



wwPDB X-ray Structure Validation Summary Report ⓘ

May 3, 2025 – 03:11 PM EDT

PDB ID : 3AD9 / pdb_00003ad9
Title : Heterotetrameric Sarcosine Oxidase from Corynebacterium sp. U-96 sarcosine-reduced form
Authors : Suzuki, H.; Moriguchi, T.; Ida, K.
Deposited on : 2010-01-15
Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 Phenix2.0rc1
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

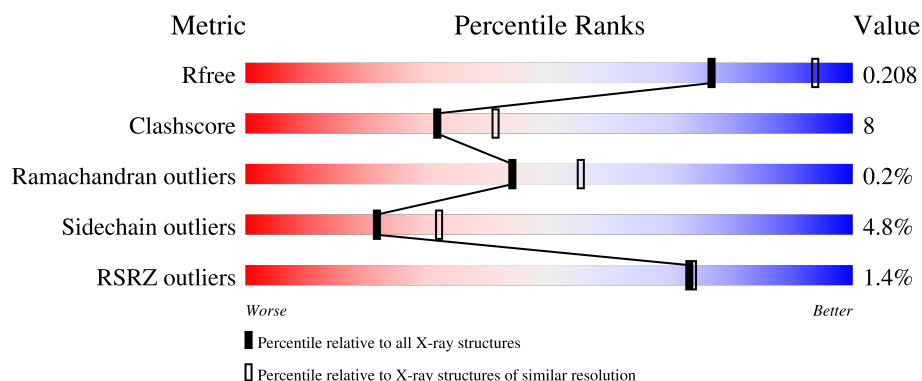
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



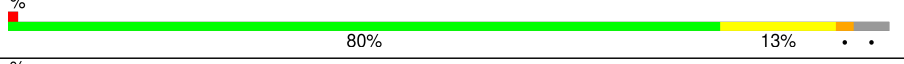

The reported resolution of this entry is 2.30 Å.

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}	164625	5963 (2.30-2.30)
Clashscore	180529	6698 (2.30-2.30)
Ramachandran outliers	177936	6640 (2.30-2.30)
Sidechain outliers	177891	6640 (2.30-2.30)
RSRZ outliers	164620	5963 (2.30-2.30)

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	964	
2	B	404	
3	C	203	
4	D	99	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	SO4	A	2559	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13569 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 SARCOSINE OXIDASE ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	963	Total	C	N	O	S	0	0	0
			7229	4507	1287	1413	22			

- Molecule 2 is a protein called SARCOSINE OXIDASE BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	397	Total	C	N	O	S	0	1	0
			3079	1965	533	571	10			

- Molecule 3 is a protein called SARCOSINE OXIDASE GAMMA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	195	Total	C	N	O	S	0	0	0
			1433	902	257	271	3			

- Molecule 4 is a protein called SARCOSINE OXIDASE DELTA SUBUNIT.

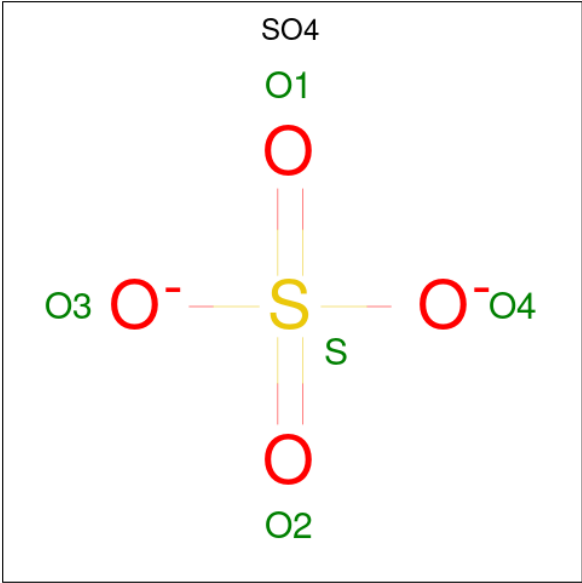
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	91	Total	C	N	O	S	0	0	0
			749	476	135	133	5			

- Molecule 5 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



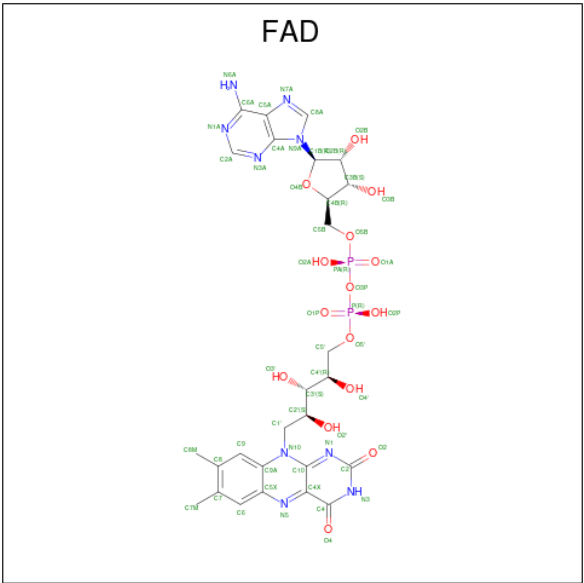
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		

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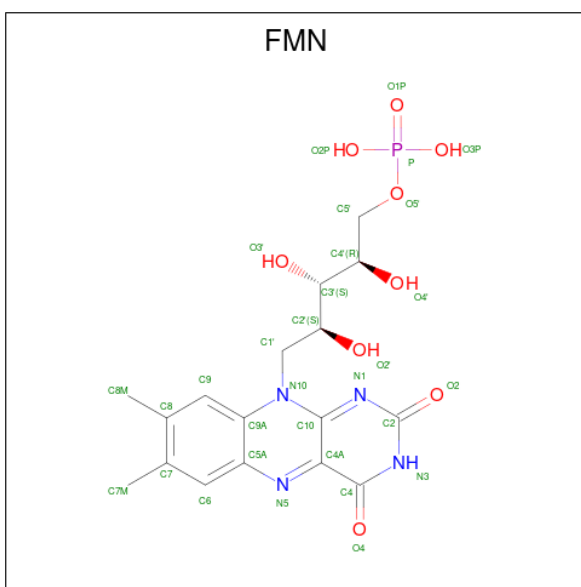
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 8 is FLAVIN MONONUCLEOTIDE (CCD ID: FMN) (formula: C₁₇H₂₁N₄O₉P).



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

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	D	1	Total Zn 1 1	0	0

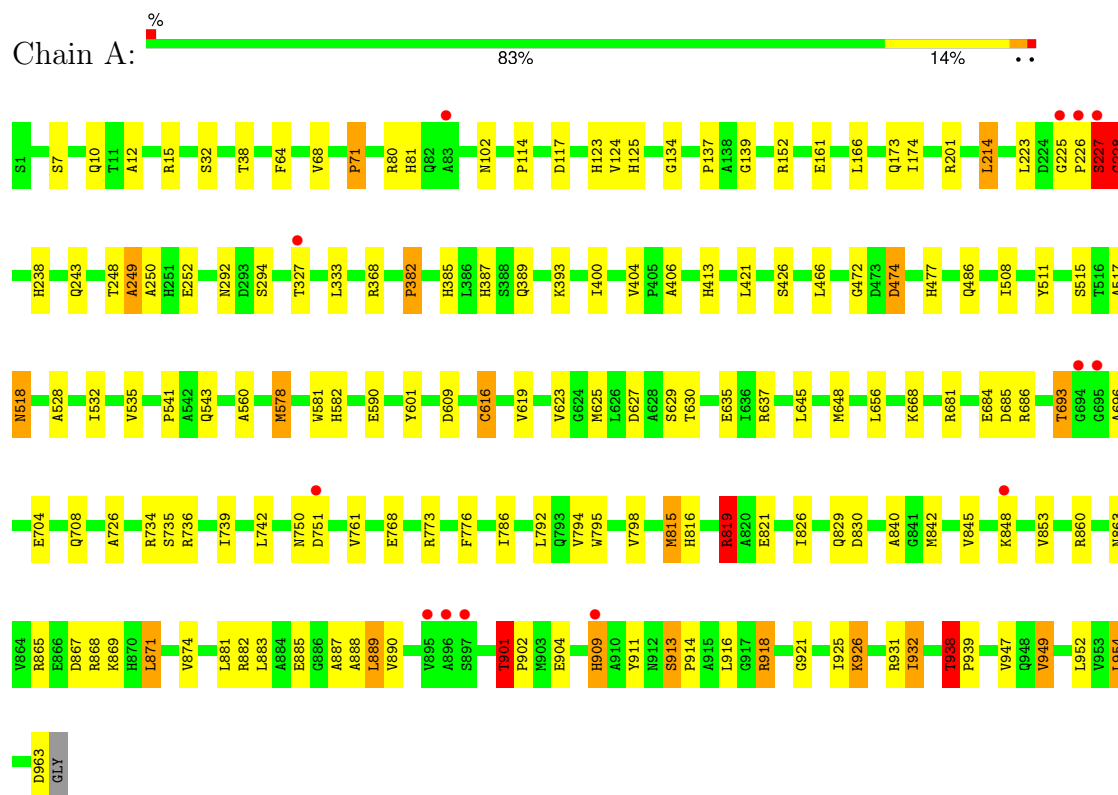
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	557	Total O 557 557	0	0
10	B	163	Total O 163 163	0	0
10	C	119	Total O 119 119	0	0
10	D	66	Total O 66 66	0	0

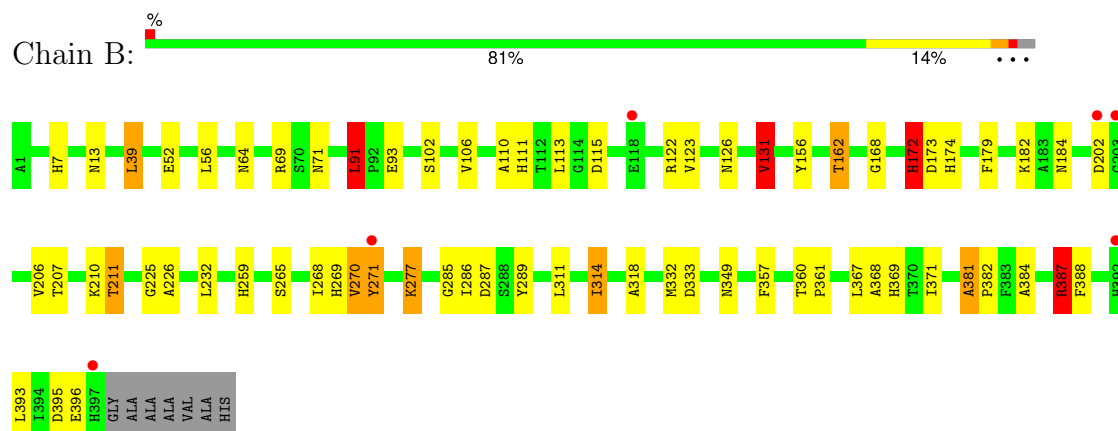
3 Residue-property plots

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.

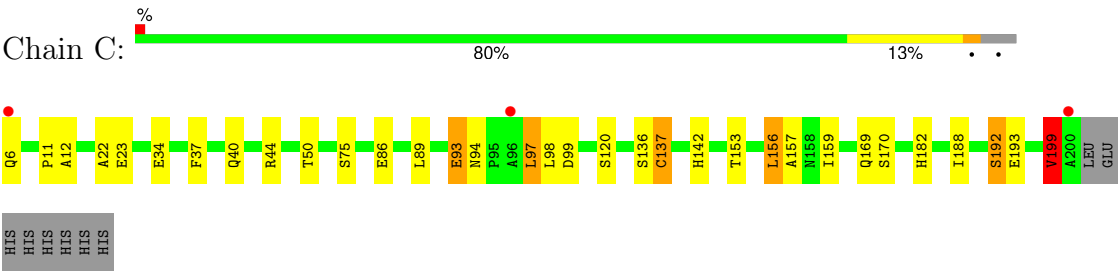
• Molecule 1: SARCOSINE OXIDASE ALPHA SUBUNIT



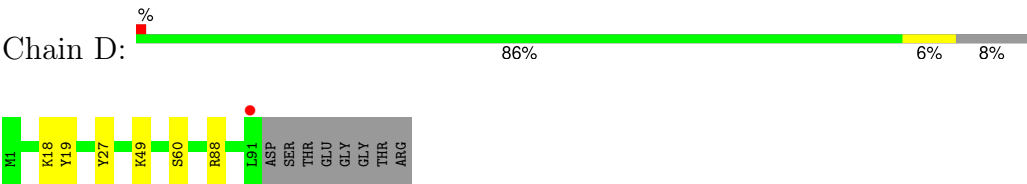
• Molecule 2: SARCOSINE OXIDASE BETA SUBUNIT



● Molecule 3: SARCOSINE OXIDASE GAMMA SUBUNIT



● Molecule 4: SARCOSINE OXIDASE DELTA SUBUNIT



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	198.56Å 198.56Å 196.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.25 – 2.30 61.25 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (61.25-2.30) 99.9 (61.25-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.18 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.169 , 0.211 0.169 , 0.208	Depositor DCC
R_{free} test set	5043 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13569	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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 i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, NAD, FAD, SO4, ZN

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	1.20	7/7361 (0.1%)	1.14	25/10017 (0.2%)
2	B	1.18	5/3160 (0.2%)	1.18	16/4300 (0.4%)
3	C	1.28	0/1461	1.20	8/1998 (0.4%)
4	D	1.08	0/772	1.10	4/1040 (0.4%)
All	All	1.20	12/12754 (0.1%)	1.15	53/17355 (0.3%)

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
2	B	0	1
3	C	0	1
All	All	0	2

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	271[A]	TYR	C-N	13.66	1.46	1.33
2	B	271[B]	TYR	C-N	13.66	1.46	1.33
1	A	249	ALA	CA-CB	-7.46	1.40	1.53
1	A	81	HIS	CA-C	-6.62	1.44	1.52
2	B	270	VAL	C-N	6.50	1.42	1.33

The worst 5 of 53 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	12	ALA	CA-C-N	-8.87	109.39	123.31
1	A	12	ALA	C-N-CA	-8.87	109.39	123.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	259	HIS	CA-C-N	7.68	127.40	119.64
2	B	259	HIS	C-N-CA	7.68	127.40	119.64
2	B	396	GLU	N-CA-C	7.64	121.20	112.57

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	268	ILE	Peptide
3	C	199	VAL	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	7229	0	7104	125	0
2	B	3079	0	3004	59	0
3	C	1433	0	1434	18	0
4	D	749	0	706	2	0
5	A	44	0	26	4	0
6	A	25	0	0	3	0
6	B	10	0	0	0	0
6	C	5	0	0	0	0
6	D	5	0	0	0	0
7	B	53	0	31	3	0
8	B	31	0	19	7	0
9	D	1	0	0	0	0
10	A	557	0	0	17	2
10	B	163	0	0	4	0
10	C	119	0	0	4	0
10	D	66	0	0	1	0
All	All	13569	0	12324	200	2

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 8.

The worst 5 of 200 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:819:ARG:HH12	1:A:909:HIS:CE1	1.38	1.41
2:B:172:HIS:NE2	8:B:406:FMN:C8M	1.79	1.41
2:B:172:HIS:NE2	8:B:406:FMN:HM82	0.96	1.28
1:A:819:ARG:NH1	1:A:909:HIS:CE1	2.11	1.16
1:A:693:THR:HG22	6:A:2559:SO4:O2	1.43	1.12

All (2) 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)
10:A:1013:HOH:O	10:A:1013:HOH:O[9_555]	1.63	0.57
10:A:1390:HOH:O	10:A:1390:HOH:O[10_665]	1.81	0.39

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	961/964 (100%)	932 (97%)	25 (3%)	4 (0%)	30	39
2	B	396/404 (98%)	382 (96%)	14 (4%)	0	100	100
3	C	193/203 (95%)	186 (96%)	7 (4%)	0	100	100
4	D	89/99 (90%)	86 (97%)	3 (3%)	0	100	100
All	All	1639/1670 (98%)	1586 (97%)	49 (3%)	4 (0%)	44	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	227	SER
1	A	228	GLY
1	A	406	ALA
1	A	71	PRO

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	747/747 (100%)	710 (95%)	37 (5%)	20	30
2	B	318/319 (100%)	305 (96%)	13 (4%)	26	39
3	C	143/151 (95%)	133 (93%)	10 (7%)	12	17
4	D	75/81 (93%)	74 (99%)	1 (1%)	65	79
All	All	1283/1298 (99%)	1222 (95%)	61 (5%)	21	32

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	874	VAL
3	C	156	LEU
1	A	949	VAL
3	C	98	LEU
3	C	199	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	111	GLN
3	C	158	ASN
1	A	518	ASN
1	A	489	GLN
3	C	182	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](#)

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 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$
6	SO4	D	2566	-	4,4,4	0.35	0	6,6,6	0.18	0
5	NAD	A	965	-	42,48,48	2.22	7 (16%)	50,73,73	1.53	7 (14%)
8	FMN	B	406	-	33,33,33	1.62	9 (27%)	48,50,50	1.83	12 (25%)
6	SO4	B	2562	-	4,4,4	0.26	0	6,6,6	0.14	0
6	SO4	A	2560	-	4,4,4	0.48	0	6,6,6	0.49	0
6	SO4	B	2561	-	4,4,4	0.34	0	6,6,6	0.63	0
6	SO4	A	2567	-	4,4,4	0.33	0	6,6,6	0.32	0
6	SO4	C	2564	-	4,4,4	0.65	0	6,6,6	0.30	0
6	SO4	A	2559	-	4,4,4	0.27	0	6,6,6	0.31	0
7	FAD	B	405	-	54,58,58	1.65	8 (14%)	71,89,89	1.55	13 (18%)
6	SO4	A	2563	-	4,4,4	0.27	0	6,6,6	0.46	0
6	SO4	A	2565	-	4,4,4	0.28	0	6,6,6	0.20	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
8	FMN	B	406	-	-	4/18/18/18	0/3/3/3
7	FAD	B	405	-	-	4/30/50/50	0/6/6/6
5	NAD	A	965	-	-	6/26/62/62	0/5/5/5

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	965	NAD	C2N-N1N	6.85	1.42	1.35
5	A	965	NAD	O7N-C7N	6.81	1.36	1.24
5	A	965	NAD	O4B-C1B	6.11	1.48	1.40
7	B	405	FAD	C4X-N5	5.71	1.43	1.30
5	A	965	NAD	PN-O3	4.89	1.64	1.59

The worst 5 of 32 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	405	FAD	C4B-O4B-C1B	-6.25	104.20	109.92
5	A	965	NAD	N3A-C2A-N1A	-4.87	122.06	128.67
8	B	406	FMN	O2P-P-O5'	4.65	118.80	106.67
8	B	406	FMN	O3P-P-O5'	-4.30	95.46	106.67
5	A	965	NAD	C4A-C5A-N7A	-3.88	105.23	109.34

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	405	FAD	C5B-O5B-PA-O2A
8	B	406	FMN	C3'-C4'-C5'-O5'
8	B	406	FMN	O4'-C4'-C5'-O5'
8	B	406	FMN	C5'-O5'-P-O2P
8	B	406	FMN	C5'-O5'-P-O3P

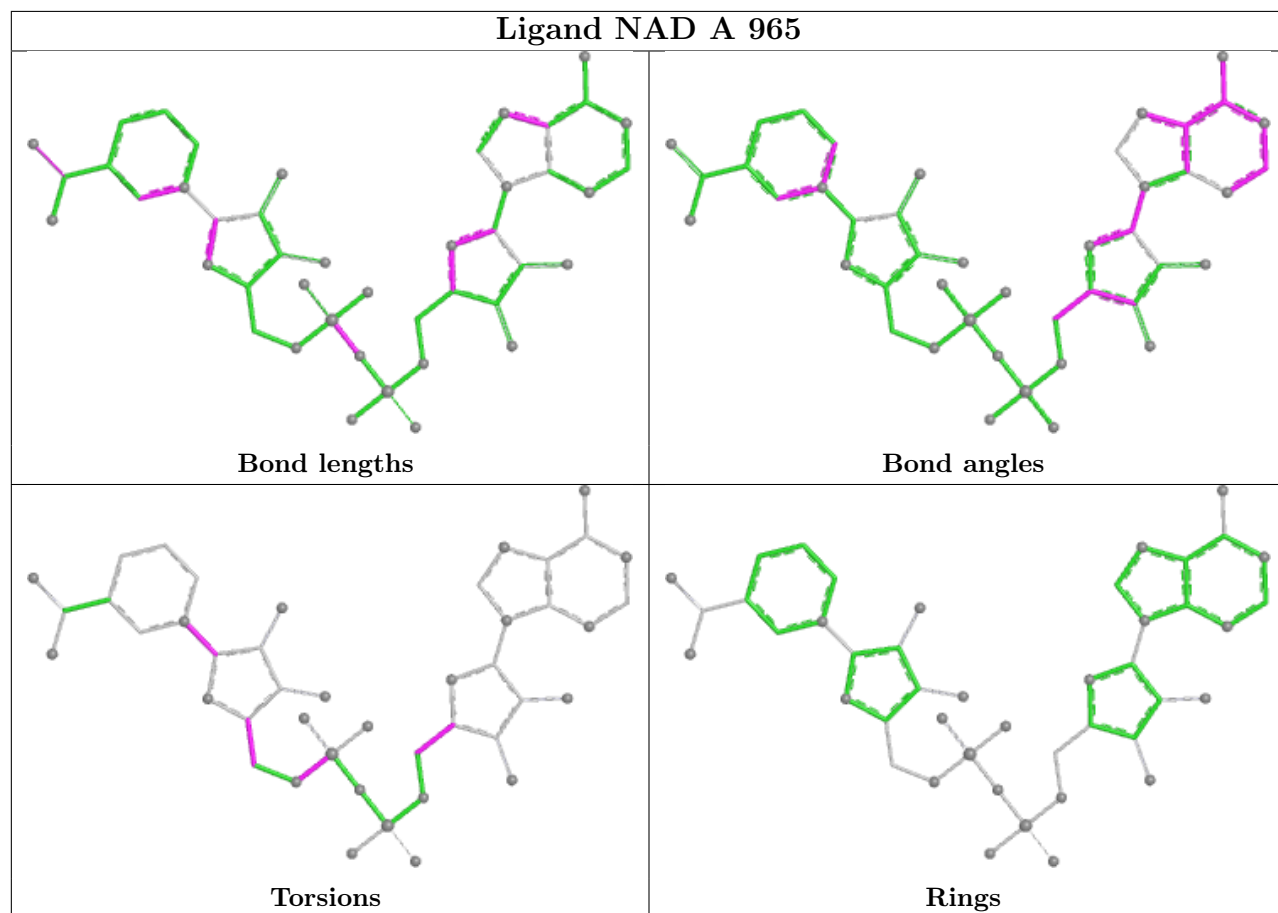
There are no ring outliers.

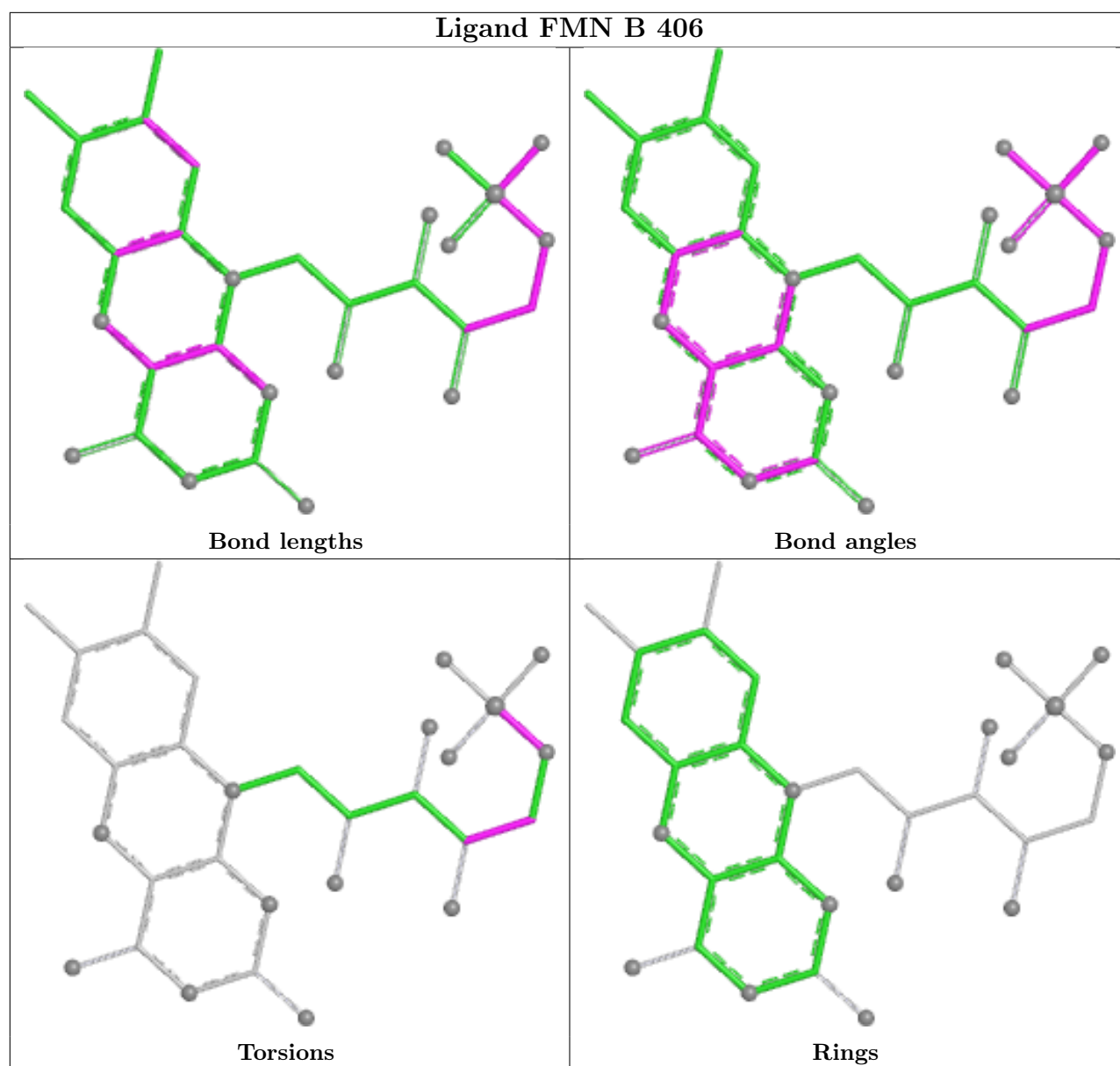
4 monomers are involved in 17 short contacts:

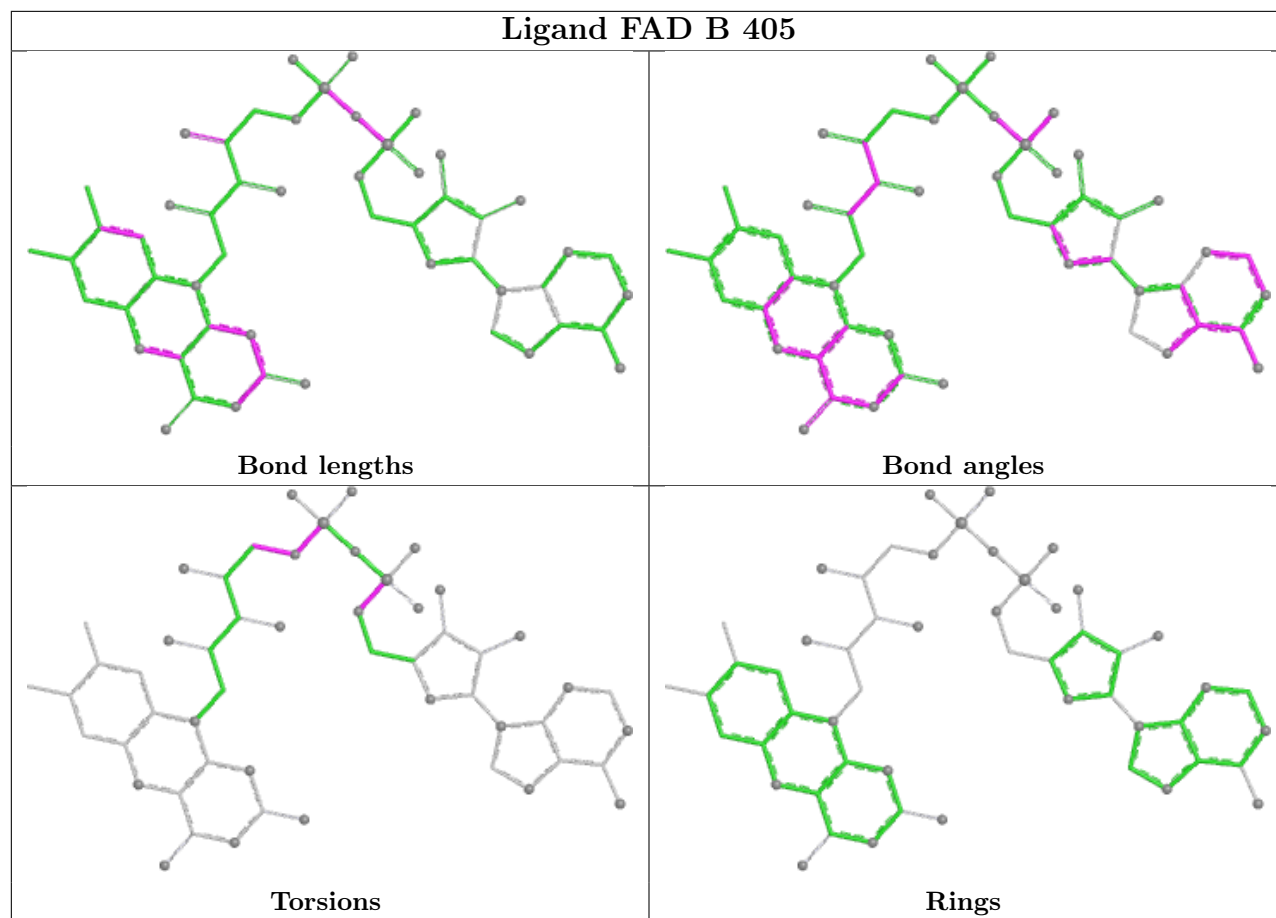
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	965	NAD	4	0
8	B	406	FMN	7	0
6	A	2559	SO4	3	0
7	B	405	FAD	3	0

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	963/964 (99%)	-0.42	13 (1%) 74 75	11, 25, 47, 67	0
2	B	397/404 (98%)	-0.41	6 (1%) 71 72	14, 28, 42, 62	1 (0%)
3	C	195/203 (96%)	-0.41	3 (1%) 71 72	14, 24, 44, 56	0
4	D	91/99 (91%)	-0.58	1 (1%) 77 78	14, 23, 43, 55	0
All	All	1646/1670 (98%)	-0.42	23 (1%) 73 74	11, 26, 46, 67	1 (0%)

The worst 5 of 23 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	200	ALA	5.0
2	B	397	HIS	4.7
1	A	327	THR	3.3
1	A	909	HIS	3.3
1	A	694	GLY	3.3

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 monosaccharides in this entry.

6.4 Ligands [i](#)

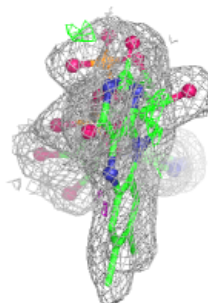
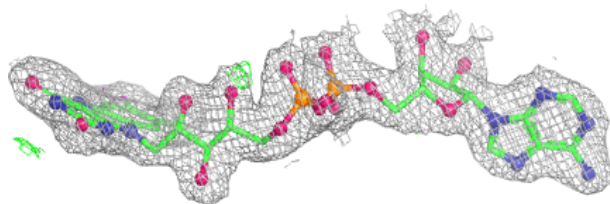
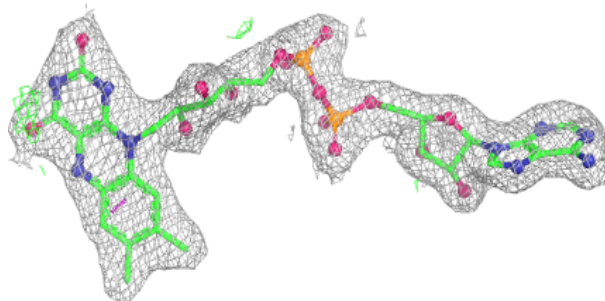
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	SO4	B	2562	5/5	0.34	0.14	156,156,156,156	0
6	SO4	A	2565	5/5	0.74	0.16	161,161,161,162	0
6	SO4	B	2561	5/5	0.86	0.16	59,59,62,63	0
6	SO4	A	2563	5/5	0.88	0.11	76,76,77,77	0
6	SO4	A	2559	5/5	0.90	0.12	80,80,80,82	0
6	SO4	D	2566	5/5	0.90	0.08	73,73,75,75	0
6	SO4	A	2560	5/5	0.94	0.11	50,51,55,55	0
6	SO4	C	2564	5/5	0.97	0.14	33,33,37,40	0
7	FAD	B	405	53/53	0.98	0.05	15,25,31,33	0
8	FMN	B	406	31/31	0.98	0.04	9,15,20,22	0
5	NAD	A	965	44/44	0.99	0.04	5,14,18,23	0
6	SO4	A	2567	5/5	0.99	0.05	34,35,36,39	0
9	ZN	D	100	1/1	1.00	0.01	21,21,21,21	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.

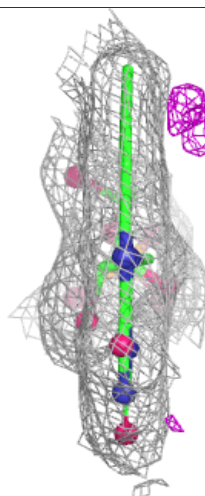
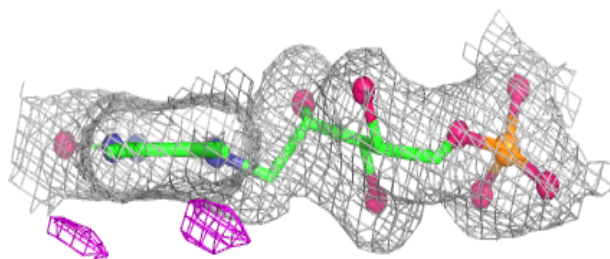
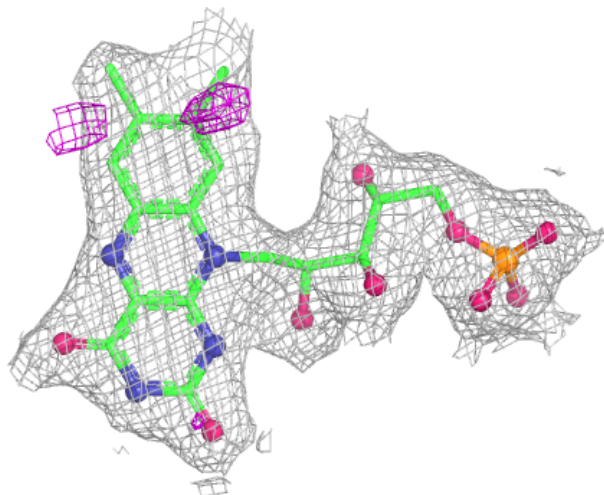
Electron density around FAD B 405:

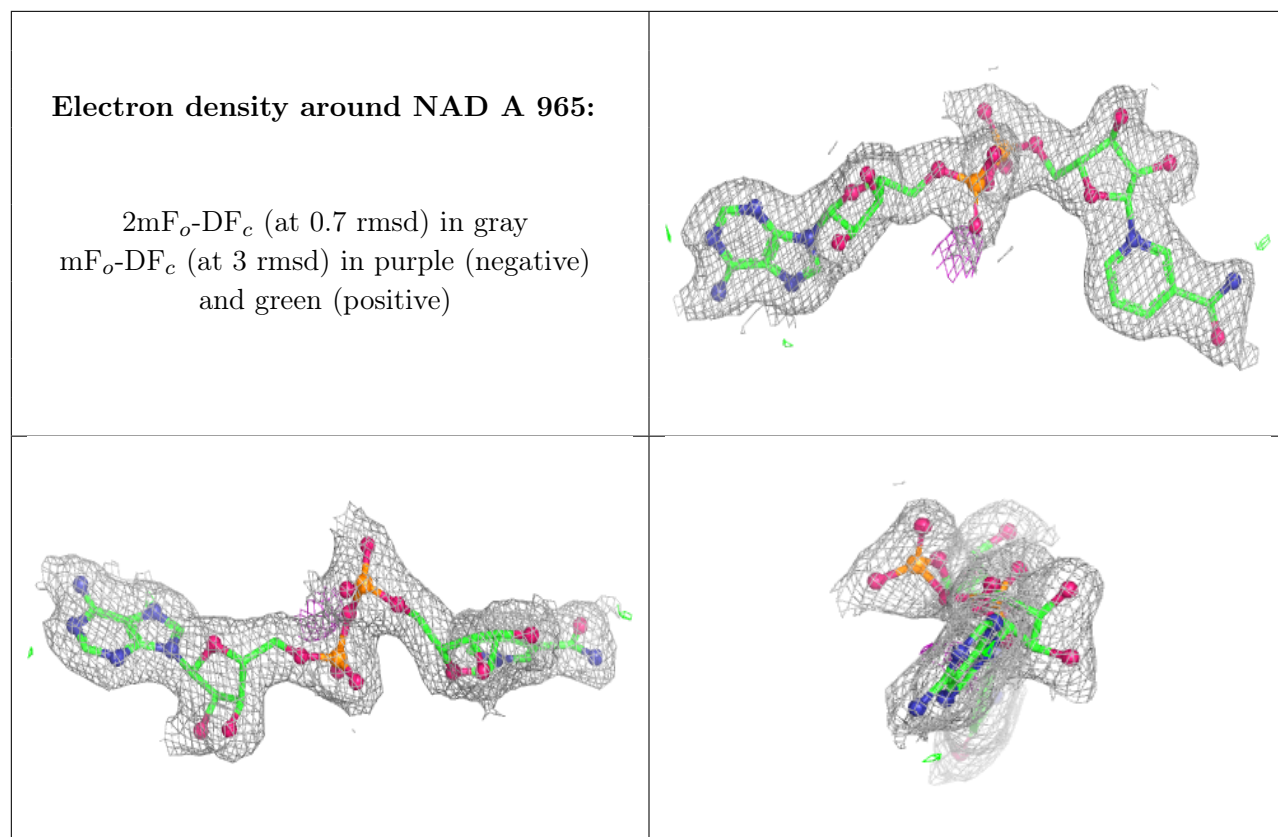
2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FMN B 406:

$2mF_o-DF_c$ (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.