Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2023 – 07:54 AM EDT

PDB ID	:	3NTD
Title	:	Structure of the Shewanella loihica PV-4 NADH-dependent persulfide reductase C531S Mutant
Authors	:	Sazinsky, M.H.; Warner, M.D.; Lukose, V.; Lee, K.H.; Crane, E.J.
Deposited on	:	2010-07-03
Resolution	:	1.99 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:

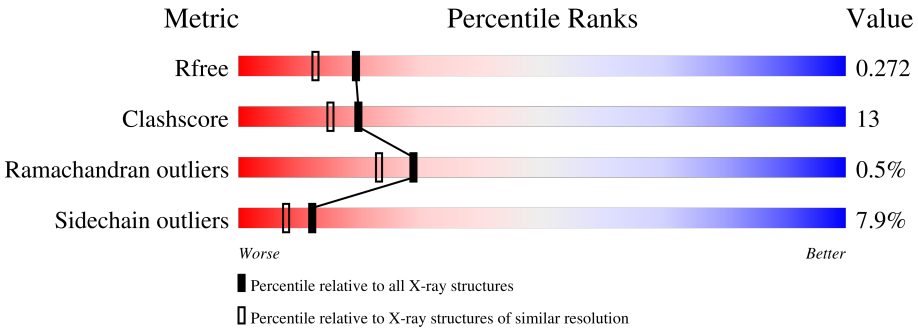
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)

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%

Mol	Chain	Length	Quality of chain
1	A	565	77% 19% .
1	B	565	75% 20% . .

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 9400 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.

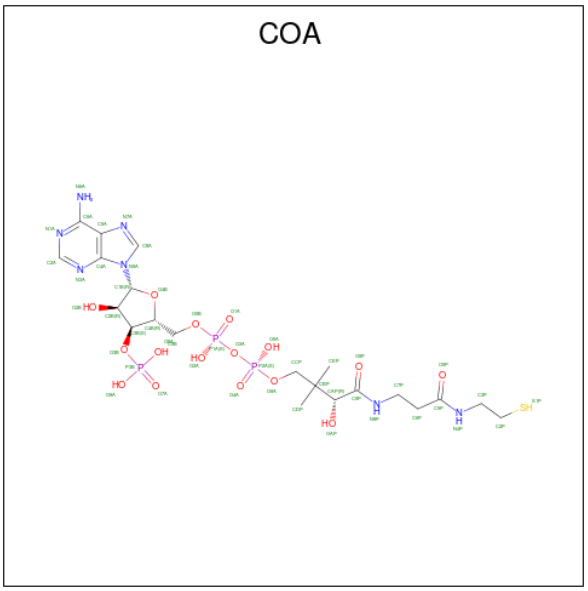
- Molecule 1 is a protein called FAD-dependent pyridine nucleotide-disulphide oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	565	Total	C	N	O	S	0	0	0
			4295	2694	762	820	19			
1	B	565	Total	C	N	O	S	0	0	0
			4294	2694	763	819	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	531	SER	CYS	engineered mutation	UNP A3QAV3
B	531	SER	CYS	engineered mutation	UNP A3QAV3

- Molecule 2 is COENZYME A (three-letter code: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

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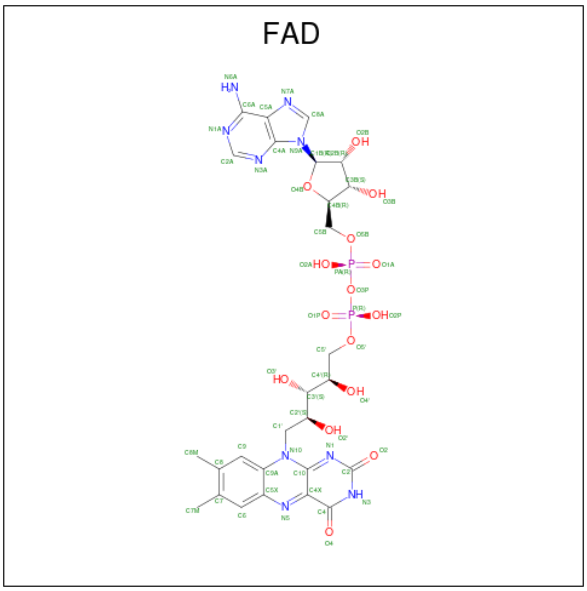
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	300	Total	O	0	0
			300	300		

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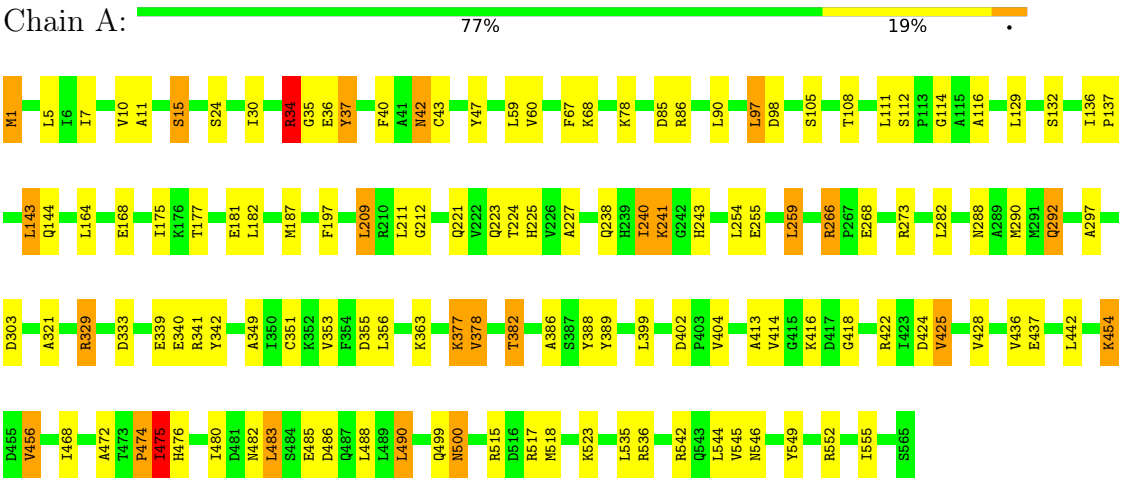
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	B	307	307	307	0	0

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. 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. 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.

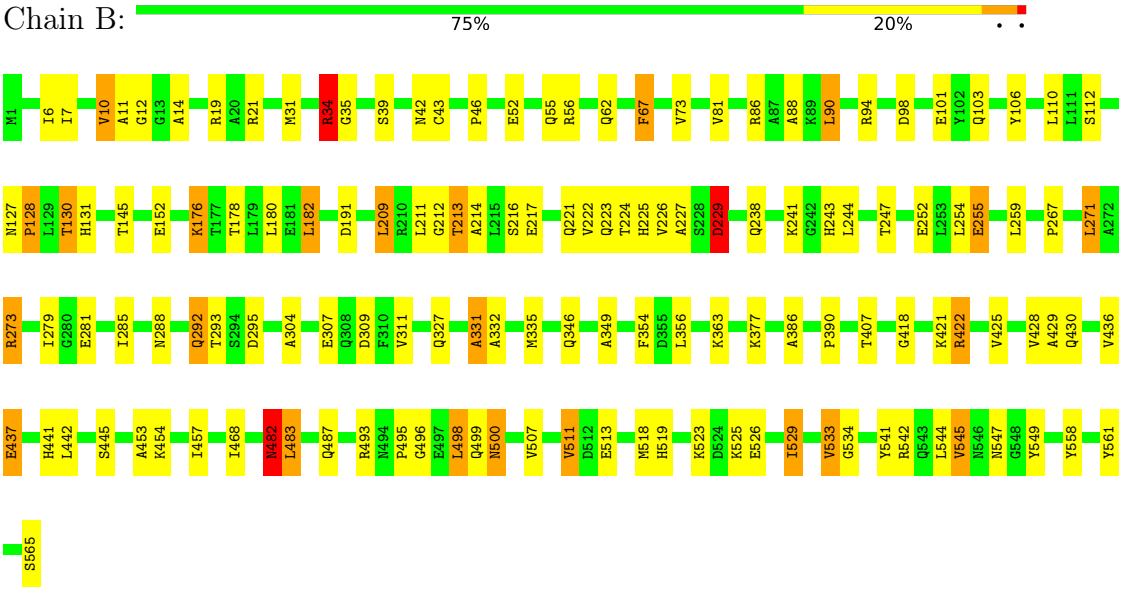
• Molecule 1: FAD-dependent pyridine nucleotide-disulphide oxidoreductase

Chain A:



• Molecule 1: FAD-dependent pyridine nucleotide-disulphide oxidoreductase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α, β, γ	133.74Å 133.74Å 79.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	115.85 – 1.99 38.37 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.3 (115.85-1.99) 99.9 (38.37-1.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	3.22 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.158 , 0.185 0.238 , 0.272	Depositor DCC
R_{free} test set	5561 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å^2)	33.0	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 38.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.56$, $< L^2 > = 0.42$	Xtriage
Estimated twinning fraction	0.000 for -h,-k,l 0.000 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
Reported twinning fraction	0.026 for H, K, L 0.323 for K, H, -L 0.651 for -h,-k,l	Depositor
Outliers	0 of 109330 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	9400	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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: CL, COA, 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	1.31	14/4365 (0.3%)	1.11	14/5910 (0.2%)
1	B	1.36	16/4364 (0.4%)	1.15	17/5908 (0.3%)
All	All	1.34	30/8729 (0.3%)	1.13	31/11818 (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
1	A	0	1

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	292	GLN	CB-CG	-7.61	1.32	1.52
1	B	453	ALA	CA-CB	7.31	1.67	1.52
1	B	558	TYR	CD1-CE1	7.24	1.50	1.39
1	B	437	GLU	CG-CD	6.88	1.62	1.51
1	B	252	GLU	CG-CD	6.43	1.61	1.51
1	B	255	GLU	CB-CG	6.22	1.64	1.52
1	A	268	GLU	CB-CG	6.19	1.64	1.52
1	B	349	ALA	CA-CB	6.06	1.65	1.52
1	B	331	ALA	CA-CB	5.90	1.64	1.52
1	A	321	ALA	CA-CB	5.70	1.64	1.52
1	B	152	GLU	CB-CG	5.68	1.62	1.52
1	B	386	ALA	CA-CB	5.68	1.64	1.52
1	B	429	ALA	CA-CB	5.68	1.64	1.52
1	A	60	VAL	CB-CG1	5.67	1.64	1.52
1	A	342	TYR	CD2-CE2	5.59	1.47	1.39
1	A	425	VAL	CB-CG1	5.54	1.64	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	472	ALA	CA-CB	5.53	1.64	1.52
1	B	21	ARG	CB-CG	5.51	1.67	1.52
1	B	354	PHE	CD2-CE2	5.47	1.50	1.39
1	A	389	TYR	CD2-CE2	5.31	1.47	1.39
1	A	413	ALA	CA-CB	5.29	1.63	1.52
1	A	47	TYR	CE1-CZ	-5.27	1.31	1.38
1	A	349	ALA	CA-CB	5.27	1.63	1.52
1	B	542	ARG	CB-CG	5.19	1.66	1.52
1	B	46	PRO	CG-CD	5.18	1.67	1.50
1	A	37	TYR	CD1-CE1	5.17	1.47	1.39
1	B	106	TYR	CD2-CE2	5.10	1.47	1.39
1	A	181	GLU	CG-CD	-5.08	1.44	1.51
1	A	353	VAL	CB-CG1	5.03	1.63	1.52
1	B	332	ALA	CA-CB	5.02	1.63	1.52

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	34	ARG	NE-CZ-NH1	11.82	126.21	120.30
1	A	34	ARG	NE-CZ-NH1	8.98	124.79	120.30
1	B	422	ARG	NE-CZ-NH1	8.20	124.40	120.30
1	B	511	VAL	CG1-CB-CG2	7.06	122.19	110.90
1	B	191	ASP	CB-CG-OD1	-6.99	112.01	118.30
1	A	34	ARG	NE-CZ-NH2	-6.91	116.85	120.30
1	B	209	LEU	CB-CG-CD2	6.68	122.35	111.00
1	B	511	VAL	CB-CA-C	-6.55	98.96	111.40
1	A	86	ARG	NE-CZ-NH2	-6.37	117.12	120.30
1	B	533	VAL	CG1-CB-CG2	6.30	120.98	110.90
1	A	483	LEU	CB-CG-CD1	5.96	121.12	111.00
1	B	19	ARG	NE-CZ-NH2	-5.91	117.34	120.30
1	B	533	VAL	CA-CB-CG2	5.88	119.72	110.90
1	A	240	ILE	CG1-CB-CG2	-5.83	98.58	111.40
1	A	259	LEU	CB-CG-CD2	5.79	120.84	111.00
1	A	86	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	B	377	LYS	CD-CE-NZ	-5.68	98.63	111.70
1	B	34	ARG	NE-CZ-NH2	-5.61	117.50	120.30
1	B	31	MET	O-C-N	-5.47	113.94	122.70
1	A	517	ARG	NE-CZ-NH2	-5.39	117.60	120.30
1	A	490	LEU	CB-CG-CD2	5.38	120.14	111.00
1	B	533	VAL	CA-CB-CG1	5.32	118.88	110.90
1	B	229	ASP	CB-CG-OD1	-5.29	113.54	118.30
1	A	552	ARG	NE-CZ-NH1	5.22	122.91	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	442	LEU	CB-CG-CD2	5.21	119.85	111.00
1	A	143	LEU	CA-CB-CG	5.20	127.26	115.30
1	A	488	LEU	CA-CB-CG	5.07	126.95	115.30
1	A	341	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	B	273	ARG	NE-CZ-NH2	-5.04	117.78	120.30
1	B	191	ASP	CB-CG-OD2	5.03	122.83	118.30
1	A	475	ILE	CB-CA-C	-5.02	101.56	111.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	474	PRO	Peptide

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	4295	0	4296	115	1
1	B	4294	0	4300	110	1
2	A	48	0	31	2	0
2	B	48	0	32	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	53	0	31	3	0
4	B	53	0	31	2	0
5	A	300	0	0	25	0
5	B	307	0	0	32	0
All	All	9400	0	8721	222	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 13.

All (222) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:229:ASP:HB3	5:B:745:HOH:O	1.41	1.18
1:B:500:ASN:HB3	5:B:803:HOH:O	1.46	1.16
1:A:266:ARG:HG3	1:A:266:ARG:HH11	1.11	1.11
1:A:474:PRO:CB	1:A:555:ILE:HD11	1.83	1.08
1:A:227:ALA:H	1:A:238:GLN:NE2	1.55	1.02
1:A:227:ALA:N	1:A:238:GLN:NE2	2.09	1.00
1:A:382:THR:HG23	1:A:456:VAL:HG22	1.44	1.00
1:A:266:ARG:HG3	1:A:266:ARG:NH1	1.68	0.99
1:A:437:GLU:HG2	5:A:618:HOH:O	1.62	0.98
1:B:55:GLN:HG2	5:B:878:HOH:O	1.65	0.96
1:B:223:GLN:HE21	1:B:225:HIS:H	1.09	0.94
1:B:346:GLN:HE21	1:B:430:GLN:HE21	1.09	0.93
1:A:85:ASP:HB3	5:A:733:HOH:O	1.67	0.92
1:A:1:MET:HB3	5:A:820:HOH:O	1.68	0.91
1:A:227:ALA:HB3	1:A:238:GLN:NE2	1.85	0.91
1:B:67:PHE:CD1	1:B:73:VAL:HG11	2.06	0.90
1:A:227:ALA:H	1:A:238:GLN:HE21	1.21	0.89
1:A:474:PRO:HB2	1:A:555:ILE:HD11	1.54	0.87
1:B:288:ASN:HD21	1:B:292:GLN:HE21	1.17	0.87
1:A:241:LYS:HE3	5:A:1106:HOH:O	1.73	0.86
1:A:418:GLY:HA2	5:A:1092:HOH:O	1.75	0.86
1:A:499:GLN:HB2	5:A:759:HOH:O	1.74	0.85
1:A:474:PRO:HB3	1:A:555:ILE:HD11	1.57	0.84
1:B:523:LYS:HE3	5:B:969:HOH:O	1.77	0.84
1:B:482:ASN:O	1:B:483:LEU:HB2	1.77	0.83
1:A:266:ARG:HH11	1:A:266:ARG:CG	1.90	0.82
1:A:241:LYS:HD3	5:A:1106:HOH:O	1.80	0.81
1:B:86:ARG:HD2	1:B:295:ASP:OD1	1.80	0.81
1:A:382:THR:HG21	1:A:456:VAL:HA	1.63	0.79
1:A:227:ALA:CB	1:A:238:GLN:HE21	1.96	0.79
5:A:814:HOH:O	1:B:441:HIS:HE1	1.66	0.78
1:B:223:GLN:NE2	1:B:225:HIS:H	1.80	0.78
1:B:243:HIS:HE1	1:B:255:GLU:HG3	1.49	0.78
1:B:127:ASN:OD1	1:B:130:THR:HG23	1.84	0.77
1:B:243:HIS:HE1	1:B:255:GLU:CG	1.98	0.77
5:A:814:HOH:O	1:B:441:HIS:CE1	2.36	0.77
1:A:227:ALA:N	1:A:238:GLN:HE22	1.82	0.77
1:A:227:ALA:CB	1:A:238:GLN:NE2	2.47	0.77
1:A:474:PRO:HB2	1:A:555:ILE:CD1	2.14	0.76
1:A:223:GLN:HE21	1:A:225:HIS:H	1.29	0.76
1:B:211:LEU:O	1:B:213:THR:HG22	1.86	0.74
1:B:496:GLY:O	1:B:499:GLN:HB2	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:436:VAL:HG21	1:A:468:ILE:CD1	2.19	0.73
1:B:519:HIS:CD2	5:B:757:HOH:O	2.41	0.72
1:A:476:HIS:CE1	1:A:555:ILE:HD13	2.25	0.71
1:A:241:LYS:CE	5:A:1106:HOH:O	2.35	0.69
1:B:67:PHE:HD1	1:B:73:VAL:HG11	1.55	0.69
1:B:67:PHE:CD1	1:B:73:VAL:CG1	2.77	0.68
1:A:227:ALA:CA	1:A:238:GLN:NE2	2.57	0.67
1:A:290:MET:HE3	1:A:339:GLU:HA	1.76	0.67
1:A:209:LEU:HD13	1:A:211:LEU:HD11	1.75	0.66
1:B:519:HIS:HE1	5:B:754:HOH:O	1.77	0.66
1:A:499:GLN:CB	5:A:759:HOH:O	2.38	0.66
1:A:288:ASN:HD21	1:A:292:GLN:HE21	1.43	0.65
1:B:94:ARG:CZ	5:B:674:HOH:O	2.44	0.65
1:B:67:PHE:HD1	1:B:73:VAL:CG1	2.09	0.65
1:A:243:HIS:HE1	1:A:255:GLU:OE2	1.79	0.65
1:A:290:MET:CE	1:A:339:GLU:HA	2.27	0.64
1:B:519:HIS:CE1	5:B:754:HOH:O	2.51	0.63
1:A:164:LEU:HD12	1:A:187:MET:HE2	1.78	0.63
1:A:436:VAL:HG21	1:A:468:ILE:HD12	1.79	0.63
1:A:116:ALA:HB3	1:A:266:ARG:HH12	1.64	0.62
1:B:243:HIS:HD2	5:B:753:HOH:O	1.82	0.62
1:A:168:GLU:HG2	5:A:922:HOH:O	1.99	0.62
1:B:55:GLN:HG3	5:B:967:HOH:O	2.00	0.61
1:A:425:VAL:HG13	1:B:428:VAL:HG21	1.82	0.61
1:A:227:ALA:N	1:A:238:GLN:HE21	1.87	0.61
1:A:290:MET:HE1	1:A:339:GLU:HG3	1.81	0.61
1:A:382:THR:CG2	1:A:456:VAL:HG22	2.26	0.61
1:A:211:LEU:HD12	1:A:211:LEU:N	2.16	0.60
1:A:187:MET:HE3	1:A:351:CYS:HB3	1.83	0.59
1:B:346:GLN:NE2	1:B:430:GLN:HE21	1.92	0.59
1:B:130:THR:HG21	1:B:259:LEU:HD23	1.85	0.59
1:B:94:ARG:HD2	5:B:839:HOH:O	2.03	0.59
1:B:436:VAL:HG21	1:B:468:ILE:HD12	1.84	0.59
1:B:130:THR:HB	1:B:259:LEU:HB3	1.85	0.58
1:A:266:ARG:NH1	1:A:266:ARG:CG	2.51	0.58
1:B:127:ASN:CG	1:B:130:THR:HG23	2.24	0.58
1:B:211:LEU:O	1:B:213:THR:CG2	2.52	0.58
1:B:223:GLN:HE21	1:B:225:HIS:N	1.90	0.57
1:A:90:LEU:HD12	1:A:105:SER:HA	1.87	0.56
1:A:355:ASP:HB2	5:A:965:HOH:O	2.04	0.56
1:B:273:ARG:HG2	1:B:279:ILE:CD1	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:VAL:CG2	1:B:468:ILE:HD12	2.36	0.56
1:B:534:GLY:HA2	5:B:767:HOH:O	2.06	0.56
1:A:223:GLN:NE2	1:A:225:HIS:H	2.02	0.56
1:A:474:PRO:CB	1:A:555:ILE:CD1	2.68	0.56
1:B:227:ALA:H	1:B:238:GLN:NE2	2.02	0.56
1:A:209:LEU:HD22	1:A:211:LEU:CD1	2.36	0.56
1:B:523:LYS:CE	5:B:969:HOH:O	2.43	0.56
1:B:498:LEU:HD23	1:B:498:LEU:H	1.71	0.56
1:A:437:GLU:CG	5:A:618:HOH:O	2.35	0.56
1:B:222:VAL:HG22	5:B:749:HOH:O	2.05	0.55
1:B:182:LEU:O	1:B:212:GLY:HA2	2.05	0.55
1:B:101:GLU:HG3	5:B:893:HOH:O	2.06	0.55
1:A:7:ILE:HD12	1:A:111:LEU:CD2	2.35	0.55
1:A:34:ARG:CD	4:A:900:FAD:O2B	2.56	0.54
1:A:424:ASP:O	1:A:428:VAL:HG23	2.08	0.54
1:B:94:ARG:CD	5:B:839:HOH:O	2.55	0.54
1:A:36:GLU:HG3	1:A:78:LYS:HE3	1.90	0.54
1:A:273:ARG:HG2	1:A:273:ARG:HH11	1.71	0.54
1:A:68:LYS:NZ	5:A:674:HOH:O	2.39	0.54
1:B:498:LEU:HD23	1:B:498:LEU:N	2.23	0.53
1:A:329:ARG:HH11	1:A:329:ARG:CG	2.22	0.53
1:B:518:MET:HE1	1:B:547:ASN:HB2	1.91	0.53
1:A:34:ARG:HD2	4:A:900:FAD:O2B	2.09	0.53
1:B:176:LYS:HG2	5:B:1025:HOH:O	2.10	0.52
1:B:243:HIS:CE1	1:B:255:GLU:HG3	2.38	0.52
1:B:425:VAL:CG1	1:B:457:ILE:HD13	2.39	0.52
1:A:11:ALA:HB2	2:A:901:COA:H32	1.92	0.51
1:A:436:VAL:CG2	1:A:468:ILE:CD1	2.86	0.51
1:A:436:VAL:HG21	1:A:468:ILE:HD11	1.91	0.51
1:A:97:LEU:HD23	1:A:98:ASP:CG	2.30	0.51
1:A:402:ASP:OD1	1:A:404:VAL:HG23	2.10	0.51
1:A:436:VAL:CG2	1:A:468:ILE:HD12	2.40	0.51
1:B:273:ARG:HG2	1:B:279:ILE:HD13	1.93	0.51
1:B:243:HIS:CE1	1:B:255:GLU:CG	2.88	0.51
1:A:378:VAL:HG12	1:A:399:LEU:HB3	1.94	0.50
1:A:518:MET:HE3	1:A:549:TYR:CE1	2.47	0.50
1:B:418:GLY:HA2	5:B:721:HOH:O	2.10	0.50
1:B:518:MET:HE3	1:B:544:LEU:HD23	1.93	0.50
1:B:418:GLY:N	5:B:721:HOH:O	2.34	0.50
1:B:541:TYR:O	1:B:545:VAL:HB	2.12	0.49
1:B:513:GLU:CG	5:B:1065:HOH:O	2.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:LYS:CD	5:A:1106:HOH:O	2.40	0.49
1:B:513:GLU:HG3	5:B:1065:HOH:O	2.13	0.49
1:B:43:CYS:SG	4:B:900:FAD:C4X	3.01	0.49
1:A:108:THR:HG23	1:A:297:ALA:O	2.12	0.48
1:A:168:GLU:CG	5:A:922:HOH:O	2.57	0.48
1:B:56:ARG:HD2	5:B:659:HOH:O	2.12	0.48
1:A:545:VAL:HG12	5:A:649:HOH:O	2.13	0.48
1:A:518:MET:HE3	1:A:544:LEU:HD23	1.95	0.48
1:A:535:LEU:HD23	1:A:536:ARG:N	2.28	0.48
1:B:131:HIS:HE1	1:B:145:THR:OG1	1.97	0.48
1:A:454:LYS:HE3	5:B:652:HOH:O	2.13	0.48
1:B:311:VAL:HG13	5:B:1060:HOH:O	2.13	0.48
1:B:493:ARG:HH12	1:B:500:ASN:HD21	1.60	0.48
1:A:288:ASN:HD21	1:A:292:GLN:HG2	1.79	0.48
1:A:7:ILE:HD12	1:A:111:LEU:HD23	1.95	0.48
1:A:209:LEU:HD22	1:A:211:LEU:HD12	1.96	0.48
1:A:288:ASN:ND2	1:A:292:GLN:HG2	2.29	0.47
1:A:211:LEU:N	1:A:211:LEU:CD1	2.77	0.47
1:B:483:LEU:HA	1:B:487:GLN:OE1	2.14	0.47
1:B:182:LEU:HG	1:B:214:ALA:HB2	1.94	0.47
1:B:518:MET:HE3	1:B:549:TYR:CE1	2.49	0.47
1:B:518:MET:CE	1:B:547:ASN:HB2	2.44	0.47
1:A:227:ALA:HB3	1:A:238:GLN:HE21	1.54	0.47
1:B:88:ALA:HB1	1:B:90:LEU:HD22	1.96	0.47
1:A:243:HIS:HD2	5:A:591:HOH:O	1.97	0.47
1:A:402:ASP:OD1	1:A:404:VAL:CG2	2.62	0.47
1:B:518:MET:HE2	1:B:547:ASN:CB	2.45	0.47
1:A:209:LEU:HD22	1:A:211:LEU:HD11	1.97	0.47
1:A:197:PHE:HB2	1:A:414:VAL:HG11	1.97	0.47
1:B:6:ILE:HG23	1:B:110:LEU:HD23	1.97	0.46
1:B:307:GLU:CD	1:B:327:GLN:HE22	2.18	0.46
1:B:418:GLY:CA	5:B:721:HOH:O	2.62	0.46
1:B:529:ILE:HD11	1:B:544:LEU:HD12	1.97	0.46
1:B:226:VAL:HA	1:B:238:GLN:HE22	1.80	0.46
1:A:114:GLY:HA2	1:A:303:ASP:HB2	1.98	0.46
1:A:209:LEU:CD1	1:A:211:LEU:HD11	2.44	0.46
1:B:34:ARG:HG2	1:B:35:GLY:N	2.30	0.46
1:B:39:SER:HB2	1:B:62:GLN:HB2	1.96	0.46
1:B:519:HIS:HD2	5:B:757:HOH:O	1.92	0.46
1:B:243:HIS:CD2	5:B:753:HOH:O	2.62	0.46
1:B:222:VAL:CG2	5:B:749:HOH:O	2.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:TYR:CD2	1:A:136:ILE:HD12	2.52	0.45
1:A:240:ILE:HD13	1:A:240:ILE:HG21	1.63	0.45
1:B:130:THR:HG22	1:B:244:LEU:HD11	1.98	0.45
1:A:182:LEU:O	1:A:212:GLY:HA2	2.16	0.45
1:B:127:ASN:ND2	1:B:130:THR:HG23	2.32	0.45
1:B:271:LEU:HB2	5:B:699:HOH:O	2.17	0.45
1:A:136:ILE:HB	1:A:137:PRO:HD3	1.99	0.45
1:A:333:ASP:OD2	1:A:340:GLU:OE1	2.35	0.45
1:B:507:VAL:HG23	5:B:758:HOH:O	2.17	0.45
1:A:386:ALA:HB1	1:A:388:TYR:CE2	2.52	0.44
1:B:454:LYS:HB2	1:B:454:LYS:HE3	1.73	0.44
1:A:243:HIS:HA	5:A:591:HOH:O	2.16	0.44
1:B:127:ASN:HD21	1:B:130:THR:CG2	2.31	0.44
1:B:331:ALA:O	1:B:335:MET:HG3	2.16	0.44
1:B:7:ILE:HG12	1:B:81:VAL:HG21	2.00	0.43
1:B:178:THR:HG21	1:B:254:LEU:HD11	2.00	0.43
1:A:112:SER:O	4:A:900:FAD:H52A	2.18	0.43
1:B:11:ALA:HB2	2:B:901:COA:H32	1.99	0.43
1:A:5:LEU:HD22	1:A:30:ILE:HB	2.00	0.43
1:A:500:ASN:HD22	1:A:500:ASN:N	2.16	0.43
1:B:309:ASP:OD2	1:B:363:LYS:HE3	2.19	0.43
1:A:416:LYS:N	5:A:821:HOH:O	2.50	0.43
1:B:43:CYS:SG	2:B:901:COA:S1P	3.16	0.43
1:B:217:GLU:HB2	1:B:247:THR:HB	2.00	0.43
1:A:475:ILE:H	1:A:475:ILE:HG13	1.54	0.43
1:A:515:ARG:N	5:A:779:HOH:O	2.44	0.43
1:B:130:THR:CG2	1:B:259:LEU:HD23	2.47	0.43
1:A:240:ILE:HG22	5:A:675:HOH:O	2.19	0.43
1:B:561:TYR:C	1:B:561:TYR:CD2	2.92	0.43
1:A:97:LEU:HD23	1:A:98:ASP:CB	2.49	0.43
1:B:223:GLN:HE21	1:B:224:THR:N	2.15	0.43
1:B:529:ILE:HD11	1:B:544:LEU:CD1	2.49	0.42
1:B:500:ASN:C	1:B:500:ASN:HD22	2.22	0.42
1:A:175:ILE:O	1:A:177:THR:HG23	2.19	0.42
1:B:12:GLY:HA3	1:B:112:SER:OG	2.19	0.42
1:B:518:MET:CE	1:B:547:ASN:CB	2.96	0.42
1:B:487:GLN:HG2	1:B:526:GLU:HG3	2.01	0.42
1:A:377:LYS:HE3	1:A:377:LYS:HB3	1.76	0.42
1:A:424:ASP:OD2	1:B:421:LYS:HD2	2.19	0.42
1:B:285:ILE:HD12	1:B:304:ALA:HB1	2.02	0.42
1:B:495:PRO:HD2	5:B:1022:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:GLN:HB2	5:A:1066:HOH:O	2.20	0.41
1:B:436:VAL:HG21	1:B:468:ILE:CD1	2.50	0.41
1:A:11:ALA:O	1:A:15:SER:HB3	2.20	0.41
1:A:34:ARG:HG2	1:A:35:GLY:N	2.35	0.41
1:A:475:ILE:CD1	1:A:480:ILE:HD12	2.50	0.41
1:B:127:ASN:HB2	1:B:128:PRO:CD	2.50	0.41
1:A:442:LEU:HD22	1:B:428:VAL:HG22	2.02	0.41
1:B:34:ARG:HD2	4:B:900:FAD:C4A	2.50	0.41
1:A:42:ASN:OD1	2:A:901:COA:H22	2.21	0.41
1:A:542:ARG:HH21	1:A:546:ASN:HD21	1.68	0.41
1:B:34:ARG:NH2	5:B:894:HOH:O	2.40	0.41
1:A:40:PHE:HA	1:A:59:LEU:O	2.21	0.41
1:A:378:VAL:HA	1:A:555:ILE:HG21	2.02	0.41
1:B:10:VAL:O	1:B:14:ALA:HB3	2.21	0.41
1:A:382:THR:HA	5:A:979:HOH:O	2.20	0.41
1:A:243:HIS:CE1	1:A:255:GLU:HB3	2.56	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)
1:A:482:ASN:OD1	1:B:482:ASN:ND2[1_554]	2.12	0.08

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	563/565 (100%)	547 (97%)	14 (2%)	2 (0%)	34	30
1	B	563/565 (100%)	544 (97%)	15 (3%)	4 (1%)	22	16
All	All	1126/1130 (100%)	1091 (97%)	29 (3%)	6 (0%)	29	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	482	ASN
1	B	10	VAL
1	B	293	THR
1	A	10	VAL
1	B	483	LEU
1	A	456	VAL

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	449/450 (100%)	415 (92%)	34 (8%)	13	8
1	B	449/450 (100%)	412 (92%)	37 (8%)	11	7
All	All	898/900 (100%)	827 (92%)	71 (8%)	12	8

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	15	SER
1	A	24	SER
1	A	34	ARG
1	A	42	ASN
1	A	43	CYS
1	A	67	PHE
1	A	97	LEU
1	A	129	LEU
1	A	132	SER
1	A	143	LEU
1	A	209	LEU
1	A	221	GLN
1	A	224	THR
1	A	241	LYS
1	A	254	LEU
1	A	259	LEU

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Mol	Chain	Res	Type
1	A	266	ARG
1	A	282	LEU
1	A	329	ARG
1	A	356	LEU
1	A	363	LYS
1	A	377	LYS
1	A	378	VAL
1	A	382	THR
1	A	422	ARG
1	A	454	LYS
1	A	475	ILE
1	A	483	LEU
1	A	485	GLU
1	A	486	ASP
1	A	490	LEU
1	A	500	ASN
1	A	523	LYS
1	B	34	ARG
1	B	42	ASN
1	B	52	GLU
1	B	67	PHE
1	B	90	LEU
1	B	98	ASP
1	B	103	GLN
1	B	128	PRO
1	B	130	THR
1	B	176	LYS
1	B	180	LEU
1	B	182	LEU
1	B	209	LEU
1	B	213	THR
1	B	216	SER
1	B	221	GLN
1	B	229	ASP
1	B	241	LYS
1	B	267	PRO
1	B	271	LEU
1	B	281	GLU
1	B	292	GLN
1	B	356	LEU
1	B	390	PRO
1	B	407	THR

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Mol	Chain	Res	Type
1	B	422	ARG
1	B	437	GLU
1	B	445	SER
1	B	482	ASN
1	B	498	LEU
1	B	500	ASN
1	B	511	VAL
1	B	525	LYS
1	B	529	ILE
1	B	533	VAL
1	B	545	VAL
1	B	565	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	171	HIS
1	A	223	GLN
1	A	238	GLN
1	A	243	HIS
1	A	292	GLN
1	A	440	GLN
1	A	441	HIS
1	A	500	ASN
1	A	543	GLN
1	A	546	ASN
1	A	547	ASN
1	B	62	GLN
1	B	131	HIS
1	B	205	GLN
1	B	223	GLN
1	B	238	GLN
1	B	243	HIS
1	B	292	GLN
1	B	370	GLN
1	B	430	GLN
1	B	440	GLN
1	B	441	HIS
1	B	482	ASN
1	B	494	ASN
1	B	500	ASN
1	B	519	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 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	FAD	A	900	-	53,58,58	1.45	6 (11%)	68,89,89	1.62	15 (22%)
4	FAD	B	900	-	53,58,58	1.61	8 (15%)	68,89,89	1.52	13 (19%)
2	COA	B	901	-	41,50,50	1.56	5 (12%)	52,75,75	2.86	12 (23%)
2	COA	A	901	-	41,50,50	1.54	4 (9%)	52,75,75	1.71	10 (19%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FAD	A	900	-	-	2/30/50/50	0/6/6/6
4	FAD	B	900	-	-	1/30/50/50	0/6/6/6
2	COA	B	901	-	-	4/44/64/64	0/3/3/3
2	COA	A	901	-	-	2/44/64/64	0/3/3/3

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	COA	O9P-C9P	6.21	1.35	1.23
4	B	900	FAD	C2A-N3A	5.35	1.40	1.32
4	B	900	FAD	C4X-N5	4.84	1.40	1.30
4	A	900	FAD	C4X-N5	4.68	1.39	1.30
2	B	901	COA	C2A-N3A	4.38	1.39	1.32
2	B	901	COA	O9P-C9P	4.12	1.31	1.23
2	B	901	COA	O5P-C5P	4.10	1.31	1.23
2	A	901	COA	C2A-N3A	3.83	1.38	1.32
4	A	900	FAD	C2A-N3A	3.78	1.38	1.32
4	A	900	FAD	O2-C2	-3.74	1.17	1.24
4	A	900	FAD	C2A-N1A	3.51	1.40	1.33
4	B	900	FAD	C10-N1	3.35	1.40	1.33
4	B	900	FAD	C2A-N1A	3.16	1.39	1.33
4	B	900	FAD	O2'-C2'	-2.99	1.37	1.43
2	B	901	COA	O4B-C1B	2.93	1.45	1.41
4	A	900	FAD	C4A-N3A	2.85	1.39	1.35
2	A	901	COA	O4B-C1B	2.84	1.45	1.41
2	A	901	COA	P1A-O2A	-2.84	1.42	1.55
4	B	900	FAD	C9A-N10	-2.67	1.36	1.41
4	A	900	FAD	C10-N10	2.50	1.42	1.37
2	B	901	COA	C6P-C5P	2.38	1.55	1.51
4	B	900	FAD	C5A-N7A	-2.18	1.31	1.39
4	B	900	FAD	O5B-C5B	-2.02	1.37	1.44

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	901	COA	C3P-N4P-C5P	-11.38	101.71	122.84
2	B	901	COA	N3A-C2A-N1A	-7.85	116.40	128.68
2	B	901	COA	C7P-C6P-C5P	-7.78	99.41	112.36
2	B	901	COA	CEP-CBP-CDP	-5.59	97.78	109.17
2	A	901	COA	CDP-CBP-CAP	5.34	118.08	108.82
2	A	901	COA	N3A-C2A-N1A	-4.87	121.07	128.68
2	B	901	COA	C6P-C5P-N4P	-4.66	108.57	116.42
4	A	900	FAD	O2-C2-N1	-4.64	114.14	121.83
4	A	900	FAD	C4-N3-C2	-4.60	117.13	125.64
4	A	900	FAD	N3A-C2A-N1A	-4.56	121.55	128.68
2	B	901	COA	CDP-CBP-CAP	4.50	116.62	108.82
2	B	901	COA	CEP-CBP-CCP	4.15	115.00	108.23
4	B	900	FAD	C5'-C4'-C3'	-3.92	104.63	112.20
4	A	900	FAD	C4X-C10-N10	3.91	122.19	116.48
4	B	900	FAD	N3A-C2A-N1A	-3.81	122.72	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	901	COA	CEP-CBP-CDP	-3.60	101.82	109.17
2	A	901	COA	C6P-C7P-N8P	-3.53	104.76	111.90
2	B	901	COA	C5B-C4B-C3B	-3.41	103.10	114.40
2	B	901	COA	C6P-C7P-N8P	-3.40	105.02	111.90
2	B	901	COA	CAP-C9P-N8P	3.34	123.22	116.58
4	A	900	FAD	O4B-C1B-C2B	-3.29	102.11	106.93
4	B	900	FAD	O4B-C4B-C3B	2.93	110.92	105.11
4	B	900	FAD	C5X-C9A-N10	2.91	120.96	117.95
2	B	901	COA	C2P-C3P-N4P	2.90	118.93	112.31
4	B	900	FAD	C9-C9A-N10	-2.84	117.99	121.84
4	B	900	FAD	C4A-C5A-N7A	-2.82	106.46	109.40
4	B	900	FAD	O2'-C2'-C1'	-2.79	103.06	109.80
4	A	900	FAD	C4-C4X-C10	2.74	121.39	116.79
2	B	901	COA	O9P-C9P-CAP	-2.68	112.90	121.06
4	B	900	FAD	O4-C4-C4X	-2.61	119.66	126.60
4	B	900	FAD	C5A-C6A-N6A	2.51	124.16	120.35
4	A	900	FAD	N3-C2-N1	2.50	124.28	119.38
4	B	900	FAD	C4X-C10-N10	2.48	120.11	116.48
4	A	900	FAD	C10-C4X-N5	-2.48	119.59	124.86
4	B	900	FAD	O3B-C3B-C4B	2.45	118.12	111.05
4	A	900	FAD	C9A-N10-C10	-2.44	116.97	120.77
2	A	901	COA	CEP-CBP-CCP	2.41	112.17	108.23
4	A	900	FAD	O4'-C4'-C5'	-2.40	104.53	109.92
2	A	901	COA	C2P-C3P-N4P	-2.40	106.83	112.31
4	B	900	FAD	O4'-C4'-C3'	2.31	114.72	109.10
4	A	900	FAD	O4-C4-C4X	-2.30	120.51	126.60
4	A	900	FAD	O3B-C3B-C2B	-2.28	104.44	111.82
2	A	901	COA	O6A-P2A-O4A	2.26	117.90	109.07
4	A	900	FAD	O2'-C2'-C3'	2.24	114.55	109.10
4	A	900	FAD	C4X-C10-N1	-2.14	119.77	124.73
2	A	901	COA	N6A-C6A-N1A	2.13	123.00	118.57
4	A	900	FAD	C4X-C4-N3	2.13	118.60	113.19
2	A	901	COA	C1B-N9A-C4A	-2.12	122.91	126.64
2	A	901	COA	O5P-C5P-C6P	2.04	125.74	122.02
4	B	900	FAD	C4X-C4-N3	2.04	118.36	113.19

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	COA	P2A-O3A-P1A-O5B
2	B	901	COA	P2A-O3A-P1A-O5B

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Mol	Chain	Res	Type	Atoms
4	A	900	FAD	PA-O3P-P-O5'
2	B	901	COA	C3B-O3B-P3B-O9A
2	B	901	COA	O5P-C5P-N4P-C3P
4	B	900	FAD	O4B-C4B-C5B-O5B
2	A	901	COA	P1A-O3A-P2A-O5A
2	B	901	COA	P1A-O3A-P2A-O4A
4	A	900	FAD	O4B-C4B-C5B-O5B

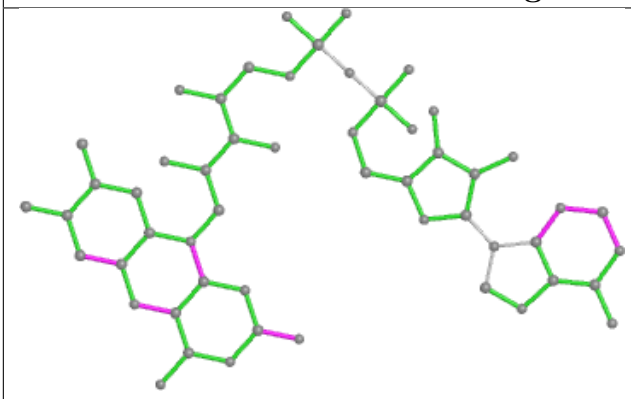
There are no ring outliers.

4 monomers are involved in 9 short contacts:

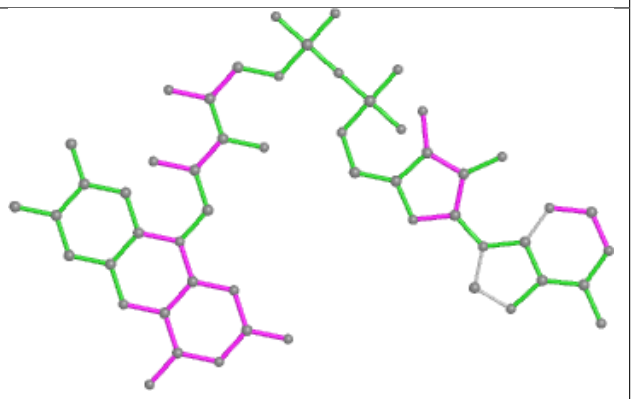
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	900	FAD	3	0
4	B	900	FAD	2	0
2	B	901	COA	2	0
2	A	901	COA	2	0

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.

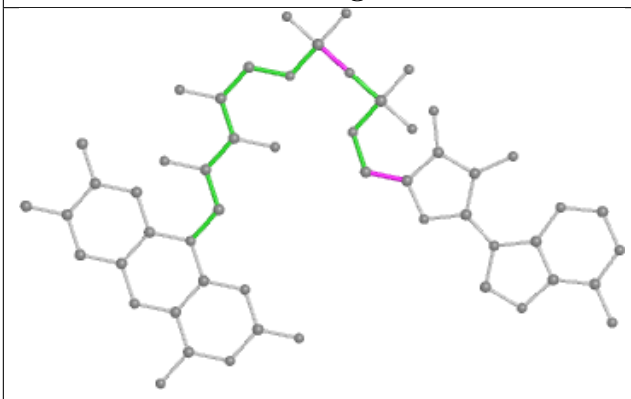
Ligand FAD A 900



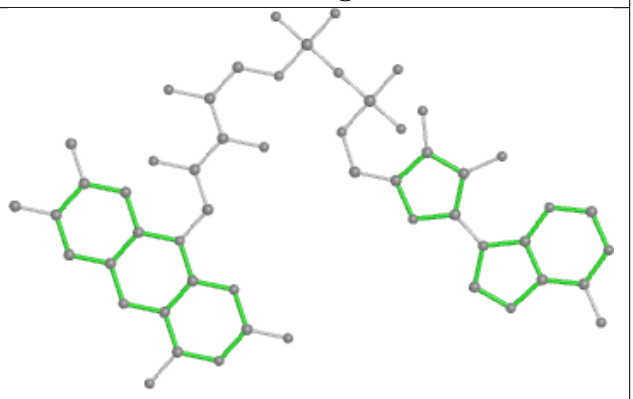
Bond lengths



Bond angles

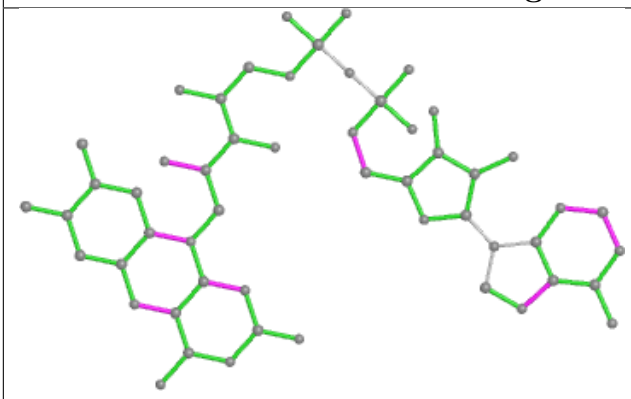


Torsions

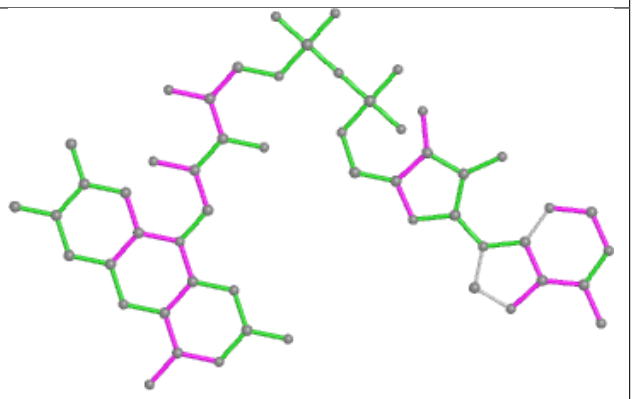


Rings

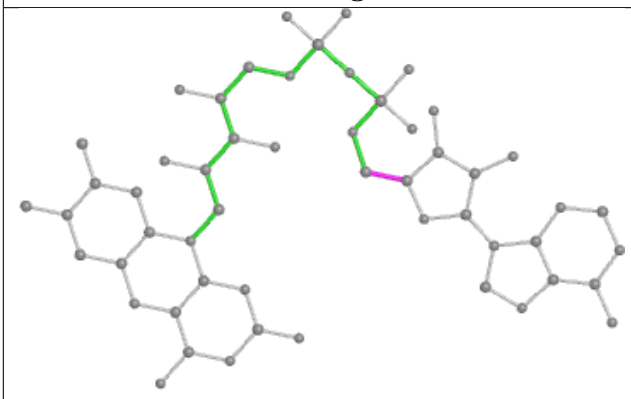
Ligand FAD B 900



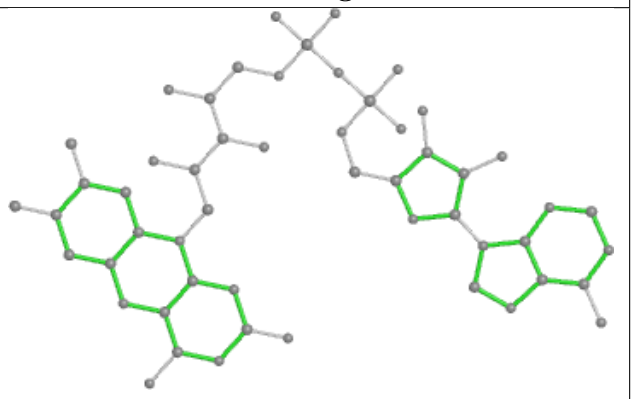
Bond lengths



Bond angles

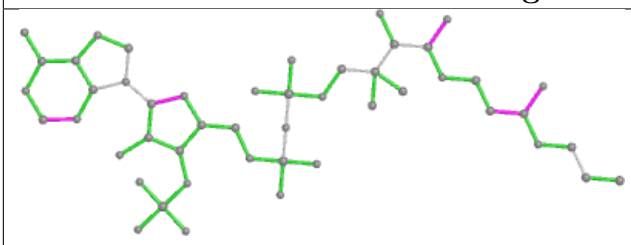


Torsions

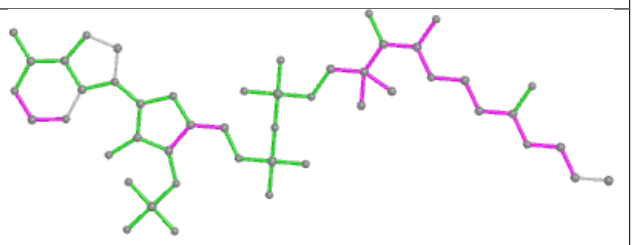


Rings

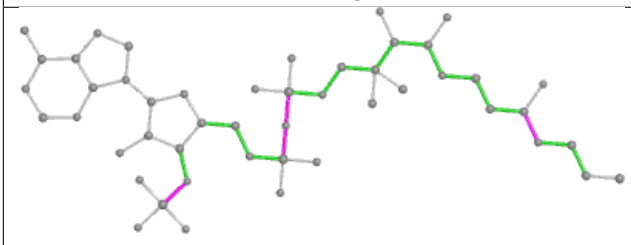
Ligand COA B 901



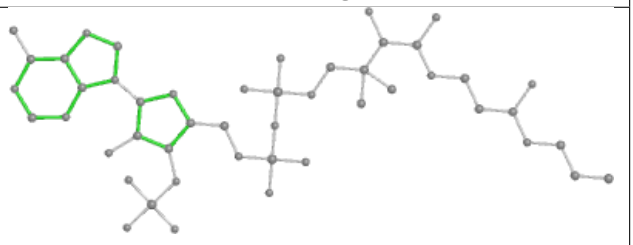
Bond lengths



Bond angles

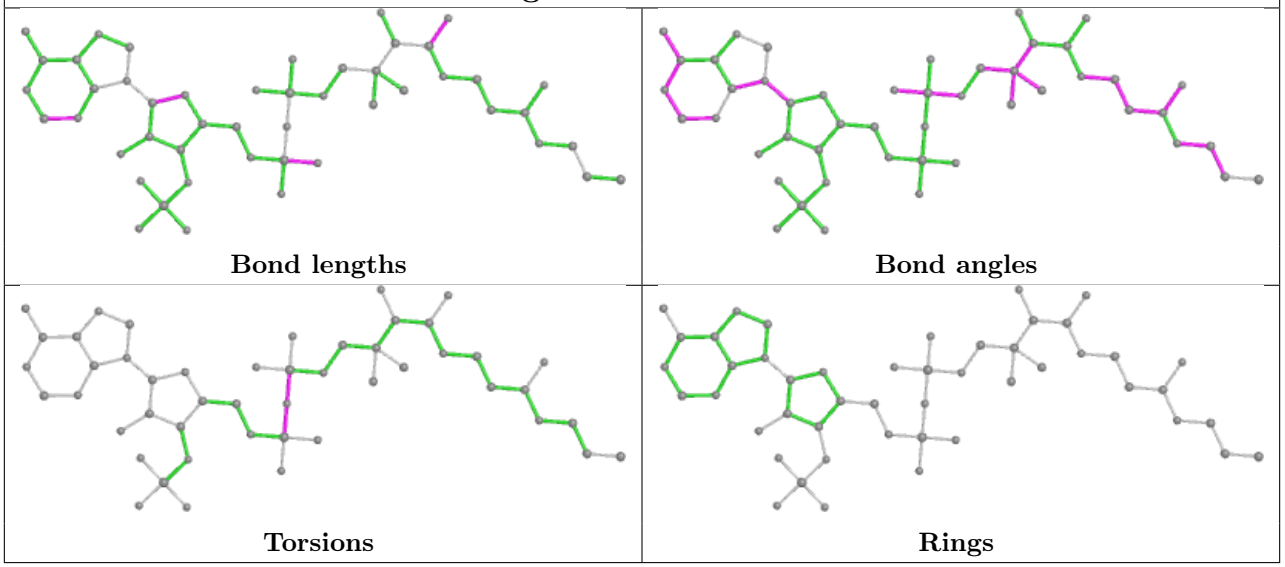


Torsions



Rings

Ligand COA A 901



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

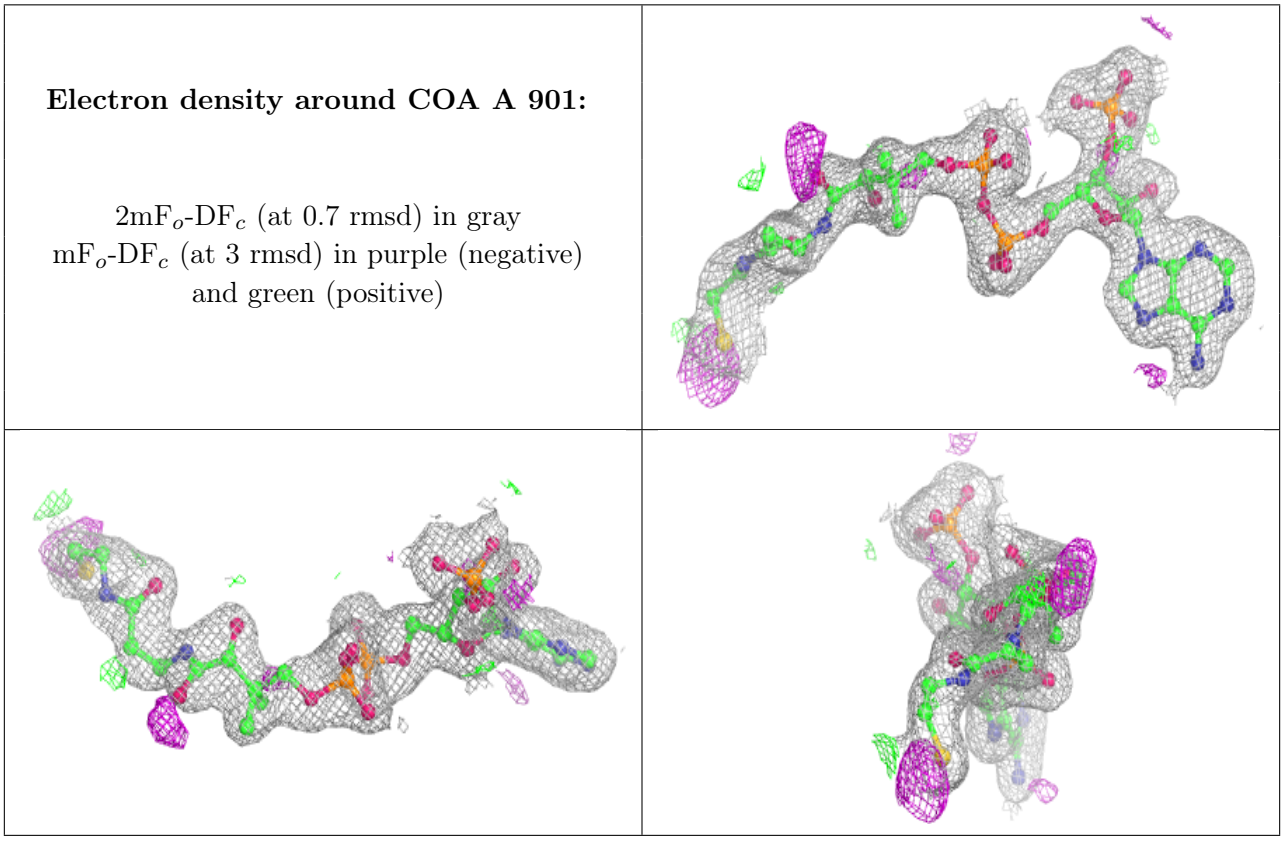
6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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.

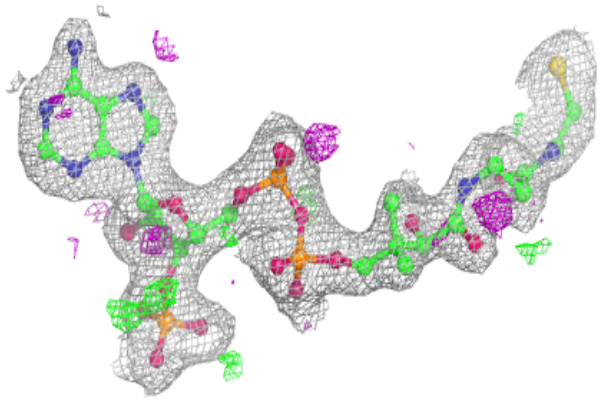
Electron density around COA A 901:

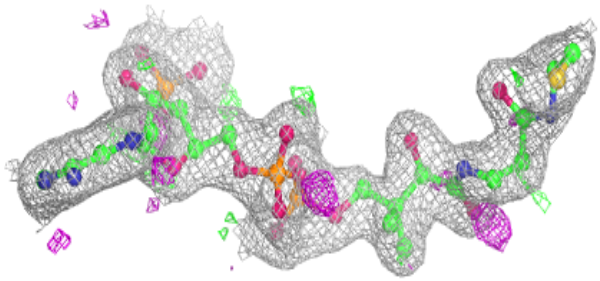
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

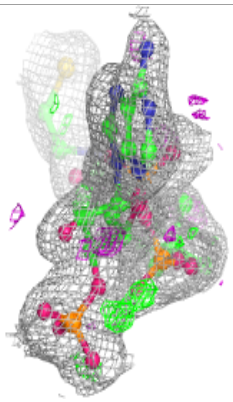


Electron density around COA B 901:

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

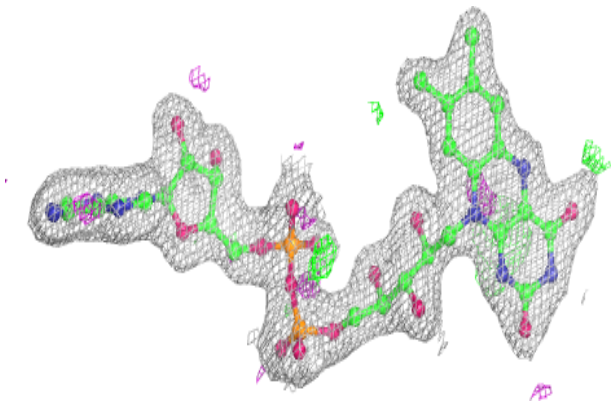


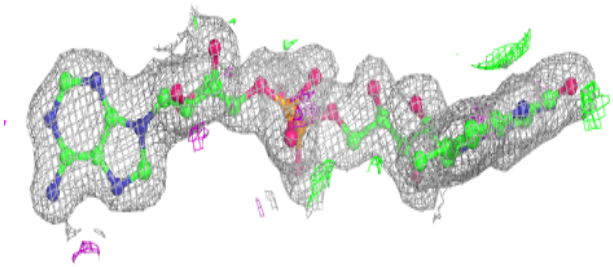


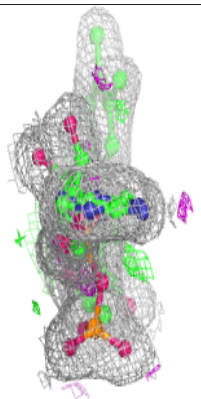


Electron density around FAD A 900:

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

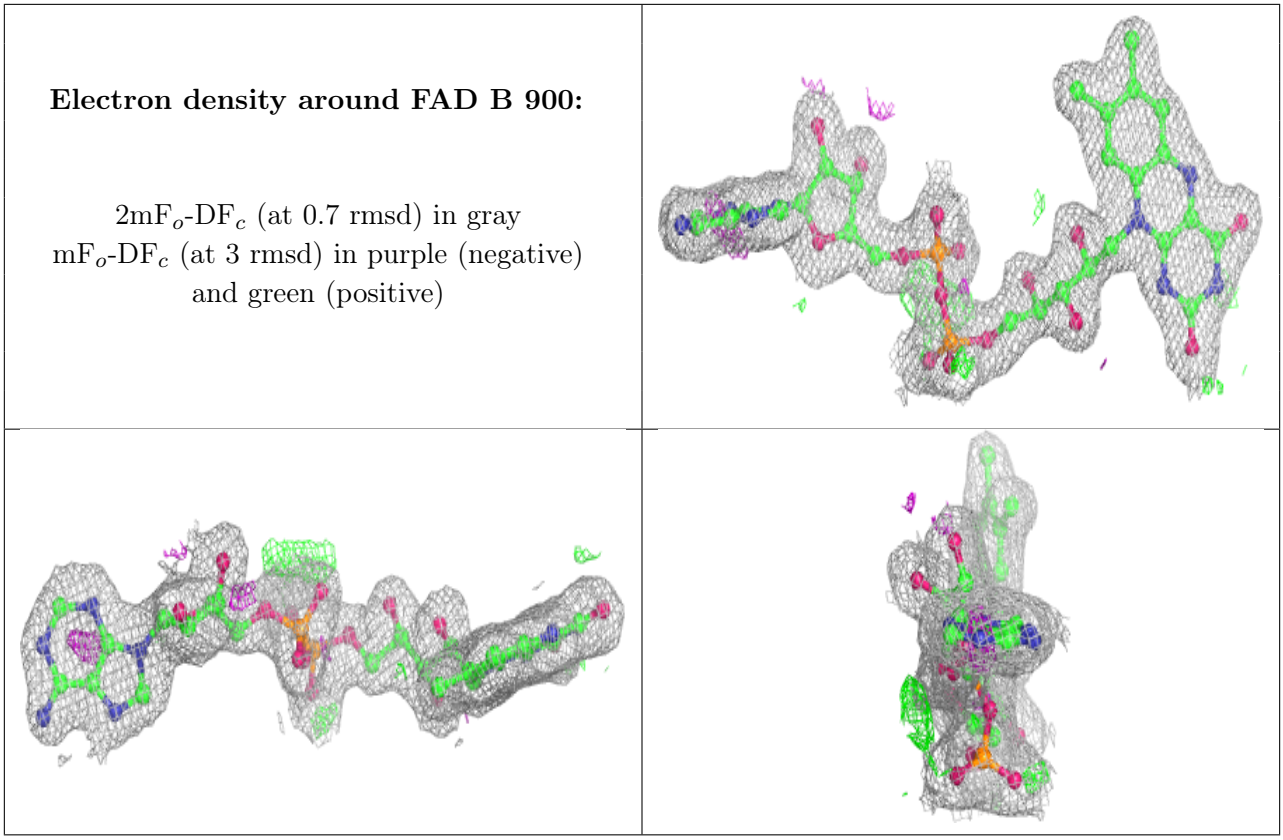






Electron density around FAD B 900:

$2mF_o$-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)

Unable to reproduce the depositors R factor - this section is therefore empty.