



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 04:41 AM UTC

PDB ID : 4FX9 / pdb_00004fx9
Title : Structure of the Pyrococcus horikoshii CoA persulfide/polysulfide reductase
Authors : Herwald, S.; Lopez, K.M.; Crane III, E.J.; Sazinsky, M.H.
Deposited on : 2012-07-02
Resolution : 2.70 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

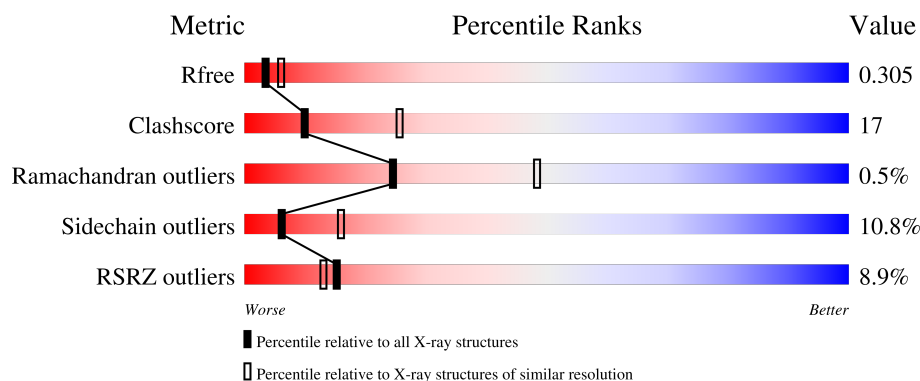
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	453	11% 62% 28% 7% .
1	B	453	6% 71% 23% . .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6918 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 Coenzyme A disulfide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	442	Total	C	N	O	S	0	0	0
			3341	2148	555	625	13			
1	B	443	Total	C	N	O	S	0	0	0
			3356	2160	563	620	13			

There are 16 discrepancies between the modelled and reference sequences:

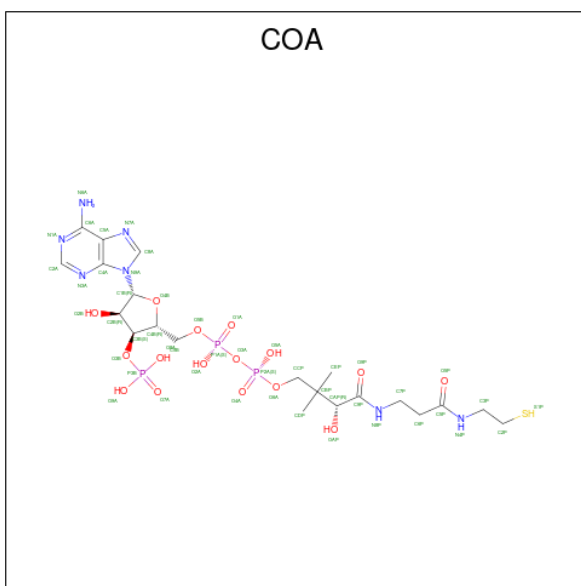
Chain	Residue	Modelled	Actual	Comment	Reference
A	446	LEU	-	expression tag	UNP O58308
A	447	GLU	-	expression tag	UNP O58308
A	448	HIS	-	expression tag	UNP O58308
A	449	HIS	-	expression tag	UNP O58308
A	450	HIS	-	expression tag	UNP O58308
A	451	HIS	-	expression tag	UNP O58308
A	452	HIS	-	expression tag	UNP O58308
A	453	HIS	-	expression tag	UNP O58308
B	446	LEU	-	expression tag	UNP O58308
B	447	GLU	-	expression tag	UNP O58308
B	448	HIS	-	expression tag	UNP O58308
B	449	HIS	-	expression tag	UNP O58308
B	450	HIS	-	expression tag	UNP O58308
B	451	HIS	-	expression tag	UNP O58308
B	452	HIS	-	expression tag	UNP O58308
B	453	HIS	-	expression tag	UNP O58308

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: $\text{C}_{21}\text{H}_{36}\text{N}_7\text{O}_{16}\text{P}_3\text{S}$).



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

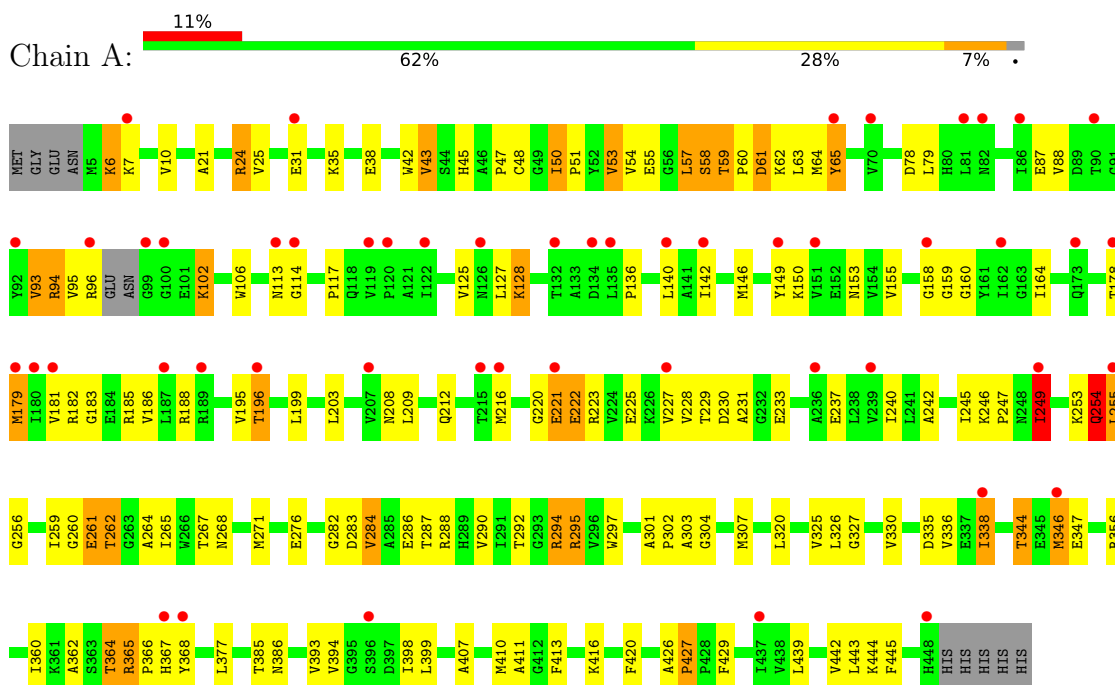
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	10	Total	O	0	0
			10	10		
4	B	9	Total	O	0	0
			9	9		

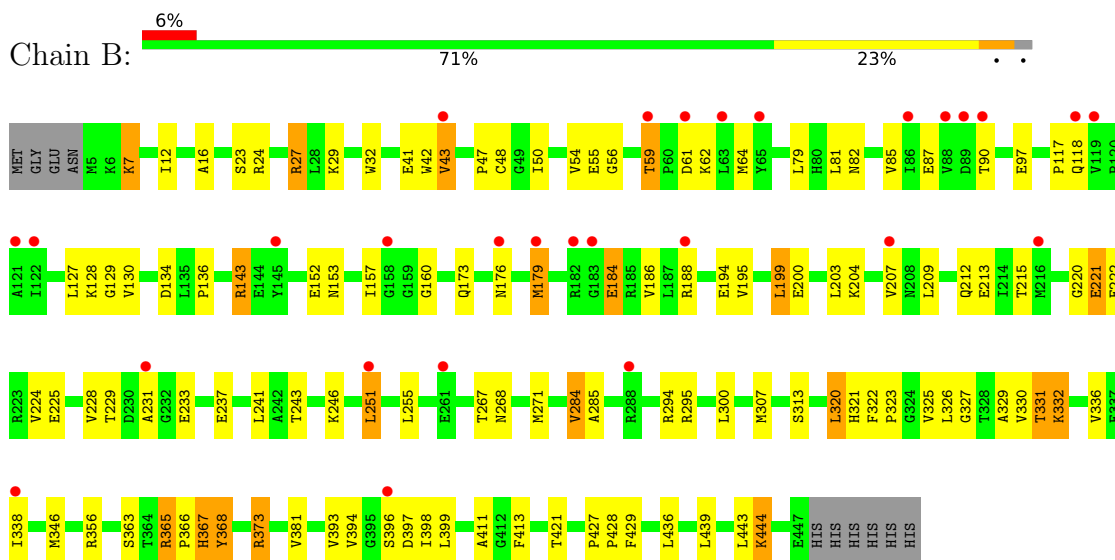
3 Residue-property plots [i](#)

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: Coenzyme A disulfide reductase



- Molecule 1: Coenzyme A disulfide reductase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	133.54Å 133.54Å 305.08Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.70 30.00 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-2.70) 99.8 (30.00-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	14.46 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.259 , 0.310 0.262 , 0.305	Depositor DCC
R_{free} test set	1480 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	56.3	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 34.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6918	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.95% 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, COA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.94	5/3410 (0.1%)	0.96	14/4642 (0.3%)
1	B	0.80	1/3426 (0.0%)	0.87	6/4657 (0.1%)
All	All	0.87	6/6836 (0.1%)	0.92	20/9299 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	50	ILE	CA-CB	-9.36	1.49	1.54
1	A	65	TYR	C-O	-6.19	1.18	1.23
1	B	160	GLY	C-O	-5.92	1.18	1.24
1	A	362	ALA	CA-CB	-5.21	1.45	1.53
1	A	65	TYR	CA-C	-5.13	1.48	1.53
1	A	365	ARG	C-O	-5.02	1.19	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	222	GLU	N-CA-C	-8.60	102.16	114.12
1	B	184	GLU	N-CA-C	8.03	120.03	111.28
1	A	50	ILE	CB-CA-C	-7.79	106.30	113.70
1	B	127	LEU	N-CA-C	-6.52	102.39	110.41
1	A	128	LYS	N-CA-C	-6.41	101.21	110.24
1	B	128	LYS	N-CA-C	-5.98	100.81	110.32
1	A	61	ASP	N-CA-C	-5.87	106.12	113.28
1	A	6	LYS	N-CA-C	-5.70	103.00	110.53
1	A	335	ASP	N-CA-C	5.68	119.68	112.87
1	A	336	VAL	N-CA-C	5.50	115.78	107.80
1	A	427	PRO	N-CA-C	5.49	117.40	110.70
1	A	136	PRO	N-CA-C	5.43	117.33	110.70
1	A	249	ILE	N-CA-CB	-5.39	105.30	112.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	VAL	CB-CA-C	-5.24	105.22	112.24
1	B	222	GLU	N-CA-C	-5.17	106.47	112.89
1	B	136	PRO	N-CA-C	5.14	116.97	110.70
1	A	427	PRO	CA-C-N	-5.13	115.60	120.83
1	A	427	PRO	C-N-CA	-5.13	115.60	120.83
1	A	102	LYS	N-CA-C	5.12	118.04	110.46
1	B	7	LYS	N-CA-C	5.06	117.20	110.53

There are no chirality outliers.

There are no planarity outliers.

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	3341	0	3291	151	0
1	B	3356	0	3334	96	0
2	A	53	0	31	1	0
2	B	53	0	31	0	0
3	A	48	0	32	6	0
3	B	48	0	32	4	0
4	A	10	0	0	0	0
4	B	9	0	0	1	0
All	All	6918	0	6751	237	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (237) 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:149:TYR:CE1	1:A:223:ARG:HD2	1.70	1.25
1:A:221:GLU:CG	1:A:222:GLU:H	1.55	1.15
1:A:195:VAL:HG13	1:A:394:VAL:CG2	1.76	1.15
1:A:195:VAL:HG13	1:A:394:VAL:HG21	1.31	1.08
1:A:221:GLU:HG2	1:A:222:GLU:H	0.94	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:LEU:HD12	1:A:255:LEU:N	1.68	1.05
1:B:443:LEU:O	1:B:444:LYS:HB2	1.52	1.04
1:A:221:GLU:HG2	1:A:222:GLU:N	1.65	1.02
1:A:221:GLU:HG2	1:A:223:ARG:H	1.28	0.98
1:A:94:ARG:HG3	1:A:94:ARG:HH11	1.26	0.98
1:B:213:GLU:CD	1:B:231:ALA:CB	2.37	0.97
1:B:213:GLU:OE2	1:B:231:ALA:HB3	1.68	0.94
1:A:199:LEU:HD11	1:A:338:ILE:HD11	1.49	0.93
1:A:199:LEU:CD1	1:A:338:ILE:CD1	2.46	0.93
1:A:255:LEU:N	1:A:255:LEU:CD1	2.30	0.92
1:B:213:GLU:HG2	1:B:231:ALA:HB2	1.54	0.90
1:B:195:VAL:HG13	1:B:394:VAL:CG2	2.02	0.90
1:A:199:LEU:HD12	1:A:338:ILE:CD1	2.01	0.89
1:A:48:CYS:SG	3:A:902:COA:S1P	2.55	0.88
1:A:195:VAL:CG1	1:A:394:VAL:HG21	2.03	0.88
1:B:195:VAL:HG13	1:B:394:VAL:HG21	1.55	0.86
1:B:179:MET:HE2	1:B:207:VAL:HG11	1.58	0.86
1:B:48:CYS:HG	3:B:902:COA:HS1	1.22	0.84
1:B:213:GLU:CD	1:B:231:ALA:HB3	1.99	0.83
1:A:195:VAL:HG13	1:A:394:VAL:HG23	1.58	0.83
1:A:195:VAL:CG1	1:A:394:VAL:CG2	2.56	0.83
1:B:284:VAL:HG22	4:B:1008:HOH:O	1.78	0.82
1:A:159:GLY:HA3	1:A:181:VAL:HG22	1.62	0.81
1:A:149:TYR:CD1	1:A:223:ARG:HD2	2.16	0.79
1:A:199:LEU:HD11	1:A:338:ILE:CD1	2.11	0.79
1:A:113:ASN:HB3	1:A:284:VAL:HG11	1.63	0.79
1:B:325:VAL:HG12	1:B:327:GLY:H	1.48	0.77
1:A:325:VAL:HG12	1:A:327:GLY:H	1.49	0.77
1:A:199:LEU:CD1	1:A:338:ILE:HD13	2.14	0.76
1:A:199:LEU:HD12	1:A:338:ILE:HD13	1.67	0.76
1:B:199:LEU:HD23	1:B:336:VAL:HG11	1.70	0.74
1:A:199:LEU:CD1	1:A:338:ILE:HD11	2.10	0.74
1:A:6:LYS:HG3	1:A:7:LYS:N	2.03	0.74
1:A:254:GLN:C	1:A:255:LEU:HD12	2.13	0.74
1:A:127:LEU:O	1:A:128:LYS:C	2.30	0.73
1:B:213:GLU:CG	1:B:231:ALA:HB2	2.19	0.72
1:B:179:MET:CE	1:B:207:VAL:HG11	2.20	0.71
1:A:94:ARG:HG3	1:A:94:ARG:NH1	2.00	0.70
1:B:43:VAL:HG22	1:B:81:LEU:HD21	1.72	0.70
1:A:21:ALA:O	1:A:25:VAL:HG23	1.91	0.70
1:A:443:LEU:O	1:A:444:LYS:CB	2.38	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:GLY:C	1:A:262:THR:H	2.00	0.69
1:B:199:LEU:HD23	1:B:336:VAL:CG1	2.23	0.69
1:A:393:VAL:HG12	1:A:398:ILE:CG2	2.22	0.69
1:B:50:ILE:O	1:B:54:VAL:HG23	1.91	0.68
1:B:43:VAL:HG11	1:B:79:LEU:HD11	1.75	0.68
1:B:157:ILE:HG22	1:B:243:THR:HG23	1.75	0.68
1:B:320:LEU:HD23	1:B:321:HIS:H	1.59	0.67
1:A:262:THR:HG22	1:A:264:ALA:H	1.59	0.67
1:A:346:MET:HG3	1:A:347:GLU:N	2.02	0.66
1:B:220:GLY:HA2	1:B:225:GLU:HG2	1.76	0.66
1:A:249:ILE:HD11	1:A:265:ILE:HG13	1.78	0.66
1:A:262:THR:HG21	1:A:297:TRP:HB2	1.77	0.65
1:B:48:CYS:SG	3:B:902:COA:S1P	2.87	0.64
1:B:157:ILE:HG22	1:B:243:THR:CG2	2.27	0.64
1:B:130:VAL:CG1	1:B:241:LEU:HD11	2.28	0.64
1:A:229:THR:C	1:A:231:ALA:H	2.05	0.64
1:B:195:VAL:CG1	1:B:394:VAL:HG21	2.27	0.63
1:A:159:GLY:CA	1:A:181:VAL:HG22	2.26	0.63
1:A:221:GLU:CG	1:A:222:GLU:N	2.29	0.63
1:A:377:LEU:HG	1:A:393:VAL:HG22	1.80	0.63
1:A:50:ILE:HB	1:A:51:PRO:HD3	1.81	0.62
1:A:199:LEU:HD23	1:A:199:LEU:O	1.99	0.62
1:B:367:HIS:O	1:B:367:HIS:ND1	2.32	0.62
1:A:262:THR:HB	1:A:286:GLU:OE2	2.00	0.61
1:B:195:VAL:HG13	1:B:394:VAL:HG23	1.81	0.61
1:A:255:LEU:CD1	1:A:255:LEU:H	2.11	0.60
1:A:57:LEU:HD23	1:A:57:LEU:N	2.16	0.60
1:B:251:LEU:HD23	1:B:251:LEU:H	1.67	0.60
1:A:58:SER:OG	1:A:59:THR:N	2.29	0.60
1:B:267:THR:O	1:B:295:ARG:NH1	2.34	0.60
1:A:45:HIS:HA	1:A:63:LEU:O	2.02	0.59
1:A:53:VAL:HG21	1:A:63:LEU:HD11	1.85	0.58
1:A:255:LEU:O	1:A:256:GLY:C	2.46	0.58
1:A:199:LEU:HD12	1:A:338:ILE:CG1	2.33	0.58
1:A:153:ASN:HB2	1:A:237:GLU:OE1	2.03	0.58
1:A:260:GLY:C	1:A:262:THR:N	2.57	0.57
1:A:60:PRO:HG3	1:A:140:LEU:HD21	1.86	0.57
1:B:271:MET:HE3	1:B:322:PHE:HB3	1.87	0.57
1:B:179:MET:HE1	1:B:203:LEU:HD22	1.85	0.57
1:B:393:VAL:CG1	1:B:398:ILE:HB	2.35	0.57
1:B:200:GLU:O	1:B:204:LYS:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:159:GLY:HA3	1:A:181:VAL:CG2	2.32	0.56
1:B:320:LEU:HD23	1:B:321:HIS:N	2.20	0.56
1:A:93:VAL:HG22	1:A:106:TRP:HD1	1.71	0.56
1:A:364:THR:HG23	1:A:364:THR:O	2.04	0.56
1:A:45:HIS:HD2	1:A:47:PRO:HD3	1.69	0.56
1:A:262:THR:CG2	1:A:297:TRP:HB2	2.35	0.56
1:B:157:ILE:HD13	1:B:215:THR:HG21	1.86	0.56
1:B:336:VAL:HG12	1:B:338:ILE:HD11	1.88	0.56
1:A:117:PRO:HA	1:A:245:ILE:HG22	1.89	0.55
1:B:393:VAL:HG12	1:B:398:ILE:CG2	2.36	0.55
1:A:292:THR:HG22	1:A:294:ARG:HB2	1.88	0.55
1:A:221:GLU:HG2	1:A:223:ARG:N	2.10	0.55
1:A:43:VAL:HG21	1:A:79:LEU:HD11	1.88	0.55
1:A:199:LEU:HD12	1:A:338:ILE:HG12	1.89	0.55
1:A:346:MET:CG	1:A:347:GLU:N	2.69	0.55
1:A:413:PHE:CE1	1:B:411:ALA:HB2	2.42	0.55
1:B:16:ALA:HB2	3:B:902:COA:H31	1.89	0.55
1:B:195:VAL:CG1	1:B:394:VAL:CG2	2.81	0.55
1:B:130:VAL:HG12	1:B:241:LEU:HD11	1.88	0.54
1:A:54:VAL:O	1:A:55:GLU:C	2.49	0.53
1:B:228:VAL:HG22	1:B:233:GLU:HG2	1.90	0.53
1:B:373:ARG:HB2	1:B:396:SER:OG	2.09	0.53
1:A:229:THR:C	1:A:231:ALA:N	2.66	0.53
1:A:181:VAL:HG12	1:A:183:GLY:H	1.73	0.53
1:A:253:LYS:O	1:A:254:GLN:C	2.50	0.52
1:B:365:ARG:HB3	1:B:366:PRO:HD2	1.91	0.52
1:B:184:GLU:H	1:B:188:ARG:HH11	1.57	0.52
1:A:268:ASN:O	1:A:271:MET:HE2	2.08	0.51
1:A:426:ALA:HB3	1:A:429:PHE:HD1	1.76	0.51
1:A:222:GLU:O	1:A:222:GLU:HG3	2.10	0.51
1:B:12:ILE:HG12	1:B:85:VAL:HG21	1.93	0.51
1:B:365:ARG:NH2	1:B:397:ASP:OD2	2.44	0.51
1:A:164:ILE:HD11	1:A:186:VAL:CG1	2.41	0.51
1:A:393:VAL:CG1	1:A:398:ILE:HB	2.40	0.51
1:A:260:GLY:O	1:A:262:THR:N	2.42	0.51
1:B:42:TRP:HA	1:B:64:MET:HE2	1.92	0.51
1:B:215:THR:HA	1:B:229:THR:HG22	1.92	0.51
1:A:114:GLY:HA2	1:A:283:ASP:HB2	1.91	0.51
1:A:158:GLY:HA3	1:A:242:ALA:O	2.11	0.51
1:A:325:VAL:HG12	1:A:326:LEU:N	2.24	0.51
1:A:60:PRO:CG	1:A:140:LEU:HD21	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:130:VAL:HG12	1:B:241:LEU:CD1	2.41	0.50
1:B:393:VAL:HG11	1:B:398:ILE:HB	1.92	0.50
1:A:10:VAL:HG11	1:A:93:VAL:HG21	1.93	0.50
1:A:185:ARG:HD2	1:A:196:THR:HG21	1.93	0.50
1:B:367:HIS:ND1	1:B:367:HIS:C	2.69	0.50
1:A:247:PRO:HB2	1:A:249:ILE:HD13	1.93	0.50
1:A:411:ALA:HB2	1:B:413:PHE:CE2	2.47	0.50
1:B:56:GLY:CA	1:B:143:ARG:HH22	2.25	0.50
1:B:325:VAL:HG12	1:B:326:LEU:N	2.26	0.50
1:A:178:THR:HG22	1:A:208:ASN:HB3	1.92	0.50
1:A:62:LYS:HA	1:B:367:HIS:CE1	2.47	0.49
1:A:267:THR:O	1:A:295:ARG:NH1	2.45	0.49
1:B:47:PRO:HD2	3:B:902:COA:H22	1.95	0.49
1:B:220:GLY:O	1:B:221:GLU:O	2.29	0.49
1:B:268:ASN:O	1:B:271:MET:HE2	2.12	0.49
1:A:393:VAL:HG12	1:A:398:ILE:HG22	1.94	0.49
1:A:93:VAL:HG22	1:A:106:TRP:CD1	2.47	0.49
1:A:53:VAL:CG2	1:A:63:LEU:HD21	2.43	0.49
1:A:60:PRO:HG3	1:A:140:LEU:CD2	2.43	0.49
1:B:23:SER:O	1:B:27:ARG:HB2	2.11	0.49
1:A:155:VAL:HG11	1:A:227:VAL:HG21	1.94	0.49
1:B:213:GLU:CG	1:B:231:ALA:CB	2.84	0.48
1:A:253:LYS:O	1:A:255:LEU:N	2.47	0.48
1:A:255:LEU:H	1:A:255:LEU:HD13	1.76	0.48
1:B:41:GLU:HG2	1:B:82:ASN:HD21	1.78	0.48
1:B:153:ASN:N	1:B:237:GLU:OE2	2.37	0.48
1:A:228:VAL:HG22	1:A:233:GLU:HG2	1.96	0.47
1:A:59:THR:OG1	1:A:61:ASP:OD1	2.30	0.47
1:A:38:GLU:HG2	1:A:42:TRP:O	2.14	0.47
1:A:149:TYR:CE1	1:A:223:ARG:CD	2.66	0.47
1:A:114:GLY:O	1:A:284:VAL:HG13	2.15	0.47
1:A:410:MET:HG3	1:B:421:THR:HG23	1.96	0.47
1:A:413:PHE:HE1	1:B:411:ALA:HB2	1.77	0.47
1:B:184:GLU:HA	1:B:212:GLN:HE21	1.80	0.47
1:A:229:THR:O	1:A:231:ALA:N	2.48	0.47
1:A:42:TRP:HA	1:A:64:MET:HE2	1.97	0.46
1:A:356:ARG:NH2	1:A:386:ASN:OD1	2.48	0.46
1:B:179:MET:HE2	1:B:179:MET:HB2	1.66	0.46
1:B:300:LEU:HD11	1:B:329:ALA:HB2	1.98	0.45
1:A:48:CYS:SG	2:A:901:FAD:C4X	3.04	0.45
1:A:292:THR:CG2	1:A:294:ARG:HB2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:ALA:CB	3:A:902:COA:H21	2.47	0.45
1:A:366:PRO:HG3	1:A:427:PRO:HA	1.98	0.45
1:B:381:VAL:HG21	1:B:443:LEU:HD13	1.98	0.45
1:A:216:MET:HE2	1:A:230:ASP:HA	1.98	0.45
1:B:129:GLY:HA3	1:B:224:VAL:HG12	1.98	0.45
1:A:160:GLY:O	1:A:164:ILE:HG13	2.16	0.45
1:A:185:ARG:HG3	1:A:188:ARG:HB3	1.99	0.45
1:B:130:VAL:HG11	1:B:241:LEU:HD11	1.98	0.45
1:A:364:THR:O	1:A:364:THR:CG2	2.57	0.44
1:A:393:VAL:CG1	1:A:398:ILE:HG22	2.47	0.44
1:A:230:ASP:OD2	1:A:230:ASP:C	2.58	0.44
1:A:393:VAL:CG1	1:A:398:ILE:CG2	2.94	0.44
1:A:360:ILE:HG21	1:A:442:VAL:HG21	1.99	0.43
1:A:429:PHE:CE1	1:B:330:VAL:HG12	2.54	0.43
1:B:59:THR:O	1:B:62:LYS:HB2	2.18	0.43
1:A:411:ALA:HB2	1:B:413:PHE:HE2	1.82	0.43
1:B:199:LEU:HG	1:B:338:ILE:HG13	2.00	0.43
1:B:427:PRO:HB2	1:B:428:PRO:HD3	2.00	0.43
1:A:45:HIS:CD2	1:A:47:PRO:HD3	2.52	0.43
1:A:95:VAL:HG12	1:A:96:ARG:N	2.33	0.43
1:A:253:LYS:C	1:A:255:LEU:N	2.74	0.43
1:B:285:ALA:HB3	1:B:307:MET:CE	2.48	0.43
1:B:325:VAL:CG1	1:B:326:LEU:N	2.82	0.43
1:B:331:THR:HG22	1:B:332:LYS:H	1.83	0.43
1:B:356:ARG:NH2	1:B:444:LYS:O	2.50	0.43
1:A:47:PRO:HD2	3:A:902:COA:S1P	2.59	0.43
1:A:128:LYS:CB	1:A:222:GLU:O	2.67	0.43
1:A:290:VAL:HB	1:A:326:LEU:HD23	2.01	0.43
1:B:322:PHE:HA	1:B:323:PRO:HD3	1.91	0.43
1:A:325:VAL:CG1	1:A:326:LEU:N	2.81	0.43
1:A:301:ALA:HB2	3:A:902:COA:H21	2.00	0.42
1:B:338:ILE:HD13	1:B:394:VAL:HG13	2.01	0.42
1:A:282:GLY:HA2	1:A:304:GLY:HA2	2.01	0.42
1:A:301:ALA:HB3	1:A:302:PRO:HD3	1.99	0.42
1:A:62:LYS:O	1:B:368:TYR:CD2	2.73	0.42
1:A:179:MET:HB3	1:A:179:MET:HE2	1.22	0.42
1:A:113:ASN:HB3	1:A:284:VAL:CG1	2.43	0.42
1:B:152:GLU:N	1:B:237:GLU:OE2	2.42	0.42
1:A:35:LYS:HG2	1:A:78:ASP:HB3	2.02	0.42
1:A:93:VAL:CG2	1:A:106:TRP:HD1	2.31	0.42
1:A:142:ILE:HG22	1:A:146:MET:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:416:LYS:HB2	1:A:445:PHE:CD1	2.56	0.41
1:A:24:ARG:HD2	1:A:24:ARG:O	2.19	0.41
1:A:420:PHE:CE1	1:B:320:LEU:HD21	2.55	0.41
1:B:118:GLN:HA	1:B:246:LYS:HE2	2.02	0.41
1:B:381:VAL:CG2	1:B:443:LEU:HD13	2.50	0.41
1:B:443:LEU:O	1:B:444:LYS:CB	2.36	0.41
1:A:426:ALA:HB3	1:A:429:PHE:CD1	2.53	0.41
1:A:385:THR:O	1:A:386:ASN:HB2	2.20	0.41
1:A:87:GLU:HA	1:A:254:GLN:HE22	1.85	0.41
1:A:344:THR:CG2	1:A:347:GLU:H	2.34	0.41
1:B:336:VAL:CG1	1:B:338:ILE:HD11	2.51	0.41
1:A:407:ALA:O	1:A:410:MET:HB2	2.21	0.41
1:B:7:LYS:HG3	1:B:32:TRP:CZ2	2.56	0.41
1:B:117:PRO:HD3	1:B:134:ASP:CG	2.46	0.41
1:A:65:TYR:HB2	3:A:902:COA:H71	2.03	0.40
1:A:303:ALA:O	1:A:307:MET:HG3	2.21	0.40
1:B:7:LYS:HG2	1:B:32:TRP:CE2	2.56	0.40
1:A:65:TYR:CD2	3:A:902:COA:H72	2.57	0.40
1:A:330:VAL:HG12	1:B:429:PHE:CE1	2.56	0.40
1:A:220:GLY:HA2	1:A:225:GLU:HG2	2.03	0.40
1:A:393:VAL:HG12	1:A:398:ILE:HG21	2.02	0.40
1:A:393:VAL:HG11	1:A:398:ILE:HB	2.04	0.40
1:B:285:ALA:HB3	1:B:307:MET:HE1	2.02	0.40
1:A:287:THR:OG1	1:A:288:ARG:N	2.54	0.40
1:A:344:THR:HG22	1:A:347:GLU:H	1.86	0.40
1:B:55:GLU:C	1:B:173:GLN:NE2	2.79	0.40
1:B:153:ASN:OD1	1:B:176:ASN:HB3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	438/453 (97%)	418 (95%)	18 (4%)	2 (0%)	24	48
1	B	441/453 (97%)	421 (96%)	18 (4%)	2 (0%)	24	48
All	All	879/906 (97%)	839 (95%)	36 (4%)	4 (0%)	24	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	221	GLU
1	A	261	GLU
1	B	444	LYS
1	A	254	GLN

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	336/368 (91%)	295 (88%)	41 (12%)	5	12
1	B	337/368 (92%)	305 (90%)	32 (10%)	8	20
All	All	673/736 (91%)	600 (89%)	73 (11%)	6	16

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	31	GLU
1	A	43	VAL
1	A	57	LEU
1	A	58	SER
1	A	59	THR
1	A	88	VAL
1	A	93	VAL
1	A	94	ARG
1	A	102	LYS
1	A	125	VAL
1	A	150	LYS

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Mol	Chain	Res	Type
1	A	179	MET
1	A	182	ARG
1	A	196	THR
1	A	203	LEU
1	A	209	LEU
1	A	212	GLN
1	A	221	GLU
1	A	240	ILE
1	A	246	LYS
1	A	249	ILE
1	A	254	GLN
1	A	255	LEU
1	A	259	ILE
1	A	261	GLU
1	A	262	THR
1	A	276	GLU
1	A	284	VAL
1	A	294	ARG
1	A	295	ARG
1	A	320	LEU
1	A	338	ILE
1	A	344	THR
1	A	346	MET
1	A	364	THR
1	A	365	ARG
1	A	367	HIS
1	A	368	TYR
1	A	399	LEU
1	A	439	LEU
1	B	24	ARG
1	B	27	ARG
1	B	29	LYS
1	B	43	VAL
1	B	59	THR
1	B	61	ASP
1	B	87	GLU
1	B	90	THR
1	B	97	GLU
1	B	143	ARG
1	B	179	MET
1	B	186	VAL
1	B	194	GLU

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Mol	Chain	Res	Type
1	B	199	LEU
1	B	209	LEU
1	B	251	LEU
1	B	255	LEU
1	B	284	VAL
1	B	294	ARG
1	B	313	SER
1	B	320	LEU
1	B	331	THR
1	B	332	LYS
1	B	346	MET
1	B	363	SER
1	B	365	ARG
1	B	367	HIS
1	B	368	TYR
1	B	373	ARG
1	B	399	LEU
1	B	436	LEU
1	B	439	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	HIS
1	A	206	HIS
1	A	254	GLN
1	A	277	ASN
1	B	82	ASN
1	B	206	HIS
1	B	212	GLN
1	B	321	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FAD	B	901	-	58,58,58	1.12	5 (8%)	85,89,89	1.55	16 (18%)
3	COA	B	902	-	47,50,50	1.72	5 (10%)	69,75,75	1.46	7 (10%)
2	FAD	A	901	-	58,58,58	1.10	5 (8%)	85,89,89	1.52	15 (17%)
3	COA	A	902	-	47,50,50	1.71	5 (10%)	69,75,75	1.42	8 (11%)

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
2	FAD	B	901	-	-	3/34/50/50	0/6/6/6
3	COA	B	902	-	-	7/48/64/64	0/3/3/3
2	FAD	A	901	-	-	1/34/50/50	0/6/6/6
3	COA	A	902	-	-	6/48/64/64	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	902	COA	O9P-C9P	9.65	1.41	1.23
3	B	902	COA	O9P-C9P	9.62	1.41	1.23
2	B	901	FAD	C4X-N5	3.81	1.39	1.30
2	A	901	FAD	C4X-N5	3.69	1.38	1.30
2	B	901	FAD	C10-N1	2.95	1.39	1.33
3	B	902	COA	C2A-N1A	2.91	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	901	FAD	C2A-N3A	2.81	1.38	1.33
3	A	902	COA	C2A-N1A	2.80	1.38	1.33
2	A	901	FAD	C10-N1	2.75	1.38	1.33
2	B	901	FAD	C2A-N1A	2.74	1.38	1.33
3	B	902	COA	C2A-N3A	2.73	1.38	1.33
2	A	901	FAD	C2A-N3A	2.73	1.38	1.33
2	A	901	FAD	C2A-N1A	2.69	1.38	1.33
3	A	902	COA	C2A-N3A	2.61	1.38	1.33
3	A	902	COA	C8A-N7A	2.21	1.35	1.31
2	A	901	FAD	C8A-N7A	2.19	1.35	1.31
2	B	901	FAD	C8A-N7A	2.18	1.35	1.31
3	A	902	COA	P2A-O3A	2.13	1.61	1.59
3	B	902	COA	C5A-N7A	-2.08	1.35	1.39
3	B	902	COA	C8A-N7A	2.03	1.35	1.31

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	FAD	N3A-C2A-N1A	-5.71	119.94	128.58
3	A	902	COA	N3A-C2A-N1A	-5.67	120.00	128.58
3	B	902	COA	N3A-C2A-N1A	-5.62	120.08	128.58
2	A	901	FAD	N3A-C2A-N1A	-5.60	120.11	128.58
3	B	902	COA	C5A-C4A-N3A	-4.78	120.13	126.72
2	B	901	FAD	C5A-C4A-N3A	-4.51	120.51	126.72
2	A	901	FAD	C5A-C4A-N3A	-4.43	120.61	126.72
3	A	902	COA	C5A-C4A-N3A	-4.28	120.83	126.72
2	B	901	FAD	N9A-C8A-N7A	-3.97	108.30	113.94
2	A	901	FAD	N9A-C8A-N7A	-3.86	108.45	113.94
3	A	902	COA	N9A-C8A-N7A	-3.77	108.58	113.94
3	B	902	COA	N3A-C4A-N9A	3.51	133.14	127.17
3	B	902	COA	C2A-N3A-C4A	3.49	120.35	111.83
2	B	901	FAD	C2A-N3A-C4A	3.49	120.35	111.83
2	B	901	FAD	C5A-N7A-C8A	3.48	108.92	103.45
3	B	902	COA	N9A-C8A-N7A	-3.41	109.10	113.94
2	A	901	FAD	C2A-N3A-C4A	3.40	120.14	111.83
2	A	901	FAD	C5A-N7A-C8A	3.40	108.80	103.45
3	A	902	COA	C2A-N3A-C4A	3.38	120.09	111.83
3	A	902	COA	C5A-N7A-C8A	3.32	108.66	103.45
3	B	902	COA	C5A-N7A-C8A	3.11	108.33	103.45
2	A	901	FAD	C4-N3-C2	-3.09	120.15	125.64
2	B	901	FAD	C4-N3-C2	-3.06	120.21	125.64
2	B	901	FAD	N3A-C4A-N9A	2.97	132.22	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	901	FAD	N3A-C4A-N9A	2.89	132.08	127.17
3	A	902	COA	N3A-C4A-N9A	2.76	131.85	127.17
2	B	901	FAD	C4X-C4-N3	2.71	120.15	113.25
2	A	901	FAD	C4X-C4-N3	2.60	119.86	113.25
2	A	901	FAD	C4X-C10-N10	2.57	120.16	116.48
2	A	901	FAD	C10-C4X-N5	-2.52	119.67	124.81
2	B	901	FAD	C4A-C5A-N7A	-2.46	107.78	110.58
2	A	901	FAD	C4A-C5A-N7A	-2.43	107.81	110.58
2	B	901	FAD	C9A-C5X-N5	-2.41	119.90	122.45
2	B	901	FAD	C10-C4X-N5	-2.40	119.91	124.81
3	A	902	COA	C4A-C5A-N7A	-2.35	107.90	110.58
2	B	901	FAD	C4X-C10-N10	2.29	119.77	116.48
2	B	901	FAD	O4-C4-C4X	-2.26	120.56	126.53
2	A	901	FAD	O4-C4-C4X	-2.24	120.62	126.53
2	B	901	FAD	C4X-C10-N1	-2.17	119.27	124.59
2	B	901	FAD	O4B-C1B-N9A	2.13	112.18	108.09
3	B	902	COA	C6A-C5A-C4A	2.13	120.08	117.18
2	B	901	FAD	C5X-C9A-N10	2.12	119.88	117.97
2	A	901	FAD	C9A-C5X-N5	-2.10	120.23	122.45
2	A	901	FAD	C4X-C10-N1	-2.06	119.53	124.59
3	A	902	COA	CDP-CBP-CAP	2.06	112.28	108.77
2	A	901	FAD	O4B-C1B-N9A	2.06	112.05	108.09

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	902	COA	C5B-O5B-P1A-O1A
3	A	902	COA	C5B-O5B-P1A-O2A
3	A	902	COA	C5B-O5B-P1A-O3A
3	B	902	COA	C5B-O5B-P1A-O2A
3	B	902	COA	C5B-O5B-P1A-O3A
3	B	902	COA	S1P-C2P-C3P-N4P
3	B	902	COA	C6P-C5P-N4P-C3P
3	B	902	COA	O5P-C5P-N4P-C3P
3	A	902	COA	P1A-O3A-P2A-O4A
2	B	901	FAD	O4B-C4B-C5B-O5B
2	B	901	FAD	O4'-C4'-C5'-O5'
3	B	902	COA	O4B-C4B-C5B-O5B
3	A	902	COA	P1A-O3A-P2A-O5A
2	A	901	FAD	O4B-C4B-C5B-O5B
2	B	901	FAD	C3B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
3	B	902	COA	O4B-C1B-N9A-C8A
3	A	902	COA	S1P-C2P-C3P-N4P

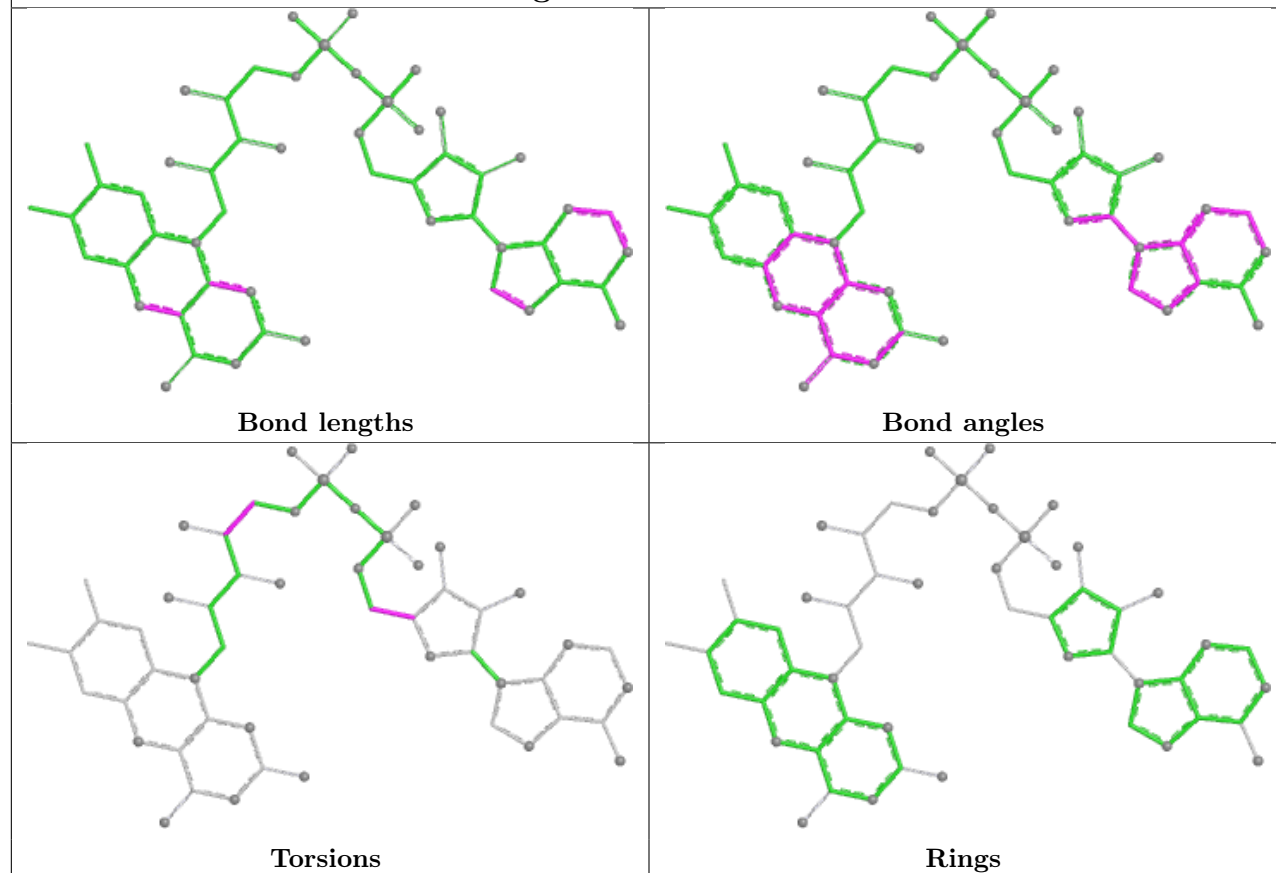
There are no ring outliers.

3 monomers are involved in 11 short contacts:

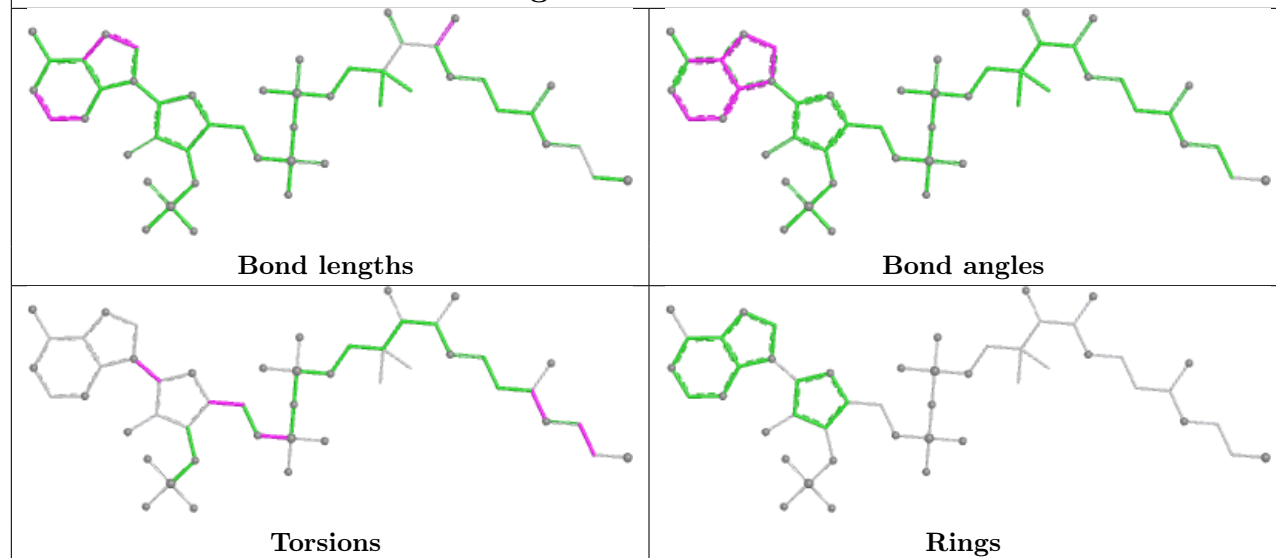
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	902	COA	4	0
2	A	901	FAD	1	0
3	A	902	COA	6	0

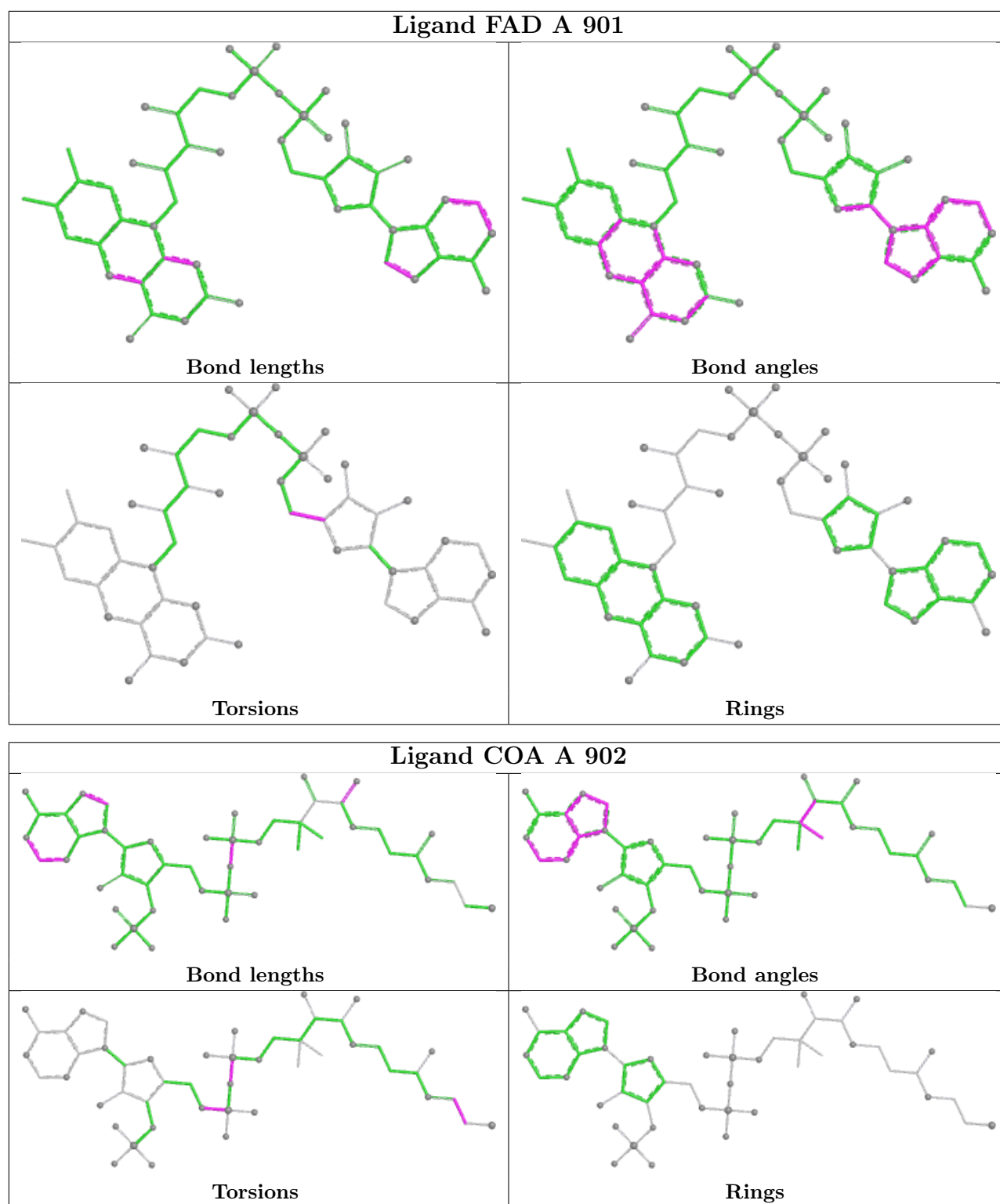
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

Ligand FAD B 901



Ligand COA B 902





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 ⓘ

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	442/453 (97%)	0.77	51 (11%) 9 8	20, 64, 94, 102	0
1	B	443/453 (97%)	0.69	28 (6%) 26 23	27, 62, 86, 97	0
All	All	885/906 (97%)	0.73	79 (8%) 15 13	20, 63, 90, 102	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	65	TYR	4.2
1	A	448	HIS	4.1
1	B	288	ARG	4.1
1	A	132	THR	3.5
1	A	179	MET	3.5
1	B	121	ALA	3.3
1	A	99	GLY	3.1
1	A	338	ILE	3.1
1	B	251	LEU	3.0
1	A	180	ILE	3.0
1	A	216	MET	3.0
1	A	158	GLY	2.8
1	A	215	THR	2.8
1	A	221	GLU	2.8
1	B	119	VAL	2.8
1	B	182	ARG	2.8
1	B	216	MET	2.8
1	B	261	GLU	2.7
1	A	119	VAL	2.7
1	B	89	ASP	2.7
1	A	65	TYR	2.6
1	A	173	GLN	2.6
1	A	81	LEU	2.6
1	A	122	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	239	VAL	2.5
1	B	61	ASP	2.5
1	A	346	MET	2.5
1	A	96	ARG	2.5
1	B	207	VAL	2.4
1	B	63	LEU	2.4
1	A	149	TYR	2.4
1	A	151	VAL	2.4
1	A	181	VAL	2.4
1	A	236	ALA	2.4
1	B	122	ILE	2.4
1	A	162	ILE	2.4
1	A	82	ASN	2.3
1	A	86	ILE	2.3
1	A	113	ASN	2.3
1	A	100	GLY	2.3
1	A	114	GLY	2.3
1	B	183	GLY	2.3
1	A	207	VAL	2.3
1	B	90	THR	2.3
1	A	249	ILE	2.2
1	B	338	ILE	2.2
1	A	196	THR	2.2
1	A	189	ARG	2.2
1	A	396	SER	2.2
1	B	158	GLY	2.2
1	A	187	LEU	2.2
1	A	7	LYS	2.2
1	A	92	TYR	2.2
1	B	179	MET	2.2
1	B	176	ASN	2.1
1	A	135	LEU	2.1
1	A	367	HIS	2.1
1	A	227	VAL	2.1
1	A	437	ILE	2.1
1	B	396	SER	2.1
1	B	231	ALA	2.1
1	A	31	GLU	2.1
1	A	178	THR	2.1
1	A	140	LEU	2.1
1	A	368	TYR	2.1
1	B	86	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	134	ASP	2.1
1	A	90	THR	2.1
1	B	59	THR	2.1
1	B	118	GLN	2.1
1	A	142	ILE	2.1
1	B	43	VAL	2.1
1	A	255	LEU	2.0
1	B	145	TYR	2.0
1	A	120	PRO	2.0
1	A	126	ASN	2.0
1	A	70	VAL	2.0
1	B	88	VAL	2.0
1	B	188	ARG	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

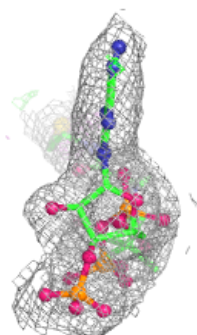
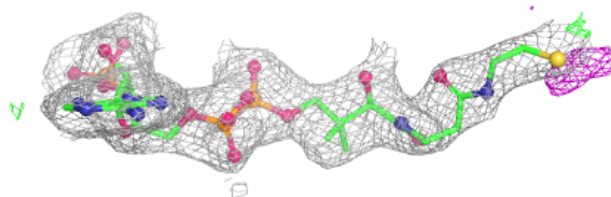
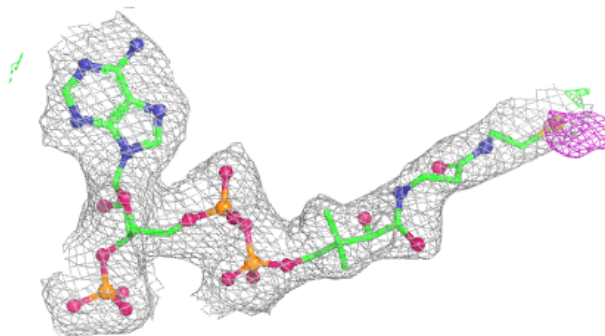
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	COA	A	902	48/48	0.92	0.09	39,51,57,61	0
3	COA	B	902	48/48	0.92	0.10	41,45,61,66	0
2	FAD	A	901	53/53	0.93	0.08	35,44,47,47	0
2	FAD	B	901	53/53	0.94	0.08	30,41,49,52	0

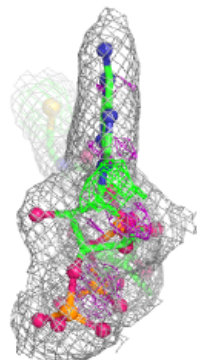
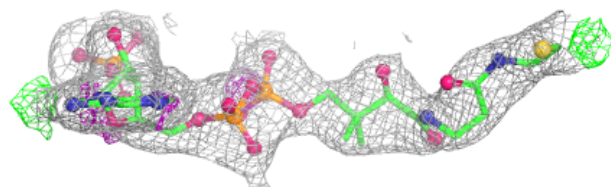
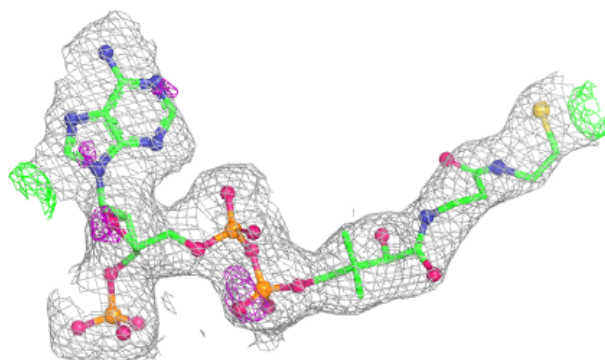
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around COA A 902:

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

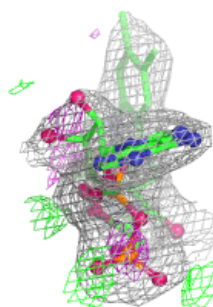
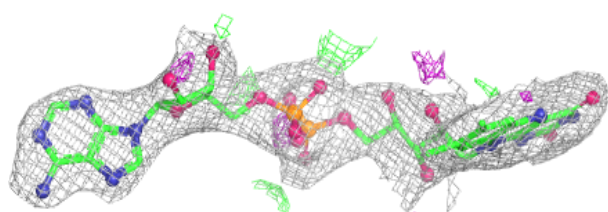
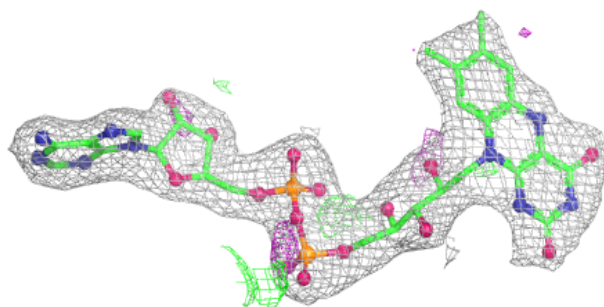
**Electron density around COA B 902:**

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

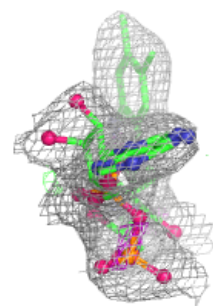
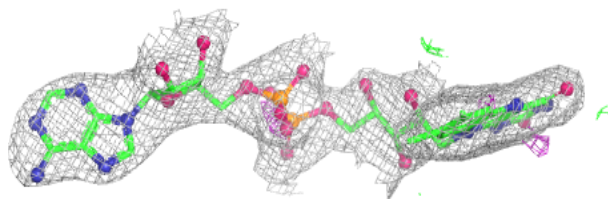
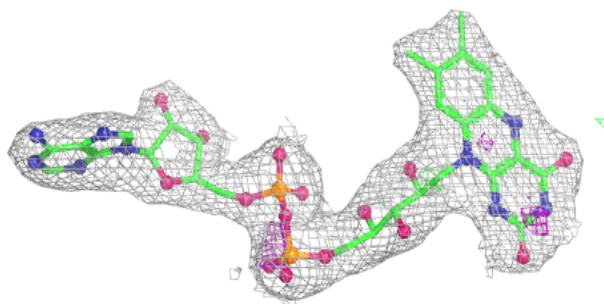


Electron density around FAD A 901:

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

**Electron density around FAD 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.