	3.65	1.49	1.40
3	B	751	FMN	C8-C7	3.64	1.49	1.40
2	A	752	FAD	C8A-N7A	3.63	1.38	1.31
4	B	753	NAP	O4D-C1D	3.60	1.45	1.40
2	A	752	FAD	C5A-C4A	3.58	1.45	1.39
4	B	753	NAP	C2N-N1N	3.57	1.38	1.35
2	A	752	FAD	C8-C7	3.51	1.49	1.40
2	B	752	FAD	C6-C7	3.45	1.44	1.39
3	A	751	FMN	C8-C7	3.44	1.49	1.40
2	A	752	FAD	C10-N1	3.42	1.40	1.33
2	B	752	FAD	C9-C9A	3.40	1.45	1.39
3	A	751	FMN	C1'-C2'	3.37	1.57	1.52
2	A	752	FAD	O2'-C2'	3.31	1.50	1.43
4	B	753	NAP	P2B-O1X	3.19	1.60	1.50
3	B	751	FMN	C6-C5A	3.19	1.44	1.40
4	B	753	NAP	C8A-N7A	3.09	1.37	1.31
2	A	752	FAD	C6-C7	3.03	1.43	1.39
4	A	753	NAP	C8A-N7A	2.93	1.37	1.31
2	B	752	FAD	O2'-C2'	2.93	1.49	1.43
4	B	753	NAP	O2D-C2D	2.91	1.50	1.43
2	B	752	FAD	PA-O5B	-2.86	1.48	1.59
4	A	753	NAP	C1B-N9A	-2.85	1.38	1.46
2	B	752	FAD	C4-N3	2.84	1.44	1.38
4	B	753	NAP	PN-O3	-2.79	1.56	1.59
2	A	752	FAD	C9-C9A	2.77	1.44	1.39
4	B	753	NAP	PA-O5B	-2.74	1.48	1.59
4	B	753	NAP	P2B-O2B	2.71	1.64	1.59
4	A	753	NAP	PA-O5B	-2.71	1.48	1.59
3	B	751	FMN	C1'-C2'	2.69	1.56	1.52
2	A	752	FAD	C4-N3	2.61	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	753	NAP	C1B-N9A	-2.53	1.39	1.46
4	B	753	NAP	C8A-N9A	2.51	1.41	1.37
2	B	752	FAD	P-O5'	-2.51	1.49	1.59
4	A	753	NAP	C8A-N9A	2.48	1.41	1.37
2	A	752	FAD	PA-O5B	-2.44	1.49	1.59
4	B	753	NAP	C5D-C4D	-2.40	1.44	1.51
4	B	753	NAP	C2A-N1A	2.39	1.38	1.33
4	B	753	NAP	PN-O5D	-2.38	1.50	1.59
2	A	752	FAD	C2A-N3A	2.38	1.38	1.33
3	B	751	FMN	C10-N1	2.34	1.37	1.33
2	B	752	FAD	C2A-N3A	2.32	1.38	1.33
3	B	751	FMN	O4'-C4'	2.32	1.48	1.43
4	B	753	NAP	C2D-C3D	2.28	1.59	1.53
3	A	751	FMN	C8M-C8	2.24	1.55	1.51
2	A	752	FAD	P-O5'	-2.18	1.50	1.59
3	A	751	FMN	C4'-C3'	2.14	1.57	1.53
3	B	751	FMN	C8M-C8	2.12	1.55	1.51
4	A	753	NAP	C2A-N1A	2.12	1.37	1.33
3	A	751	FMN	C10-N1	2.10	1.37	1.33
4	B	753	NAP	C5A-C6A	2.09	1.46	1.41
3	B	751	FMN	C4'-C3'	2.06	1.57	1.53

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	753	NAP	N9A-C8A-N7A	5.98	122.42	113.94
4	B	753	NAP	C4A-N9A-C8A	-5.74	99.71	105.74
4	A	753	NAP	N9A-C8A-N7A	5.71	122.03	113.94
2	B	752	FAD	C2A-N1A-C6A	5.59	127.92	118.73
4	A	753	NAP	C4A-N9A-C8A	-5.48	99.98	105.74
2	A	752	FAD	C2A-N1A-C6A	5.44	127.67	118.73
4	B	753	NAP	C5A-C4A-N3A	-4.90	119.97	126.72
4	B	753	NAP	C2D-C3D-C4D	4.71	111.71	102.61
4	A	753	NAP	C5A-C4A-N3A	-4.63	120.34	126.72
2	A	752	FAD	N3A-C2A-N1A	-4.36	121.98	128.58
2	B	752	FAD	N3A-C2A-N1A	-4.36	121.99	128.58
4	A	753	NAP	C5A-N7A-C8A	-4.18	96.89	103.45
4	B	753	NAP	C5A-N7A-C8A	-4.15	96.93	103.45
3	B	751	FMN	C9A-C5A-N5	3.99	126.69	122.45
3	A	751	FMN	C9A-C5A-N5	3.99	126.69	122.45
4	A	753	NAP	C1B-N9A-C8A	3.82	135.56	127.09
4	B	753	NAP	C1B-N9A-C8A	3.53	134.93	127.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	752	FAD	O2-C2-N3	3.16	124.64	118.58
3	B	751	FMN	O4-C4-N3	-3.12	114.25	120.11
4	A	753	NAP	N3A-C2A-N1A	-3.11	123.88	128.58
4	B	753	NAP	N3A-C2A-N1A	-3.07	123.94	128.58
2	A	752	FAD	O2P-P-O3P	3.04	115.50	107.27
2	B	752	FAD	C1B-N9A-C8A	2.99	133.74	127.09
2	A	752	FAD	C2'-C1'-N10	-2.99	96.07	110.20
3	A	751	FMN	O4-C4-N3	-2.96	114.55	120.11
4	B	753	NAP	N3A-C4A-N9A	2.85	132.01	127.17
2	B	752	FAD	O2-C2-N3	2.84	124.04	118.58
3	B	751	FMN	C9-C9A-N10	2.79	125.61	121.85
2	B	752	FAD	O2P-P-O3P	2.78	114.79	107.27
3	A	751	FMN	C9-C9A-N10	2.75	125.55	121.85
2	B	752	FAD	C2'-C1'-N10	-2.74	97.26	110.20
2	A	752	FAD	C1B-N9A-C8A	2.73	133.15	127.09
3	B	751	FMN	C4'-C3'-C2'	-2.72	109.04	113.57
3	B	751	FMN	C5A-C9A-N10	-2.70	115.53	117.97
4	A	753	NAP	O2N-PN-O3	2.68	113.63	104.64
3	A	751	FMN	C5A-C9A-N10	-2.66	115.56	117.97
4	A	753	NAP	O2A-PA-O3	2.55	114.17	107.27
4	A	753	NAP	N3A-C4A-N9A	2.55	131.50	127.17
4	B	753	NAP	C2A-N1A-C6A	2.53	122.89	118.73
2	A	752	FAD	C8M-C8-C9	-2.49	115.18	119.57
4	A	753	NAP	C2A-N1A-C6A	2.48	122.80	118.73
2	B	752	FAD	C10-N1-C2	2.48	122.21	116.85
2	B	752	FAD	C5X-C9A-N10	-2.46	115.75	117.97
2	B	752	FAD	C9-C9A-N10	2.44	125.14	121.85
2	A	752	FAD	C9A-C5X-N5	2.43	125.04	122.45
2	A	752	FAD	C5X-C9A-N10	-2.42	115.78	117.97
4	B	753	NAP	C4D-O4D-C1D	2.40	112.12	109.92
2	A	752	FAD	O5'-C5'-C4'	2.38	115.72	109.36
2	A	752	FAD	C10-N1-C2	2.38	122.00	116.85
2	B	752	FAD	C8M-C8-C9	-2.37	115.40	119.57
4	B	753	NAP	O3D-C3D-C4D	-2.37	104.28	111.08
2	B	752	FAD	O5'-C5'-C4'	2.34	115.62	109.36
2	B	752	FAD	C4A-N9A-C1B	-2.33	121.19	126.63
4	B	753	NAP	O2A-PA-O3	2.29	113.45	107.27
4	B	753	NAP	C6N-N1N-C2N	-2.29	119.93	121.88
2	B	752	FAD	O2A-PA-O3P	2.28	113.45	107.27
2	B	752	FAD	C9A-C5X-N5	2.26	124.86	122.45
2	A	752	FAD	C9-C9A-N10	2.22	124.84	121.85
2	A	752	FAD	C4A-N9A-C1B	-2.19	121.51	126.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	751	FMN	O3P-P-O5'	-2.17	101.01	106.67
4	A	753	NAP	C5A-C4A-N9A	2.16	108.16	105.81
3	B	751	FMN	O4'-C4'-C3'	-2.14	104.24	109.25
3	B	751	FMN	C6-C5A-N5	-2.10	114.96	118.44
2	A	752	FAD	O2'-C2'-C3'	-2.09	104.34	109.25
3	A	751	FMN	O4'-C4'-C3'	-2.07	104.40	109.25
4	A	753	NAP	C4A-C5A-N7A	2.06	112.93	110.58
4	B	753	NAP	C4A-C5A-N7A	2.05	112.92	110.58
4	B	753	NAP	C5A-C4A-N9A	2.03	108.02	105.81
3	B	751	FMN	C7M-C7-C6	-2.02	116.01	119.57
2	A	752	FAD	O2-C2-N1	-2.01	118.46	121.80
3	B	751	FMN	C8M-C8-C7	2.01	124.86	120.76
4	B	753	NAP	O4B-C1B-C2B	-2.00	103.14	106.59

There are no chirality outliers.

All (23) torsion outliers are listed below:

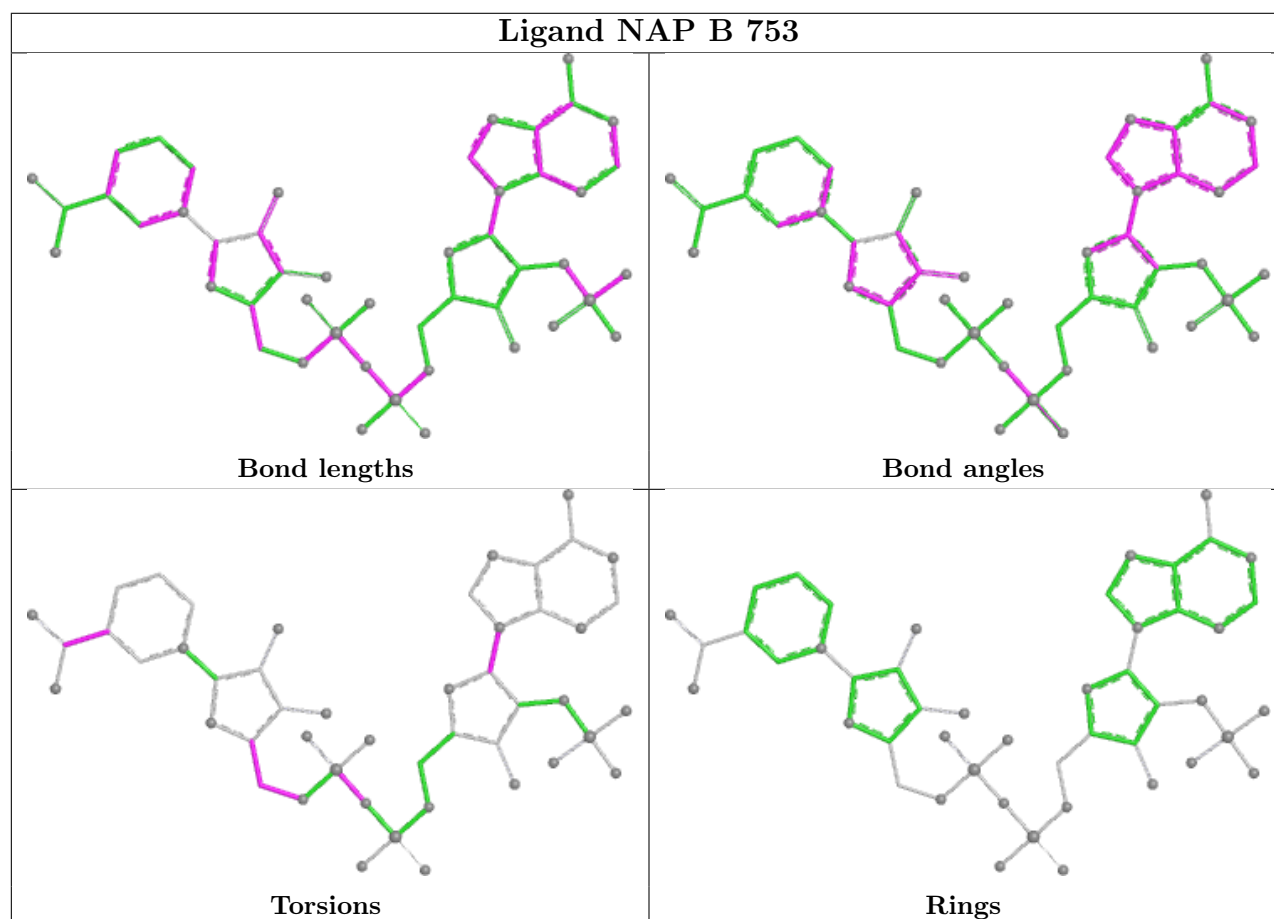
Mol	Chain	Res	Type	Atoms
4	B	753	NAP	C2N-C3N-C7N-O7N
4	B	753	NAP	C2N-C3N-C7N-N7N
4	B	753	NAP	C4N-C3N-C7N-N7N
4	B	753	NAP	C4N-C3N-C7N-O7N
4	B	753	NAP	O4D-C4D-C5D-O5D
4	B	753	NAP	C3D-C4D-C5D-O5D
4	A	753	NAP	C2B-C1B-N9A-C8A
4	B	753	NAP	C2B-C1B-N9A-C8A
2	A	752	FAD	PA-O3P-P-O2P
2	B	752	FAD	PA-O3P-P-O2P
4	B	753	NAP	PA-O3-PN-O1N
4	A	753	NAP	C5B-O5B-PA-O1A
4	A	753	NAP	C2B-C1B-N9A-C4A
4	B	753	NAP	C2B-C1B-N9A-C4A
4	B	753	NAP	C4D-C5D-O5D-PN
2	B	752	FAD	O2'-C2'-C3'-C4'
2	B	752	FAD	O2'-C2'-C3'-O3'
4	B	753	NAP	O4B-C1B-N9A-C8A
2	B	752	FAD	PA-O3P-P-O1P
4	A	753	NAP	O4B-C1B-N9A-C8A
4	A	753	NAP	C2B-O2B-P2B-O3X
2	A	752	FAD	PA-O3P-P-O1P
4	B	753	NAP	PA-O3-PN-O2N

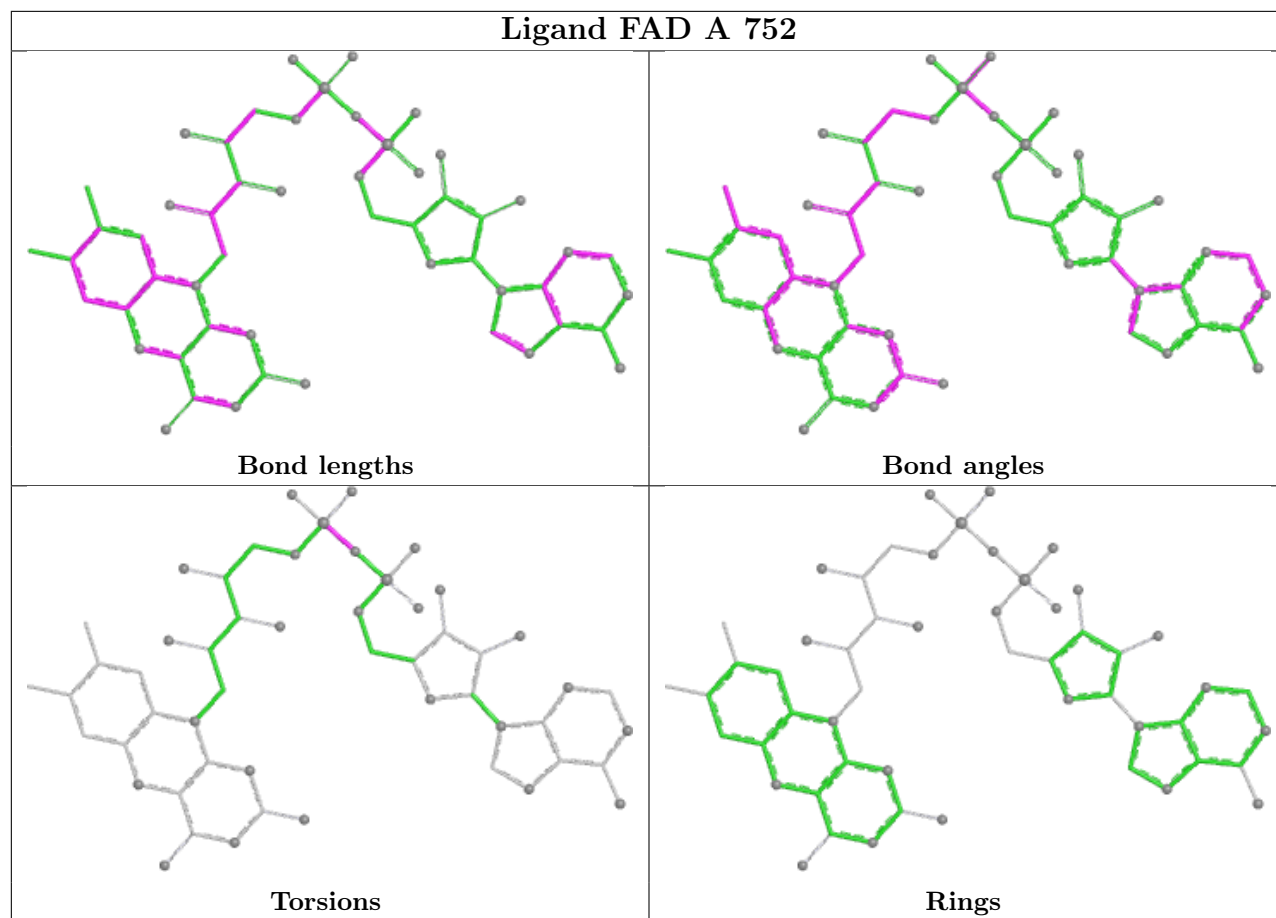
There are no ring outliers.

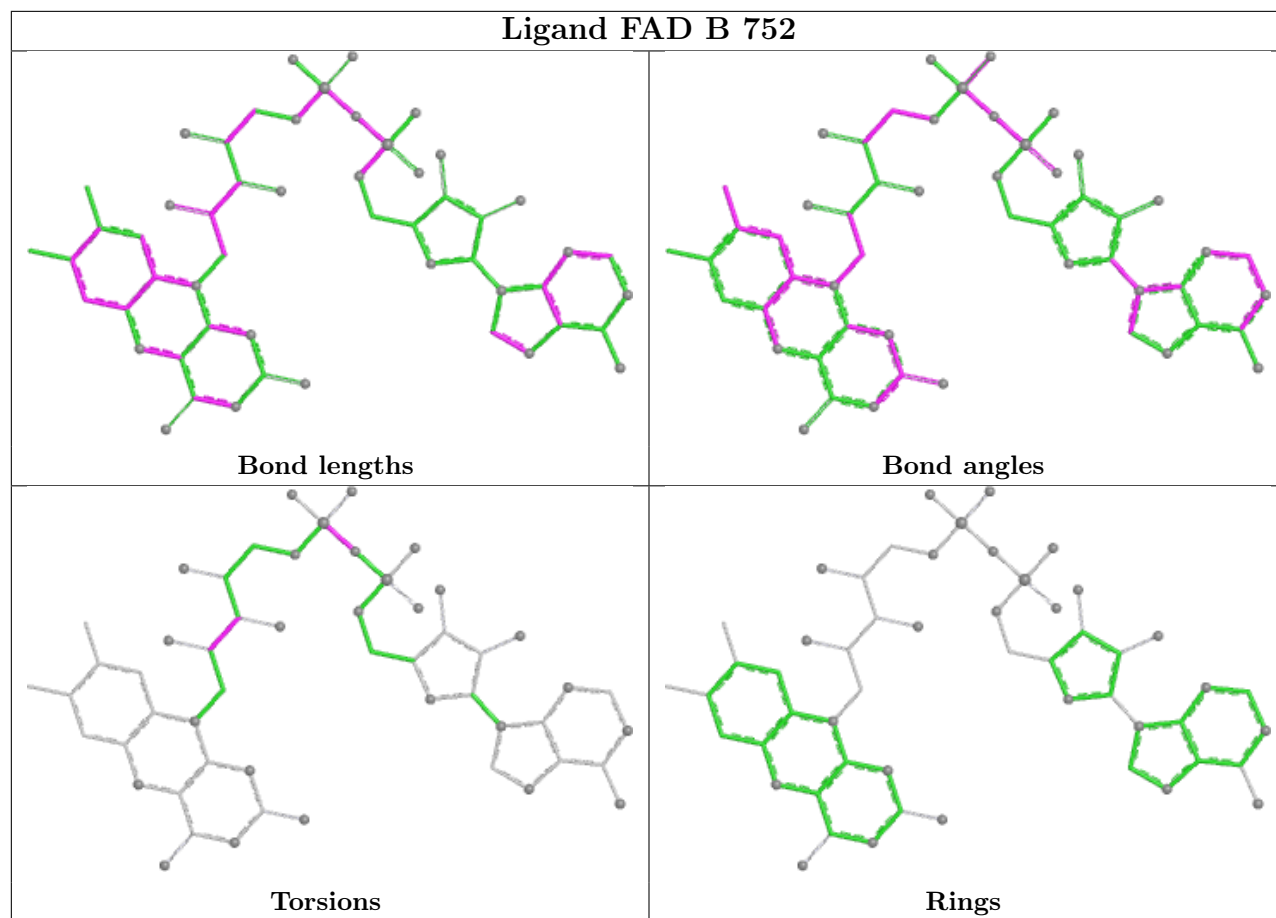
2 monomers are involved in 2 short contacts:

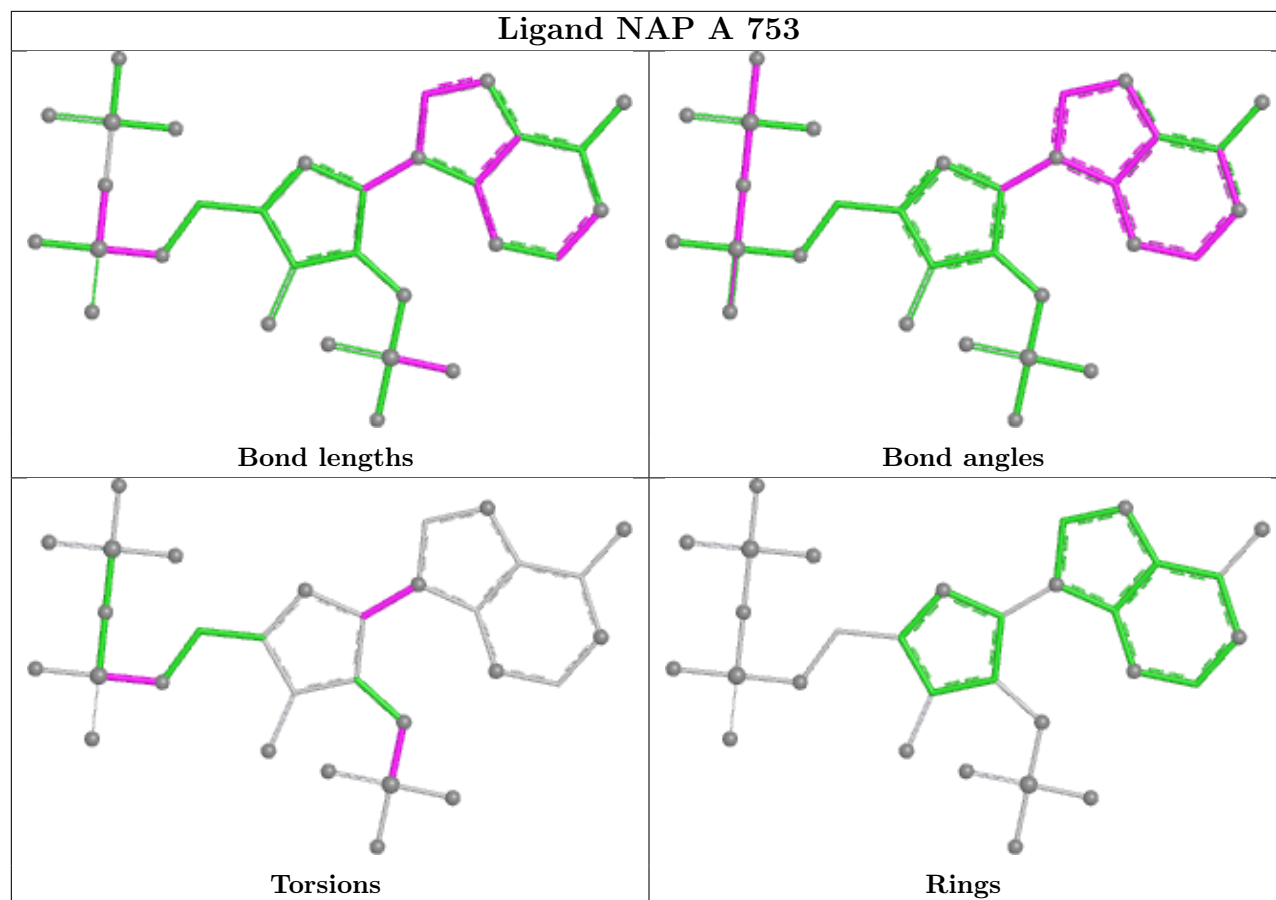
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	753	NAP	1	0
3	B	751	FMN	1	0

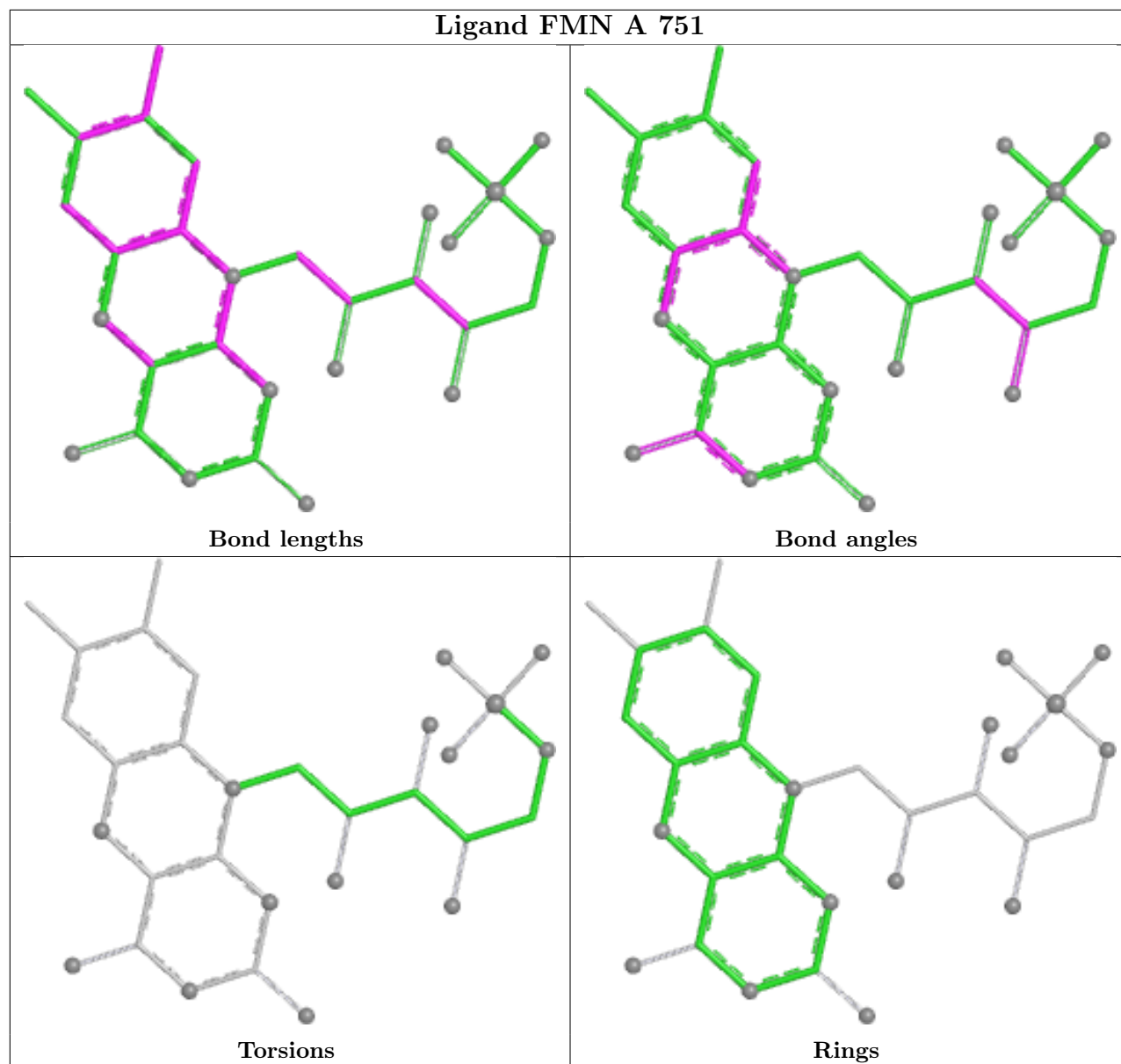
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

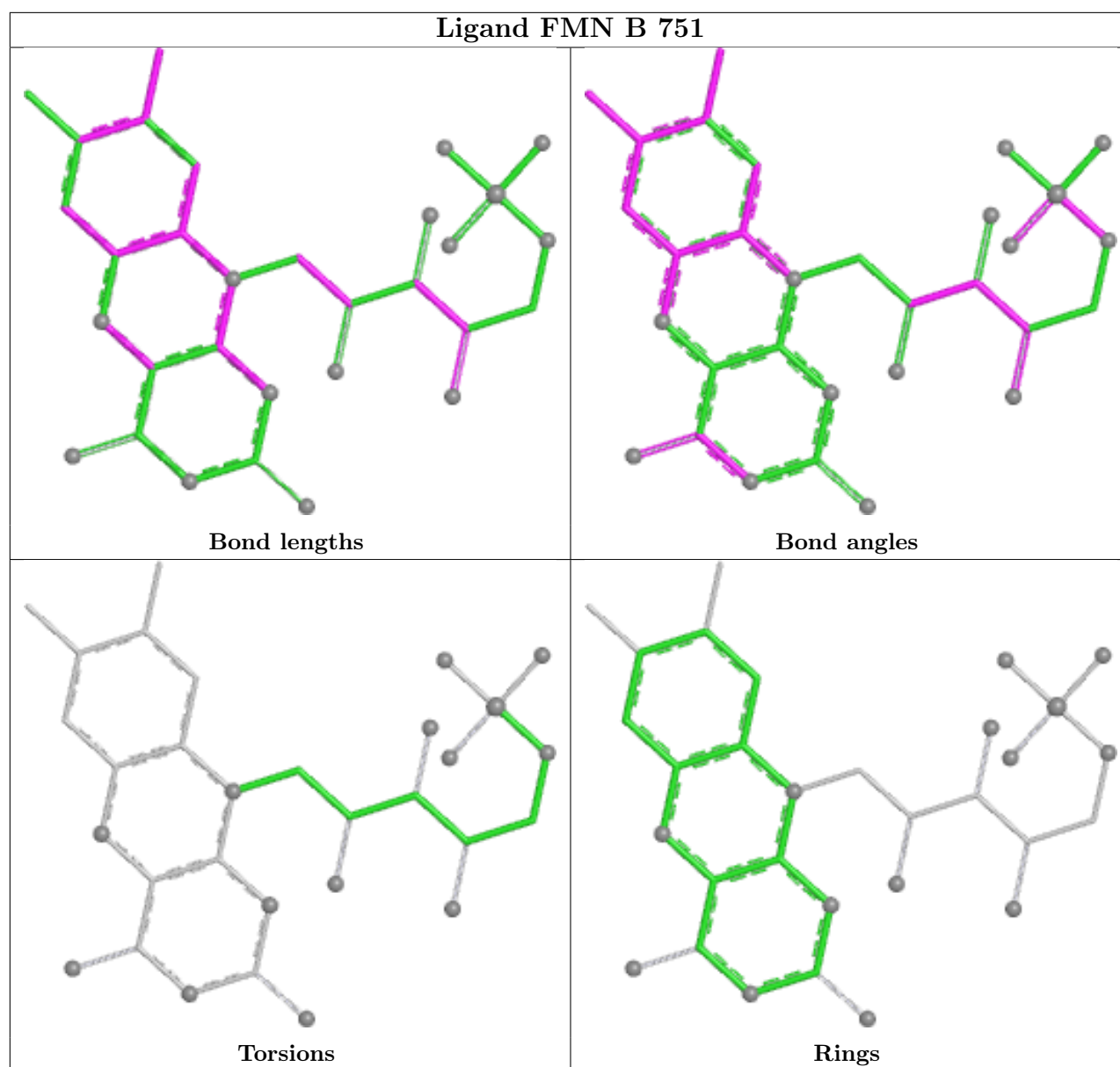












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	607/618 (98%)	0.29	31 (5%)	33 32	13, 23, 40, 57	0
1	B	595/618 (96%)	1.80	220 (36%)	1 0	22, 42, 53, 58	0
All	All	1202/1236 (97%)	1.03	251 (20%)	2 2	13, 33, 51, 58	0

All (251) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	143	TYR	6.3
1	B	503	VAL	6.1
1	B	255	ASP	5.8
1	B	163	ASP	5.5
1	B	245	ILE	5.2
1	B	299	THR	5.1
1	B	178	ASN	5.0
1	B	177	GLY	5.0
1	B	402	GLN	4.9
1	A	366	CYS	4.8
1	B	525	PHE	4.7
1	B	280	ASP	4.7
1	B	181	TYR	4.4
1	B	512	VAL	4.0
1	A	242	GLU	4.0
1	B	144	GLY	4.0
1	B	277	PRO	4.0
1	A	655	ALA	3.9
1	B	239	THR	3.8
1	B	342	GLY	3.8
1	A	503	VAL	3.8
1	B	545	GLY	3.8
1	B	213	GLY	3.7
1	B	556	GLN	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	240	GLY	3.7
1	B	491	GLY	3.7
1	A	243	SER	3.6
1	B	332	ALA	3.6
1	B	210	ASP	3.6
1	B	467	VAL	3.6
1	B	281	ALA	3.5
1	B	74	GLU	3.5
1	B	300	GLU	3.5
1	A	245	ILE	3.5
1	A	486	GLY	3.5
1	B	298	GLY	3.5
1	B	326	TYR	3.5
1	A	239	THR	3.5
1	B	291	THR	3.5
1	B	494	THR	3.5
1	B	261	VAL	3.4
1	B	254	THR	3.4
1	B	381	THR	3.4
1	B	338	GLY	3.4
1	A	299	THR	3.4
1	B	340	ILE	3.4
1	B	240	GLY	3.4
1	B	350	SER	3.4
1	B	359	ASN	3.4
1	B	399	PRO	3.4
1	B	310	ILE	3.3
1	B	432	ALA	3.3
1	B	397	SER	3.3
1	B	454	LEU	3.3
1	B	197	GLN	3.3
1	B	357	GLU	3.2
1	B	333	LEU	3.2
1	B	141	ALA	3.2
1	B	278	PRO	3.2
1	B	502	PRO	3.2
1	B	146	GLY	3.1
1	B	495	ASN	3.1
1	B	499	ALA	3.1
1	B	597	ALA	3.1
1	B	679	TRP	3.1
1	B	176	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	481	TYR	3.1
1	B	477	VAL	3.1
1	B	256	ILE	3.0
1	B	337	LEU	3.0
1	B	422	TRP	3.0
1	B	184	PHE	3.0
1	B	345	LEU	3.0
1	B	323	VAL	3.0
1	B	279	PHE	3.0
1	B	412	SER	3.0
1	B	511	LEU	2.9
1	B	70	SER	2.9
1	B	285	PHE	2.9
1	B	514	MET	2.9
1	A	297	GLN	2.9
1	B	179	LYS	2.9
1	B	588	ASP	2.9
1	B	418	LEU	2.9
1	B	247	GLN	2.9
1	B	341	LEU	2.9
1	B	348	VAL	2.8
1	A	604	HIS	2.8
1	B	107	ARG	2.8
1	B	393	ALA	2.8
1	B	411	SER	2.8
1	B	131	ILE	2.8
1	B	584	GLN	2.8
1	A	106	HIS	2.8
1	B	606	VAL	2.8
1	B	248	TYR	2.8
1	B	161	GLU	2.8
1	B	96	GLU	2.8
1	B	540	VAL	2.8
1	B	552	TRP	2.8
1	B	510	ALA	2.7
1	A	484	LYS	2.7
1	A	300	GLU	2.7
1	A	603	SER	2.7
1	B	358	SER	2.7
1	B	246	ARG	2.7
1	B	252	VAL	2.7
1	B	238	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	288	ALA	2.7
1	B	456	ALA	2.7
1	B	611	LEU	2.7
1	B	380	ILE	2.7
1	B	640	ARG	2.7
1	B	274	ASN	2.7
1	B	470	ASN	2.7
1	B	325	VAL	2.7
1	B	324	ALA	2.7
1	B	383	PRO	2.7
1	A	580	GLU	2.7
1	A	298	GLY	2.6
1	B	229	ALA	2.6
1	B	493	ALA	2.6
1	B	142	THR	2.6
1	B	180	THR	2.6
1	A	244	SER	2.6
1	B	264	GLY	2.6
1	A	485	ALA	2.6
1	B	262	TYR	2.6
1	B	390	TYR	2.6
1	B	576	TYR	2.6
1	B	365	PRO	2.6
1	B	408	MET	2.6
1	B	583	ALA	2.6
1	B	114	SER	2.6
1	B	636	ARG	2.6
1	B	73	VAL	2.6
1	B	343	ALA	2.6
1	B	479	VAL	2.6
1	B	198	LEU	2.5
1	B	437	CYS	2.5
1	B	257	ASP	2.5
1	B	272	TYR	2.5
1	B	539	GLY	2.5
1	B	635	ALA	2.5
1	B	366	CYS	2.4
1	A	441	ARG	2.4
1	B	92	GLY	2.4
1	B	406	ARG	2.4
1	B	680	SER	2.4
1	B	355	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	214	ASN	2.4
1	B	561	GLY	2.4
1	B	225	GLN	2.4
1	B	265	GLU	2.4
1	B	250	LEU	2.4
1	A	241	GLU	2.4
1	B	251	VAL	2.4
1	B	334	VAL	2.4
1	B	267	GLY	2.4
1	A	418	LEU	2.4
1	B	315	ILE	2.4
1	B	461	ILE	2.4
1	B	570	ARG	2.4
1	B	426	ALA	2.4
1	B	414	GLU	2.4
1	A	407	LYS	2.4
1	B	509	ARG	2.3
1	B	587	ARG	2.3
1	B	592	THR	2.3
1	B	152	ALA	2.3
1	B	353	ASN	2.3
1	B	415	GLY	2.3
1	B	490	LYS	2.3
1	B	598	PHE	2.3
1	B	283	ASN	2.3
1	B	382	ASN	2.3
1	B	590	ALA	2.3
1	B	603	SER	2.3
1	B	560	VAL	2.3
1	A	273	GLU	2.3
1	B	354	LEU	2.3
1	B	528	THR	2.3
1	B	220	ILE	2.3
1	B	268	ARG	2.3
1	B	398	GLU	2.3
1	B	193	LYS	2.3
1	B	670	THR	2.3
1	B	259	ALA	2.2
1	B	513	PRO	2.2
1	B	150	ASP	2.2
1	B	444	ILE	2.2
1	B	608	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	678	VAL	2.2
1	A	640	ARG	2.2
1	B	91	THR	2.2
1	B	159	LEU	2.2
1	B	80	GLY	2.2
1	B	199	GLY	2.2
1	B	270	LYS	2.2
1	B	221	THR	2.2
1	A	70	SER	2.2
1	B	599	SER	2.2
1	B	571	ARG	2.2
1	B	249	GLU	2.2
1	B	162	THR	2.2
1	B	413	GLY	2.2
1	B	148	PRO	2.2
1	B	524	PRO	2.2
1	B	655	ALA	2.2
1	B	360	LYS	2.2
1	B	666	LYS	2.2
1	B	297	GLN	2.1
1	B	423	VAL	2.1
1	B	71	SER	2.1
1	A	406	ARG	2.1
1	B	269	LEU	2.1
1	B	303	LEU	2.1
1	B	351	LEU	2.1
1	B	367	PRO	2.1
1	B	404	LEU	2.1
1	B	438	PRO	2.1
1	B	120	TYR	2.1
1	B	547	ILE	2.1
1	B	537	GLY	2.1
1	B	75	LYS	2.1
1	B	286	LEU	2.1
1	B	90	GLN	2.1
1	B	328	ALA	2.1
1	B	376	TYR	2.1
1	B	474	ILE	2.1
1	A	107	ARG	2.1
1	B	517	ARG	2.1
1	B	492	GLU	2.1
1	B	349	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	183	HIS	2.1
1	A	274	ASN	2.1
1	B	158	TRP	2.1
1	B	186	ALA	2.1
1	B	174	PHE	2.0
1	B	127	SER	2.0
1	B	433	ILE	2.0
1	B	375	THR	2.0
1	B	557	GLY	2.0
1	B	420	LEU	2.0
1	B	496	TRP	2.0
1	B	194	ARG	2.0
1	A	232	GLU	2.0
1	B	501	GLU	2.0
1	B	149	THR	2.0
1	B	529	THR	2.0
1	B	190	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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
3	FMN	B	751	31/31	0.81	0.16	36,45,47,48	0
4	NAP	B	753	48/48	0.82	0.15	32,49,64,65	0
2	FAD	B	752	53/53	0.83	0.15	33,38,57,58	0
4	NAP	A	753	31/48	0.94	0.09	21,25,40,44	0
2	FAD	A	752	53/53	0.96	0.07	12,16,28,32	0
3	FMN	A	751	31/31	0.97	0.06	17,20,22,23	0

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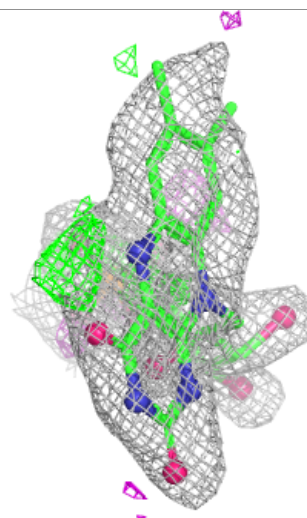
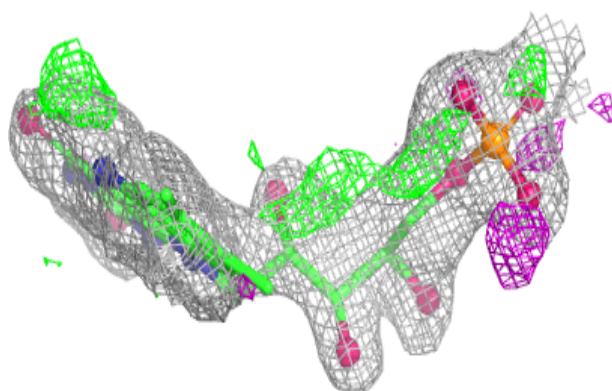
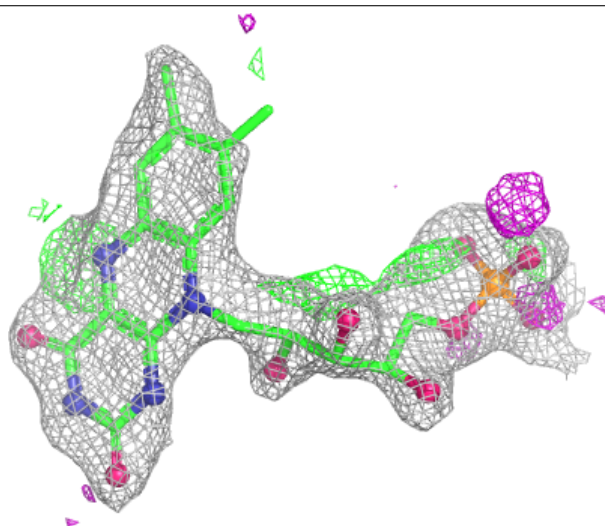
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CA	A	762	1/1	0.98	0.05	28,28,28,28	0

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

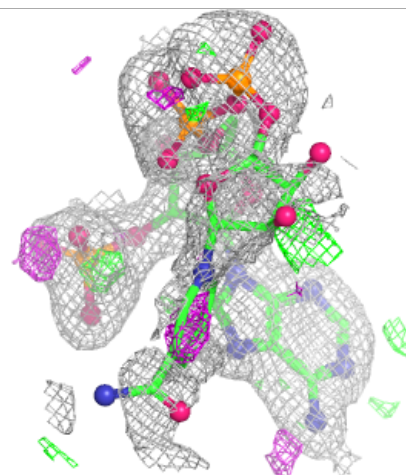
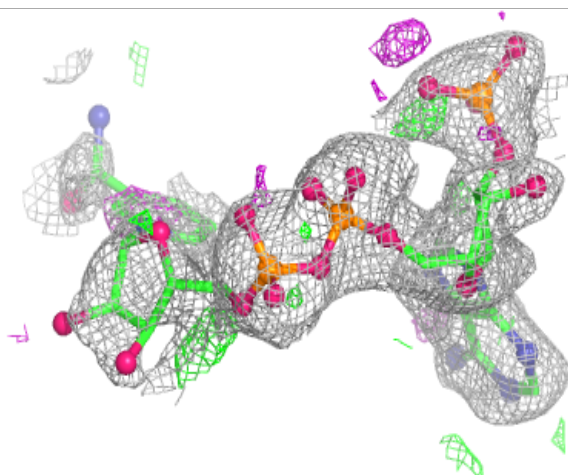
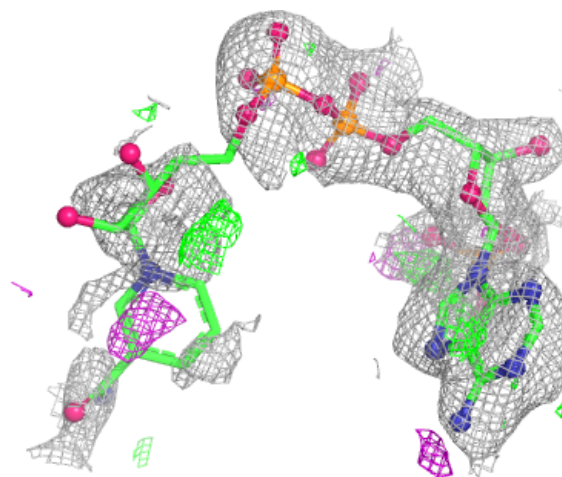
Electron density around FMN B 751:

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



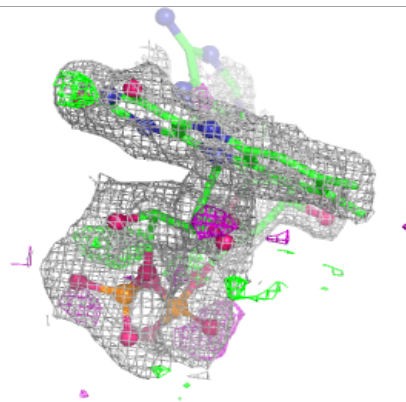
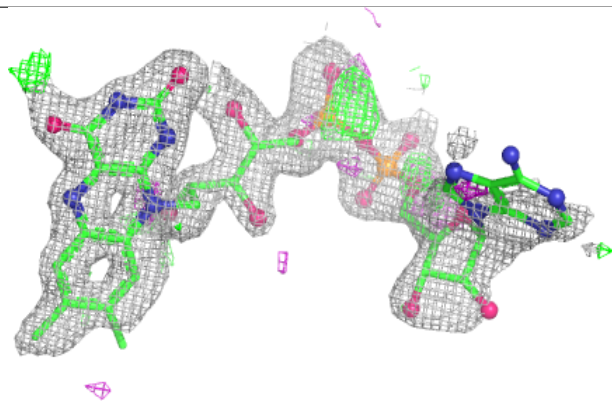
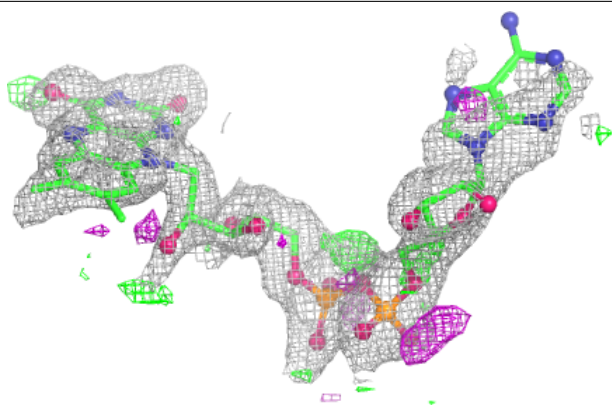
Electron density around NAP B 753:

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



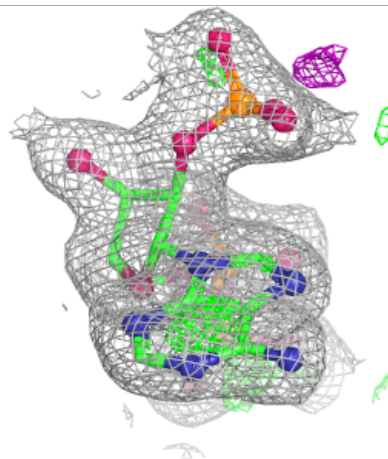
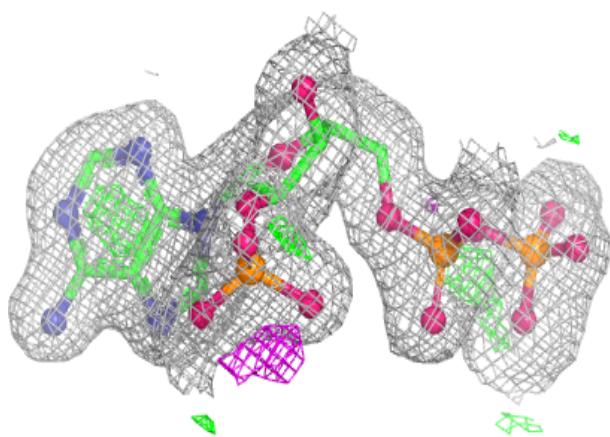
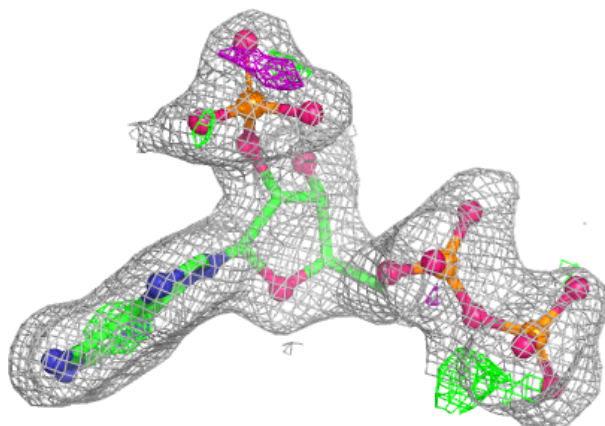
Electron density around FAD B 752:

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



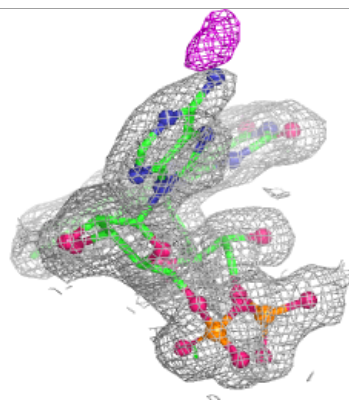
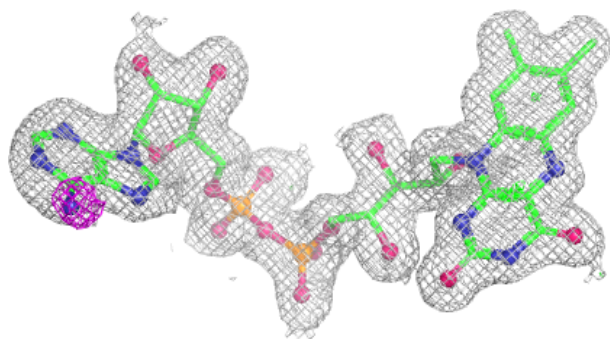
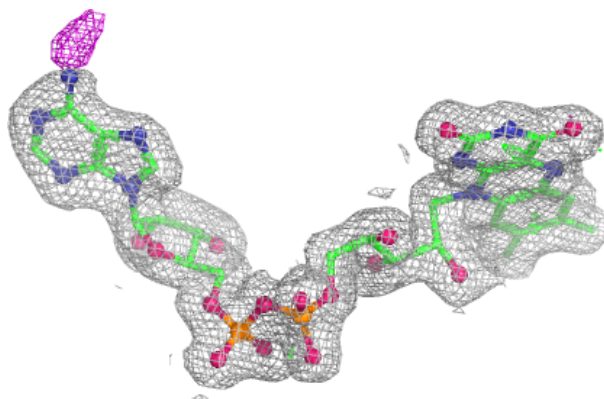
Electron density around NAP A 753:

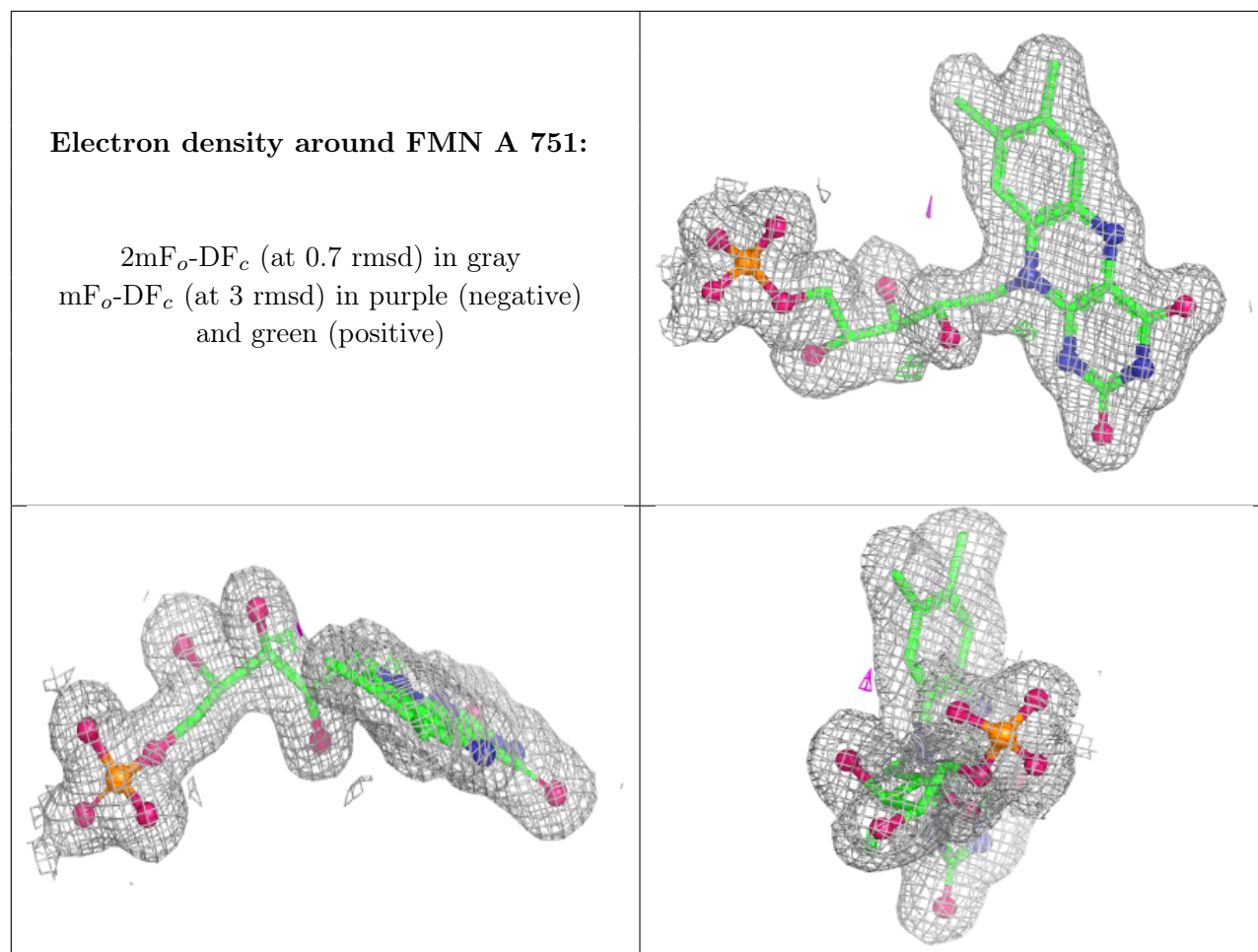
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FAD A 752:

$2mF_o - DF_c$ (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.