



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:29 AM UTC

PDB ID : 3AE4 / pdb_00003ae4
Title : Crystal structure of porcine heart mitochondrial complex II bound with 2-Iodo-N-methyl-benzamide
Authors : Harada, S.; Sasaki, T.; Shindo, M.; Kido, Y.; Inaoka, D.K.; Omori, J.; Osanai, A.; Sakamoto, K.; Mao, J.; Matsuoka, S.; Inoue, M.; Honma, T.; Tanaka, A.; Kita, K.
Deposited on : 2010-02-04
Resolution : 2.91 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](#) ①) 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

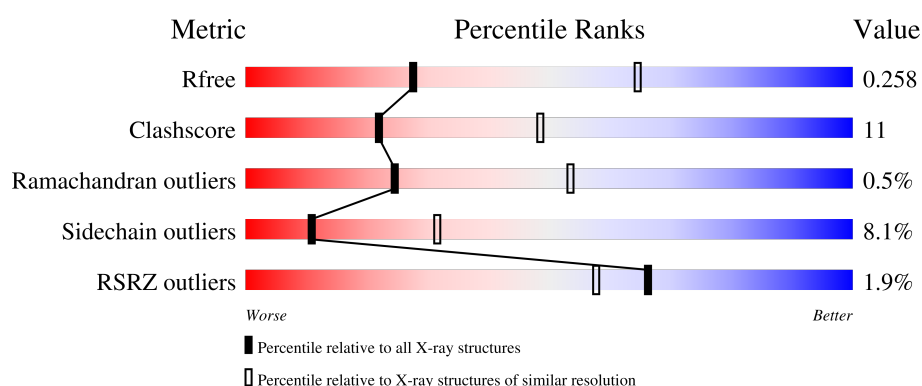
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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	2995 (2.94-2.90)
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)
RSRZ outliers	180081	2995 (2.94-2.90)

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	622	2% 73% 22% • •
2	B	252	% 66% 25% • 5%
3	C	140	3% 81% 16% • •
4	D	103	3% 88% 10% • •

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 8657 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 Succinate dehydrogenase [ubiquinone] flavoprotein subunit, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	613	Total	C	N	O	S	0	0	0
			4729	2954	848	895	32			

- Molecule 2 is a protein called Succinate dehydrogenase [ubiquinone] iron-sulfur subunit, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	239	Total	C	N	O	S	0	0	0
			1922	1214	326	360	22			

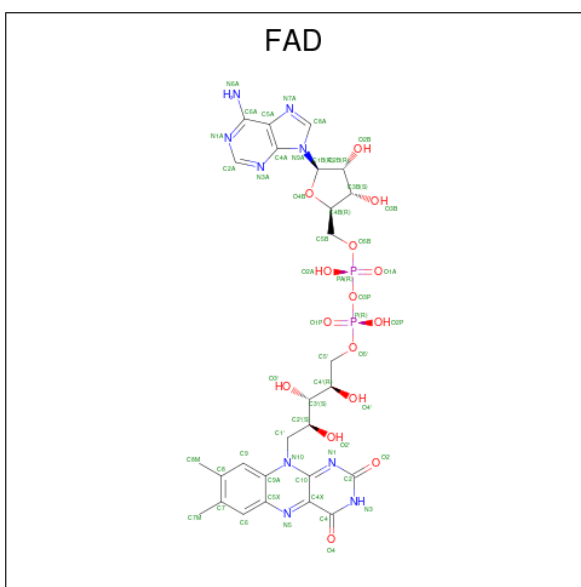
- Molecule 3 is a protein called Succinate dehydrogenase cytochrome b560 subunit, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	138	Total	C	N	O	S	0	0	0
			1064	695	179	183	7			

- Molecule 4 is a protein called Succinate dehydrogenase [ubiquinone] cytochrome b small subunit, mitochondrial.

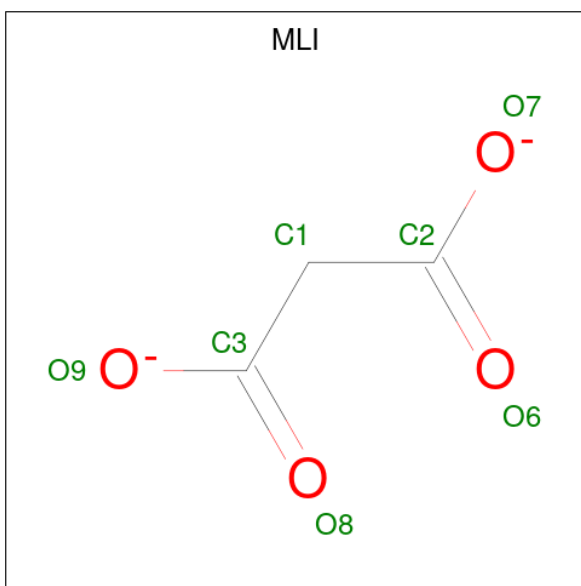
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	102	Total	C	N	O	S	0	0	0
			765	499	128	133	5			

- Molecule 5 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

- Molecule 6 is MALONATE ION (CCD ID: MLI) (formula: $\text{C}_3\text{H}_2\text{O}_4$).



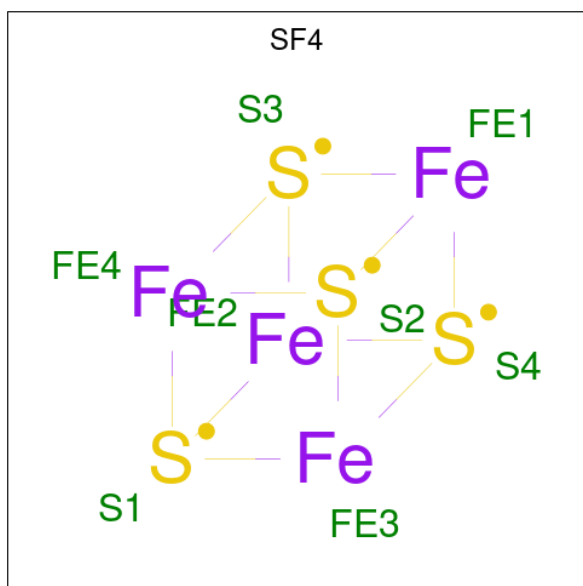
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	3	4		

- Molecule 7 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



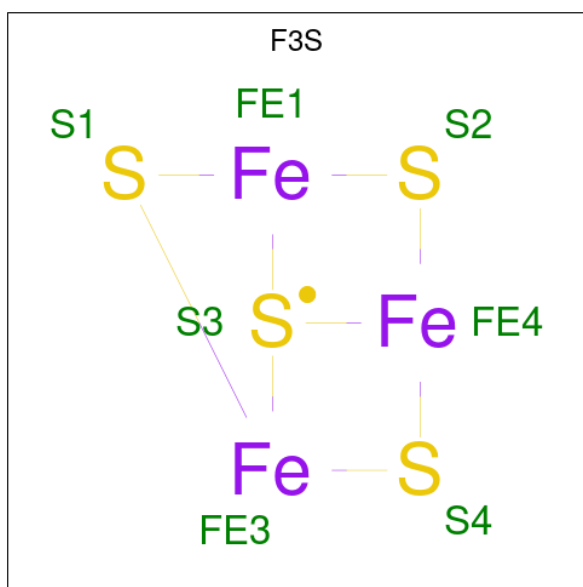
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 8 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe_4S_4).



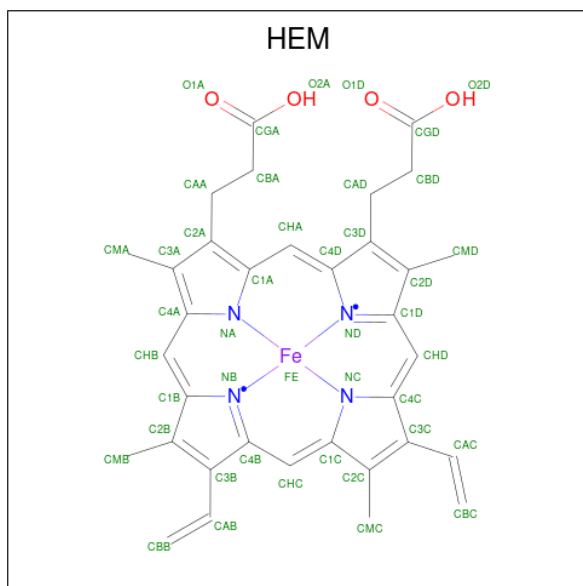
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	Fe	S	0	0
			8	4	4		

- Molecule 9 is FE3-S4 CLUSTER (CCD ID: F3S) (formula: Fe_3S_4).



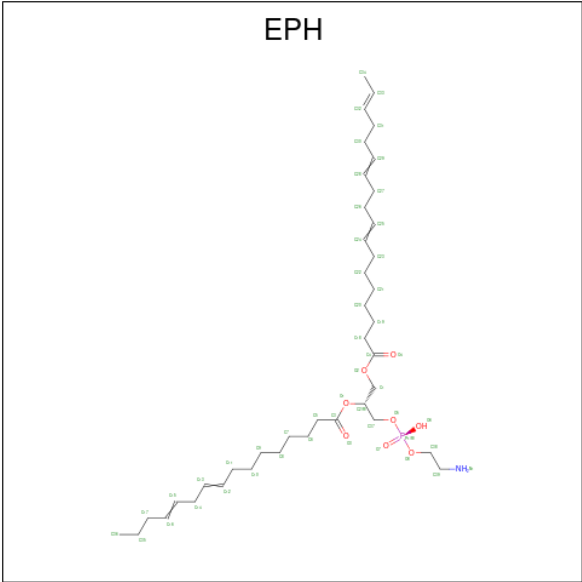
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 10 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



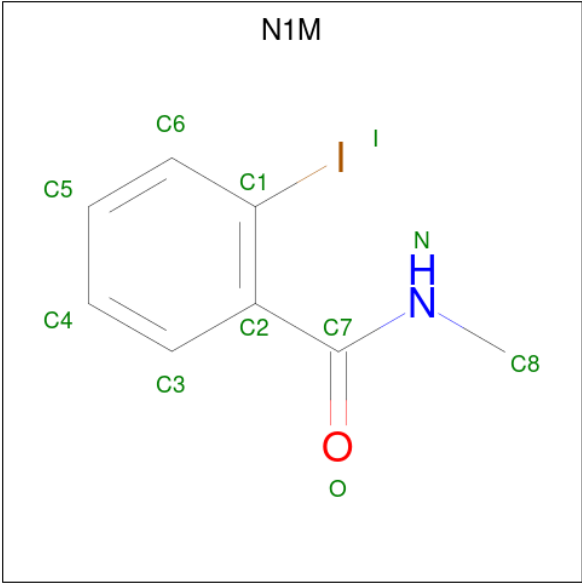
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
10	C	1	Total	C	Fe	N	O	
			43	34	1	4	4	

- Molecule 11 is L-ALPHA-PHOSPHATIDYL-BETA-OLEOYL-GAMMA-PALMITOYL-PHOSPHATIDYLETHANOLAMINE (CCD ID: EPH) (formula: $C_{39}H_{68}NO_8P$).

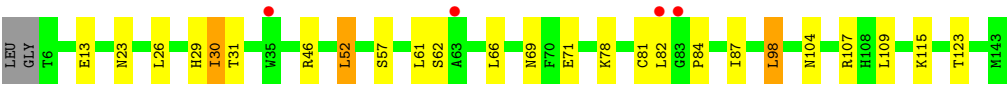


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
11	D	1	Total	C	N	O	P	0	0
			44	34	1	8	1		

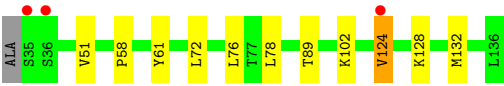
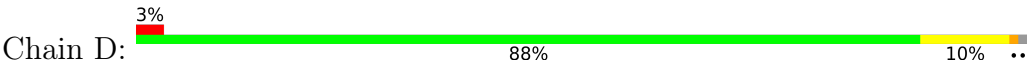
- Molecule 12 is 2-iodo-N-methylbenzamide (CCD ID: N1M) (formula: C₈H₈INO).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
12	D	1	Total	C	I	N	O	0	0
			11	8	1	1	1		



● Molecule 4: Succinate dehydrogenase [ubiquinone] cytochrome b small subunit, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	71.56Å 84.22Å 294.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.24 – 2.91 48.24 – 2.91	Depositor EDS
% Data completeness (in resolution range)	85.0 (48.24-2.91) 84.9 (48.24-2.91)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.19 (at 2.91Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.222 , 0.257 0.220 , 0.258	Depositor DCC
R_{free} test set	1682 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	67.2	Xtriage
Anisotropy	0.857	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 55.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8657	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: F3S, N1M, EPH, MLI, HEM, SF4, FAD, FES

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	0.47	0/4828	0.85	3/6531 (0.0%)
2	B	0.47	0/1964	0.86	4/2648 (0.2%)
3	C	0.46	0/1091	0.85	0/1483
4	D	0.45	0/784	0.89	0/1066
All	All	0.47	0/8667	0.85	7/11728 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	402	ALA	N-CA-C	-7.26	104.56	113.19
2	B	102	LEU	CA-C-N	6.33	126.35	119.89
2	B	102	LEU	C-N-CA	6.33	126.35	119.89
1	A	358	ILE	CB-CA-C	-6.30	105.08	110.94
2	B	69	ILE	CB-CA-C	-5.63	105.13	110.70
2	B	67	GLU	N-CA-C	-5.51	107.80	114.75
1	A	358	ILE	N-CA-C	5.30	113.19	108.63

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	4729	0	4618	116	0
2	B	1922	0	1900	47	0
3	C	1064	0	1104	16	0
4	D	765	0	773	8	0
5	A	53	0	31	14	0
6	A	7	0	2	1	0
7	B	4	0	0	0	0
8	B	8	0	0	0	0
9	B	7	0	0	0	0
10	C	43	0	30	3	0
11	D	44	0	53	2	0
12	D	11	0	8	0	0
All	All	8657	0	8519	182	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 11.

All (182) 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:57:HIS:NE2	5:A:700:FAD:HM82	1.04	1.35
1:A:57:HIS:CE1	5:A:700:FAD:HM82	1.88	1.08
2:B:102:LEU:HB3	2:B:166:THR:HG21	1.34	1.08
1:A:509:THR:HG23	1:A:512:ARG:HH12	1.30	0.97
1:A:61:ALA:HB3	1:A:155:GLY:HA3	1.49	0.93
1:A:150:VAL:H	1:A:154:THR:HG22	1.35	0.90
1:A:59:VAL:HG23	1:A:159:LEU:HD23	1.61	0.82
3:C:52:LEU:HB3	10:C:1305:HEM:HAC	1.62	0.80
1:A:179:ALA:N	5:A:700:FAD:N1A	2.31	0.77
1:A:254:HIS:HD2	1:A:256:THR:H	1.33	0.76
1:A:415:LEU:HG	5:A:700:FAD:C2	2.15	0.75
1:A:353:VAL:HG12	1:A:358:ILE:HD11	1.70	0.73
1:A:29:GLY:H	1:A:58:THR:HG21	1.52	0.71
1:A:373:THR:HG22	1:A:374:ASN:O	1.89	0.71
2:B:27:LYS:HB3	2:B:28:PRO:CD	2.21	0.70
4:D:72:LEU:O	4:D:76:LEU:HB2	1.91	0.69
1:A:361:LEU:HD12	1:A:362:PRO:HD2	1.73	0.69
1:A:418:LEU:HD11	5:A:700:FAD:H4'	1.75	0.68
1:A:254:HIS:CD2	1:A:256:THR:H	2.12	0.68
1:A:561:GLU:OE2	1:A:581:ARG:HD3	1.94	0.67
2:B:129:LEU:HD11	2:B:195:ARG:HB2	1.75	0.67
1:A:246:GLN:NE2	1:A:600:ARG:HE	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:THR:O	1:A:59:VAL:HG21	1.94	0.67
1:A:504:LEU:HD21	1:A:530:MET:HG3	1.76	0.66
3:C:29:HIS:HD2	3:C:30:ILE:HG13	1.60	0.66
1:A:405:HIS:ND1	1:A:409:ARG:HG3	2.10	0.66
3:C:71:GLU:HG2	4:D:132:MET:HE1	1.78	0.66
2:B:35:ILE:HD11	2:B:51:ILE:HG22	1.78	0.65
3:C:46:ARG:HE	10:C:1305:HEM:CGD	2.10	0.65
1:A:258:ILE:HD11	1:A:265:ILE:HG13	1.80	0.63
1:A:265:ILE:HD12	1:A:358:ILE:HG21	1.81	0.63
1:A:246:GLN:HA	1:A:584:THR:O	1.98	0.62
2:B:166:THR:O	2:B:166:THR:HG22	1.99	0.62
1:A:61:ALA:HA	5:A:700:FAD:N5	2.15	0.61
1:A:239:THR:HG23	1:A:588:VAL:HG13	1.82	0.61
1:A:150:VAL:N	1:A:154:THR:HG22	2.12	0.61
1:A:563:ASP:H	1:A:571:GLN:NE2	1.99	0.61
1:A:264:LEU:HD13	1:A:365:HIS:CE1	2.36	0.60
1:A:258:ILE:HG22	1:A:260:GLY:H	1.67	0.60
1:A:116:MET:HA	1:A:161:THR:HG21	1.84	0.59
1:A:181:ASP:HA	1:A:237:MET:HG2	1.83	0.59
1:A:166:SER:HB2	1:A:173:TYR:OH	2.03	0.58
1:A:563:ASP:H	1:A:571:GLN:HE22	1.52	0.58
1:A:401:CYS:C	1:A:403:SER:H	2.12	0.57
2:B:102:LEU:CB	2:B:166:THR:HG21	2.22	0.57
1:A:84:THR:HA	1:A:410:LEU:HD22	1.86	0.57
1:A:118:PHE:HA	1:A:150:VAL:HG22	1.86	0.57
1:A:265:ILE:CD1	1:A:358:ILE:HG21	2.35	0.57
1:A:485:ARG:HH21	1:A:550:HIS:HD2	1.53	0.56
3:C:29:HIS:CD2	3:C:30:ILE:HG13	2.40	0.56
1:A:190:ARG:HD2	1:A:440:PRO:HB2	1.87	0.56
1:A:409:ARG:NH2	6:A:701:MLI:O9	2.37	0.56
1:A:129:ARG:NH1	1:A:130:ALA:H	2.04	0.56
2:B:69:ILE:O	2:B:159:ILE:HD11	2.05	0.55
1:A:246:GLN:HE22	1:A:600:ARG:HE	1.51	0.55
2:B:21:PRO:HG3	2:B:108:ILE:O	2.06	0.55
3:C:52:LEU:HD21	3:C:98:LEU:HA	1.89	0.54
1:A:581:ARG:HH22	1:A:604:ASP:CG	2.15	0.54
1:A:91:LEU:HA	1:A:602:VAL:HG11	1.89	0.54
2:B:230:ASN:HD22	2:B:230:ASN:C	2.15	0.54
1:A:566:LYS:HB2	1:A:567:PRO:HD2	1.89	0.53
1:A:414:SER:O	1:A:418:LEU:HD13	2.09	0.53
2:B:203:LEU:HD11	2:B:241:MET:HE2	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:516:TRP:HB3	2:B:60:THR:HG21	1.90	0.53
2:B:27:LYS:HE2	2:B:28:PRO:HD3	1.91	0.53
1:A:157:SER:O	1:A:161:THR:HG23	2.08	0.53
1:A:239:THR:CG2	1:A:588:VAL:HG13	2.39	0.53
3:C:69:ASN:ND2	3:C:71:GLU:HB2	2.24	0.52
1:A:497:ILE:HG22	1:A:534:LEU:HD13	1.91	0.52
1:A:254:HIS:HD2	1:A:256:THR:N	2.07	0.52
3:C:69:ASN:HD21	3:C:71:GLU:HB2	1.75	0.52
1:A:127:TYR:OH	1:A:129:ARG:HD3	2.10	0.51
1:A:246:GLN:HE22	1:A:600:ARG:HH21	1.59	0.51
2:B:65:CYS:O	2:B:66:ARG:HG3	2.10	0.51
2:B:21:PRO:HD3	2:B:109:LYS:HG3	1.92	0.51
1:A:61:ALA:HA	5:A:700:FAD:C5X	2.41	0.51
2:B:178:TYR:CD1	2:B:214:ARG:HB2	2.46	0.51
2:B:208:ASP:HB2	2:B:209:PRO:CD	2.41	0.51
2:B:27:LYS:HB3	2:B:28:PRO:HD3	1.93	0.51
2:B:52:LYS:HA	2:B:56:ASP:O	2.10	0.51
2:B:74:ALA:O	2:B:102:LEU:HG	2.11	0.51
2:B:189:ARG:NH2	2:B:193:ASP:OD1	2.44	0.51
1:A:374:ASN:HB3	1:A:376:LYS:H	1.76	0.50
1:A:104:ALA:HB3	1:A:105:PRO:HD3	1.93	0.50
1:A:10:SER:C	1:A:12:GLN:H	2.19	0.49
1:A:27:ALA:H	5:A:700:FAD:H4B	1.75	0.49
1:A:66:ASN:ND2	1:A:411:GLY:O	2.41	0.49
3:C:57:SER:O	3:C:61:LEU:HB2	2.13	0.49
3:C:104:ASN:HD22	3:C:107:ARG:HE	1.61	0.49
1:A:48:VAL:HG12	5:A:700:FAD:H2A	1.95	0.49
1:A:568:ILE:HG13	1:A:569:GLN:N	2.27	0.49
1:A:61:ALA:HA	5:A:700:FAD:C6	2.43	0.48
1:A:113:ASN:HD22	2:B:138:GLY:H	1.60	0.48
4:D:61:TYR:HA	11:D:1306:EPH:H52	1.95	0.48
1:A:130:ALA:HB2	1:A:145:HIS:CD2	2.48	0.48
2:B:155:LEU:HD23	2:B:185:MET:HE2	1.95	0.48
1:A:29:GLY:N	1:A:58:THR:HG21	2.24	0.48
2:B:14:PHE:O	2:B:32:THR:HA	2.13	0.48
1:A:258:ILE:HD12	1:A:263:CYS:O	2.13	0.48
2:B:166:THR:O	2:B:166:THR:CG2	2.62	0.47
1:A:64:GLY:HA2	1:A:154:THR:HG21	1.95	0.47
1:A:298:ARG:HH22	1:A:409:ARG:NH2	2.13	0.47
1:A:342:SER:OG	1:A:353:VAL:HG21	2.14	0.47
2:B:129:LEU:HD11	2:B:195:ARG:CB	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:109:LEU:HD13	4:D:51:VAL:HG12	1.97	0.47
1:A:57:HIS:C	1:A:59:VAL:N	2.73	0.47
1:A:117:PRO:O	1:A:150:VAL:HG13	2.14	0.47
1:A:249:GLU:HB2	1:A:540:ALA:HB2	1.97	0.46
1:A:113:ASN:ND2	2:B:138:GLY:H	2.13	0.46
4:D:124:VAL:HG13	4:D:128:LYS:HB3	1.97	0.46
1:A:114:TYR:CD2	1:A:165:ARG:HD2	2.50	0.46
1:A:405:HIS:CG	1:A:409:ARG:HG3	2.50	0.46
1:A:414:SER:O	1:A:417:ASP:HB3	2.14	0.46
1:A:415:LEU:HG	5:A:700:FAD:N3	2.28	0.46
1:A:554:ASP:HB2	1:A:555:PHE:CD2	2.51	0.46
2:B:69:ILE:O	2:B:159:ILE:CD1	2.64	0.45
1:A:353:VAL:CG1	1:A:358:ILE:HD11	2.44	0.45
3:C:62:SER:HB2	3:C:66:LEU:HD12	1.98	0.45
1:A:267:GLU:OE1	1:A:298:ARG:HD3	2.17	0.45
2:B:64:SER:HB3	2:B:70:CYS:SG	2.56	0.45
2:B:188:TYR:OH	2:B:237:GLU:HG2	2.17	0.45
1:A:57:HIS:O	1:A:59:VAL:N	2.51	0.44
2:B:230:ASN:ND2	2:B:233:LYS:H	2.15	0.44
1:A:23:VAL:HG23	1:A:207:ALA:HB3	1.99	0.44
1:A:57:HIS:CE1	1:A:226:SER:HA	2.52	0.44
1:A:58:THR:HG22	5:A:700:FAD:O4'	2.17	0.44
5:A:700:FAD:H9	5:A:700:FAD:H1'1	1.84	0.44
1:A:457:LEU:O	1:A:509:THR:HG21	2.17	0.44
1:A:577:GLN:H	1:A:577:GLN:CD	2.26	0.44
1:A:270:ARG:HD2	1:A:294:ASP:O	2.18	0.44
4:D:128:LYS:O	4:D:132:MET:HG3	2.17	0.44
1:A:556:LYS:HE2	1:A:556:LYS:HB3	1.85	0.44
2:B:185:MET:HE3	2:B:231:PRO:HA	2.00	0.44
4:D:58:PRO:HB3	11:D:1306:EPH:H28	2.00	0.44
1:A:299:ASP:O	1:A:303:ARG:HB2	2.17	0.44
2:B:103:PRO:O	2:B:166:THR:HG23	2.17	0.44
1:A:60:ALA:HB1	5:A:700:FAD:HM71	1.99	0.43
4:D:89:THR:HG23	4:D:102:LYS:NZ	2.33	0.43
1:A:174:PHE:HB3	1:A:177:TYR:CD1	2.53	0.43
2:B:44:VAL:HG11	2:B:77:ILE:HD12	2.00	0.43
1:A:401:CYS:C	1:A:403:SER:N	2.75	0.43
1:A:182:LEU:HD23	1:A:184:MET:HE1	1.99	0.43
1:A:461:ASN:ND2	1:A:508:LYS:HE3	2.34	0.43
1:A:464:ILE:O	1:A:508:LYS:N	2.51	0.43
1:A:98:HIS:C	1:A:98:HIS:CD2	2.97	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:172:TRP:HZ3	3:C:30:ILE:HG23	1.83	0.43
2:B:164:CYS:HB2	2:B:221:CYS:HB2	2.01	0.43
1:A:131:PHE:CD1	1:A:411:GLY:HA3	2.54	0.43
1:A:272:GLU:HB3	1:A:333:GLN:HE21	1.84	0.43
1:A:129:ARG:HA	1:A:129:ARG:HD2	1.54	0.42
1:A:200:GLY:HA3	1:A:514:MET:HB3	2.01	0.42
2:B:132:LYS:HG2	2:B:197:ASP:HB3	2.01	0.42
2:B:208:ASP:HB2	2:B:209:PRO:HD3	2.01	0.42
2:B:27:LYS:HB3	2:B:28:PRO:HD2	2.00	0.42
2:B:67:GLU:HB3	2:B:69:ILE:HD11	2.00	0.42
2:B:45:LEU:O	2:B:49:ILE:HG12	2.20	0.42
1:A:182:LEU:HD13	1:A:237:MET:HE2	2.02	0.42
1:A:501:TYR:HA	1:A:504:LEU:HD23	2.01	0.42
2:B:165:SER:HA	2:B:181:PRO:HD2	2.02	0.42
1:A:88:SER:HB2	1:A:406:GLY:HA3	2.02	0.41
1:A:374:ASN:HD22	1:A:374:ASN:HA	1.67	0.41
1:A:52:PHE:CD1	1:A:53:PRO:HD2	2.56	0.41
2:B:198:PHE:HD2	2:B:201:GLU:HG3	1.86	0.41
1:A:130:ALA:O	1:A:271:GLY:HA3	2.21	0.41
1:A:416:LEU:O	1:A:417:ASP:C	2.64	0.41
3:C:46:ARG:HG2	10:C:1305:HEM:O2D	2.21	0.41
2:B:47:ALA:O	2:B:51:ILE:HG23	2.21	0.41
1:A:162:LEU:HD23	1:A:162:LEU:HA	1.96	0.41
1:A:231:THR:HA	1:A:529:LEU:HD21	2.03	0.41
1:A:493:GLY:O	1:A:497:ILE:HG12	2.20	0.41
1:A:29:GLY:HA3	1:A:418:LEU:HD23	2.03	0.41
1:A:152:ASP:HB2	1:A:339:PRO:HD2	2.03	0.41
1:A:521:VAL:HA	1:A:524:LEU:HD12	2.02	0.41
2:B:20:ASP:HA	2:B:21:PRO:HD3	1.88	0.41
1:A:114:TYR:CE2	1:A:165:ARG:HD2	2.56	0.41
1:A:556:LYS:H	1:A:556:LYS:HD2	1.86	0.41
2:B:154:GLY:N	2:B:157:GLU:OE1	2.54	0.41
3:C:104:ASN:ND2	3:C:107:ARG:HE	2.19	0.41
1:A:443:LYS:HB2	1:A:446:ALA:HB2	2.02	0.40
2:B:108:ILE:HD12	2:B:112:VAL:HG12	2.02	0.40
3:C:78:LYS:HA	3:C:82:LEU:HD21	2.03	0.40
2:B:16:ILE:HG21	2:B:59:LEU:HD11	2.04	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	611/622 (98%)	573 (94%)	36 (6%)	2 (0%)	36	64
2	B	237/252 (94%)	225 (95%)	11 (5%)	1 (0%)	30	58
3	C	136/140 (97%)	129 (95%)	5 (4%)	2 (2%)	8	27
4	D	100/103 (97%)	95 (95%)	5 (5%)	0	100	100
All	All	1084/1117 (97%)	1022 (94%)	57 (5%)	5 (0%)	24	53

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	84	PRO
1	A	260	GLY
1	A	11	THR
2	B	175	GLY
3	C	81	CYS

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	499/506 (99%)	457 (92%)	42 (8%)	10	30
2	B	214/220 (97%)	195 (91%)	19 (9%)	9	27
3	C	117/118 (99%)	107 (92%)	10 (8%)	10	29
4	D	76/76 (100%)	74 (97%)	2 (3%)	40	72
All	All	906/920 (98%)	833 (92%)	73 (8%)	11	31

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

Mol	Chain	Res	Type
1	A	11	THR
1	A	54	THR
1	A	58	THR
1	A	59	VAL
1	A	72	MET
1	A	74	GLU
1	A	129	ARG
1	A	165	ARG
1	A	167	LEU
1	A	175	VAL
1	A	182	LEU
1	A	220	THR
1	A	264	LEU
1	A	277	ILE
1	A	294	ASP
1	A	306	THR
1	A	324	LEU
1	A	329	LEU
1	A	341	ILE
1	A	352	ASP
1	A	368	MET
1	A	374	ASN
1	A	403	SER
1	A	410	LEU
1	A	419	VAL
1	A	452	MET
1	A	461	ASN
1	A	463	THR
1	A	485	ARG
1	A	488	SER
1	A	504	LEU
1	A	514	MET
1	A	534	LEU
1	A	541	GLU
1	A	554	ASP
1	A	556	LYS
1	A	561	GLU
1	A	569	GLN
1	A	572	GLN
1	A	588	VAL
1	A	597	LEU
1	A	605	LYS

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Mol	Chain	Res	Type
2	B	49	ILE
2	B	51	ILE
2	B	66	ARG
2	B	69	ILE
2	B	86	THR
2	B	87	ARG
2	B	88	ARG
2	B	96	VAL
2	B	105	MET
2	B	134	GLU
2	B	146	ILE
2	B	159	ILE
2	B	189	ARG
2	B	192	ILE
2	B	201	GLU
2	B	214	ARG
2	B	220	ASN
2	B	224	THR
2	B	235	ILE
3	C	13	GLU
3	C	23	ASN
3	C	26	LEU
3	C	30	ILE
3	C	31	THR
3	C	52	LEU
3	C	87	ILE
3	C	98	LEU
3	C	115	LYS
3	C	123	THR
4	D	78	LEU
4	D	124	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	ASN
1	A	143	GLN
1	A	145	HIS
1	A	160	HIS
1	A	246	GLN
1	A	254	HIS
1	A	280	GLN

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Mol	Chain	Res	Type
1	A	333	GLN
1	A	374	ASN
1	A	384	ASN
1	A	408	ASN
1	A	453	ASN
1	A	461	ASN
1	A	474	GLN
1	A	527	GLN
1	A	550	HIS
1	A	571	GLN
1	A	572	GLN
1	A	579	HIS
2	B	38	ASN
2	B	121	GLN
2	B	144	GLN
2	B	174	ASN
2	B	186	GLN
2	B	220	ASN
2	B	230	ASN
3	C	23	ASN
3	C	29	HIS
3	C	104	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	$\# Z > 2$
7	FES	B	302	2	0,4,4	-	-	-		
9	F3S	B	304	2	0,9,9	-	-	-		
5	FAD	A	700	1	58,58,58	1.12	6 (10%)	85,89,89	1.63	16 (18%)
11	EPH	D	1306	-	43,43,48	1.77	8 (18%)	46,48,53	1.25	3 (6%)
8	SF4	B	303	2	0,12,12	-	-	-		
12	N1M	D	1201	-	11,11,11	0.90	1 (9%)	14,14,14	0.92	0
10	HEM	C	1305	4,3	50,50,50	1.86	9 (18%)	67,82,82	1.17	1 (1%)
6	MLI	A	701	-	6,6,6	1.19	0	7,7,7	0.84	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
7	FES	B	302	2	-	-	0/1/1/1
9	F3S	B	304	2	-	-	0/3/3/3
5	FAD	A	700	1	-	6/34/50/50	0/6/6/6
11	EPH	D	1306	-	-	28/47/47/52	-
8	SF4	B	303	2	-	-	0/6/5/5
12	N1M	D	1201	-	-	2/6/6/6	0/1/1/1
10	HEM	C	1305	4,3	-	7/14/54/54	-
6	MLI	A	701	-	-	2/4/4/4	-

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	C	1305	HEM	C3D-C2D	7.76	1.53	1.36
11	D	1306	EPH	P1-O6	5.21	1.79	1.55
10	C	1305	HEM	FE-NB	4.58	2.09	1.94
11	D	1306	EPH	O1-C3	4.44	1.46	1.34
11	D	1306	EPH	O2-C4	4.37	1.46	1.33
10	C	1305	HEM	FE-NA	4.17	2.08	1.95
11	D	1306	EPH	C29-C28	3.82	1.53	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	D	1306	EPH	C25-C24	3.79	1.53	1.31
5	A	700	FAD	C4X-N5	3.77	1.38	1.30
11	D	1306	EPH	C13-C12	3.77	1.53	1.31
10	C	1305	HEM	FE-ND	3.44	2.05	1.94
11	D	1306	EPH	C15-C16	3.23	1.53	1.29
10	C	1305	HEM	CAC-C3C	2.94	1.55	1.47
5	A	700	FAD	C10-N1	2.79	1.38	1.33
5	A	700	FAD	C2A-N1A	2.68	1.38	1.33
12	D	1201	N1M	C2-C1	2.65	1.49	1.39
10	C	1305	HEM	CAB-C3B	2.63	1.54	1.47
11	D	1306	EPH	P1-O7	2.60	1.59	1.50
5	A	700	FAD	C2A-N3A	2.58	1.38	1.33
10	C	1305	HEM	CMB-C2B	2.26	1.55	1.50
10	C	1305	HEM	CMA-C3A	2.23	1.55	1.50
10	C	1305	HEM	CMC-C2C	2.16	1.55	1.50
5	A	700	FAD	C8A-N7A	2.10	1.35	1.31
5	A	700	FAD	C5A-N7A	-2.03	1.35	1.39

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	700	FAD	N3A-C2A-N1A	-6.17	119.24	128.58
10	C	1305	HEM	C4D-ND-C1D	5.21	111.38	105.21
11	D	1306	EPH	O1-C3-C5	4.41	121.03	111.48
5	A	700	FAD	C5A-C4A-N3A	-4.32	120.77	126.72
5	A	700	FAD	N9A-C8A-N7A	-3.79	108.56	113.94
5	A	700	FAD	C5A-N7A-C8A	3.49	108.94	103.45
5	A	700	FAD	C2A-N3A-C4A	3.42	120.18	111.83
5	A	700	FAD	C4-N3-C2	-3.21	119.94	125.64
5	A	700	FAD	C10-C4X-N5	-2.84	119.00	124.81
11	D	1306	EPH	O2-C4-C18	2.78	120.32	111.83
5	A	700	FAD	C4X-C10-N10	2.78	120.46	116.48
5	A	700	FAD	C4X-C4-N3	2.77	120.30	113.25
5	A	700	FAD	N3A-C4A-N9A	2.77	131.87	127.17
5	A	700	FAD	C4A-C5A-N7A	-2.52	107.70	110.58
5	A	700	FAD	O2P-P-O3P	2.50	114.03	107.27
5	A	700	FAD	O4-C4-C4X	-2.48	120.00	126.53
11	D	1306	EPH	C30-C29-C28	-2.42	111.86	126.42
5	A	700	FAD	C4X-C10-N1	-2.42	118.67	124.59
5	A	700	FAD	C10-N1-C2	2.28	121.78	116.85
5	A	700	FAD	C9A-C5X-N5	-2.26	120.06	122.45

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	C	1305	HEM	C2B-C3B-CAB-CBB
10	C	1305	HEM	C4B-C3B-CAB-CBB
10	C	1305	HEM	C2C-C3C-CAC-CBC
10	C	1305	HEM	C4C-C3C-CAC-CBC
11	D	1306	EPH	C25-C26-C27-C28
11	D	1306	EPH	C37-O5-P1-O7
11	D	1306	EPH	C5-C3-O1-C2
11	D	1306	EPH	C38-O8-P1-O5
11	D	1306	EPH	C38-O8-P1-O6
11	D	1306	EPH	C38-O8-P1-O7
11	D	1306	EPH	O8-C38-C39-N1
12	D	1201	N1M	O-C7-N-C8
12	D	1201	N1M	C2-C7-N-C8
11	D	1306	EPH	O3-C3-O1-C2
11	D	1306	EPH	C18-C4-O2-C1
11	D	1306	EPH	O4-C4-O2-C1
11	D	1306	EPH	C27-C28-C29-C30
5	A	700	FAD	O3'-C3'-C4'-C5'
5	A	700	FAD	C2'-C3'-C4'-C5'
10	C	1305	HEM	C3D-CAD-CBD-CGD
11	D	1306	EPH	C11-C12-C13-C14
5	A	700	FAD	C2'-C3'-C4'-O4'
11	D	1306	EPH	C6-C7-C8-C9
11	D	1306	EPH	C20-C21-C22-C23
11	D	1306	EPH	C5-C6-C7-C8
11	D	1306	EPH	C23-C24-C25-C26
11	D	1306	EPH	C19-C20-C21-C22
11	D	1306	EPH	C11-C10-C9-C8
5	A	700	FAD	O3'-C3'-C4'-O4'
11	D	1306	EPH	C4-C18-C19-C20
11	D	1306	EPH	C18-C19-C20-C21
11	D	1306	EPH	C12-C13-C14-C15
5	A	700	FAD	PA-O3P-P-O5'
11	D	1306	EPH	C39-C38-O8-P1
11	D	1306	EPH	C37-O5-P1-O8
11	D	1306	EPH	C2-C37-O5-P1
11	D	1306	EPH	O2-C1-C2-O1
10	C	1305	HEM	CAA-CBA-CGA-O1A
10	C	1305	HEM	CAA-CBA-CGA-O2A
6	A	701	MLI	C2-C1-C3-O8

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Mol	Chain	Res	Type	Atoms
11	D	1306	EPH	O1-C3-C5-C6
5	A	700	FAD	O4B-C4B-C5B-O5B
6	A	701	MLI	C2-C1-C3-O9
11	D	1306	EPH	O3-C3-C5-C6
11	D	1306	EPH	C19-C18-C4-O2

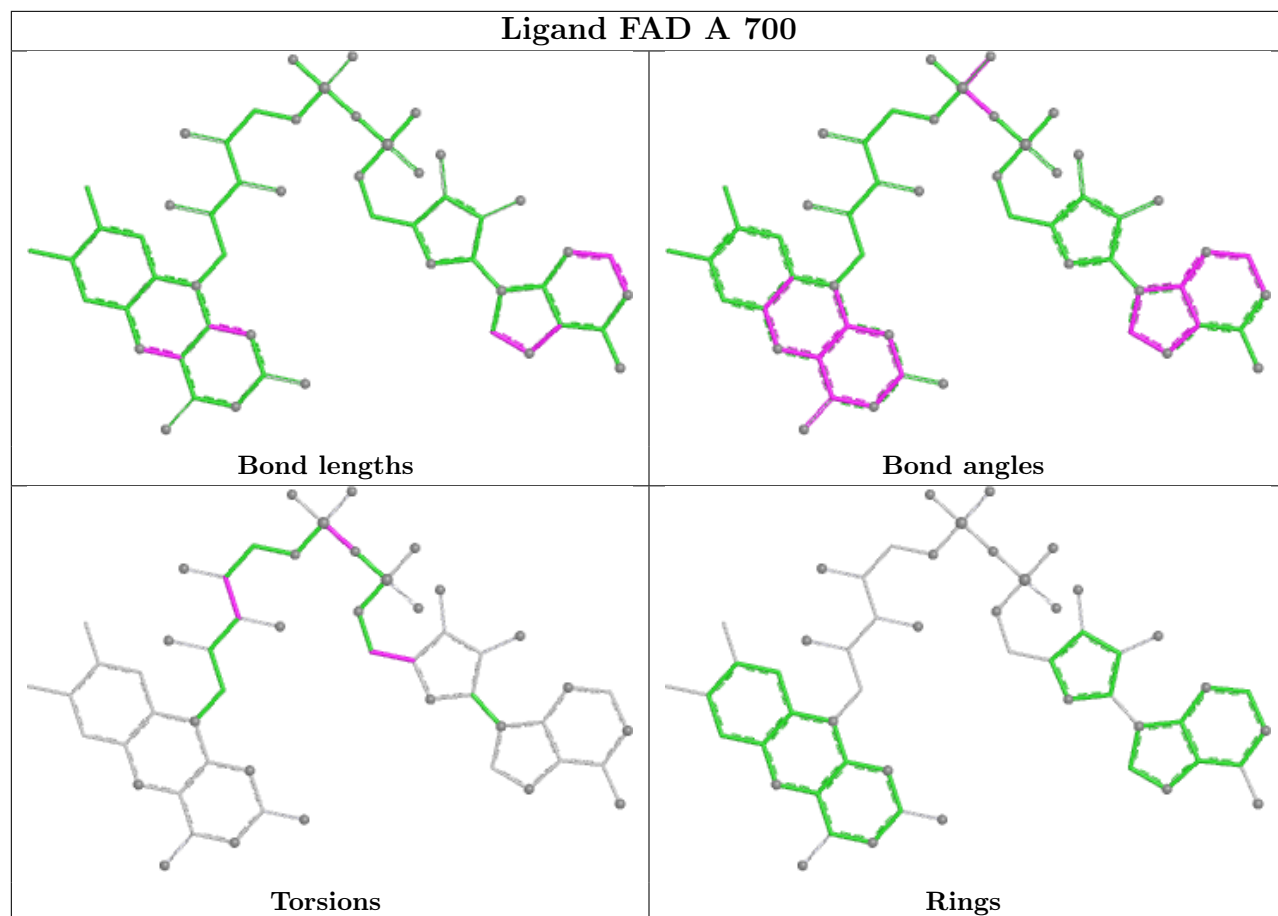
There are no ring outliers.

4 monomers are involved in 20 short contacts:

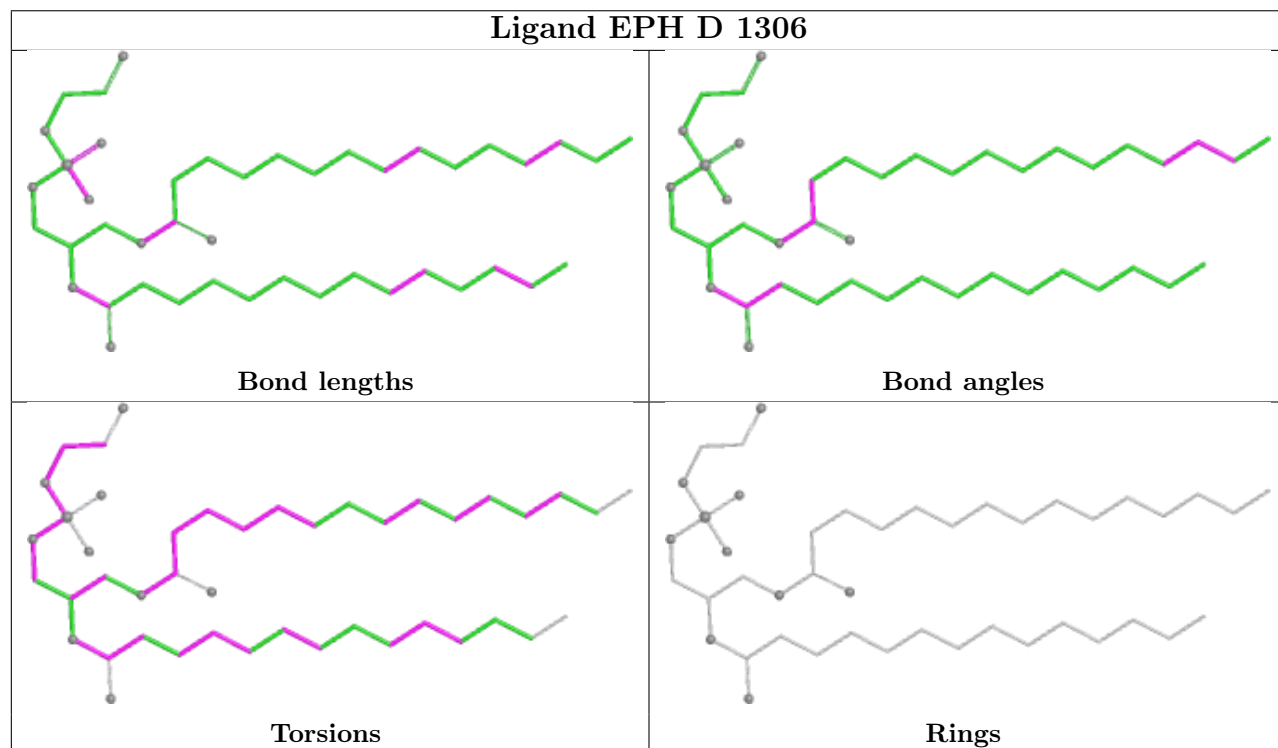
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	700	FAD	14	0
11	D	1306	EPH	2	0
10	C	1305	HEM	3	0
6	A	701	MLI	1	0

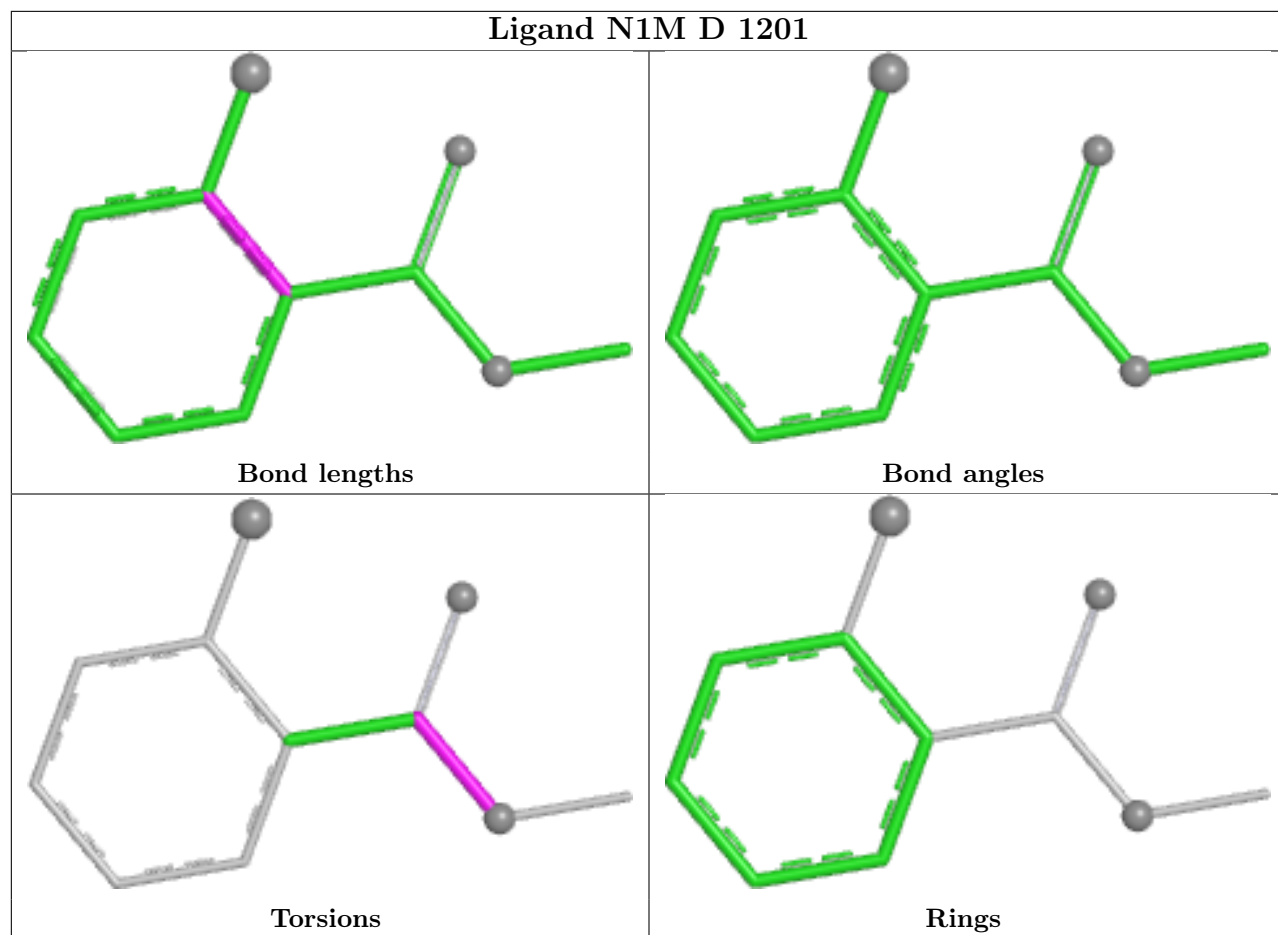
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

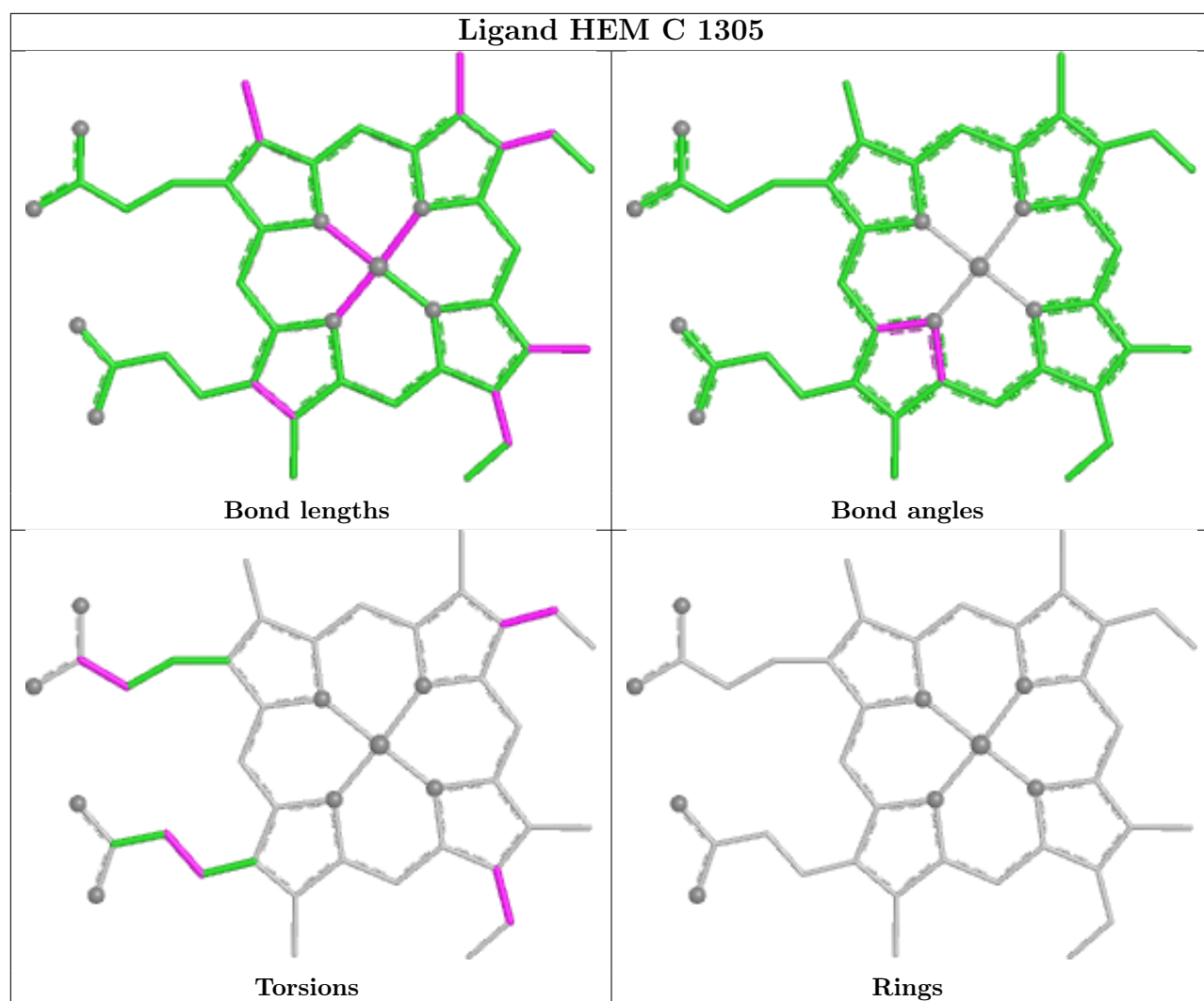
Ligand FAD A 700



Ligand EPH D 1306







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

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	613/622 (98%)	0.19	11 (1%) 67 59	56, 75, 106, 131	0
2	B	239/252 (94%)	0.07	3 (1%) 75 67	56, 70, 105, 126	0
3	C	138/140 (98%)	0.23	4 (2%) 53 44	56, 79, 126, 138	0
4	D	102/103 (99%)	0.15	3 (2%) 53 44	61, 79, 99, 115	0
All	All	1092/1117 (97%)	0.16	21 (1%) 66 58	56, 75, 107, 138	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	9	PRO	4.0
3	C	82	LEU	3.6
1	A	129	ARG	3.5
4	D	35	SER	3.3
3	C	83	GLY	3.1
1	A	406	GLY	3.0
3	C	63	ALA	3.0
4	D	36	SER	2.9
1	A	449	GLU	2.8
1	A	442	ILE	2.8
1	A	404	VAL	2.6
1	A	10	SER	2.6
1	A	405	HIS	2.3
1	A	299	ASP	2.3
1	A	568	ILE	2.3
4	D	124	VAL	2.2
2	B	223	GLY	2.2
1	A	37	GLY	2.1
2	B	35	ILE	2.1
1	A	327	HIS	2.0
3	C	35	TRP	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](#)

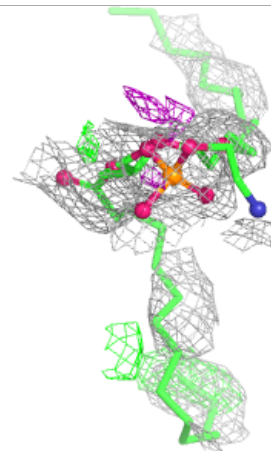
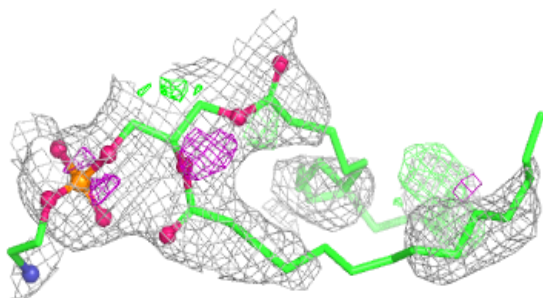
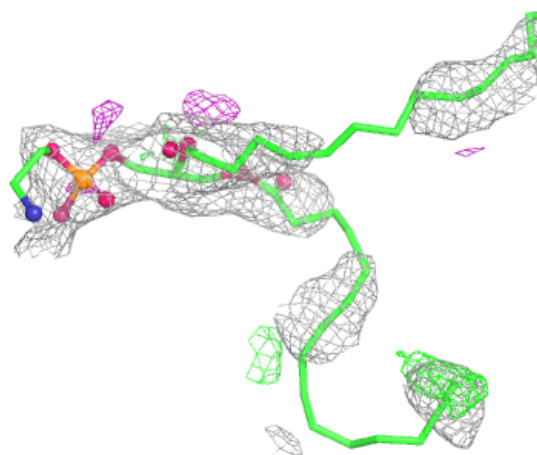
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
11	EPH	D	1306	44/49	0.83	0.23	106,118,123,123	0
5	FAD	A	700	53/53	0.86	0.15	47,60,75,75	0
6	MLI	A	701	7/7	0.91	0.12	79,81,82,82	0
10	HEM	C	1305	43/43	0.97	0.10	62,64,70,72	0
7	FES	B	302	4/4	0.98	0.05	61,62,64,64	0
9	F3S	B	304	7/7	0.98	0.06	66,67,70,70	0
12	N1M	D	1201	11/11	0.98	0.11	80,80,81,84	0
8	SF4	B	303	8/8	0.99	0.03	59,59,60,60	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.

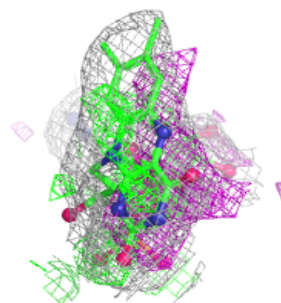
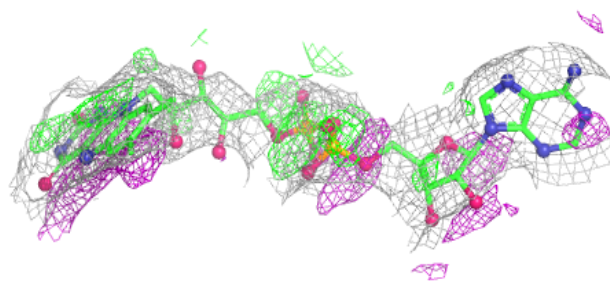
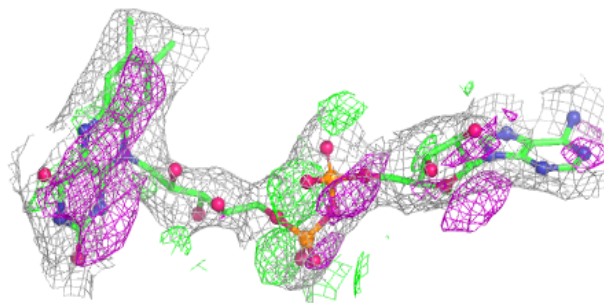
Electron density around EPH D 1306:

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



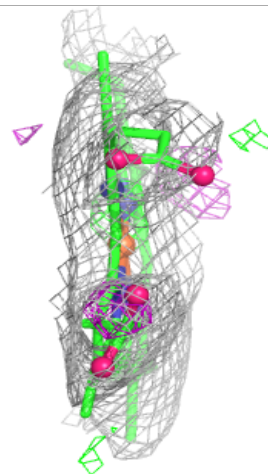
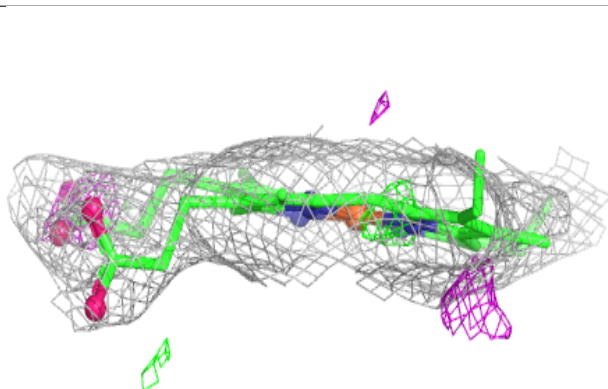
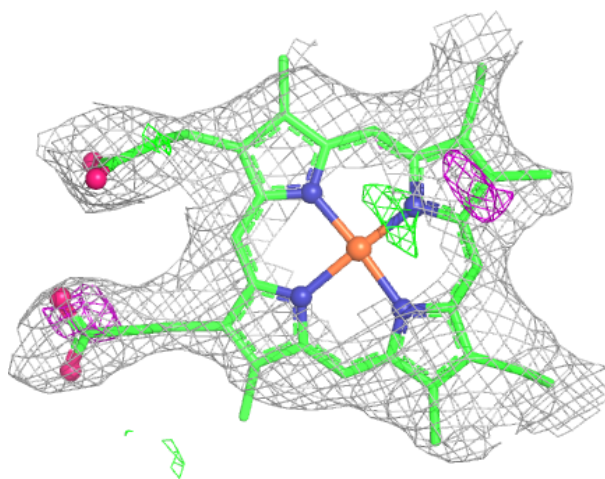
Electron density around FAD A 700:

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



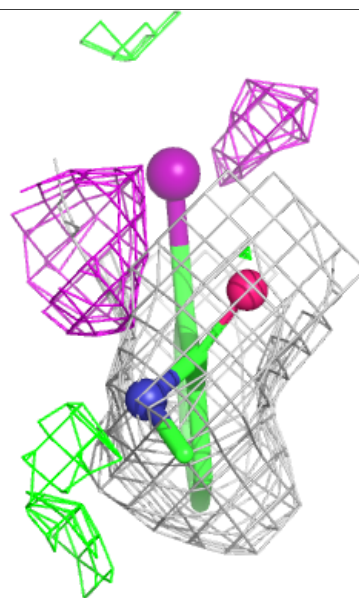
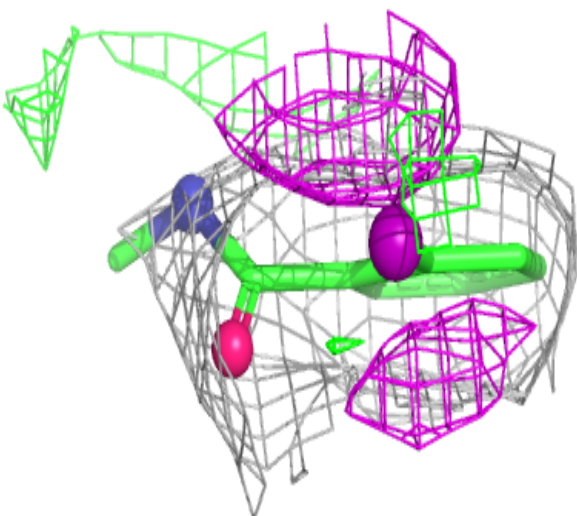
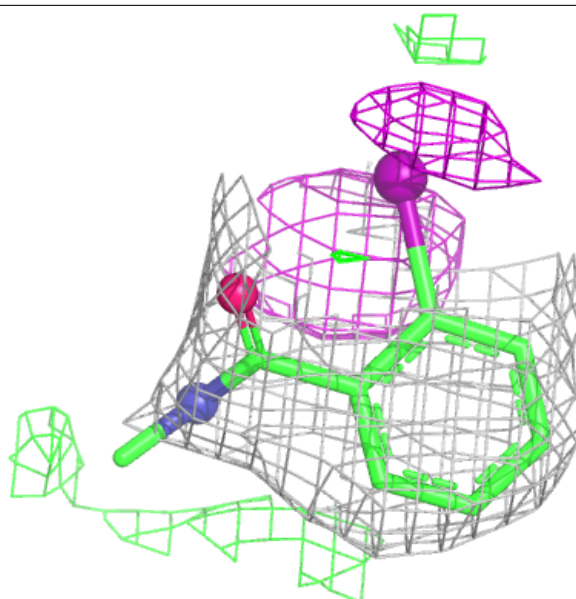
Electron density around HEM C 1305:

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



Electron density around N1M D 1201:

$2mF_o-DF_c$ (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.