



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:13 AM UTC

PDB ID : 5JCL / pdb_00005jcl
Title : Structure and catalytic mechanism of monodehydroascorbate reductase, MD-HAR, from *Oryza sativa* L. japonica
Authors : Park, A.K.; Kim, H.W.
Deposited on : 2016-04-15
Resolution : 1.80 Å(reported)

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

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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

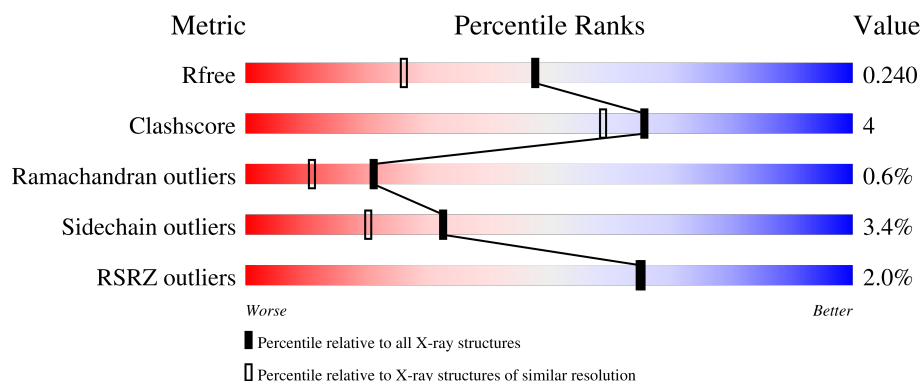
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

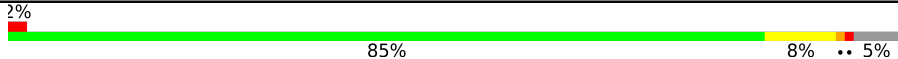
The reported resolution of this entry is 1.80 Å.

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	451	
1	B	451	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7575 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 Os09g0567300 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	429	Total	C	N	O	S	0	0	0
			3253	2097	531	619	6			
1	B	428	Total	C	N	O	S	0	0	0
			3240	2088	531	615	6			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	HIS	-	expression tag	UNP Q652L6
A	-14	HIS	-	expression tag	UNP Q652L6
A	-13	HIS	-	expression tag	UNP Q652L6
A	-12	HIS	-	expression tag	UNP Q652L6
A	-11	HIS	-	expression tag	UNP Q652L6
A	-10	HIS	-	expression tag	UNP Q652L6
A	-9	ALA	-	expression tag	UNP Q652L6
A	-8	SER	-	expression tag	UNP Q652L6
A	-7	GLU	-	expression tag	UNP Q652L6
A	-6	ASN	-	expression tag	UNP Q652L6
A	-5	LEU	-	expression tag	UNP Q652L6
A	-4	TYR	-	expression tag	UNP Q652L6
A	-3	PHE	-	expression tag	UNP Q652L6
A	-2	GLN	-	expression tag	UNP Q652L6
A	-1	GLY	-	expression tag	UNP Q652L6
A	0	ALA	-	expression tag	UNP Q652L6
A	1	MET	-	expression tag	UNP Q652L6
A	2	ALA	-	expression tag	UNP Q652L6
A	3	SER	-	expression tag	UNP Q652L6
A	196	ALA	GLU	engineered mutation	UNP Q652L6
B	-15	HIS	-	expression tag	UNP Q652L6
B	-14	HIS	-	expression tag	UNP Q652L6
B	-13	HIS	-	expression tag	UNP Q652L6
B	-12	HIS	-	expression tag	UNP Q652L6
B	-11	HIS	-	expression tag	UNP Q652L6

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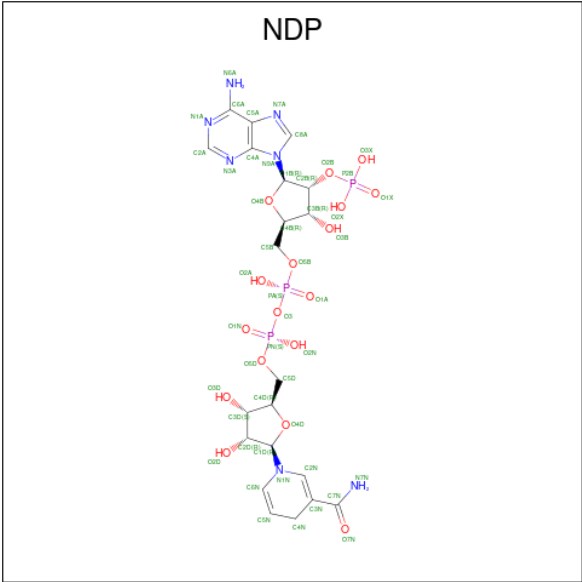
Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	HIS	-	expression tag	UNP Q652L6
B	-9	ALA	-	expression tag	UNP Q652L6
B	-8	SER	-	expression tag	UNP Q652L6
B	-7	GLU	-	expression tag	UNP Q652L6
B	-6	ASN	-	expression tag	UNP Q652L6
B	-5	LEU	-	expression tag	UNP Q652L6
B	-4	TYR	-	expression tag	UNP Q652L6
B	-3	PHE	-	expression tag	UNP Q652L6
B	-2	GLN	-	expression tag	UNP Q652L6
B	-1	GLY	-	expression tag	UNP Q652L6
B	0	ALA	-	expression tag	UNP Q652L6
B	1	MET	-	expression tag	UNP Q652L6
B	2	ALA	-	expression tag	UNP Q652L6
B	3	SER	-	expression tag	UNP Q652L6
B	196	ALA	GLU	engineered mutation	UNP Q652L6

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



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 NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

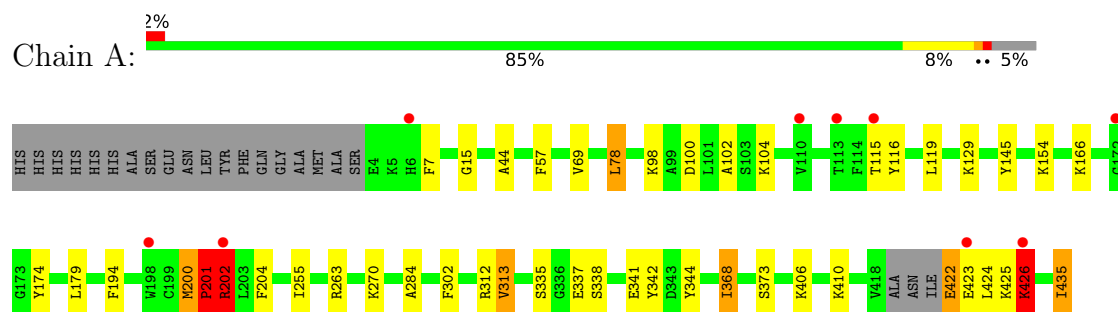
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	435	Total	O	0	0
			435	435		
4	B	445	Total	O	0	0
			445	445		

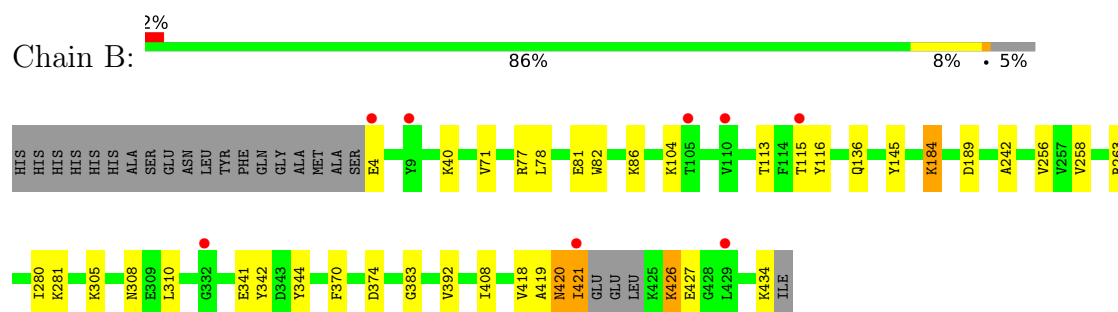
3 Residue-property plots [i](#)

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: Os09g0567300 protein



• Molecule 1: Os09g0567300 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.25Å 85.52Å 131.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80 50.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.5 (50.00-1.80) 99.5 (50.00-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.42 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.182 , 0.234 0.192 , 0.240	Depositor DCC
R_{free} test set	4126 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7575	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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, NDP

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	1.24	4/3322 (0.1%)	1.20	11/4494 (0.2%)
1	B	1.18	2/3309 (0.1%)	1.10	5/4477 (0.1%)
All	All	1.21	6/6631 (0.1%)	1.15	16/8971 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	200	MET	C-O	7.87	1.27	1.23
1	A	179	LEU	C-O	-5.45	1.17	1.24
1	A	302	PHE	C-N	5.43	1.38	1.33
1	B	242	ALA	CA-C	-5.35	1.45	1.52
1	A	119	LEU	N-CA	5.17	1.52	1.46
1	B	256	VAL	CA-C	-5.13	1.46	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	PRO	N-CA-C	11.55	129.59	113.53
1	A	44	ALA	CA-C-N	7.89	129.71	119.84
1	A	44	ALA	C-N-CA	7.89	129.71	119.84
1	B	408	ILE	N-CA-C	7.83	118.61	110.62
1	A	201	PRO	CA-C-N	6.94	134.80	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	PRO	C-N-CA	6.94	134.80	121.54
1	B	419	ALA	N-CA-C	-6.43	98.92	109.40
1	A	194	PHE	N-CA-C	5.86	113.49	108.22
1	A	200	MET	CA-C-N	5.61	125.27	119.05
1	A	200	MET	C-N-CA	5.61	125.27	119.05
1	A	57	PHE	CA-C-N	-5.24	114.26	119.56
1	A	57	PHE	C-N-CA	-5.24	114.26	119.56
1	A	313	VAL	CB-CA-C	5.06	118.29	110.50
1	B	280	ILE	N-CA-C	-5.05	101.30	108.58
1	B	374	ASP	N-CA-C	-5.03	102.24	109.24
1	B	184	LYS	N-CA-C	-5.01	105.90	111.36

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	201	PRO	Peptide
1	A	422	GLU	Peptide
1	A	425	LYS	Peptide
1	A	426	LYS	Peptide

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	3253	0	3256	28	0
1	B	3240	0	3244	19	0
2	A	53	0	31	1	0
2	B	53	0	31	0	0
3	A	48	0	26	1	0
3	B	48	0	26	3	0
4	A	435	0	0	8	0
4	B	445	0	0	6	0
All	All	7575	0	6614	50	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (50) 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:136:GLN:HG2	4:B:866:HOH:O	1.55	1.03
1:B:71:VAL:CG1	1:B:77:ARG:HG2	2.12	0.79
3:B:501:NDP:O2B	4:B:601:HOH:O	2.08	0.71
1:A:100:ASP:OD1	4:A:601:HOH:O	2.14	0.64
1:B:71:VAL:HG11	1:B:77:ARG:HD3	1.79	0.63
1:A:200:MET:SD	1:A:202:ARG:NH2	2.73	0.62
3:B:501:NDP:P2B	4:B:601:HOH:O	2.59	0.59
1:A:312:ARG:O	4:A:602:HOH:O	2.16	0.59
1:A:102:ALA:N	4:A:609:HOH:O	2.37	0.57
1:A:263:ARG:NE	4:A:606:HOH:O	2.35	0.56
1:A:7:PHE:O	1:A:116:TYR:HA	2.07	0.54
1:A:200:MET:C	1:A:202:ARG:HB2	2.33	0.53
1:A:200:MET:HB2	1:A:204:PHE:CD2	2.43	0.53
1:A:154:LYS:NZ	4:A:615:HOH:O	2.42	0.52
1:B:71:VAL:HG13	1:B:77:ARG:HG2	1.92	0.51
1:B:421:ILE:O	1:B:426:LYS:HG3	2.11	0.50
1:A:201:PRO:N	1:A:202:ARG:HB2	2.26	0.50
1:A:69:VAL:HG23	1:A:78:LEU:HD22	1.93	0.50
1:A:104:LYS:O	1:A:115:THR:HG23	2.12	0.49
1:A:424:LEU:C	1:A:426:LYS:HB2	2.38	0.48
3:B:501:NDP:O1X	4:B:602:HOH:O	2.20	0.48
1:A:284:ALA:HB1	1:A:341:GLU:HB2	1.96	0.47
1:B:71:VAL:HG12	1:B:77:ARG:HG2	1.96	0.47
1:B:420:ASN:O	1:B:421:ILE:HG23	2.15	0.47
1:B:145:TYR:CE1	1:B:258:VAL:HB	2.49	0.47
1:B:421:ILE:O	1:B:421:ILE:HD12	2.15	0.47
1:B:392:VAL:HG23	1:B:418:VAL:CG2	2.44	0.46
1:B:104:LYS:HA	1:B:116:TYR:CZ	2.51	0.46
1:B:370:PHE:CZ	1:B:383:GLY:HA3	2.51	0.45
1:A:422:GLU:N	4:A:628:HOH:O	2.50	0.45
1:B:281:LYS:HE3	1:B:310:LEU:HD11	1.99	0.45
1:B:4:GLU:CD	1:B:113:THR:OG1	2.60	0.44
4:A:654:HOH:O	1:B:263:ARG:HG3	2.17	0.44
1:B:71:VAL:HG11	1:B:77:ARG:CD	2.48	0.43
1:A:200:MET:N	1:A:201:PRO:CD	2.81	0.43
1:A:342:TYR:CE2	1:A:344:TYR:HB2	2.55	0.42
1:B:82:TRP:CZ2	1:B:86:LYS:HD2	2.54	0.42
1:A:368:ILE:HD13	1:A:368:ILE:HA	1.84	0.42
1:A:174:TYR:CZ	1:A:200:MET:HE1	2.55	0.42
1:A:166:LYS:NZ	4:A:632:HOH:O	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:435:ILE:HD13	1:A:435:ILE:HA	1.85	0.42
1:A:410:LYS:HD3	1:A:435:ILE:HG22	2.02	0.42
1:A:255:ILE:HG21	1:A:255:ILE:HD13	1.85	0.41
1:A:174:TYR:HB3	3:A:501:NDP:H42N	2.02	0.41
1:A:129:LYS:HA	1:A:145:TYR:CE1	2.55	0.41
1:A:15:GLY:HA3	2:A:500:FAD:O5B	2.20	0.41
1:A:406:LYS:HE3	4:B:992:HOH:O	2.21	0.41
1:B:4:GLU:N	4:B:625:HOH:O	2.53	0.41
1:B:342:TYR:CE2	1:B:344:TYR:HB2	2.56	0.41
1:A:200:MET:HE3	1:A:200:MET:HB3	1.77	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	425/451 (94%)	404 (95%)	18 (4%)	3 (1%)	18	8
1	B	424/451 (94%)	406 (96%)	16 (4%)	2 (0%)	24	14
All	All	849/902 (94%)	810 (95%)	34 (4%)	5 (1%)	21	11

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	420	ASN
1	A	202	ARG
1	B	427	GLU
1	A	426	LYS
1	A	337	GLU

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	335/352 (95%)	324 (97%)	11 (3%)	33	21
1	B	333/352 (95%)	321 (96%)	12 (4%)	31	18
All	All	668/704 (95%)	645 (97%)	23 (3%)	32	20

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	LEU
1	A	98	LYS
1	A	202	ARG
1	A	270	LYS
1	A	313	VAL
1	A	335	SER
1	A	338	SER
1	A	368	ILE
1	A	373	SER
1	A	423	GLU
1	A	435	ILE
1	B	40	LYS
1	B	78	LEU
1	B	81	GLU
1	B	115	THR
1	B	184	LYS
1	B	189	ASP
1	B	305	LYS
1	B	308	ASN
1	B	341	GLU
1	B	421	ILE
1	B	426	LYS
1	B	434	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	141	ASN
1	A	186	ASN
1	A	430	GLN
1	B	318	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NDP	A	501	-	51,52,52	1.16	4 (7%)	71,80,80	1.88	18 (25%)
3	NDP	B	501	-	51,52,52	1.43	7 (13%)	71,80,80	2.04	18 (25%)
2	FAD	A	500	-	58,58,58	2.09	15 (25%)	85,89,89	1.92	30 (35%)
2	FAD	B	502	-	58,58,58	1.65	12 (20%)	85,89,89	1.74	21 (24%)

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
3	NDP	A	501	-	-	6/34/77/77	0/5/5/5
3	NDP	B	501	-	-	8/34/77/77	0/5/5/5
2	FAD	A	500	-	-	1/34/50/50	0/6/6/6
2	FAD	B	502	-	-	2/34/50/50	0/6/6/6

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	FAD	PA-O3P	-6.35	1.52	1.59
2	B	502	FAD	C1'-C2'	6.09	1.61	1.52
2	A	500	FAD	P-O3P	5.31	1.65	1.59
2	A	500	FAD	C1'-C2'	4.67	1.59	1.52
3	B	501	NDP	C5A-C4A	4.17	1.46	1.39
2	B	502	FAD	C9A-C5X	4.00	1.47	1.41
2	A	500	FAD	C5A-C4A	3.79	1.45	1.39
2	A	500	FAD	C4-N3	-3.68	1.32	1.38
3	B	501	NDP	C4A-N9A	-3.54	1.30	1.37
2	A	500	FAD	C5A-N7A	-3.52	1.32	1.39
2	B	502	FAD	C4X-N5	3.48	1.38	1.30
3	A	501	NDP	C5A-C4A	3.12	1.44	1.39
2	A	500	FAD	C9A-C5X	3.09	1.46	1.41
2	B	502	FAD	C4-N3	-2.97	1.33	1.38
2	A	500	FAD	C8A-N7A	2.92	1.37	1.31
2	A	500	FAD	O4B-C4B	2.91	1.51	1.45
2	B	502	FAD	C4A-N9A	-2.90	1.31	1.37
2	A	500	FAD	C8A-N9A	-2.89	1.32	1.37
2	B	502	FAD	C2-N3	-2.83	1.32	1.39
2	B	502	FAD	C5A-N7A	-2.82	1.33	1.39
3	B	501	NDP	P2B-O2B	2.78	1.64	1.59
3	B	501	NDP	C2A-N3A	-2.73	1.29	1.33
2	B	502	FAD	P-O3P	2.69	1.62	1.59
3	A	501	NDP	C2N-C3N	2.63	1.42	1.35
2	A	500	FAD	C5A-C6A	2.59	1.48	1.41
2	A	500	FAD	C9-C9A	2.39	1.43	1.39
2	A	500	FAD	C10-N10	2.38	1.42	1.37
2	B	502	FAD	C8-C7	2.35	1.46	1.40
2	B	502	FAD	C5A-C4A	2.32	1.43	1.39
2	A	500	FAD	O2B-C2B	-2.26	1.37	1.43
2	B	502	FAD	C8A-N7A	2.22	1.36	1.31
3	A	501	NDP	C5A-C6A	2.18	1.47	1.41
3	B	501	NDP	C2N-C3N	2.17	1.41	1.35
3	A	501	NDP	PN-O3	2.17	1.61	1.59
3	B	501	NDP	C2A-N1A	2.16	1.37	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	502	FAD	C8A-N9A	-2.13	1.33	1.37
3	B	501	NDP	C3B-C2B	2.12	1.57	1.53
2	A	500	FAD	P-O2P	-2.08	1.45	1.55

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	NDP	C4A-N9A-C8A	6.73	112.81	105.74
3	B	501	NDP	N3A-C4A-N9A	6.22	137.75	127.17
3	A	501	NDP	N3A-C4A-N9A	5.82	137.06	127.17
3	B	501	NDP	C5A-C4A-N3A	-5.38	119.31	126.72
3	B	501	NDP	N3A-C2A-N1A	-5.26	120.62	128.58
3	A	501	NDP	C4A-N9A-C8A	5.17	111.16	105.74
2	A	500	FAD	N3A-C2A-N1A	-4.97	121.05	128.58
2	A	500	FAD	C4A-N9A-C8A	4.82	110.80	105.74
3	A	501	NDP	C5A-C4A-N3A	-4.81	120.09	126.72
3	B	501	NDP	C2A-N3A-C4A	4.46	122.71	111.83
2	B	502	FAD	C4A-N9A-C8A	4.38	110.34	105.74
2	B	502	FAD	N3A-C2A-N1A	-4.31	122.05	128.58
2	A	500	FAD	C9A-C5X-N5	-4.14	118.06	122.45
2	B	502	FAD	C9A-C5X-N5	-3.92	118.30	122.45
2	B	502	FAD	N3A-C4A-N9A	3.87	133.74	127.17
2	B	502	FAD	C4-C4X-N5	3.81	123.47	118.21
2	A	500	FAD	C2A-N1A-C6A	3.69	124.79	118.73
2	A	500	FAD	O4B-C1B-N9A	-3.55	101.27	108.09
2	B	502	FAD	C5A-C4A-N3A	-3.54	121.84	126.72
3	A	501	NDP	N3A-C2A-N1A	-3.52	123.25	128.58
3	A	501	NDP	C6N-N1N-C2N	3.47	123.03	119.32
3	B	501	NDP	N9A-C8A-N7A	-3.42	109.08	113.94
2	A	500	FAD	N3A-C4A-N9A	3.38	132.92	127.17
2	B	502	FAD	O4-C4-C4X	-3.32	117.76	126.53
3	A	501	NDP	N9A-C8A-N7A	-3.17	109.43	113.94
3	A	501	NDP	C5A-C6A-N6A	-3.08	115.66	123.29
3	A	501	NDP	C2A-N3A-C4A	3.05	119.28	111.83
3	B	501	NDP	C1D-N1N-C2N	-3.02	116.17	121.14
2	A	500	FAD	O4-C4-C4X	-2.99	118.65	126.53
2	A	500	FAD	C5X-N5-C4X	2.92	122.82	118.09
2	A	500	FAD	C4B-O4B-C1B	-2.91	103.04	109.47
2	B	502	FAD	C3B-C2B-C1B	-2.87	96.03	101.46
2	A	500	FAD	C4A-N9A-C1B	-2.86	119.94	126.63
3	B	501	NDP	C2A-N1A-C6A	2.77	123.28	118.73
3	B	501	NDP	C5A-C4A-N9A	-2.75	102.81	105.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	FAD	C4A-N9A-C1B	-2.72	120.27	126.63
2	A	500	FAD	C4-C4X-N5	2.72	121.97	118.21
3	A	501	NDP	C5A-C4A-N9A	-2.72	102.85	105.81
3	A	501	NDP	C4B-O4B-C1B	-2.67	103.56	109.47
2	A	500	FAD	O2P-P-O1P	2.67	124.84	112.44
2	A	500	FAD	N9A-C8A-N7A	-2.66	110.16	113.94
2	A	500	FAD	C3B-C2B-C1B	-2.62	96.51	101.46
2	A	500	FAD	O4B-C1B-C2B	-2.62	101.01	106.62
2	A	500	FAD	C1'-N10-C9A	2.62	125.72	120.63
3	B	501	NDP	O4D-C1D-C2D	2.62	112.21	106.62
2	A	500	FAD	C2A-N3A-C4A	2.61	118.21	111.83
2	B	502	FAD	C2A-N3A-C4A	2.60	118.18	111.83
2	A	500	FAD	C5X-C9A-N10	2.59	120.31	117.97
3	B	501	NDP	C6N-N1N-C2N	2.59	122.09	119.32
2	B	502	FAD	C2A-N1A-C6A	2.59	122.98	118.73
2	A	500	FAD	O2A-PA-O3P	2.57	114.21	107.27
2	B	502	FAD	N9A-C8A-N7A	-2.56	110.31	113.94
2	A	500	FAD	O2P-P-O3P	2.52	114.10	107.27
2	A	500	FAD	N6A-C6A-N1A	2.52	123.99	118.38
2	A	500	FAD	O2-C2-N1	-2.47	117.70	121.80
3	A	501	NDP	O3X-P2B-O2B	-2.47	96.23	105.85
3	A	501	NDP	C1D-N1N-C2N	-2.44	117.11	121.14
2	B	502	FAD	C8M-C8-C7	-2.43	115.79	120.76
2	A	500	FAD	C6-C5X-N5	2.41	122.45	118.44
3	A	501	NDP	N6A-C6A-N1A	2.41	123.74	118.38
2	A	500	FAD	C5A-C4A-N3A	-2.38	123.44	126.72
3	B	501	NDP	C3B-C2B-C1B	-2.38	98.25	102.81
2	B	502	FAD	C4X-C4-N3	2.38	119.31	113.25
2	A	500	FAD	O2A-PA-O5B	-2.36	96.87	107.57
2	A	500	FAD	O4-C4-N3	2.36	124.54	120.11
3	A	501	NDP	C4A-N9A-C1B	-2.35	121.13	126.63
3	A	501	NDP	C3B-C2B-C1B	-2.33	98.35	102.81
2	A	500	FAD	O3B-C3B-C2B	2.30	119.19	111.82
3	B	501	NDP	C5A-N7A-C8A	2.30	107.06	103.45
2	B	502	FAD	N6A-C6A-N1A	2.27	123.44	118.38
2	B	502	FAD	C5A-C6A-N6A	-2.25	117.72	123.29
2	A	500	FAD	C5A-C4A-N9A	-2.24	103.37	105.81
3	B	501	NDP	O4B-C1B-C2B	-2.23	102.74	106.59
2	B	502	FAD	C5X-C9A-N10	2.14	119.90	117.97
3	B	501	NDP	O2N-PN-O3	2.13	113.02	107.27
2	B	502	FAD	C1'-N10-C9A	2.12	124.75	120.63
3	A	501	NDP	C5A-N7A-C8A	2.12	106.78	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	NDP	C4A-C5A-N7A	-2.09	108.19	110.58
2	B	502	FAD	O5B-PA-O1A	-2.09	100.64	108.94
2	A	500	FAD	C5A-C6A-N6A	-2.07	118.15	123.29
3	A	501	NDP	C3N-C2N-N1N	-2.07	120.17	123.20
2	A	500	FAD	C5'-C4'-C3'	2.07	116.11	112.22
3	B	501	NDP	C4D-O4D-C1D	-2.06	104.92	109.47
2	B	502	FAD	O2'-C2'-C1'	2.06	118.66	110.20
3	B	501	NDP	C6A-C5A-N7A	2.06	136.05	132.09
3	A	501	NDP	O3-PN-O1N	-2.05	104.54	110.70
2	B	502	FAD	O3P-PA-O1A	2.03	116.80	110.70

There are no chirality outliers.

All (17) torsion outliers are listed below:

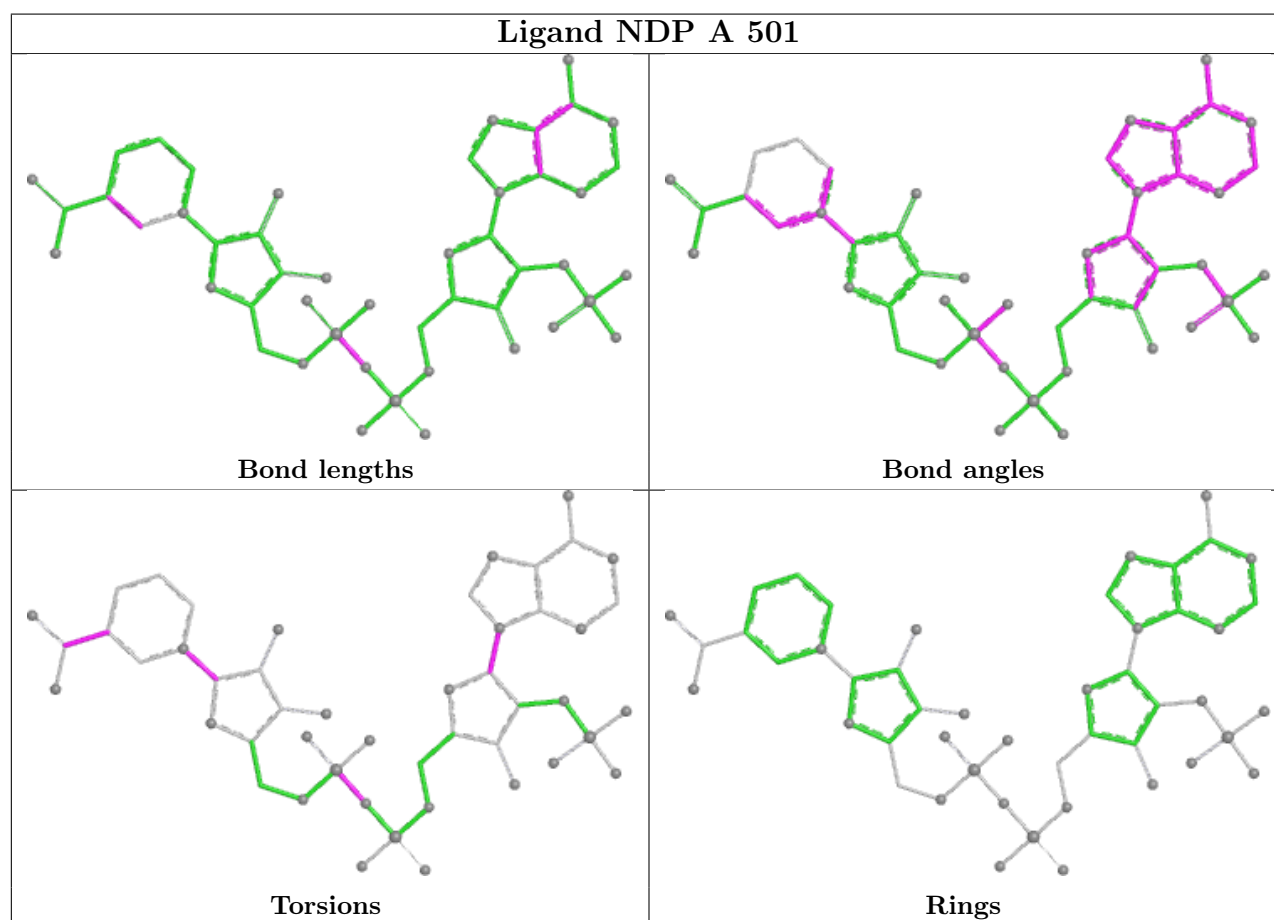
Mol	Chain	Res	Type	Atoms
2	B	502	FAD	N10-C1'-C2'-O2'
3	A	501	NDP	C2B-C1B-N9A-C8A
3	B	501	NDP	C2B-C1B-N9A-C8A
3	B	501	NDP	C2B-C1B-N9A-C4A
3	A	501	NDP	O4D-C1D-N1N-C2N
3	B	501	NDP	O4D-C1D-N1N-C2N
3	A	501	NDP	C2B-C1B-N9A-C4A
3	B	501	NDP	C3B-C2B-O2B-P2B
3	B	501	NDP	C1B-C2B-O2B-P2B
3	A	501	NDP	PA-O3-PN-O1N
3	B	501	NDP	PA-O3-PN-O1N
3	A	501	NDP	PA-O3-PN-O2N
3	B	501	NDP	PA-O3-PN-O2N
3	A	501	NDP	C2N-C3N-C7N-N7N
2	B	502	FAD	O4B-C4B-C5B-O5B
2	A	500	FAD	O4B-C4B-C5B-O5B
3	B	501	NDP	O4B-C1B-N9A-C8A

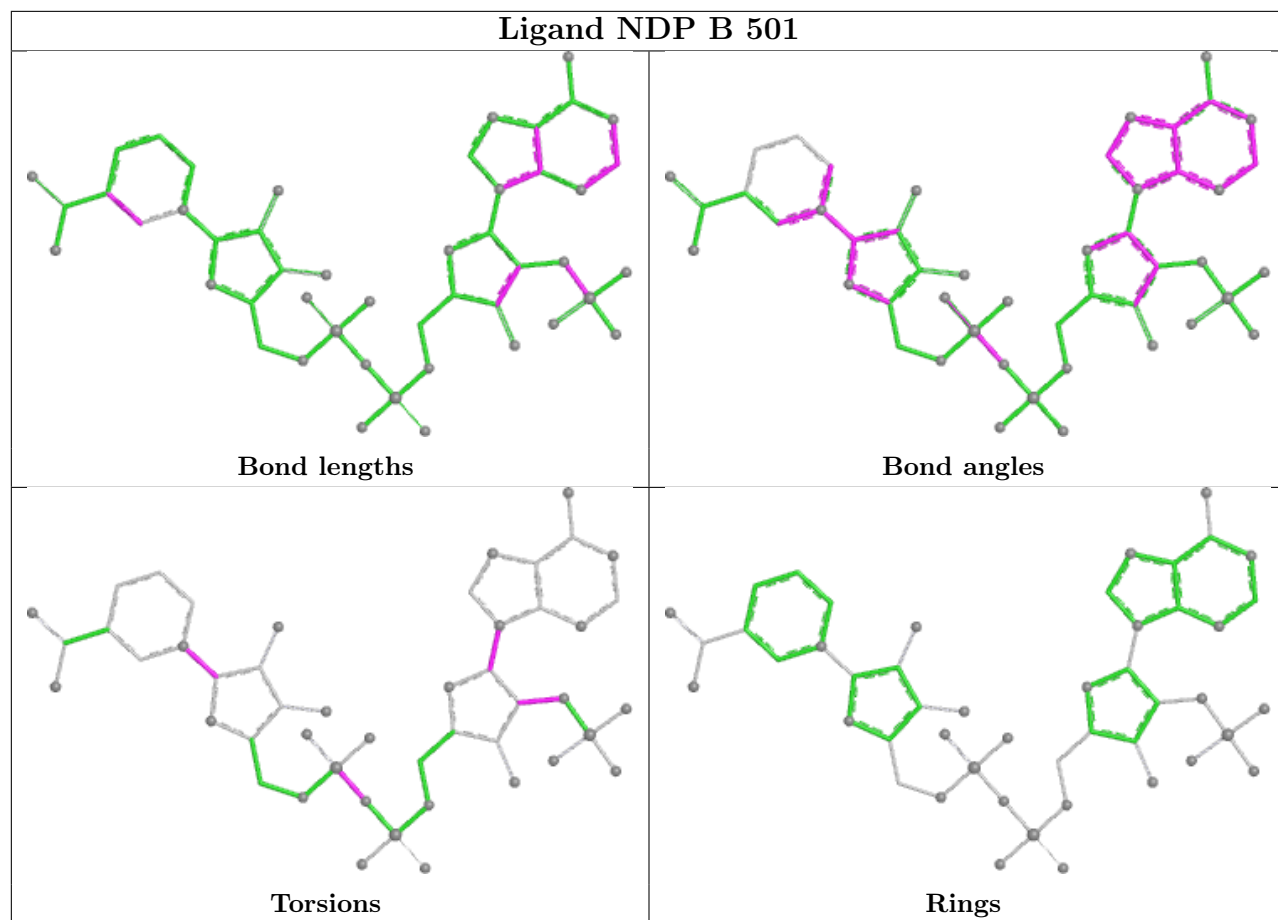
There are no ring outliers.

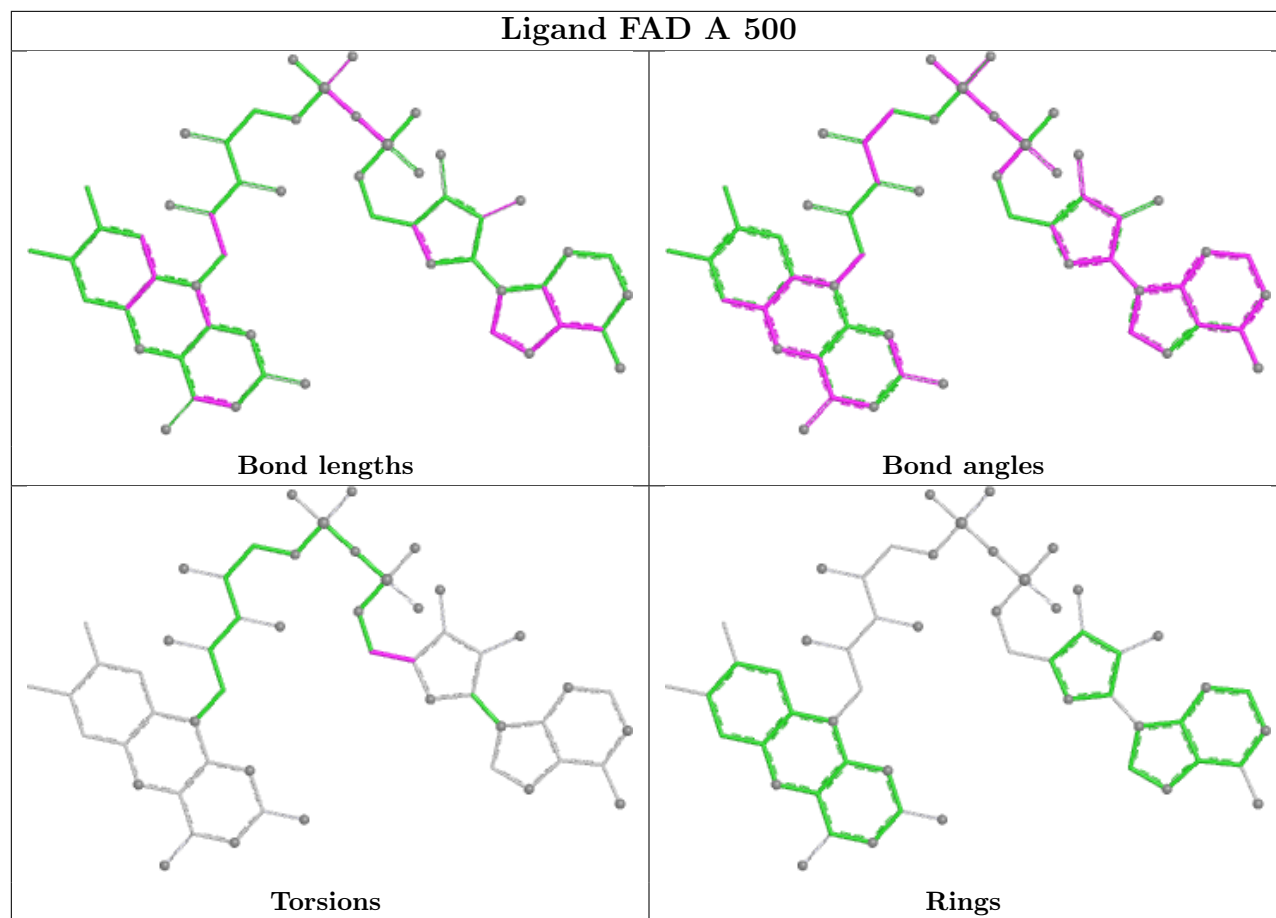
3 monomers are involved in 5 short contacts:

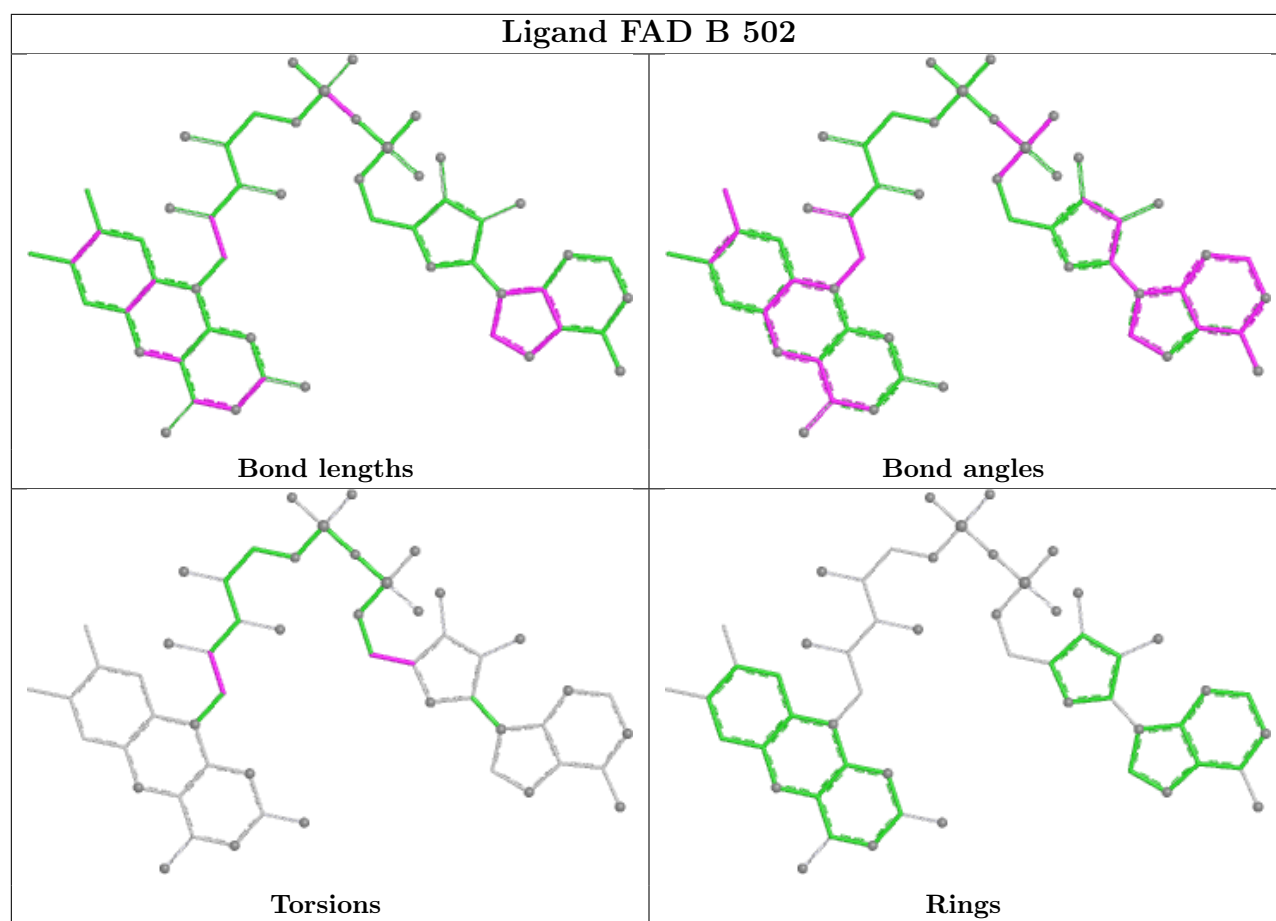
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	NDP	1	0
3	B	501	NDP	3	0
2	A	500	FAD	1	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	429/451 (95%)	0.26	9 (2%) 63 64	20, 34, 59, 103	0
1	B	428/451 (94%)	0.14	8 (1%) 66 66	22, 34, 60, 88	0
All	All	857/902 (95%)	0.20	17 (1%) 65 65	20, 34, 60, 103	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	421	ILE	5.6
1	B	332	GLY	4.1
1	B	115	THR	2.8
1	A	202	ARG	2.5
1	A	198	TRP	2.5
1	A	426	LYS	2.4
1	A	115	THR	2.4
1	A	6	HIS	2.4
1	A	423	GLU	2.3
1	B	4	GLU	2.3
1	A	113	THR	2.3
1	B	105	THR	2.3
1	A	110	VAL	2.1
1	B	9	TYR	2.1
1	B	110	VAL	2.1
1	A	172	GLY	2.1
1	B	429	LEU	2.1

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](#)

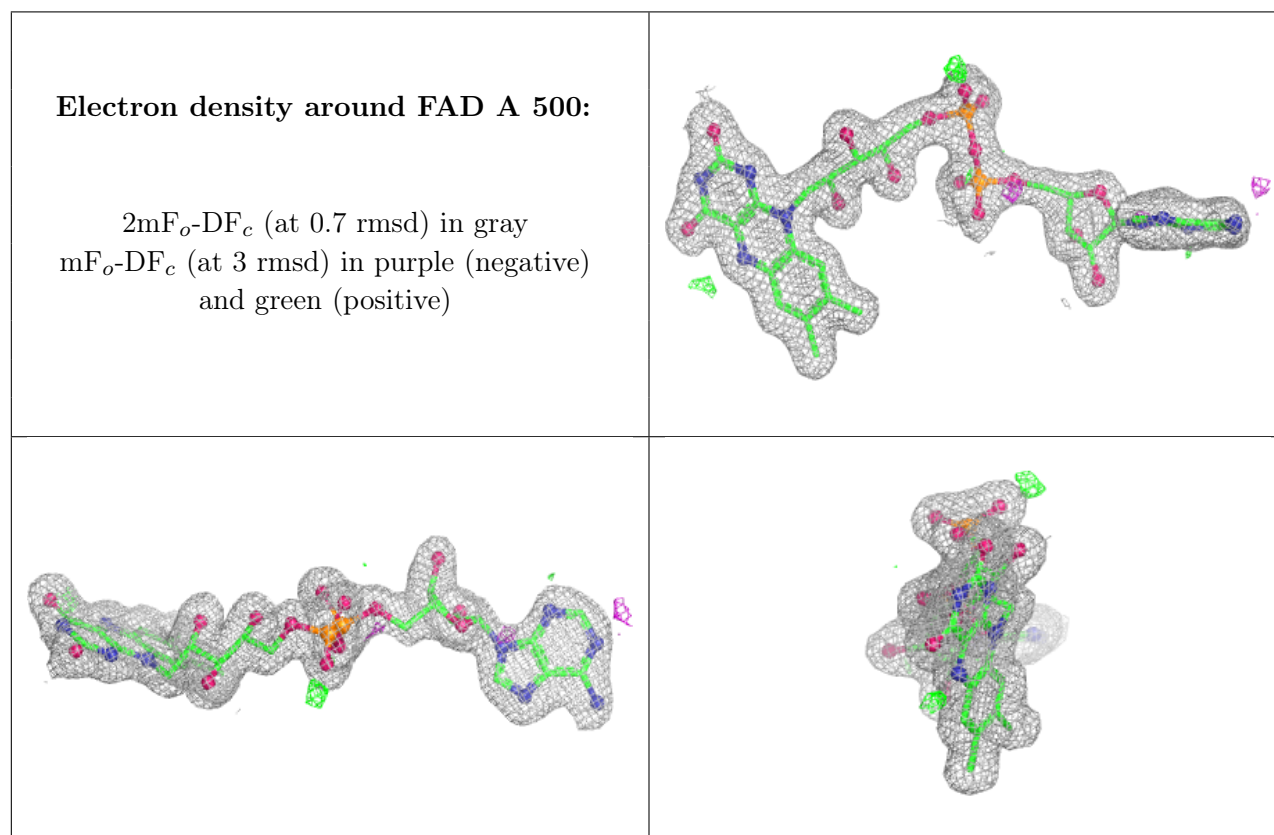
There are no oligosaccharides in this entry.

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.

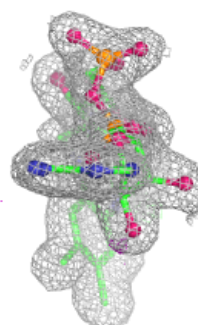
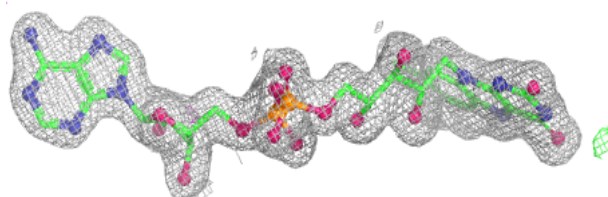
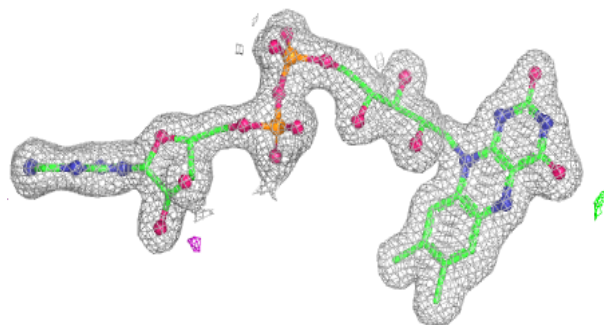
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FAD	A	500	53/53	0.97	0.06	20,24,32,39	0
2	FAD	B	502	53/53	0.97	0.06	20,25,36,40	0
3	NDP	A	501	48/48	0.97	0.06	19,28,51,60	0
3	NDP	B	501	48/48	0.97	0.06	20,25,51,61	0

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.

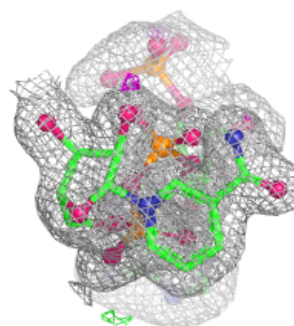
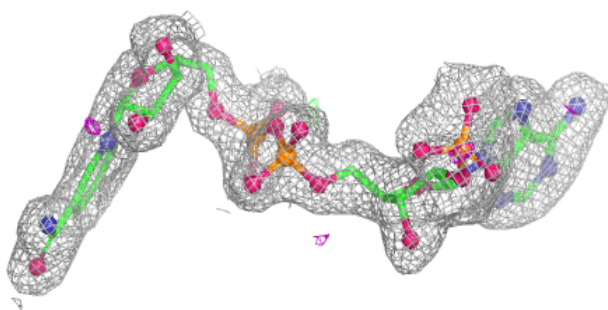
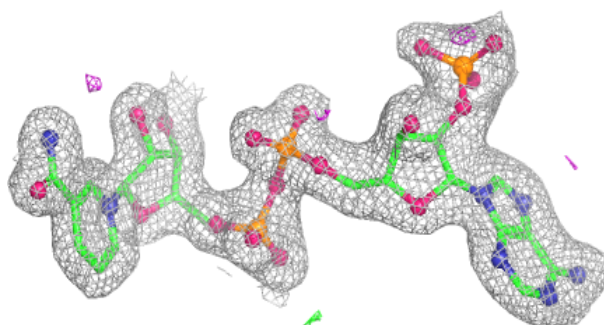


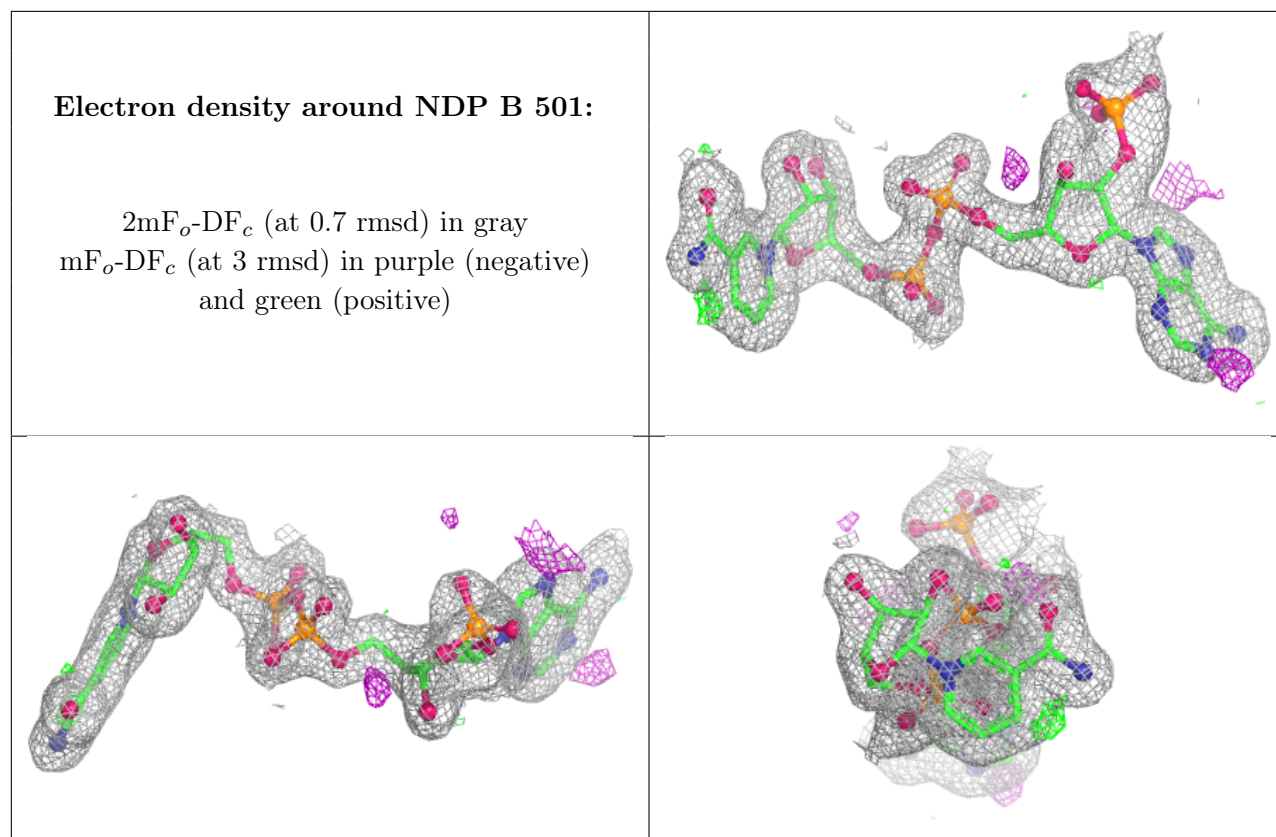
Electron density around FAD B 502:

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

**Electron density around NDP A 501:**

$2mF_o-DF_c$ (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.