Full wwPDB X-ray Structure Validation Report (i)

Oct 20, 2022 – 07:27 pm BST

PDB ID	:	2CDU
Title	:	The Crystal Structure of Water-forming NAD(P)H Oxidase from Lactobacillus sanfranciscensis
Authors	:	Lountos, G.T.; Jiang, R.; Wellborn, W.B.; Thaler, T.L.; Bommarius, A.S.; Orville, A.M.
Deposited on	:	2006-01-28
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 (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 (i)) were used in the production of this report:

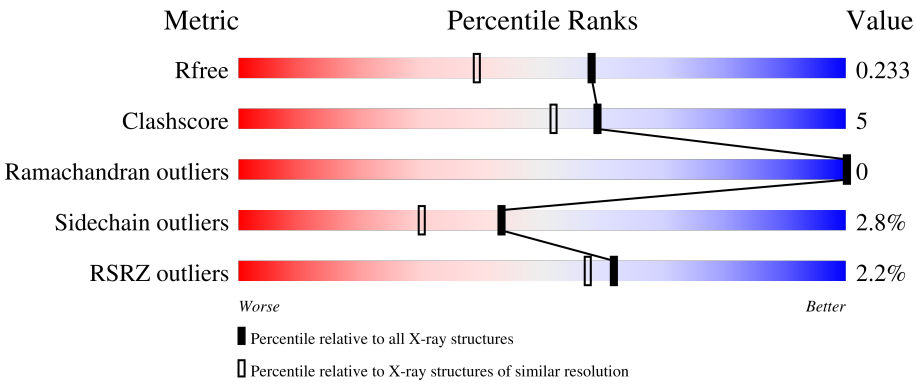
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

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 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (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 <=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	452	2% 87% 11% •
1	B	452	2% 89% 9% ••

2 Entry composition (i)

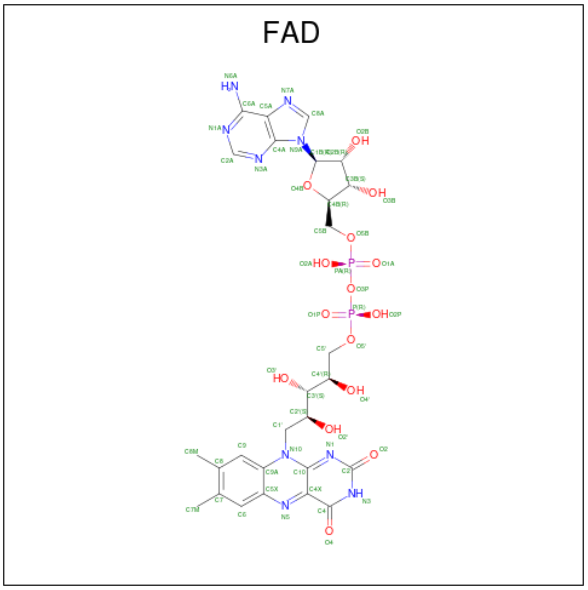
There are 4 unique types of molecules in this entry. The entry contains 7846 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 OXIDASE.

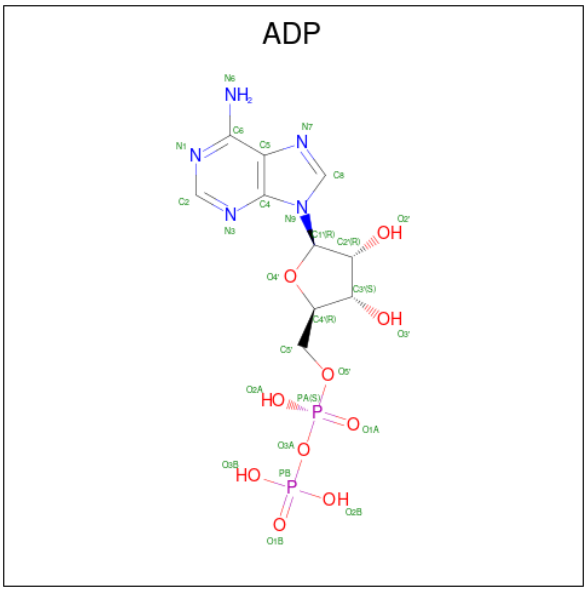
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	451	3493	2197	587	692	17	0	4	0
1	B	449	3485	2191	585	691	18	0	8	0

- 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	53	27	9	15	2	0	0

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
3	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 4 is water.

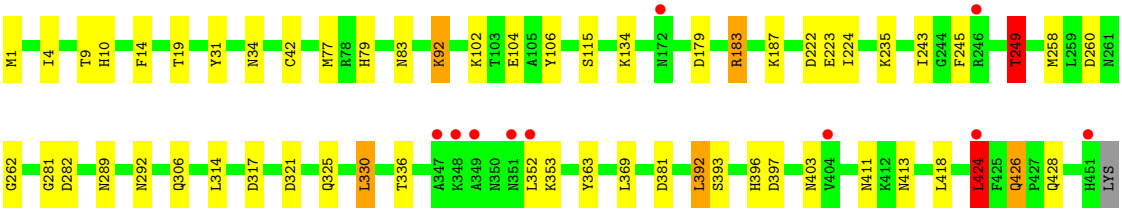
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	382	Total	O	0	0
			382	382		
4	B	326	Total	O	0	0
			326	326		

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.

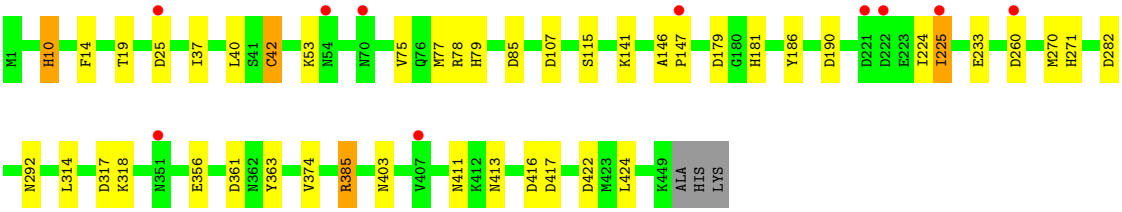
- Molecule 1: NADPH OXIDASE

Chain A:



- Molecule 1: NADPH OXIDASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	59.64Å 92.64Å 163.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	81.65 – 1.80 40.30 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.1 (81.65-1.80) 96.0 (40.30-1.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.56 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.178 , 0.223 0.189 , 0.233	Depositor DCC
R_{free} test set	4054 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	23.8	Xtriage
Anisotropy	0.613	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7846	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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: ADP, CSX, 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	0.66	0/3555	0.89	14/4818 (0.3%)
1	B	0.64	0/3572	0.84	10/4840 (0.2%)
All	All	0.65	0/7127	0.86	24/9658 (0.2%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	183	ARG	NE-CZ-NH2	-9.39	115.60	120.30
1	A	183	ARG	NE-CZ-NH1	9.32	124.96	120.30
1	B	385	ARG	NE-CZ-NH2	-9.31	115.65	120.30
1	A	392	LEU	CA-CB-CG	8.66	135.21	115.30
1	A	330	LEU	CA-CB-CG	8.40	134.61	115.30
1	A	282	ASP	CB-CG-OD2	8.03	125.53	118.30
1	B	416	ASP	CB-CG-OD2	6.91	124.52	118.30
1	A	381	ASP	CB-CG-OD2	6.76	124.38	118.30
1	B	385	ARG	NE-CZ-NH1	6.65	123.62	120.30
1	B	179	ASP	CB-CG-OD2	5.94	123.65	118.30
1	B	417	ASP	CB-CG-OD2	5.77	123.49	118.30
1	A	260	ASP	CB-CG-OD2	5.73	123.46	118.30
1	A	424	LEU	CA-CB-CG	5.61	128.19	115.30
1	A	249	THR	N-CA-CB	-5.59	99.67	110.30
1	A	317	ASP	CB-CG-OD2	5.44	123.19	118.30
1	A	222	ASP	CB-CG-OD2	5.42	123.17	118.30
1	B	260	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	85	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	392	LEU	CB-CG-CD2	5.23	119.89	111.00
1	A	179	ASP	CB-CG-OD2	5.22	123.00	118.30
1	B	361	ASP	CB-CG-OD2	5.09	122.88	118.30
1	B	107	ASP	CB-CG-OD2	5.09	122.88	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	78	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	A	321	ASP	CB-CG-OD1	5.03	122.82	118.30

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	3493	0	3464	43	0
1	B	3485	0	3458	30	0
2	A	53	0	31	1	0
2	B	53	0	31	2	0
3	A	27	0	12	1	0
3	B	27	0	12	0	0
4	A	382	0	0	12	1
4	B	326	0	0	6	1
All	All	7846	0	7008	67	1

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 5.

All (67) 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:289:ASN:HD22	1:A:325:GLN:HE22	1.09	0.99
1:A:249:THR:HG21	1:A:262:GLY:O	1.68	0.93
1:A:411:ASN:HD21	1:B:413:ASN:HD21	1.20	0.85
1:A:363:TYR:CZ	1:A:396[B]:HIS:HD2	1.96	0.84
1:A:289:ASN:HD22	1:A:325:GLN:NE2	1.80	0.80
1:A:413:ASN:HD21	1:B:411:ASN:HD21	1.29	0.78
1:B:25:ASP:OD1	4:B:2014:HOH:O	2.02	0.77
1:B:318:LYS:NZ	4:B:2224:HOH:O	2.21	0.74
1:A:223:GLU:OE2	1:A:235:LYS:NZ	2.21	0.73
1:A:363:TYR:CE1	1:A:396[B]:HIS:HD2	2.06	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:ILE:HG23	1:B:77[A]:MET:HE2	1.70	0.72
1:B:270:MET:SD	4:B:2194:HOH:O	2.48	0.72
1:B:10[A]:HIS:ND1	4:B:2006:HOH:O	2.23	0.71
1:A:10[A]:HIS:ND1	4:A:2006:HOH:O	2.24	0.69
1:A:403:ASN:HD21	1:B:424:LEU:H	1.38	0.68
1:A:10[A]:HIS:CE1	4:A:2006:HOH:O	2.46	0.67
1:A:428[B]:GLN:OE1	4:A:2352:HOH:O	2.14	0.66
1:A:281:GLY:H	1:A:306:GLN:HE21	1.43	0.65
1:A:34:ASN:ND2	4:A:2024:HOH:O	2.30	0.64
1:A:187:LYS:NZ	4:A:2173:HOH:O	2.29	0.64
1:B:225:ILE:HD12	1:B:233:GLU:HB3	1.81	0.61
1:A:31:TYR:OH	1:A:104:GLU:OE2	2.09	0.60
1:A:363:TYR:CE1	1:A:396[B]:HIS:CD2	2.88	0.59
1:A:396[B]:HIS:ND1	1:A:397:ASP:N	2.50	0.59
1:B:42[B]:CSX:OD	2:B:500:FAD:C9A	2.51	0.58
1:B:53:LYS:NZ	4:B:2037:HOH:O	2.37	0.57
1:A:42[B]:CSX:OD	2:A:500:FAD:C9A	2.53	0.56
1:A:424:LEU:H	1:B:403:ASN:HD21	1.53	0.56
1:A:34:ASN:CG	4:A:2026:HOH:O	2.44	0.56
1:A:183:ARG:HD3	4:A:2168:HOH:O	2.06	0.55
1:A:281:GLY:H	1:A:306:GLN:NE2	2.05	0.54
1:A:249:THR:HG23	1:A:258:MET:CE	2.38	0.53
1:B:40:LEU:HD22	1:B:42[A]:CSX:OD	2.08	0.53
1:A:363:TYR:CZ	1:A:396[B]:HIS:CD2	2.87	0.53
1:A:83:ASN:HB3	1:A:92:LYS:HB2	1.90	0.53
1:B:37:ILE:HG23	1:B:77[A]:MET:CE	2.39	0.52
1:A:79:HIS:HE1	4:A:2093:HOH:O	1.92	0.51
1:B:146:ALA:HB3	1:B:147:PRO:HD3	1.92	0.51
1:B:19:THR:HG23	1:B:314:LEU:HD22	1.95	0.49
1:B:271:HIS:HE1	1:B:317:ASP:OD1	1.96	0.48
1:A:134:LYS:HG3	1:A:245:PHE:CE2	2.48	0.48
1:A:249:THR:HG22	4:A:2227:HOH:O	2.13	0.47
1:A:77:MET:O	1:A:79:HIS:HD2	1.98	0.47
1:B:75:VAL:CG1	1:B:77[B]:MET:HE3	2.45	0.46
1:A:243:ILE:HA	3:A:501:ADP:H5'2	1.98	0.45
1:B:190:ASP:OD2	1:B:356:GLU:OE2	2.35	0.45
1:A:9:THR:HG22	4:A:2034:HOH:O	2.16	0.45
1:A:10[B]:HIS:CE1	4:A:2373:HOH:O	2.70	0.45
1:A:4:ILE:HD12	1:A:106:TYR:HB3	1.99	0.45
1:A:411:ASN:ND2	1:B:413:ASN:HD21	2.00	0.44
1:B:225:ILE:CD1	1:B:233:GLU:HB3	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:THR:HG23	1:A:258:MET:HE1	1.99	0.43
1:A:115:SER:HB2	1:A:245:PHE:HB3	1.99	0.43
1:B:10[A]:HIS:CD2	4:B:2005:HOH:O	2.72	0.43
1:B:115:SER:HB3	1:B:282:ASP:HB3	2.00	0.43
1:A:336:THR:O	1:A:393:SER:HA	2.19	0.42
1:B:363:TYR:HB3	1:B:374[B]:VAL:CG2	2.49	0.42
1:B:75:VAL:CG1	1:B:77[B]:MET:CE	2.98	0.41
1:A:325:GLN:HB2	1:B:422:ASP:OD2	2.20	0.41
1:B:42[B]:CSX:OD	2:B:500:FAD:N10	2.53	0.41
1:A:19:THR:HG23	1:A:314:LEU:HD22	2.02	0.41
1:B:19:THR:HG23	1:B:314:LEU:CD2	2.51	0.41
1:A:292:ASN:HB2	4:A:2256:HOH:O	2.22	0.40
1:B:181:HIS:CD2	1:B:186:TYR:CD2	3.10	0.40
1:A:289:ASN:HB2	1:A:325:GLN:HE21	1.86	0.40

All (1) 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)
4:A:2121:HOH:O	4:B:2074:HOH:O[2_554]	2.11	0.09

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	451/452 (100%)	437 (97%)	14 (3%)	0	100	100
1	B	453/452 (100%)	441 (97%)	12 (3%)	0	100	100
All	All	904/904 (100%)	878 (97%)	26 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	384/383 (100%)	370 (96%)	14 (4%)	35	20
1	B	387/383 (101%)	379 (98%)	8 (2%)	53	42
All	All	771/766 (101%)	749 (97%)	22 (3%)	43	29

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	14	PHE
1	A	92	LYS
1	A	102	LYS
1	A	224	ILE
1	A	249	THR
1	A	330	LEU
1	A	352	LEU
1	A	353	LYS
1	A	369	LEU
1	A	392	LEU
1	A	418	LEU
1	A	424	LEU
1	A	426	GLN
1	B	10[A]	HIS
1	B	10[B]	HIS
1	B	14	PHE
1	B	141	LYS
1	B	224	ILE
1	B	225	ILE
1	B	292	ASN
1	B	385	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	ASN
1	A	79	HIS
1	A	271	HIS
1	A	306	GLN
1	A	325	GLN
1	A	403	ASN
1	A	411	ASN
1	A	426	GLN
1	A	451	HIS
1	B	79	HIS
1	B	271	HIS
1	B	292	ASN
1	B	403	ASN
1	B	411	ASN
1	B	428	GLN
1	B	446	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)

4 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	CSX	B	42[A]	-	3,6,7	0.60	0	1,6,8	2.66	1 (100%)
1	CSX	A	42[A]	-	3,6,7	0.66	0	1,6,8	1.67	0
1	CSX	B	42[B]	-	3,6,7	0.60	0	1,6,8	2.66	1 (100%)
1	CSX	A	42[B]	-	3,6,7	0.66	0	1,6,8	1.67	0

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
1	CSX	B	42[A]	-	-	1/1/5/7	-
1	CSX	A	42[A]	-	-	1/1/5/7	-
1	CSX	B	42[B]	-	-	1/1/5/7	-
1	CSX	A	42[B]	-	-	1/1/5/7	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	42[A]	CSX	CA-CB-SG	-2.66	107.55	113.36
1	B	42[B]	CSX	CA-CB-SG	-2.66	107.55	113.36

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	42[A]	CSX	N-CA-CB-SG
1	A	42[B]	CSX	N-CA-CB-SG
1	B	42[A]	CSX	N-CA-CB-SG
1	B	42[B]	CSX	N-CA-CB-SG

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	42[A]	CSX	1	0
1	B	42[B]	CSX	2	0
1	A	42[B]	CSX	1	0

5.5 Carbohydrates (i)

There are no monosaccharides 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	ADP	A	501	-	24,29,29	1.23	2 (8%)	29,45,45	1.51	5 (17%)
3	ADP	B	501	-	24,29,29	1.22	3 (12%)	29,45,45	1.54	4 (13%)
2	FAD	B	500	-	53,58,58	1.38	6 (11%)	68,89,89	1.38	9 (13%)
2	FAD	A	500	-	53,58,58	1.24	3 (5%)	68,89,89	1.49	10 (14%)

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	ADP	A	501	-	-	8/12/32/32	0/3/3/3
3	ADP	B	501	-	-	3/12/32/32	0/3/3/3
2	FAD	B	500	-	-	2/30/50/50	0/6/6/6
2	FAD	A	500	-	-	2/30/50/50	0/6/6/6

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	FAD	C4X-N5	4.83	1.40	1.30
2	B	500	FAD	C4X-N5	4.75	1.40	1.30
3	B	501	ADP	C2-N3	3.74	1.38	1.32
3	A	501	ADP	C2-N3	3.62	1.37	1.32
2	B	500	FAD	C10-N1	3.33	1.40	1.33
2	A	500	FAD	O4B-C1B	3.10	1.45	1.41
2	B	500	FAD	C5'-C4'	2.68	1.55	1.51
2	B	500	FAD	P-O5'	2.40	1.69	1.59
3	A	501	ADP	PB-O3B	2.37	1.64	1.54
3	B	501	ADP	C2-N1	2.37	1.38	1.33
2	B	500	FAD	O5'-C5'	-2.31	1.35	1.44
3	B	501	ADP	PB-O3B	2.28	1.63	1.54
2	A	500	FAD	C10-N1	2.11	1.37	1.33
2	B	500	FAD	C7M-C7	2.02	1.55	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	501	ADP	N3-C2-N1	-5.67	119.81	128.68
3	A	501	ADP	N3-C2-N1	-5.22	120.53	128.68
2	B	500	FAD	N3A-C2A-N1A	-4.64	121.42	128.68
2	A	500	FAD	N3A-C2A-N1A	-4.20	122.11	128.68
2	A	500	FAD	O4B-C1B-C2B	-3.49	101.82	106.93
2	A	500	FAD	C9A-C5X-N5	-3.43	118.70	122.43
2	B	500	FAD	C4X-C10-N10	3.31	121.32	116.48
3	A	501	ADP	O2B-PB-O3A	3.19	115.34	104.64
2	A	500	FAD	C4X-C10-N10	2.97	120.82	116.48
2	A	500	FAD	C10-C4X-N5	-2.84	118.84	124.86
3	B	501	ADP	C1'-N9-C4	-2.83	121.66	126.64
3	B	501	ADP	O2B-PB-O3A	2.73	113.77	104.64
2	A	500	FAD	C10-N1-C2	2.72	122.35	116.90
2	B	500	FAD	C10-C4X-N5	-2.69	119.15	124.86
3	A	501	ADP	C1'-N9-C4	-2.65	121.98	126.64
2	A	500	FAD	C4-C4X-N5	2.61	121.95	118.23
2	A	500	FAD	C5X-C9A-N10	2.54	120.57	117.95
2	B	500	FAD	C4X-C10-N1	-2.45	119.06	124.73
3	A	501	ADP	PA-O3A-PB	-2.44	124.44	132.83
2	A	500	FAD	C4X-C4-N3	2.36	119.19	113.19
2	B	500	FAD	O2P-P-O1P	2.33	123.74	112.24
3	B	501	ADP	PA-O3A-PB	-2.28	124.99	132.83
2	B	500	FAD	O4B-C1B-C2B	-2.26	103.62	106.93
2	B	500	FAD	C4-C4X-C10	2.23	120.53	116.79
2	A	500	FAD	C4X-C10-N1	-2.13	119.78	124.73
3	A	501	ADP	C2'-C3'-C4'	2.11	106.74	102.64
2	B	500	FAD	C4X-C4-N3	2.08	118.48	113.19
2	B	500	FAD	C9A-C5X-N5	-2.06	120.19	122.43

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	ADP	C5'-O5'-PA-O1A
3	A	501	ADP	C5'-O5'-PA-O2A
3	A	501	ADP	C5'-O5'-PA-O3A
3	B	501	ADP	C3'-C4'-C5'-O5'
3	B	501	ADP	O4'-C4'-C5'-O5'
2	A	500	FAD	PA-O3P-P-O5'
2	B	500	FAD	PA-O3P-P-O5'
3	A	501	ADP	PA-O3A-PB-O1B
3	A	501	ADP	O4'-C4'-C5'-O5'
2	B	500	FAD	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
3	A	501	ADP	PA-O3A-PB-O2B
3	A	501	ADP	PA-O3A-PB-O3B
2	A	500	FAD	O4B-C4B-C5B-O5B
3	B	501	ADP	C5'-O5'-PA-O1A
3	A	501	ADP	C3'-C4'-C5'-O5'

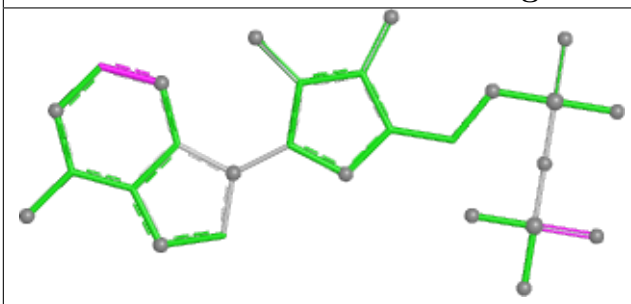
There are no ring outliers.

3 monomers are involved in 4 short contacts:

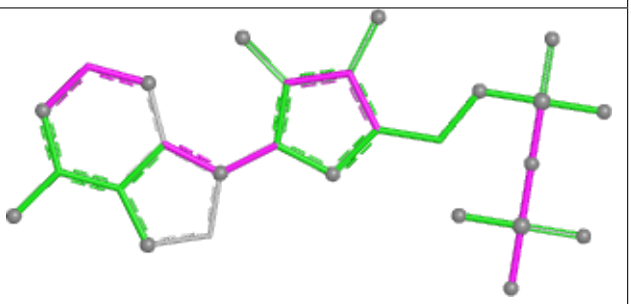
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	501	ADP	1	0
2	B	500	FAD	2	0
2	A	500	FAD	1	0

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 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.

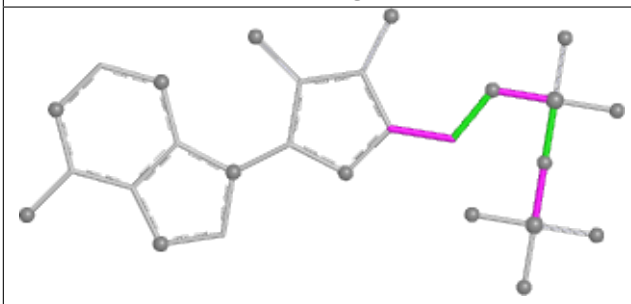
Ligand ADP A 501



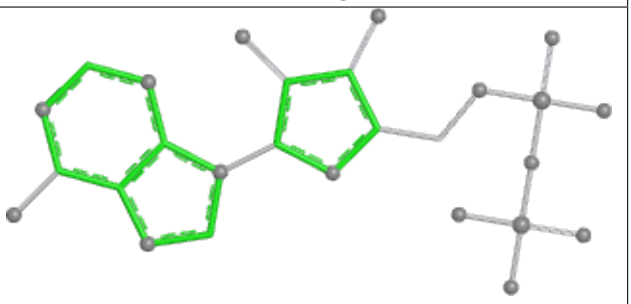
Bond lengths



Bond angles

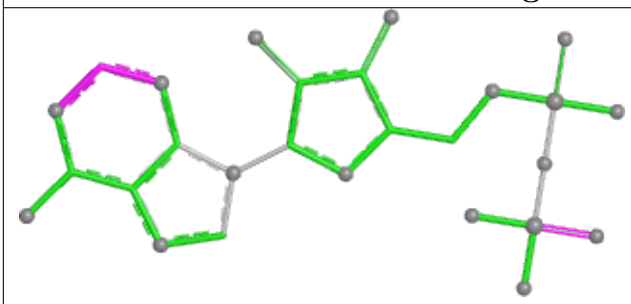


Torsions

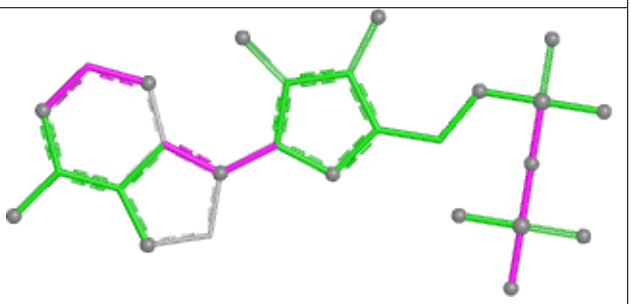


Rings

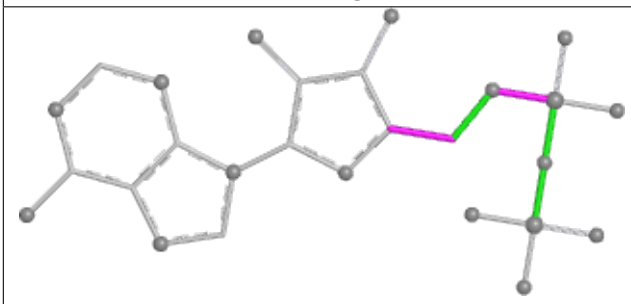
Ligand ADP B 501



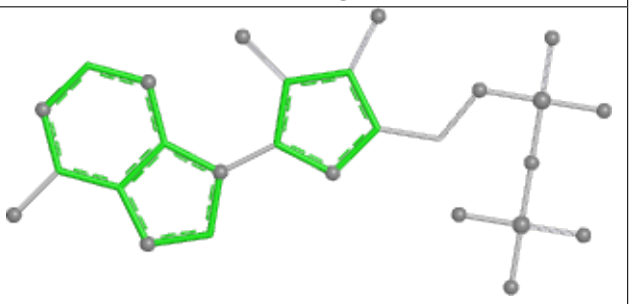
Bond lengths



Bond angles

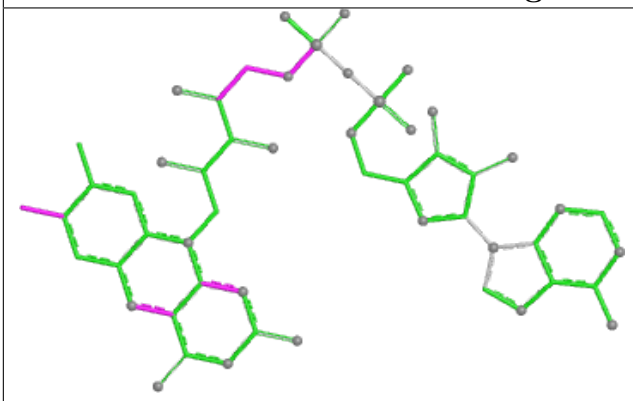


Torsions

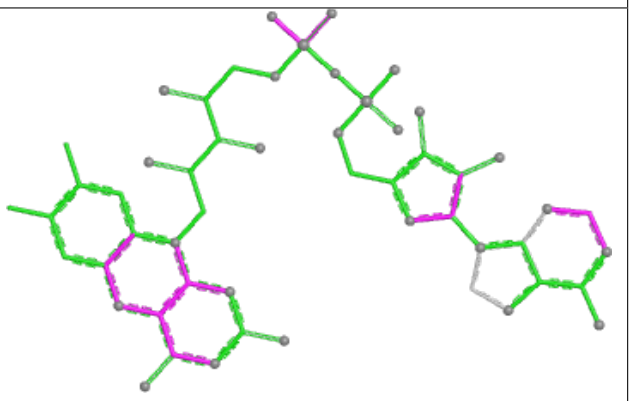


Rings

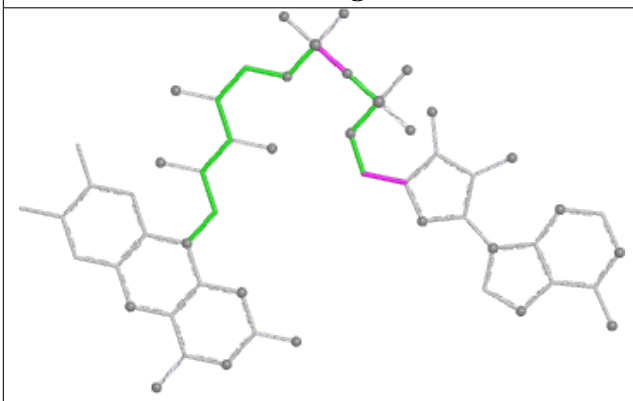
Ligand FAD B 500



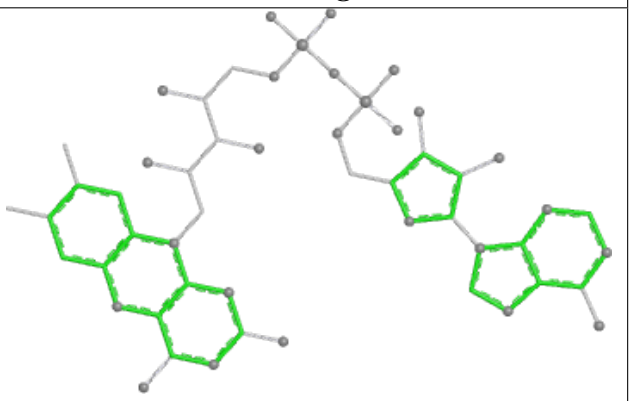
Bond lengths



Bond angles

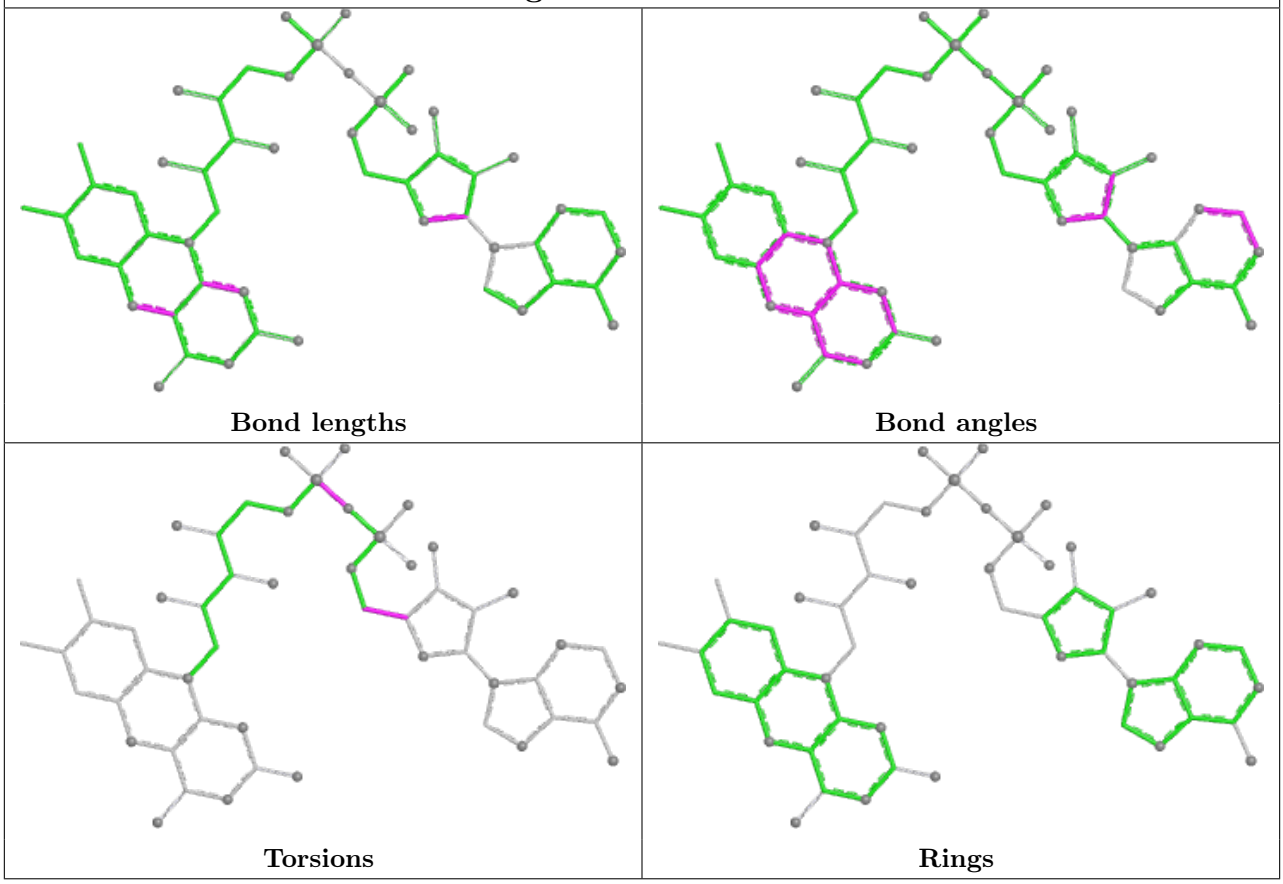


Torsions



Rings

Ligand FAD A 500



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	42:CSX	C	43:GLY	N	1.65
1	B	41:SER	C	42:CSX	N	1.60

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	450/452 (99%)	-0.14	10 (2%)	62	57	16, 24, 37, 56	0
1	B	448/452 (99%)	-0.00	10 (2%)	62	57	15, 28, 42, 54	0
All	All	898/904 (99%)	-0.07	20 (2%)	62	57	15, 26, 41, 56	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	451	HIS	5.5
1	B	221	ASP	5.2
1	B	25	ASP	3.9
1	A	348	LYS	3.4
1	B	147	PRO	3.3
1	A	351	ASN	3.1
1	A	349	ALA	2.9
1	B	222	ASP	2.8
1	B	351	ASN	2.7
1	A	347	ALA	2.5
1	B	260	ASP	2.4
1	A	172	ASN	2.3
1	B	54	ASN	2.3
1	A	246	ARG	2.3
1	B	70	ASN	2.3
1	A	424	LEU	2.3
1	A	404	VAL	2.2
1	B	407	VAL	2.2
1	A	352	LEU	2.1
1	B	225	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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(Å^2)	Q<0.9
1	CSX	B	42[A]	7/8	0.95	0.14	18,25,25,26	1
1	CSX	B	42[B]	7/8	0.95	0.14	21,25,25,26	1
1	CSX	A	42[A]	7/8	0.98	0.12	18,19,20,22	1
1	CSX	A	42[B]	7/8	0.98	0.12	18,19,21,22	1

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

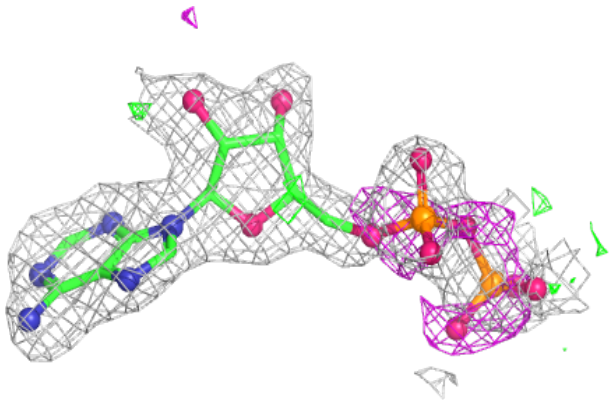
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

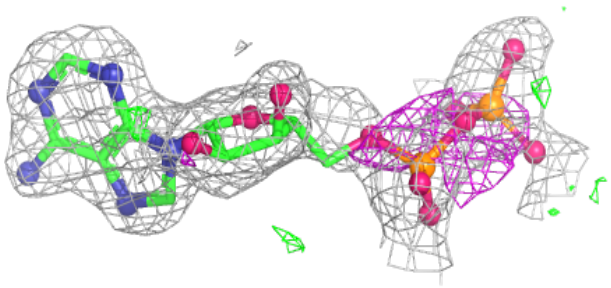
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ADP	B	501	27/27	0.85	0.18	37,43,59,61	0
2	FAD	B	500	53/53	0.86	0.16	20,24,27,31	9
2	FAD	A	500	53/53	0.90	0.15	15,20,23,23	9
3	ADP	A	501	27/27	0.92	0.09	22,30,48,49	0

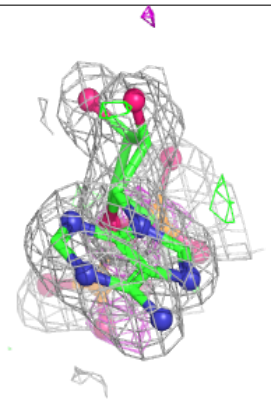
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP B 501:

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

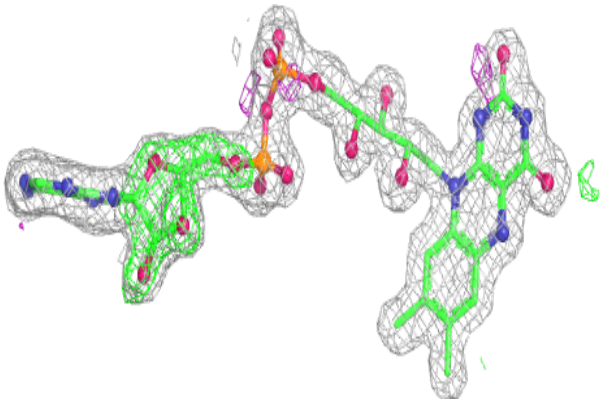


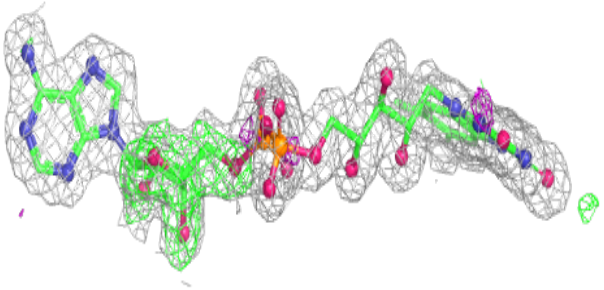


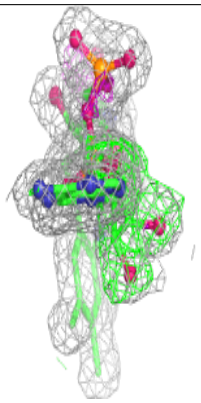


Electron density around FAD B 500:

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

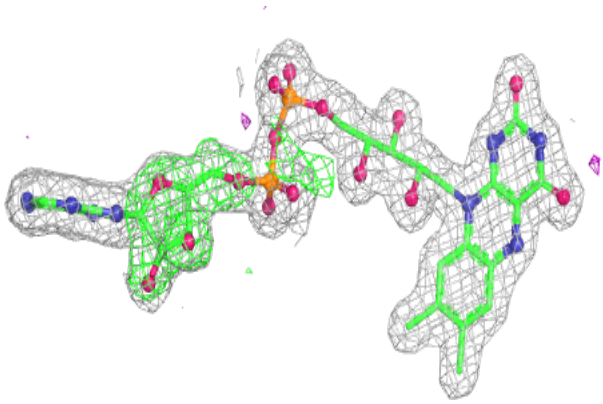


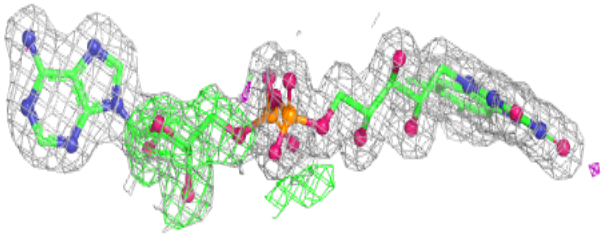


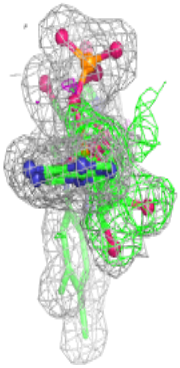


Electron density around FAD A 500:

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

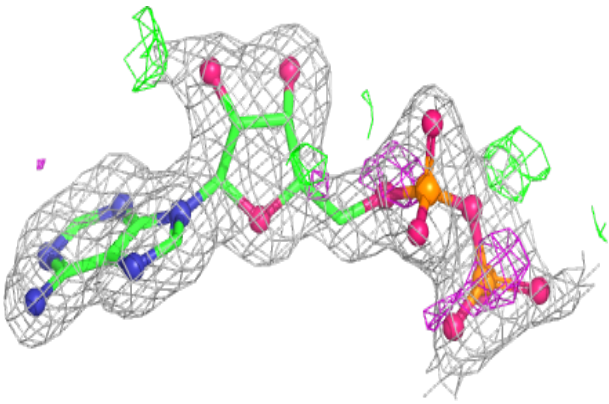


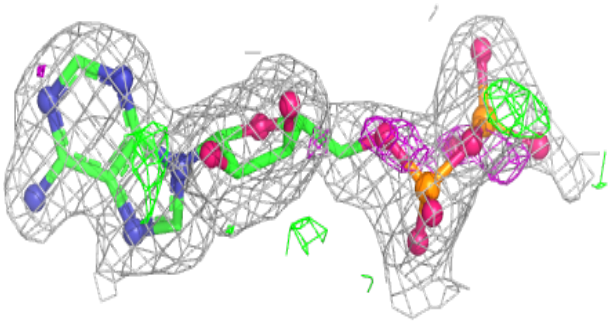


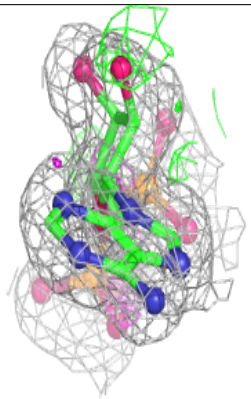


Electron density around ADP A 501:

$2mF_o$-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.