



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 25, 2023 – 04:00 PM EDT

PDB ID : 3AD8
Title : Heterotetrameric Sarcosine Oxidase from *Corynebacterium* sp. U-96 in complex with pyrrole 2-carboxylate
Authors : Suzuki, H.; Moriguchi, T.; Ida, K.
Deposited on : 2010-01-15
Resolution : 2.20 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

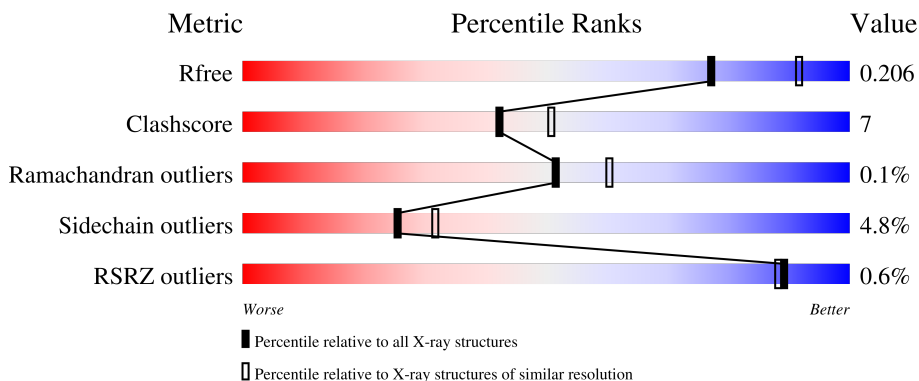
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="flex-grow: 1; position: relative;"> <div style="position: absolute; top: -10px; left: 0; right: 0; text-align: center;">%</div> <div style="background-color: green; width: 87%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: yellow; width: 11%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: orange; width: 2%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: red; width: 2%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: grey; width: 0%; height: 15px; margin-bottom: 2px;"></div> </div> <div style="margin-left: 10px;"> <p>87% 11% .</p> </div> </div>
2	B	404	<div style="display: flex; align-items: center;"> <div style="width: 10px; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="flex-grow: 1; position: relative;"> <div style="position: absolute; top: -10px; left: 0; right: 0; text-align: center;">%</div> <div style="background-color: green; width: 85%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: yellow; width: 12%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: orange; width: 3%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: red; width: 0%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: grey; width: 0%; height: 15px; margin-bottom: 2px;"></div> </div> <div style="margin-left: 10px;"> <p>85% 12% .</p> </div> </div>
3	C	203	<div style="display: flex; align-items: center;"> <div style="flex-grow: 1; position: relative;"> <div style="position: absolute; top: -10px; left: 0; right: 0; text-align: center;">%</div> <div style="background-color: green; width: 80%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: yellow; width: 13%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: orange; width: 7%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: red; width: 0%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: grey; width: 0%; height: 15px; margin-bottom: 2px;"></div> </div> <div style="margin-left: 10px;"> <p>80% 13% . .</p> </div> </div>
4	D	99	<div style="display: flex; align-items: center;"> <div style="flex-grow: 1; position: relative;"> <div style="position: absolute; top: -10px; left: 0; right: 0; text-align: center;">%</div> <div style="background-color: green; width: 87%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: yellow; width: 5%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: orange; width: 8%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: red; width: 0%; height: 15px; margin-bottom: 2px;"></div> <div style="background-color: grey; width: 0%; height: 15px; margin-bottom: 2px;"></div> </div> <div style="margin-left: 10px;"> <p>87% 5% 8%</p> </div> </div>

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	2509	-	-	X	-
8	FMN	B	406	-	-	X	-

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 13736 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
			Total	C	N	O	S			
1	A	963	7229	4507	1287	1413	22	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	404	3108	1981	541	576	10	0	0	0

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

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

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

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

- Molecule 5 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



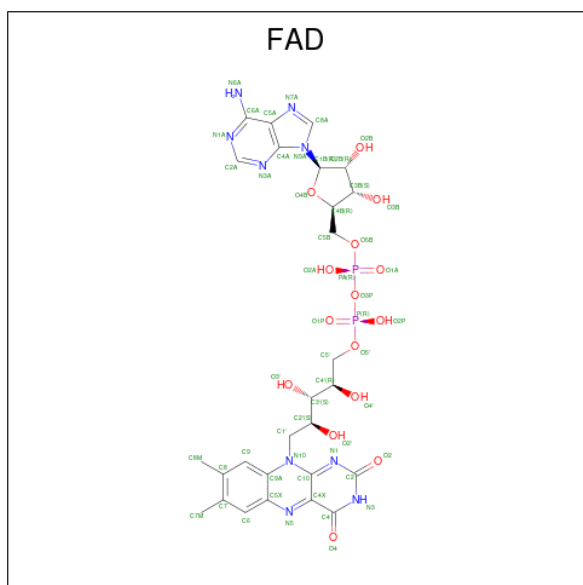
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
6	A	1	5	4	1	0	0
6	A	1	5	4	1	0	0
6	A	1	5	4	1	0	0

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

- Molecule 7 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



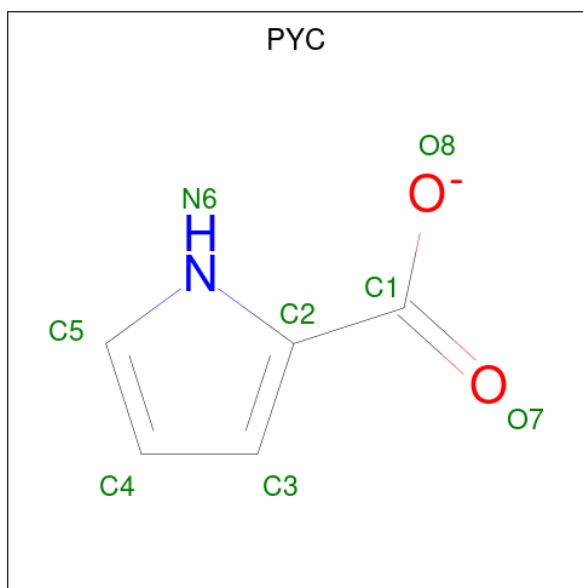
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
7	B	1	53	27	9	15	2	0	0

- Molecule 8 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



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

- Molecule 9 is PYRROLE-2-CARBOXYLATE (three-letter code: PYC) (formula: $C_5H_4NO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
9	B	1	8	5	1	2	0	0

- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

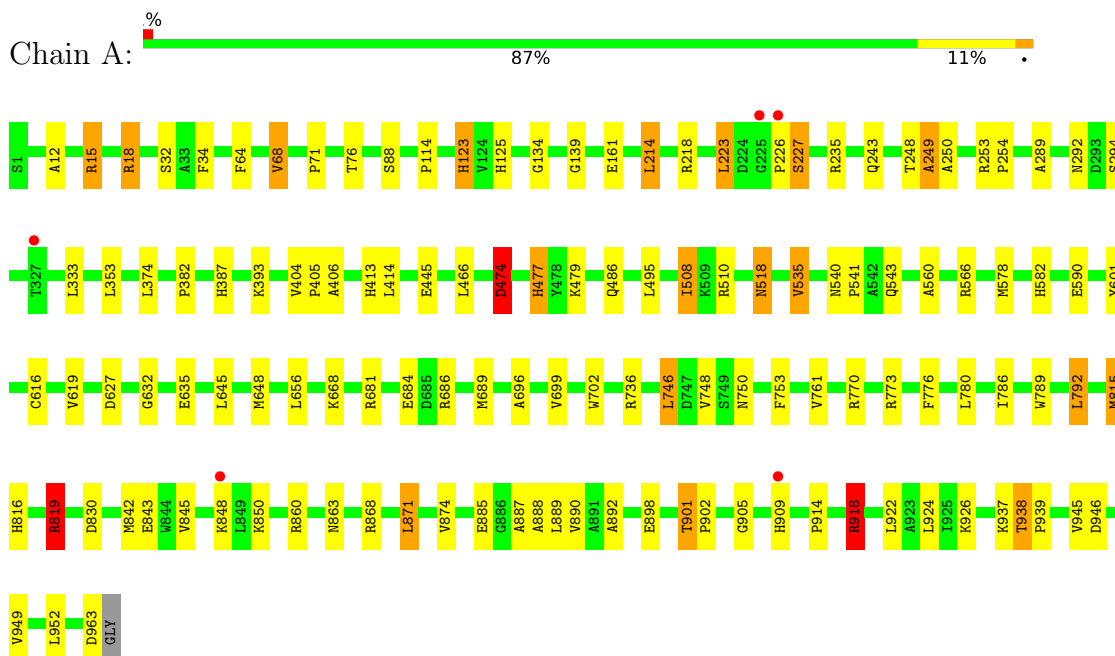
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	631	Total 631	O 631	0	0
11	B	189	Total 189	O 189	0	0
11	C	133	Total 133	O 133	0	0
11	D	67	Total 67	O 67	0	0

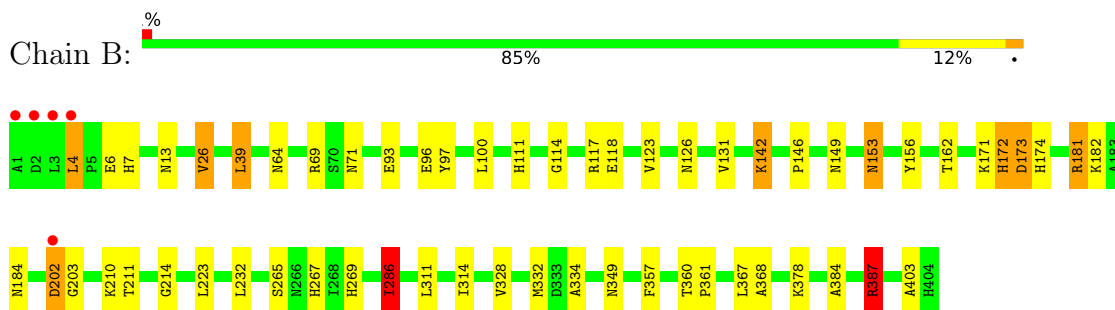
3 Residue-property plots [i](#)

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

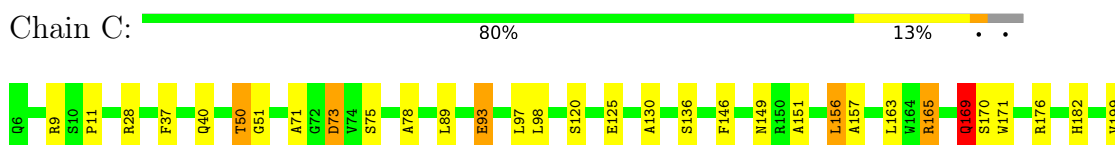
- Molecule 1: SARCOSINE OXIDASE ALPHA SUBUNIT



- Molecule 2: SARCOSINE OXIDASE BETA SUBUNIT




- Molecule 3: SARCOSINE OXIDASE GAMMA SUBUNIT



A200
LEU
GLU
HIS
HIS
HIS
HIS
HIS
HIS

● Molecule 4: SARCOSINE OXIDASE DELTA SUBUNIT

Chain D:  87% 5% 8%

H1
I4
R12
F63
E77
F78
K79
L91
ASP
SER
THR
GLU
GLY
GLY
THR
ARG

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	198.80Å 198.80Å 196.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.31 – 2.20 61.31 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (61.31-2.20) 100.0 (61.31-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.18 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.161 , 0.206 0.161 , 0.206	Depositor DCC
R_{free} test set	5799 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtrriage
Anisotropy	0.054	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13736	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.27% 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: FAD, NAD, FMN, SO4, PYC, 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	1.04	5/7361 (0.1%)	0.93	19/10017 (0.2%)
2	B	1.01	2/3189 (0.1%)	0.91	6/4340 (0.1%)
3	C	1.11	3/1461 (0.2%)	1.00	6/1998 (0.3%)
4	D	1.03	1/772 (0.1%)	0.88	1/1040 (0.1%)
All	All	1.04	11/12783 (0.1%)	0.93	32/17395 (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	1
3	C	0	1
All	All	0	2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	477	HIS	CB-CG	9.30	1.66	1.50
1	A	477	HIS	CA-CB	8.05	1.71	1.53
1	A	477	HIS	CA-C	6.84	1.70	1.52
4	D	53	PHE	CE1-CZ	6.38	1.49	1.37
1	A	249	ALA	CA-CB	-5.99	1.39	1.52
3	C	130	ALA	CA-CB	5.88	1.64	1.52
3	C	125	GLU	CB-CG	-5.67	1.41	1.52
1	A	445	GLU	CG-CD	5.64	1.60	1.51
3	C	169	GLN	CB-CG	5.41	1.67	1.52
2	B	387	ARG	CG-CD	5.31	1.65	1.51
2	B	334	ALA	CA-CB	5.17	1.63	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	387	ARG	NE-CZ-NH1	-14.34	113.13	120.30
1	A	15	ARG	NE-CZ-NH2	-12.63	113.98	120.30
1	A	819	ARG	NE-CZ-NH2	-12.61	113.99	120.30
1	A	819	ARG	NE-CZ-NH1	10.36	125.48	120.30
3	C	176	ARG	NE-CZ-NH2	-8.83	115.88	120.30
2	B	387	ARG	NE-CZ-NH2	8.37	124.49	120.30
1	A	477	HIS	CB-CA-C	8.22	126.83	110.40
1	A	918	ARG	NE-CZ-NH1	8.10	124.35	120.30
3	C	176	ARG	NE-CZ-NH1	8.08	124.34	120.30
3	C	165	ARG	NE-CZ-NH1	8.05	124.33	120.30
1	A	770	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	A	15	ARG	NE-CZ-NH1	6.93	123.77	120.30
2	B	117	ARG	NE-CZ-NH2	-6.77	116.91	120.30
1	A	214	LEU	CA-CB-CG	-6.67	99.95	115.30
3	C	28	ARG	NE-CZ-NH2	-6.57	117.02	120.30
1	A	918	ARG	NE-CZ-NH2	-6.45	117.08	120.30
1	A	819	ARG	CG-CD-NE	-6.43	98.30	111.80
3	C	28	ARG	NE-CZ-NH1	6.39	123.49	120.30
1	A	214	LEU	CB-CG-CD1	6.09	121.36	111.00
2	B	181	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	A	819	ARG	CD-NE-CZ	6.01	132.01	123.60
1	A	566	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	A	235	ARG	NE-CZ-NH2	-5.73	117.44	120.30
1	A	12	ALA	C-N-CA	-5.72	110.29	122.30
1	A	474	ASP	CB-CG-OD2	-5.68	113.19	118.30
2	B	286	ILE	CG1-CB-CG2	5.47	123.44	111.40
1	A	18	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	819	ARG	CB-CG-CD	5.28	125.33	111.60
4	D	12	ARG	NE-CZ-NH1	5.11	122.86	120.30
3	C	200	ALA	N-CA-C	5.11	124.80	111.00
2	B	173	ASP	CB-CG-OD1	5.09	122.88	118.30
1	A	68	VAL	CG1-CB-CG2	5.03	118.95	110.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	227	SER	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	101	1
2	B	3108	0	3035	54	0
3	C	1433	0	1434	17	0
4	D	749	0	706	3	0
5	A	44	0	26	4	0
6	A	35	0	0	3	0
6	B	15	0	0	0	0
6	C	5	0	0	0	0
6	D	5	0	0	0	0
7	B	53	0	31	4	0
8	B	31	0	18	9	0
9	B	8	0	4	0	0
10	D	1	0	0	0	0
11	A	631	0	0	13	1
11	B	189	0	0	5	0
11	C	133	0	0	2	0
11	D	67	0	0	0	0
All	All	13736	0	12358	173	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 7.

All (173) 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)
2:B:172:HIS:NE2	8:B:406:FMN:C8M	1.71	1.50
1:A:830:ASP:OD1	1:A:909:HIS:CE1	1.88	1.26
1:A:696:ALA:HB3	6:A:2509:SO4:O3	1.41	1.17
2:B:384:ALA:O	2:B:387:ARG:HD3	1.44	1.12
2:B:387:ARG:HH11	2:B:387:ARG:HG3	0.98	1.11
1:A:830:ASP:CG	1:A:909:HIS:HE1	1.56	1.09
2:B:387:ARG:HH11	2:B:387:ARG:CG	1.66	1.06
1:A:885:GLU:HG2	1:A:909:HIS:HA	1.37	1.05
1:A:696:ALA:HB1	11:A:1461:HOH:O	1.57	1.04
2:B:387:ARG:HG3	2:B:387:ARG:NH1	1.76	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:736:ARG:NH2	1:A:963:ASP:O	1.99	0.95
1:A:696:ALA:CB	6:A:2509:SO4:O3	2.14	0.95
3:C:169:GLN:HE21	3:C:171:TRP:HE1	1.13	0.94
1:A:830:ASP:OD1	1:A:909:HIS:HE1	1.36	0.92
1:A:226:PRO:HB2	11:A:1256:HOH:O	1.70	0.90
3:C:169:GLN:NE2	3:C:171:TRP:HE1	1.72	0.88
1:A:15:ARG:CD	1:A:161:GLU:OE2	2.22	0.87
3:C:37:PHE:O	3:C:93:GLU:HG2	1.75	0.87
2:B:100:LEU:HD12	2:B:171:LYS:HD2	1.57	0.87
1:A:689:MET:HE2	1:A:699:VAL:HG11	1.56	0.87
1:A:15:ARG:HD2	1:A:161:GLU:OE2	1.75	0.85
1:A:868:ARG:O	1:A:926:LYS:O	1.95	0.84
1:A:689:MET:CE	1:A:699:VAL:HG11	2.08	0.83
1:A:819:ARG:HH22	1:A:909:HIS:CE1	1.98	0.81
1:A:830:ASP:CG	1:A:909:HIS:CE1	2.44	0.79
1:A:888:ALA:O	1:A:938:THR:HG23	1.83	0.79
2:B:172:HIS:HD1	2:B:173:ASP:H	1.32	0.78
1:A:249:ALA:HB1	11:A:978:HOH:O	1.83	0.78
1:A:830:ASP:OD2	1:A:909:HIS:HE1	1.66	0.78
1:A:510:ARG:HH21	8:B:406:FMN:H5'1	1.50	0.76
2:B:172:HIS:CD2	8:B:406:FMN:C8M	2.68	0.76
2:B:387:ARG:CG	2:B:387:ARG:NH1	2.37	0.75
1:A:892:ALA:HA	1:A:937:LYS:HD2	1.68	0.75
3:C:50:THR:HG22	11:C:482:HOH:O	1.87	0.74
2:B:269:HIS:CE1	2:B:403:ALA:H	2.06	0.73
1:A:292:ASN:HD22	1:A:294:SER:H	1.34	0.73
1:A:736:ARG:NH1	11:A:1486:HOH:O	2.21	0.72
1:A:816:HIS:O	1:A:819:ARG:HD3	1.89	0.72
2:B:202:ASP:OD1	2:B:203:GLY:N	2.22	0.72
1:A:909:HIS:NE2	1:A:922:LEU:HD12	2.05	0.72
1:A:901:THR:CG2	11:A:1118:HOH:O	2.39	0.71
1:A:15:ARG:HD3	1:A:161:GLU:OE2	1.89	0.70
1:A:510:ARG:HH21	8:B:406:FMN:C5'	2.04	0.70
1:A:918:ARG:HH11	1:A:918:ARG:HG3	1.57	0.70
2:B:153:ASN:H	2:B:153:ASN:HD22	1.38	0.70
1:A:696:ALA:HB3	6:A:2509:SO4:S	2.31	0.69
2:B:93:GLU:HG2	11:B:655:HOH:O	1.92	0.69
3:C:73:ASP:OD2	3:C:75:SER:HB2	1.92	0.69
1:A:292:ASN:ND2	1:A:294:SER:H	1.90	0.68
1:A:860:ARG:HH21	1:A:863:ASN:HD21	1.42	0.67
1:A:819:ARG:NH2	1:A:909:HIS:NE2	2.42	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:HIS:HD2	11:A:1191:HOH:O	1.77	0.66
1:A:736:ARG:HG2	1:A:748:VAL:HG23	1.78	0.65
1:A:937:LYS:NZ	1:A:946:ASP:OD2	2.29	0.65
3:C:37:PHE:O	3:C:93:GLU:CG	2.44	0.65
2:B:210:LYS:HE3	2:B:211:THR:O	1.97	0.65
2:B:149:ASN:HD21	2:B:153:ASN:HD21	1.45	0.64
1:A:830:ASP:OD2	1:A:909:HIS:CE1	2.48	0.64
1:A:243:GLN:OE1	1:A:413:HIS:HE1	1.80	0.64
2:B:269:HIS:HE1	2:B:403:ALA:H	1.46	0.64
1:A:901:THR:HG23	11:A:1118:HOH:O	1.98	0.63
1:A:223:LEU:HD23	1:A:226:PRO:HB3	1.80	0.63
2:B:171:LYS:HE2	2:B:174:HIS:CD2	2.34	0.62
1:A:632:GLY:HA3	1:A:696:ALA:HB2	1.81	0.61
3:C:157:ALA:O	3:C:182:HIS:HE1	1.84	0.60
1:A:750:ASN:HD22	1:A:773:ARG:HH12	1.49	0.60
2:B:142:LYS:HE2	2:B:146:PRO:O	2.02	0.59
1:A:945:VAL:HG13	11:A:1127:HOH:O	2.02	0.59
1:A:816:HIS:HA	1:A:819:ARG:HD2	1.85	0.59
2:B:172:HIS:HD1	2:B:173:ASP:N	1.98	0.58
1:A:842:MET:O	1:A:845:VAL:HG12	2.03	0.58
11:A:1227:HOH:O	3:C:182:HIS:HD2	1.86	0.58
1:A:582:HIS:CD2	1:A:601:TYR:OH	2.57	0.58
1:A:382:PRO:HG2	1:A:404:VAL:HG12	1.86	0.58
2:B:111:HIS:HD2	2:B:156:TYR:O	1.87	0.58
2:B:114:GLY:O	2:B:118:GLU:HG2	2.03	0.58
1:A:635:GLU:OE2	1:A:686:ARG:NH1	2.32	0.57
1:A:474:ASP:OD2	1:A:477:HIS:CD2	2.58	0.56
1:A:540:ASN:O	1:A:543:GLN:HB2	2.05	0.56
2:B:7:HIS:HE1	11:B:426:HOH:O	1.86	0.56
2:B:172:HIS:CE1	8:B:406:FMN:C8M	2.76	0.56
2:B:172:HIS:NE2	8:B:406:FMN:C8	2.63	0.55
1:A:510:ARG:NH2	8:B:406:FMN:C5'	2.70	0.55
1:A:249:ALA:HB2	5:A:965:NAD:O3	2.07	0.55
1:A:888:ALA:O	1:A:938:THR:CG2	2.51	0.55
1:A:750:ASN:ND2	1:A:773:ARG:HH12	2.06	0.54
2:B:64:ASN:HA	7:B:405:FAD:C6	2.37	0.54
1:A:746:LEU:HD13	1:A:748:VAL:HG12	1.89	0.54
2:B:39:LEU:HD13	2:B:368:ALA:HA	1.91	0.53
1:A:510:ARG:NH2	8:B:406:FMN:H5'1	2.22	0.53
2:B:286:ILE:HD11	2:B:328:VAL:HG21	1.90	0.53
2:B:360:THR:HB	2:B:361:PRO:HD3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ALA:HB2	5:A:965:NAD:PA	2.48	0.52
2:B:123:VAL:HG23	2:B:162:THR:HG22	1.92	0.52
1:A:945:VAL:CG1	11:A:1127:HOH:O	2.56	0.52
1:A:736:ARG:HG2	1:A:748:VAL:CG2	2.40	0.52
2:B:4:LEU:HD22	2:B:96:GLU:HA	1.91	0.52
3:C:11:PRO:HG2	3:C:120:SER:HB3	1.91	0.52
1:A:249:ALA:CB	5:A:965:NAD:O2A	2.58	0.52
2:B:118:GLU:OE2	11:B:802:HOH:O	2.19	0.51
2:B:26:VAL:HG13	2:B:223:LEU:HD23	1.92	0.51
1:A:114:PRO:HA	2:B:314:ILE:HG13	1.91	0.51
1:A:495:LEU:HD21	1:A:535:VAL:HG22	1.92	0.51
1:A:508:ILE:HD13	1:A:508:ILE:N	2.26	0.51
2:B:126:ASN:HB3	2:B:131:VAL:HG22	1.93	0.51
1:A:248:THR:O	1:A:249:ALA:HB3	2.12	0.50
1:A:486:GLN:HE22	1:A:518:ASN:ND2	2.10	0.49
1:A:871:LEU:HD13	1:A:952:LEU:HD11	1.93	0.49
4:D:77:GLU:CD	4:D:79:LYS:NZ	2.66	0.49
1:A:582:HIS:HE1	1:A:627:ASP:OD2	1.96	0.48
1:A:249:ALA:HB2	5:A:965:NAD:O2A	2.14	0.48
2:B:13:ASN:ND2	2:B:184:ASN:HD21	2.13	0.47
2:B:265:SER:O	2:B:269:HIS:HD2	1.98	0.47
1:A:689:MET:HE1	1:A:699:VAL:HG11	1.90	0.47
1:A:938:THR:HG22	1:A:939:PRO:HD2	1.97	0.47
1:A:736:ARG:HB2	1:A:780:LEU:HD21	1.98	0.46
1:A:64:PHE:HB3	1:A:71:PRO:HD2	1.97	0.46
3:C:169:GLN:NE2	3:C:171:TRP:NE1	2.53	0.46
1:A:901:THR:HG22	11:A:1118:HOH:O	2.10	0.46
2:B:111:HIS:CE1	2:B:265:SER:OG	2.69	0.46
3:C:40:GLN:HA	3:C:89:LEU:O	2.14	0.46
1:A:689:MET:HE2	1:A:699:VAL:CG1	2.37	0.46
2:B:265:SER:O	2:B:269:HIS:HA	2.16	0.46
1:A:510:ARG:NH2	8:B:406:FMN:H5'2	2.31	0.46
2:B:349:ASN:HB2	2:B:367:LEU:HD22	1.98	0.46
1:A:253:ARG:HB2	1:A:254:PRO:HD2	1.98	0.45
1:A:289:ALA:HB2	1:A:374:LEU:HD11	1.97	0.45
1:A:405:PRO:HB2	1:A:414:LEU:HD13	1.98	0.45
1:A:905:GLY:HA3	1:A:924:LEU:O	2.16	0.45
3:C:136:SER:OG	3:C:156:LEU:HD22	2.17	0.45
1:A:898:GLU:HG3	11:A:1298:HOH:O	2.16	0.45
2:B:69:ARG:HD2	2:B:71:ASN:OD1	2.14	0.45
1:A:815:MET:HE2	1:A:816:HIS:HD2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:118:GLU:OE1	11:B:809:HOH:O	2.20	0.45
1:A:901:THR:HA	1:A:902:PRO:HD3	1.77	0.45
1:A:387:HIS:HD2	11:A:968:HOH:O	1.98	0.45
1:A:486:GLN:HE22	1:A:518:ASN:HD22	1.64	0.45
2:B:96:GLU:O	2:B:181:ARG:NH2	2.50	0.45
4:D:77:GLU:OE2	4:D:79:LYS:NZ	2.49	0.44
1:A:123:HIS:HE1	1:A:560:ALA:O	2.01	0.44
2:B:210:LYS:NZ	2:B:214:GLY:HA2	2.33	0.44
3:C:51:GLY:HA3	11:C:223:HOH:O	2.17	0.44
1:A:668:LYS:HD3	1:A:702:TRP:CH2	2.53	0.44
1:A:819:ARG:NH2	1:A:830:ASP:OD2	2.51	0.44
1:A:134:GLY:O	1:A:139:GLY:HA3	2.18	0.44
1:A:860:ARG:NH2	1:A:863:ASN:HD21	2.13	0.44
2:B:153:ASN:HD22	2:B:153:ASN:N	2.06	0.43
1:A:387:HIS:HE1	1:A:393:LYS:O	2.01	0.43
1:A:479:LYS:HB3	1:A:479:LYS:HE2	1.87	0.43
4:D:77:GLU:CD	4:D:79:LYS:HZ2	2.22	0.43
1:A:249:ALA:HB1	1:A:250:ALA:H	1.66	0.43
1:A:648:MET:HA	1:A:648:MET:HE2	2.01	0.43
1:A:689:MET:CE	1:A:699:VAL:CG1	2.88	0.43
2:B:149:ASN:HD21	2:B:153:ASN:ND2	2.13	0.43
1:A:616:CYS:HB2	1:A:914:PRO:HG2	2.01	0.43
2:B:39:LEU:HD13	2:B:368:ALA:CA	2.49	0.42
1:A:776:PHE:CE2	1:A:819:ARG:HG3	2.53	0.42
2:B:111:HIS:HE1	2:B:265:SER:OG	2.02	0.42
2:B:182:LYS:HA	2:B:182:LYS:HD3	1.80	0.42
3:C:71:ALA:O	3:C:78:ALA:HA	2.20	0.42
3:C:151:ALA:HA	3:C:163:LEU:O	2.20	0.42
2:B:6:GLU:HB2	11:B:790:HOH:O	2.20	0.41
2:B:210:LYS:HD2	2:B:211:THR:H	1.85	0.41
2:B:357:PHE:HB3	7:B:405:FAD:C2	2.50	0.41
2:B:64:ASN:HA	7:B:405:FAD:C5X	2.51	0.41
1:A:18:ARG:HA	1:A:34:PHE:CD1	2.55	0.41
1:A:792:LEU:HD12	3:C:9:ARG:HD3	2.03	0.41
2:B:64:ASN:HB2	7:B:405:FAD:C4X	2.50	0.41
1:A:887:ALA:HB1	1:A:938:THR:HG21	2.03	0.41
1:A:76:THR:HA	1:A:88:SER:HA	2.02	0.40
1:A:656:LEU:HD23	1:A:681:ARG:HB2	2.04	0.40
2:B:332:MET:CE	2:B:332:MET:HA	2.52	0.40
3:C:149:ASN:N	3:C:165:ARG:O	2.50	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:HIS:NE2	1:A:477:HIS:NE2[11_555]	1.27	0.93
11:A:1080:HOH:O	11:A:1080:HOH:O[9_555]	1.63	0.57

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	961/964 (100%)	941 (98%)	18 (2%)	2 (0%)	47	55
2	B	402/404 (100%)	387 (96%)	15 (4%)	0	100	100
3	C	193/203 (95%)	187 (97%)	6 (3%)	0	100	100
4	D	89/99 (90%)	87 (98%)	2 (2%)	0	100	100
All	All	1645/1670 (98%)	1602 (97%)	41 (2%)	2 (0%)	51	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	227	SER
1	A	406	ALA

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%)	709 (95%)	38 (5%)	24	29

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	319/319 (100%)	305 (96%)	14 (4%)	28	35
3	C	143/151 (95%)	134 (94%)	9 (6%)	18	20
4	D	75/81 (93%)	74 (99%)	1 (1%)	69	81
All	All	1284/1298 (99%)	1222 (95%)	62 (5%)	25	32

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	SER
1	A	68	VAL
1	A	123	HIS
1	A	214	LEU
1	A	218	ARG
1	A	223	LEU
1	A	333	LEU
1	A	353	LEU
1	A	466	LEU
1	A	474	ASP
1	A	508	ILE
1	A	518	ASN
1	A	535	VAL
1	A	541	PRO
1	A	578	MET
1	A	590	GLU
1	A	619	VAL
1	A	645	LEU
1	A	684	GLU
1	A	746	LEU
1	A	753	PHE
1	A	761	VAL
1	A	786	ILE
1	A	789	TRP
1	A	792	LEU
1	A	815	MET
1	A	819	ARG
1	A	843	GLU
1	A	848	LYS
1	A	850	LYS
1	A	871	LEU
1	A	874	VAL
1	A	889	LEU

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Mol	Chain	Res	Type
1	A	890	VAL
1	A	901	THR
1	A	918	ARG
1	A	938	THR
1	A	949	VAL
2	B	4	LEU
2	B	26	VAL
2	B	39	LEU
2	B	97	TYR
2	B	142	LYS
2	B	153	ASN
2	B	172	HIS
2	B	202	ASP
2	B	232	LEU
2	B	267	HIS
2	B	286	ILE
2	B	311	LEU
2	B	378	LYS
2	B	387	ARG
3	C	50	THR
3	C	73	ASP
3	C	93	GLU
3	C	97	LEU
3	C	98	LEU
3	C	146	PHE
3	C	156	LEU
3	C	169	GLN
3	C	170	SER
4	D	4	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	82	GLN
1	A	123	HIS
1	A	125	HIS
1	A	181	GLN
1	A	238	HIS
1	A	292	ASN
1	A	387	HIS
1	A	413	HIS

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Mol	Chain	Res	Type
1	A	477	HIS
1	A	518	ASN
1	A	540	ASN
1	A	582	HIS
1	A	750	ASN
1	A	816	HIS
1	A	863	ASN
1	A	912	ASN
2	B	7	HIS
2	B	13	ASN
2	B	14	ASN
2	B	111	HIS
2	B	153	ASN
2	B	267	HIS
2	B	269	HIS
2	B	344	GLN
2	B	369	HIS
3	C	6	GLN
3	C	158	ASN
3	C	169	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 1 is monoatomic - leaving 16 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
5	NAD	A	965	-	42,48,48	1.65	5 (11%)	50,73,73	1.60	7 (14%)
7	FAD	B	405	-	53,58,58	1.55	6 (11%)	68,89,89	1.39	12 (17%)
6	SO4	A	2502	-	4,4,4	0.23	0	6,6,6	0.29	0
6	SO4	A	2504	-	4,4,4	0.18	0	6,6,6	0.34	0
6	SO4	A	2507	-	4,4,4	0.18	0	6,6,6	0.46	0
6	SO4	A	2509	-	4,4,4	0.36	0	6,6,6	0.45	0
6	SO4	A	2500	-	4,4,4	0.15	0	6,6,6	0.27	0
6	SO4	D	2503	-	4,4,4	0.16	0	6,6,6	0.19	0
8	FMN	B	406	-	33,33,33	2.07	9 (27%)	48,50,50	2.35	11 (22%)
6	SO4	A	2501	-	4,4,4	0.29	0	6,6,6	0.59	0
6	SO4	C	2508	-	4,4,4	0.60	0	6,6,6	0.28	0
9	PYC	B	801	-	8,8,8	1.39	0	8,10,10	1.44	1 (12%)
6	SO4	B	2510	-	4,4,4	0.44	0	6,6,6	0.65	0
6	SO4	B	2511	-	4,4,4	0.18	0	6,6,6	0.45	0
6	SO4	A	2506	-	4,4,4	0.26	0	6,6,6	0.31	0
6	SO4	B	2505	-	4,4,4	0.16	0	6,6,6	0.25	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
5	NAD	A	965	-	-	5/26/62/62	0/5/5/5
9	PYC	B	801	-	-	0/2/4/4	0/1/1/1
8	FMN	B	406	-	-	6/18/18/18	0/3/3/3
7	FAD	B	405	-	-	3/30/50/50	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	405	FAD	C4X-N5	7.15	1.44	1.30
5	A	965	NAD	O7N-C7N	6.35	1.36	1.24
8	B	406	FMN	O5'-C5'	-6.24	1.20	1.44
5	A	965	NAD	C2N-N1N	4.82	1.40	1.35
8	B	406	FMN	C4A-N5	3.91	1.38	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	965	NAD	C2B-C1B	-3.29	1.48	1.53
7	B	405	FAD	C10-N1	3.15	1.39	1.33
8	B	406	FMN	C8M-C8	3.01	1.57	1.51
8	B	406	FMN	C9A-C5A	-2.94	1.36	1.41
8	B	406	FMN	C10-N1	2.90	1.39	1.33
8	B	406	FMN	C1'-C2'	2.89	1.56	1.52
7	B	405	FAD	C9-C8	2.79	1.43	1.39
7	B	405	FAD	C10-N10	2.48	1.42	1.37
5	A	965	NAD	C5B-C4B	2.46	1.59	1.51
8	B	406	FMN	C4'-C3'	2.40	1.58	1.53
5	A	965	NAD	C5A-N7A	-2.38	1.31	1.39
8	B	406	FMN	C4A-C10	-2.38	1.37	1.44
8	B	406	FMN	C7M-C7	2.21	1.55	1.51
7	B	405	FAD	O3B-C3B	2.11	1.47	1.43
7	B	405	FAD	C6-C5X	2.04	1.43	1.40

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	406	FMN	C5'-C4'-C3'	-9.29	94.25	112.20
8	B	406	FMN	O4'-C4'-C5'	-6.14	96.13	109.92
5	A	965	NAD	O4B-C1B-C2B	-5.69	98.61	106.93
8	B	406	FMN	O5'-P-O1P	5.21	121.09	106.47
5	A	965	NAD	N3A-C2A-N1A	-4.11	122.25	128.68
8	B	406	FMN	C9A-C5A-N5	-3.65	118.47	122.43
7	B	405	FAD	C4-N3-C2	-3.54	119.10	125.64
5	A	965	NAD	C5B-C4B-C3B	-3.47	102.17	115.18
8	B	406	FMN	C4'-C3'-C2'	3.47	120.58	113.36
7	B	405	FAD	C9A-C5X-N5	-3.39	118.75	122.43
7	B	405	FAD	N3A-C2A-N1A	-3.34	123.45	128.68
8	B	406	FMN	O5'-C5'-C4'	3.10	117.63	109.36
7	B	405	FAD	O4B-C1B-C2B	-3.08	102.42	106.93
5	A	965	NAD	C4A-C5A-N7A	-3.05	106.22	109.40
7	B	405	FAD	C4X-C4-N3	2.89	120.53	113.19
8	B	406	FMN	O2P-P-O5'	-2.74	99.45	106.73
8	B	406	FMN	O4-C4-C4A	-2.73	119.36	126.60
5	A	965	NAD	O4D-C1D-C2D	-2.70	102.98	106.93
7	B	405	FAD	C10-C4X-N5	-2.68	119.17	124.86
5	A	965	NAD	C2N-C3N-C4N	2.45	121.03	118.26
5	A	965	NAD	C6N-N1N-C2N	-2.43	119.76	121.97
7	B	405	FAD	C6-C5X-C9A	2.43	122.37	118.94
7	B	405	FAD	C5X-N5-C4X	2.34	121.96	118.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	406	FMN	C5A-C9A-N10	2.29	120.32	117.95
9	B	801	PYC	O7-C1-C2	-2.28	116.61	121.24
8	B	406	FMN	P-O5'-C5'	2.28	124.58	118.30
7	B	405	FAD	C4A-C5A-N7A	-2.22	107.09	109.40
7	B	405	FAD	C4-C4X-N5	2.21	121.38	118.23
7	B	405	FAD	O4-C4-C4X	-2.16	120.88	126.60
8	B	406	FMN	C4-N3-C2	-2.14	121.69	125.64
7	B	405	FAD	C5'-C4'-C3'	-2.07	108.21	112.20

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	406	FMN	C1'-C2'-C3'-C4'
8	B	406	FMN	C2'-C3'-C4'-C5'
8	B	406	FMN	C3'-C4'-C5'-O5'
8	B	406	FMN	O4'-C4'-C5'-O5'
8	B	406	FMN	O3'-C3'-C4'-C5'
8	B	406	FMN	O2'-C2'-C3'-C4'
7	B	405	FAD	C4'-C5'-O5'-P
5	A	965	NAD	PA-O3-PN-O5D
5	A	965	NAD	O4D-C4D-C5D-O5D
5	A	965	NAD	O4B-C4B-C5B-O5B
5	A	965	NAD	C5D-O5D-PN-O1N
7	B	405	FAD	C5'-O5'-P-O1P
5	A	965	NAD	C3D-C4D-C5D-O5D
7	B	405	FAD	O4B-C4B-C5B-O5B

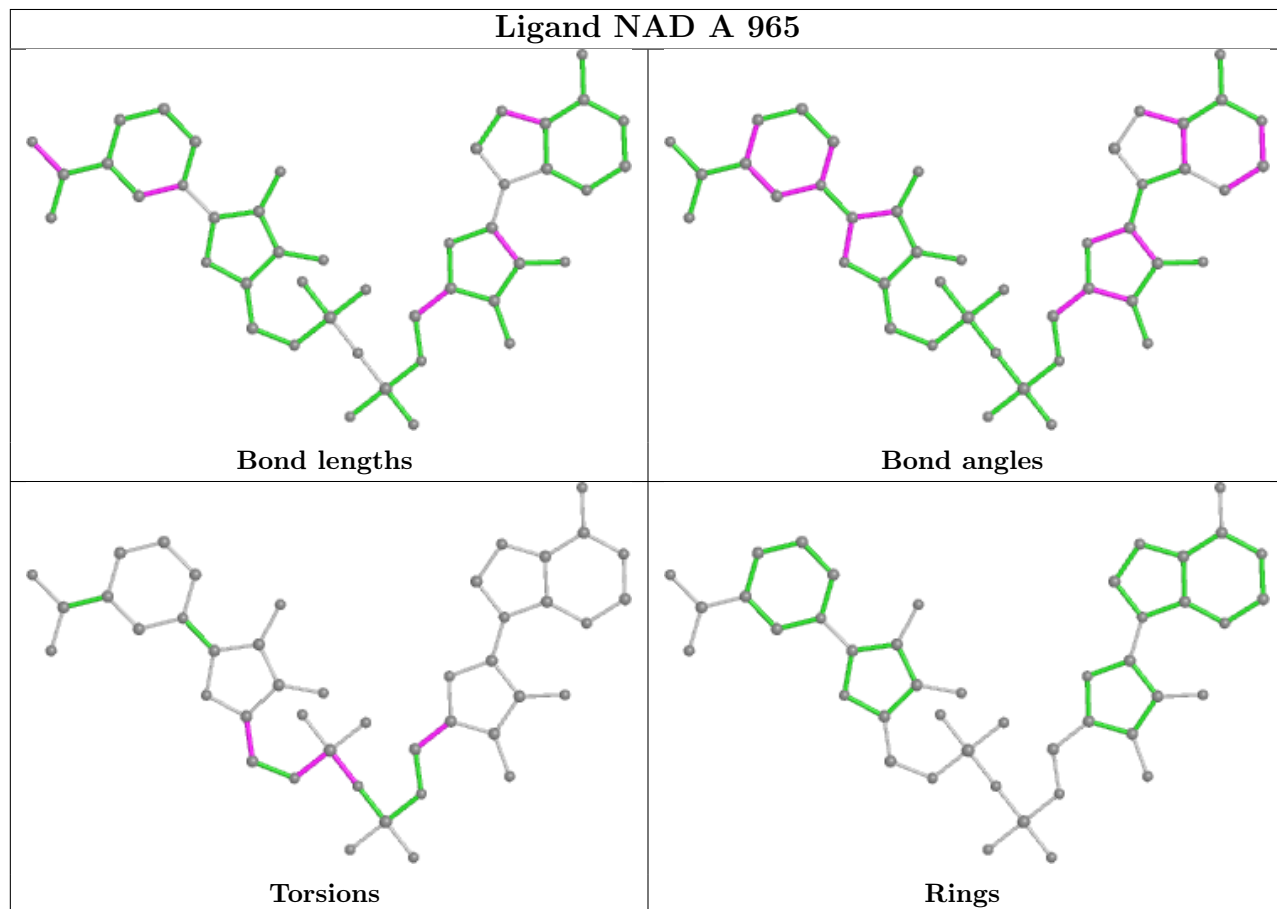
There are no ring outliers.

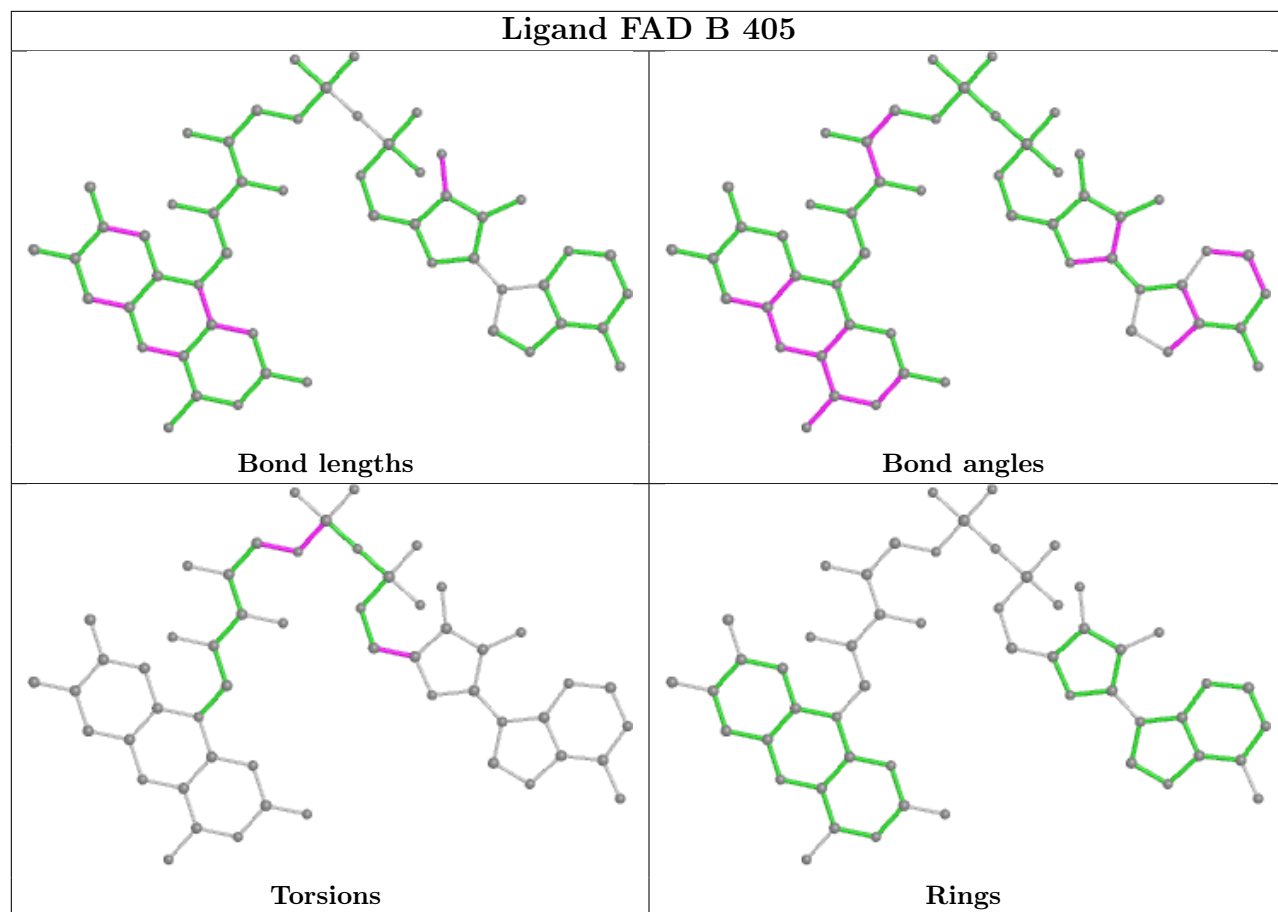
4 monomers are involved in 20 short contacts:

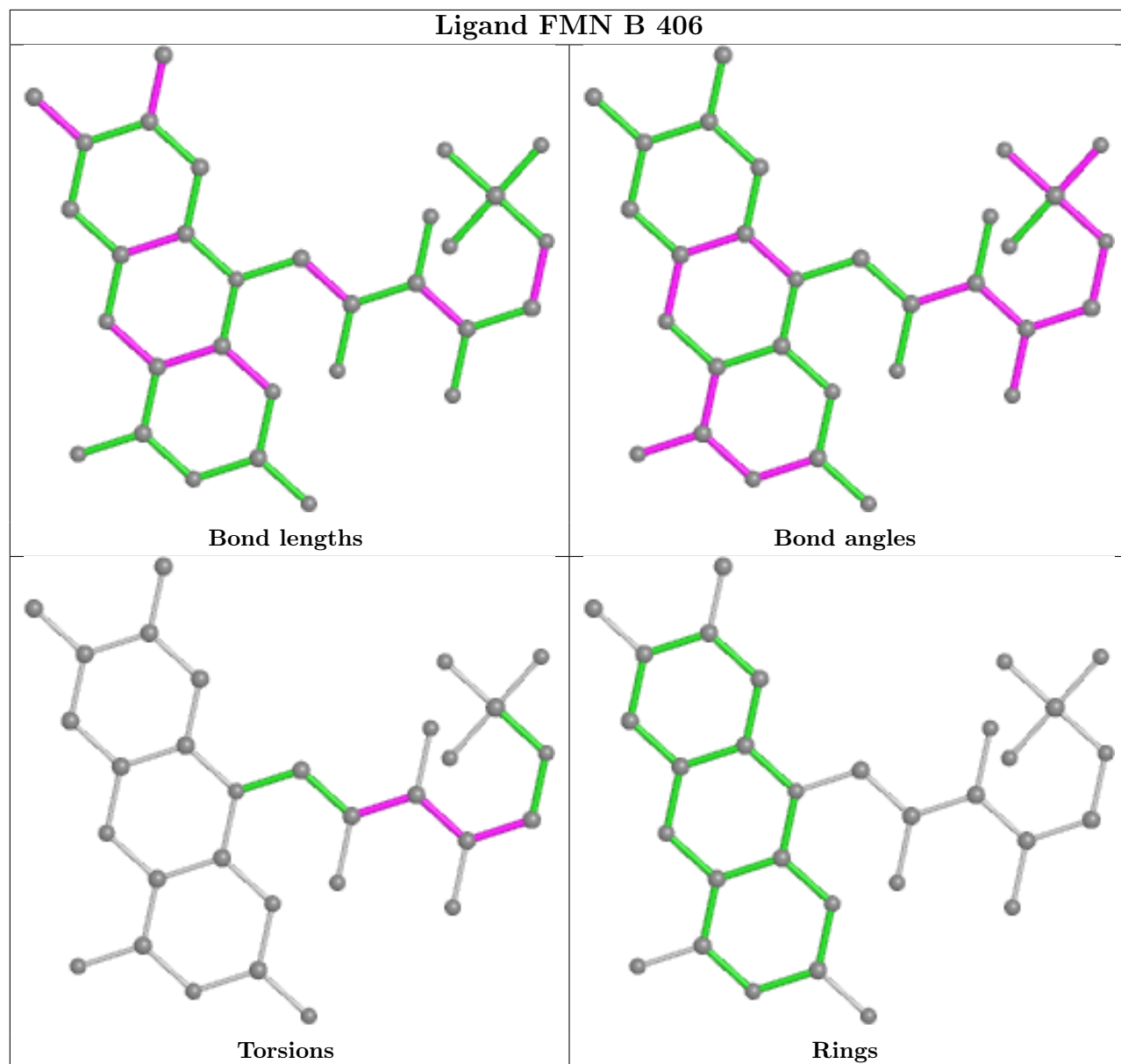
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	965	NAD	4	0
7	B	405	FAD	4	0
6	A	2509	SO4	3	0
8	B	406	FMN	9	0

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%)	-0.62	5 (0%) 91 90	11, 24, 41, 59	0
2	B	404/404 (100%)	-0.55	5 (1%) 79 77	13, 25, 38, 84	0
3	C	195/203 (96%)	-0.74	0 100 100	14, 22, 42, 49	0
4	D	91/99 (91%)	-0.80	0 100 100	15, 22, 42, 54	0
All	All	1653/1670 (98%)	-0.63	10 (0%) 89 88	11, 24, 41, 84	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	2	ASP	8.4
2	B	1	ALA	7.5
2	B	3	LEU	7.3
1	A	909	HIS	3.8
1	A	226	PRO	2.8
2	B	202	ASP	2.7
2	B	4	LEU	2.4
1	A	327	THR	2.3
1	A	848	LYS	2.1
1	A	225	GLY	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 monosaccharides in this entry.

6.4 Ligands

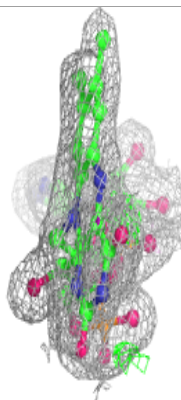
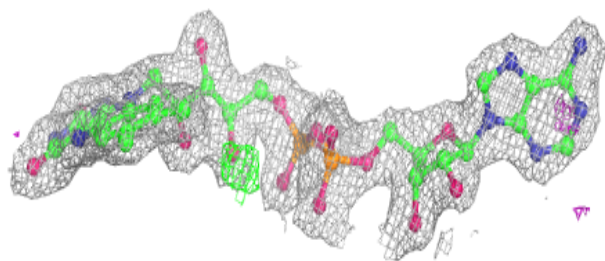
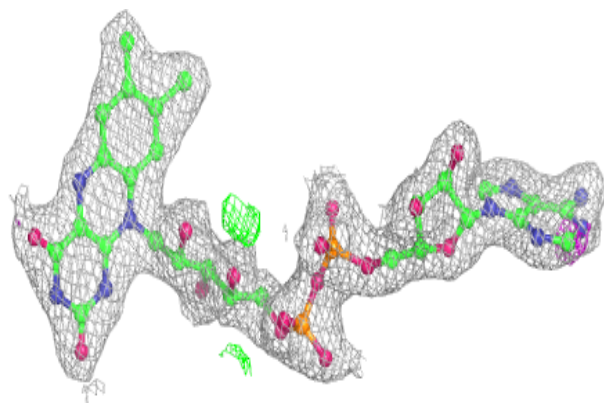
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	SO4	B	2511	5/5	0.90	0.11	73,73,75,76	0
6	SO4	A	2506	5/5	0.93	0.27	65,65,67,68	0
6	SO4	B	2510	5/5	0.94	0.23	52,53,57,57	0
6	SO4	D	2503	5/5	0.94	0.13	82,82,83,83	0
6	SO4	A	2502	5/5	0.95	0.29	76,76,77,78	0
6	SO4	A	2507	5/5	0.96	0.20	59,60,62,64	0
9	PYC	B	801	8/8	0.96	0.14	44,44,46,48	0
6	SO4	A	2501	5/5	0.97	0.12	46,47,50,51	0
7	FAD	B	405	53/53	0.97	0.10	11,22,27,28	0
6	SO4	B	2505	5/5	0.97	0.20	56,57,58,60	0
8	FMN	B	406	31/31	0.98	0.09	11,15,19,23	0
6	SO4	A	2509	5/5	0.98	0.13	44,46,49,52	0
6	SO4	A	2500	5/5	0.99	0.13	65,66,67,68	0
6	SO4	C	2508	5/5	0.99	0.13	30,33,35,35	0
5	NAD	A	965	44/44	0.99	0.09	8,14,19,25	0
6	SO4	A	2504	5/5	1.00	0.08	29,30,35,37	0
10	ZN	D	100	1/1	1.00	0.08	19,19,19,19	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.

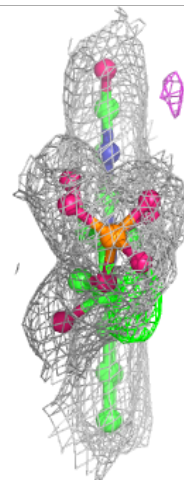
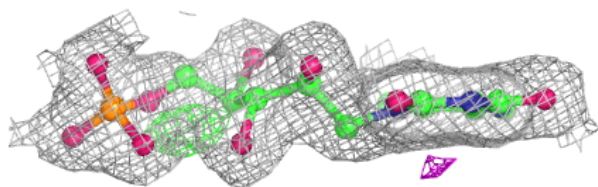
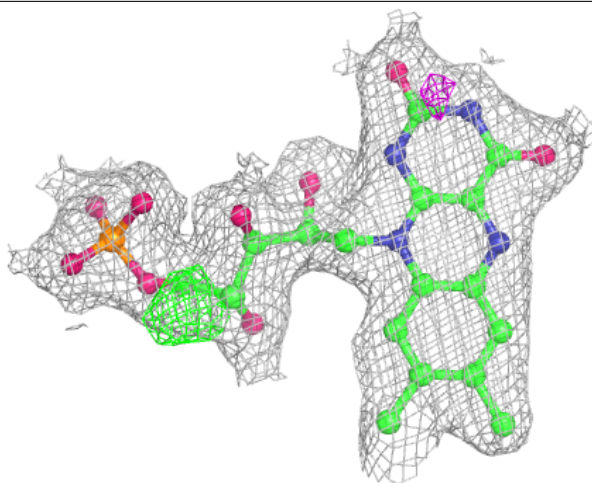
Electron density around FAD B 405:

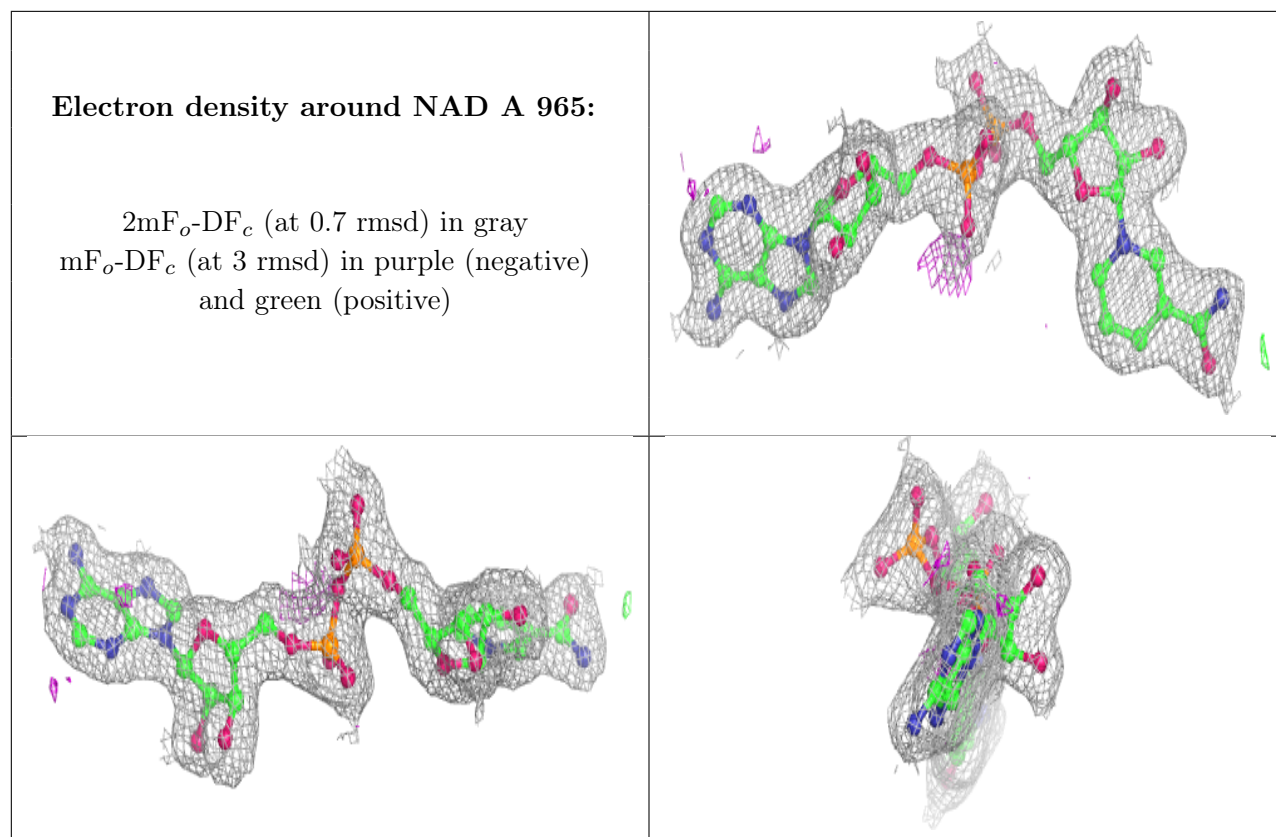
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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.