



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 11:31 PM BST

PDB ID : 3AYF
Title : Crystal structure of nitric oxide reductase
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Deposited on : 2011-05-06
Resolution : 2.50 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 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 : 2.13.1
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) : 2.13.1

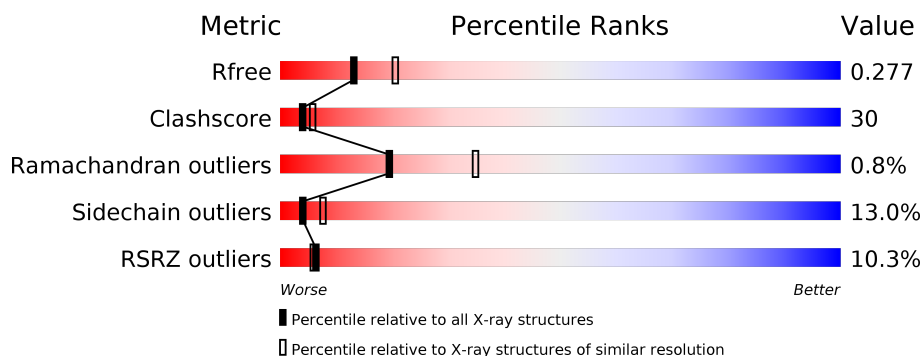
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on 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	800	<div> <div>10%</div> <div>49%</div> <div>37%</div> <div>8%</div> <div>6%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	EPE	A	808	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6455 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.

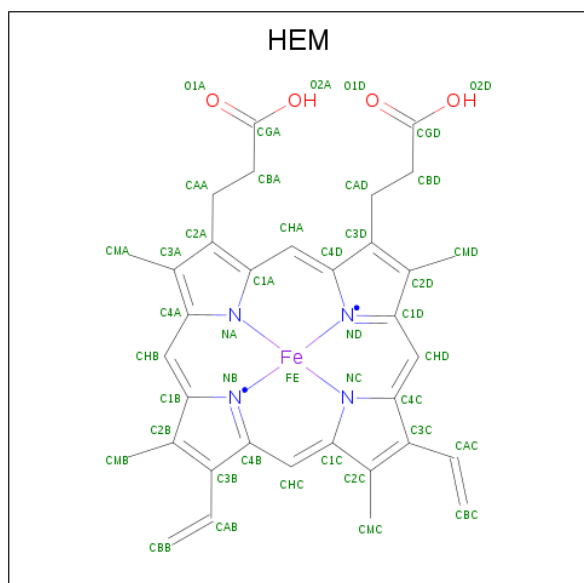
- Molecule 1 is a protein called Nitric oxide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	754	6078	4055	966	1034	23	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	793	LEU	-	expression tag	UNP B3Y963
A	794	GLU	-	expression tag	UNP B3Y963
A	795	HIS	-	expression tag	UNP B3Y963
A	796	HIS	-	expression tag	UNP B3Y963
A	797	HIS	-	expression tag	UNP B3Y963
A	798	HIS	-	expression tag	UNP B3Y963
A	799	HIS	-	expression tag	UNP B3Y963
A	800	HIS	-	expression tag	UNP B3Y963

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

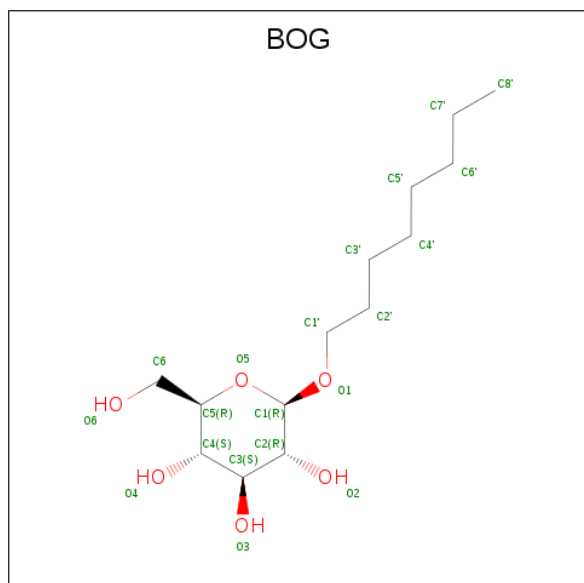
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn		
			1	1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

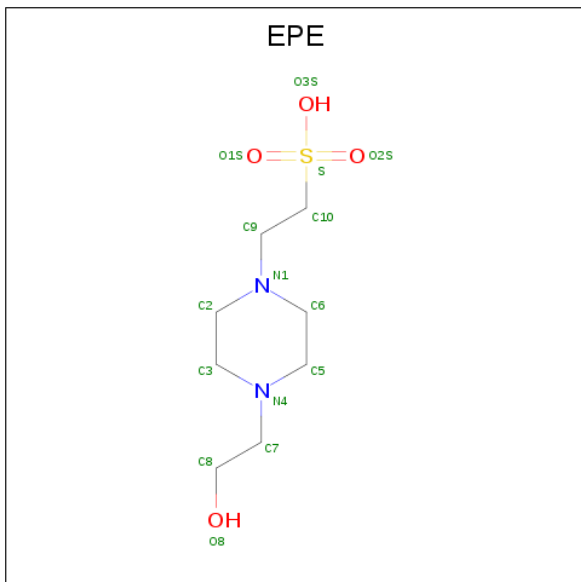
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca		
			1	1	0	0

- Molecule 5 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: C₁₄H₂₈O₆).



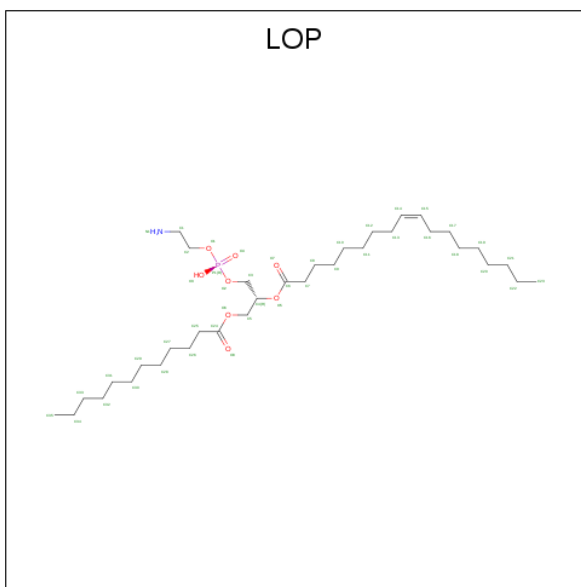
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O		
			20	14	6	0	0
5	A	1	Total	C	O		
			20	14	6	0	0
5	A	1	Total	C	O		
			20	14	6	0	0

- Molecule 6 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 7 is (1R)-2-{[(R)-(2-AMINOETHOXY)(HYDROXY)PHOSPHORYL]OXY}-1-[(DODECANOYLOXY)METHYL]ETHYL (9Z)-OCTADEC-9-ENOATE (three-letter code: LOP) (formula: $C_{35}H_{68}NO_8P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	N	O	P	0	0
			45	35	1	8	1		
7	A	1	Total	C	N	O	P	0	0
			45	35	1	8	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	124	Total	O	0	0
			124	124		

E777

LEU
PRO
VAL
ALA
ASN
HIS
TRP
MET
LYS
ASP
ARG
LEU
LYS
ASN
SER
LEU
GLU
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	110.41Å 149.65Å 151.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.78 – 2.50 43.53 – 2.51	Depositor EDS
% Data completeness (in resolution range)	98.9 (19.78-2.50) 99.0 (43.53-2.51)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.32 (at 2.51Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.244 , 0.282 0.241 , 0.277	Depositor DCC
R_{free} test set	2127 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.384	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 65.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6455	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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: ZN, CA, LOP, HEM, EPE, BOG

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	0.44	0/6262	0.65	2/8517 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	220	LEU	CA-CB-CG	5.51	127.96	115.30
1	A	602	GLY	N-CA-C	-5.08	100.40	113.10

There are no chirality outliers.

There are no planarity outliers.

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	6078	0	6095	368	0
2	A	86	0	60	3	0
3	A	1	0	0	0	0
4	A	1	0	0	0	0
5	A	60	0	84	4	0
6	A	15	0	18	0	0
7	A	90	0	134	31	0
8	A	124	0	0	9	0
All	All	6455	0	6391	384	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 30.

All (384) 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:198:ARG:HH11	1:A:198:ARG:HG3	0.97	1.12
1:A:172:ARG:HG3	1:A:207:ILE:HD11	1.33	1.04
1:A:293:LEU:H	1:A:293:LEU:HD23	1.28	0.99
1:A:465:ASP:HA	7:A:809:LOP:H11	1.47	0.97
1:A:126:LEU:H	1:A:126:LEU:HD22	1.33	0.93
1:A:56:ARG:HB3	1:A:62:VAL:HA	1.49	0.93
1:A:198:ARG:NH1	1:A:198:ARG:HG3	1.72	0.90
1:A:56:ARG:HA	1:A:63:LEU:HD22	1.54	0.88
1:A:585:LEU:HA	1:A:588:LEU:HD22	1.55	0.87
1:A:198:ARG:HH12	1:A:203:ASP:HA	1.41	0.85
1:A:75:VAL:CG2	1:A:210:ILE:HG12	2.07	0.85
1:A:742:ARG:HD2	1:A:746:ASP:OD2	1.77	0.85
7:A:810:LOP:H112	7:A:810:LOP:H272	1.59	0.84
1:A:172:ARG:HA	1:A:207:ILE:HD12	1.61	0.82
1:A:422:LEU:O	1:A:436:ILE:HG23	1.81	0.81
1:A:543:TYR:O	1:A:547:THR:HG23	1.81	0.81
1:A:293:LEU:H	1:A:293:LEU:CD2	1.94	0.81
1:A:341:TRP:O	1:A:345:ILE:HG22	1.81	0.80
7:A:810:LOP:H292	7:A:810:LOP:H132	1.64	0.79
1:A:624:VAL:HA	1:A:628:VAL:HG13	1.65	0.78
1:A:255:GLY:O	1:A:259:THR:HG23	1.84	0.78
1:A:198:ARG:HH11	1:A:198:ARG:CG	1.87	0.77
1:A:55:VAL:HG13	1:A:64:MET:HB3	1.67	0.77
1:A:104:LYS:O	1:A:108:GLU:HG2	1.85	0.76
7:A:810:LOP:O7	7:A:810:LOP:H91	1.82	0.76
1:A:300:PRO:HD2	1:A:382:GLU:HG3	1.69	0.75
1:A:318:VAL:HG23	1:A:358:LEU:HD13	1.69	0.74
1:A:614:PHE:O	1:A:618:THR:HG22	1.88	0.74
1:A:164:LEU:O	1:A:168:ARG:HG3	1.88	0.73
1:A:126:LEU:HB3	1:A:130:GLU:HG2	1.71	0.73
1:A:517:VAL:O	1:A:521:VAL:HG12	1.89	0.73
1:A:175:PHE:HD1	1:A:187:LEU:HD11	1.52	0.72
1:A:104:LYS:HE2	1:A:108:GLU:OE2	1.90	0.72
1:A:585:LEU:O	1:A:588:LEU:HB2	1.90	0.71
1:A:735:VAL:HG12	1:A:739:LEU:HD22	1.72	0.70
1:A:624:VAL:HG11	7:A:810:LOP:H211	1.73	0.70
1:A:44:ILE:HG12	1:A:567:PRO:HD2	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:ARG:HD3	1:A:208:GLN:NE2	2.06	0.70
1:A:210:ILE:HG22	1:A:211:ALA:N	2.07	0.69
7:A:810:LOP:H112	7:A:810:LOP:H251	1.73	0.69
7:A:810:LOP:O6	7:A:810:LOP:C6	2.40	0.69
1:A:51:ARG:HD2	1:A:495:ASN:ND2	2.08	0.69
1:A:299:THR:HB	1:A:382:GLU:HG2	1.74	0.69
1:A:466:LYS:HB2	7:A:809:LOP:O3	1.92	0.69
1:A:302:GLN:HE22	1:A:377:LEU:HD13	1.58	0.68
1:A:380:GLY:O	1:A:777:GLU:HG2	1.92	0.68
1:A:188:ILE:HG21	1:A:207:ILE:HG22	1.74	0.68
1:A:199:ALA:HB2	1:A:418:ASN:HB3	1.75	0.68
1:A:207:ILE:HG23	8:A:905:HOH:O	1.94	0.68
1:A:198:ARG:NH1	1:A:203:ASP:HA	2.07	0.67
1:A:216:TRP:O	1:A:220:LEU:HD13	1.94	0.67
1:A:194:PRO:O	1:A:198:ARG:HD3	1.94	0.67
1:A:375:ALA:HB3	1:A:376:PRO:HD3	1.76	0.66
1:A:110:MET:HG2	1:A:159:ALA:O	1.94	0.66
1:A:363:ILE:HG22	1:A:515:PHE:CE1	2.30	0.66
1:A:334:ASP:HA	1:A:343:TYR:CD2	2.30	0.65
1:A:168:ARG:HD3	1:A:208:GLN:HE21	1.61	0.65
1:A:363:ILE:HG22	1:A:515:PHE:CZ	2.31	0.65
1:A:75:VAL:HG22	1:A:210:ILE:HG12	1.79	0.65
1:A:321:MET:HE1	1:A:750:LEU:HD11	1.79	0.65
1:A:370:MET:HE1	1:A:526:LEU:HD21	1.79	0.65
1:A:217:THR:HA	1:A:220:LEU:CD1	2.27	0.65
1:A:225:ARG:HE	1:A:232:TYR:HA	1.62	0.65
1:A:93:TYR:HE2	1:A:649:PRO:HG3	1.61	0.64
1:A:157:THR:O	1:A:161:VAL:HG23	1.97	0.64
1:A:135:ARG:O	1:A:139:ILE:HG12	1.97	0.64
1:A:621:TRP:CZ2	7:A:810:LOP:H192	2.33	0.64
1:A:648:THR:HB	1:A:649:PRO:HD3	1.80	0.63
7:A:809:LOP:O4	7:A:809:LOP:H12	1.97	0.63
1:A:393:LEU:HD22	1:A:397:LEU:HD21	1.80	0.63
1:A:65:THR:HG23	1:A:68:THR:H	1.63	0.63
1:A:142:MET:HE3	1:A:242:ALA:HA	1.80	0.63
1:A:293:LEU:N	1:A:293:LEU:HD23	2.07	0.63
1:A:331:THR:HG21	1:A:740:LEU:HD13	1.80	0.62
1:A:400:LEU:HB2	1:A:440:ILE:HD11	1.81	0.62
1:A:217:THR:HA	1:A:220:LEU:HD11	1.80	0.62
1:A:198:ARG:HH12	1:A:203:ASP:CA	2.12	0.62
1:A:302:GLN:NE2	1:A:377:LEU:HD13	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:347:PRO:HD2	1:A:350:ILE:HD12	1.81	0.62
1:A:436:ILE:O	1:A:440:ILE:HG23	2.00	0.61
1:A:771:THR:HG22	1:A:772:HIS:CD2	2.35	0.61
1:A:107:THR:O	1:A:111:GLN:HG3	2.00	0.61
1:A:225:ARG:NE	1:A:232:TYR:HA	2.16	0.61
1:A:274:ARG:HG3	1:A:274:ARG:O	2.00	0.61
1:A:55:VAL:CG1	1:A:64:MET:HB3	2.31	0.61
7:A:809:LOP:O3	7:A:809:LOP:H4	2.00	0.61
1:A:588:LEU:HD23	1:A:622:ASN:HB3	1.81	0.61
1:A:607:PRO:O	1:A:673:ARG:HD2	2.01	0.61
1:A:298:LEU:CD1	1:A:303:VAL:HG13	2.31	0.61
1:A:581:GLU:O	1:A:584:PRO:HD2	2.00	0.61
7:A:810:LOP:C11	7:A:810:LOP:H251	2.31	0.61
1:A:480:VAL:HB	1:A:481:PRO:CD	2.31	0.60
1:A:569:VAL:HG23	5:A:807:BOG:H3'1	1.83	0.60
1:A:298:LEU:HD12	1:A:303:VAL:HG13	1.84	0.60
1:A:393:LEU:O	1:A:397:LEU:HD22	2.02	0.60
1:A:633:ILE:HD12	2:A:802:HEM:HBD2	1.82	0.60
1:A:309:PHE:HA	1:A:312:VAL:HG13	1.82	0.60
1:A:365:THR:HB	8:A:990:HOH:O	2.01	0.60
1:A:516:GLU:HG2	1:A:549:LEU:HD21	1.83	0.60
1:A:172:ARG:NH2	1:A:208:GLN:OE1	2.31	0.60
1:A:34:SER:OG	7:A:809:LOP:H233	2.00	0.60
1:A:56:ARG:HB2	1:A:61:GLU:O	2.02	0.60
1:A:402:GLY:O	1:A:406:ILE:HG12	2.01	0.59
1:A:691:TRP:CE3	1:A:695:ILE:HD11	2.36	0.59
1:A:179:ASP:OD1	1:A:183:LEU:HB3	2.02	0.59
1:A:172:ARG:HG3	1:A:207:ILE:CD1	2.21	0.59
1:A:304:LYS:HG2	1:A:391:ASP:OD1	2.02	0.59
1:A:86:THR:HG23	1:A:92:SER:O	2.03	0.59
1:A:646:PHE:C	1:A:649:PRO:HD2	2.23	0.59
1:A:93:TYR:CE2	1:A:649:PRO:HG3	2.37	0.59
1:A:694:ASN:HB3	7:A:810:LOP:H312	1.84	0.58
1:A:113:TYR:HB2	1:A:162:TYR:CD2	2.38	0.58
1:A:164:LEU:O	1:A:164:LEU:HD22	2.02	0.58
1:A:101:GLU:O	1:A:105:VAL:HG23	2.04	0.57
1:A:692:MET:CE	1:A:748:ILE:HG23	2.34	0.57
1:A:151:THR:O	1:A:153:VAL:N	2.34	0.57
1:A:187:LEU:HD13	1:A:188:ILE:H	1.69	0.57
1:A:299:THR:HB	1:A:382:GLU:CG	2.33	0.57
1:A:704:THR:HG21	1:A:742:ARG:HG3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:ASP:O	1:A:349:ASN:HA	2.04	0.57
1:A:220:LEU:HD22	1:A:221:SER:N	2.19	0.57
1:A:429:GLU:O	1:A:430:TYR:HB2	2.04	0.57
1:A:291:ILE:CG2	1:A:607:PRO:HG3	2.35	0.57
1:A:148:ASN:O	1:A:151:THR:O	2.23	0.57
1:A:421:PHE:CE1	1:A:492:PRO:HD3	2.40	0.56
1:A:156:LEU:H	1:A:156:LEU:HD23	1.71	0.56
1:A:198:ARG:NH1	1:A:202:ALA:O	2.38	0.56
1:A:395:TRP:HA	1:A:398:VAL:HG12	1.86	0.56
1:A:663:PHE:HB3	8:A:922:HOH:O	2.06	0.56
1:A:691:TRP:HE3	1:A:695:ILE:HD11	1.71	0.56
1:A:433:LEU:HD13	1:A:438:GLN:CG	2.35	0.56
1:A:675:ILE:CG2	1:A:767:LEU:HD22	2.36	0.56
1:A:181:TRP:CG	1:A:727:PRO:HB3	2.41	0.55
1:A:511:VAL:O	1:A:515:PHE:CD2	2.58	0.55
1:A:774:GLU:N	1:A:774:GLU:OE1	2.39	0.55
1:A:62:VAL:HG23	8:A:983:HOH:O	2.07	0.55
1:A:139:ILE:HD13	1:A:142:MET:HE3	1.89	0.55
1:A:156:LEU:N	1:A:156:LEU:HD23	2.21	0.55
1:A:142:MET:CE	1:A:242:ALA:HA	2.37	0.55
1:A:299:THR:CB	1:A:382:GLU:HG2	2.37	0.55
1:A:395:TRP:HA	1:A:398:VAL:CG1	2.36	0.55
1:A:439:ILE:O	1:A:442:VAL:HG22	2.07	0.55
1:A:454:ARG:HH11	1:A:454:ARG:HG2	1.72	0.54
1:A:207:ILE:HG13	1:A:207:ILE:O	2.08	0.54
1:A:139:ILE:HD11	1:A:241:ASP:O	2.08	0.54
1:A:264:PHE:O	1:A:268:ILE:HG12	2.08	0.54
1:A:400:LEU:HG	1:A:440:ILE:CG1	2.38	0.54
1:A:688:PHE:O	1:A:692:MET:HG2	2.08	0.54
1:A:744:VAL:HB	1:A:745:PRO:CD	2.36	0.54
1:A:53:LYS:HE2	1:A:54:GLU:HG3	1.90	0.54
1:A:682:ASN:ND2	1:A:684:LYS:HB3	2.23	0.54
1:A:166:LYS:N	1:A:166:LYS:HE2	2.23	0.54
1:A:175:PHE:CE2	1:A:210:ILE:HG21	2.43	0.53
1:A:203:ASP:HB3	8:A:904:HOH:O	2.07	0.53
1:A:378:VAL:HG11	1:A:456:VAL:HG13	1.89	0.53
1:A:682:ASN:HD21	1:A:684:LYS:HB3	1.73	0.53
1:A:483:PHE:CZ	7:A:809:LOP:H222	2.43	0.53
1:A:539:VAL:HG22	8:A:984:HOH:O	2.08	0.53
1:A:339:ILE:HG22	1:A:340:ASN:N	2.23	0.53
1:A:187:LEU:HD13	1:A:188:ILE:N	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:LEU:CG	1:A:440:ILE:HD11	2.39	0.53
1:A:685:TRP:CE3	1:A:685:TRP:HA	2.43	0.53
1:A:758:VAL:O	1:A:762:LYS:HG2	2.09	0.53
1:A:534:THR:OG1	1:A:537:SER:HB3	2.09	0.53
1:A:113:TYR:CD2	1:A:159:ALA:HA	2.44	0.53
1:A:400:LEU:HG	1:A:440:ILE:HD11	1.91	0.52
1:A:569:VAL:CG2	5:A:807:BOG:H3'1	2.39	0.52
1:A:131:LYS:HD3	1:A:131:LYS:O	2.10	0.52
1:A:760:ALA:O	1:A:763:ALA:HB3	2.09	0.52
1:A:367:TRP:CZ3	1:A:667:VAL:HG21	2.45	0.51
1:A:378:VAL:HB	1:A:456:VAL:HG22	1.91	0.51
1:A:291:ILE:HG21	1:A:607:PRO:HG3	1.91	0.51
1:A:771:THR:HG22	1:A:772:HIS:HD2	1.74	0.51
1:A:203:ASP:O	1:A:204:SER:HB2	2.10	0.51
1:A:395:TRP:O	1:A:399:VAL:HG23	2.09	0.51
1:A:515:PHE:HB2	1:A:660:TYR:CE1	2.44	0.51
1:A:172:ARG:HA	1:A:207:ILE:CD1	2.36	0.51
1:A:393:LEU:HD22	1:A:397:LEU:CD2	2.40	0.51
1:A:203:ASP:O	1:A:204:SER:CB	2.58	0.51
1:A:331:THR:CG2	1:A:740:LEU:HD13	2.41	0.51
1:A:434:GLY:O	1:A:437:TRP:HB2	2.11	0.51
1:A:480:VAL:HB	1:A:481:PRO:HD3	1.92	0.51
1:A:534:THR:OG1	1:A:537:SER:CB	2.59	0.51
1:A:138:VAL:HG22	1:A:142:MET:HE2	1.91	0.51
1:A:55:VAL:HA	1:A:154:LEU:O	2.11	0.51
1:A:699:GLY:O	1:A:703:ILE:HB	2.10	0.51
1:A:76:PHE:HB2	1:A:210:ILE:HD11	1.93	0.51
1:A:201:VAL:CG2	1:A:209:GLN:HE22	2.24	0.50
1:A:64:MET:HE1	1:A:212:ASP:HA	1.91	0.50
1:A:19:ASN:O	1:A:23:LYS:HD2	2.11	0.50
1:A:382:GLU:HB2	1:A:773:GLY:O	2.11	0.50
1:A:65:THR:HG23	1:A:68:THR:N	2.25	0.50
1:A:583:ILE:HB	1:A:584:PRO:HD3	1.94	0.50
1:A:616:ILE:HD13	7:A:810:LOP:H82	1.93	0.50
1:A:384:LYS:O	1:A:454:ARG:NH2	2.44	0.50
1:A:52:PRO:HA	1:A:152:ASP:OD1	2.11	0.50
1:A:56:ARG:HA	1:A:63:LEU:CD2	2.35	0.50
1:A:462:ARG:HG2	1:A:462:ARG:HH11	1.77	0.50
1:A:53:LYS:O	1:A:65:THR:HA	2.12	0.50
1:A:440:ILE:HD12	1:A:440:ILE:O	2.11	0.50
1:A:88:LEU:HA	1:A:236:TRP:HZ3	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ILE:CG2	1:A:207:ILE:HG22	2.40	0.50
1:A:209:GLN:O	1:A:212:ASP:HB2	2.12	0.50
1:A:217:THR:O	1:A:220:LEU:HD22	2.11	0.50
1:A:25:ILE:HD13	5:A:806:BOG:H6'1	1.92	0.50
1:A:429:GLU:HG2	1:A:430:TYR:CD2	2.47	0.50
1:A:53:LYS:O	1:A:53:LYS:HD2	2.11	0.50
1:A:23:LYS:NZ	7:A:809:LOP:O4	2.37	0.50
1:A:518:PHE:O	1:A:522:VAL:HG23	2.12	0.49
1:A:309:PHE:HB3	1:A:668:LEU:HG	1.93	0.49
1:A:302:GLN:HE22	1:A:377:LEU:CD1	2.23	0.49
1:A:384:LYS:O	1:A:385:LYS:HB2	2.12	0.49
1:A:55:VAL:HG22	1:A:64:MET:HB2	1.94	0.49
1:A:65:THR:H	1:A:68:THR:HG1	1.58	0.49
1:A:599:MET:HE2	7:A:810:LOP:H71	1.94	0.49
1:A:162:TYR:O	1:A:165:GLU:HB3	2.13	0.49
1:A:441:LEU:HD22	1:A:445:MET:HE2	1.95	0.49
1:A:56:ARG:HB2	1:A:61:GLU:C	2.33	0.49
1:A:759:PHE:O	1:A:763:ALA:HB2	2.13	0.49
1:A:79:TYR:CE2	1:A:188:ILE:HG12	2.47	0.49
1:A:248:PHE:O	1:A:251:VAL:HG13	2.13	0.49
1:A:339:ILE:C	1:A:340:ASN:HD22	2.17	0.48
1:A:615:LEU:O	1:A:618:THR:HG23	2.13	0.48
1:A:56:ARG:CD	1:A:62:VAL:HG12	2.43	0.48
1:A:163:GLY:O	1:A:167:VAL:HG23	2.12	0.48
1:A:240:GLU:HB2	8:A:918:HOH:O	2.13	0.48
1:A:475:TYR:CZ	7:A:809:LOP:H281	2.48	0.48
1:A:483:PHE:HZ	7:A:809:LOP:H222	1.78	0.48
1:A:441:LEU:HD22	1:A:445:MET:CE	2.43	0.48
1:A:55:VAL:HB	1:A:154:LEU:HD12	1.96	0.48
1:A:505:TRP:O	1:A:509:LEU:HB2	2.12	0.48
1:A:126:LEU:HD23	1:A:131:LYS:HG2	1.96	0.48
1:A:203:ASP:O	1:A:203:ASP:CG	2.52	0.48
1:A:647:LEU:HG	1:A:709:GLY:HA3	1.94	0.48
1:A:293:LEU:HA	1:A:296:GLN:HG3	1.95	0.47
1:A:184:LYS:NZ	1:A:344:ASP:O	2.46	0.47
1:A:675:ILE:O	1:A:768:ARG:HD3	2.14	0.47
1:A:199:ALA:HB2	1:A:418:ASN:CB	2.44	0.47
1:A:265:ILE:O	1:A:269:LEU:HB2	2.14	0.47
1:A:305:ALA:O	1:A:308:TYR:HB2	2.14	0.47
1:A:751:ILE:O	1:A:754:VAL:HG22	2.15	0.47
1:A:675:ILE:HG22	1:A:767:LEU:HD22	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:599:MET:CE	7:A:810:LOP:H71	2.43	0.47
1:A:515:PHE:CB	1:A:660:TYR:CE1	2.98	0.47
1:A:280:GLN:N	1:A:537:SER:OG	2.46	0.47
1:A:26:LEU:HA	1:A:547:THR:HG22	1.97	0.47
1:A:586:THR:C	1:A:588:LEU:H	2.18	0.47
1:A:635:LEU:HB2	8:A:901:HOH:O	2.15	0.47
1:A:474:PHE:O	1:A:478:ILE:CG2	2.61	0.47
1:A:75:VAL:CG2	1:A:210:ILE:CG1	2.85	0.47
1:A:139:ILE:CD1	1:A:241:ASP:O	2.63	0.47
1:A:757:LEU:HD22	1:A:761:ILE:HG13	1.96	0.47
7:A:809:LOP:O7	7:A:809:LOP:O6	2.33	0.47
1:A:126:LEU:H	1:A:126:LEU:CD2	2.10	0.47
1:A:77:GLN:HB3	1:A:427:GLY:HA3	1.97	0.47
1:A:298:LEU:HB2	1:A:768:ARG:HH21	1.80	0.47
7:A:810:LOP:H181	7:A:810:LOP:C33	2.44	0.47
1:A:445:MET:O	1:A:448:TRP:HB3	2.15	0.46
1:A:379:GLY:HA3	1:A:455:GLY:O	2.16	0.46
1:A:70:ILE:HG23	1:A:493:ASP:HA	1.97	0.46
1:A:378:VAL:CB	1:A:456:VAL:HG22	2.46	0.46
1:A:400:LEU:CB	1:A:440:ILE:HD11	2.45	0.46
1:A:643:HIS:HA	2:A:802:HEM:O1D	2.16	0.46
1:A:498:MET:O	1:A:501:PHE:HB3	2.16	0.46
1:A:73:GLN:HA	1:A:213:PHE:CD2	2.51	0.46
1:A:510:TRP:CZ2	1:A:514:ILE:HD11	2.51	0.46
1:A:70:ILE:HD13	1:A:493:ASP:HB2	1.98	0.46
1:A:279:MET:HA	1:A:537:SER:OG	2.15	0.46
7:A:810:LOP:H181	7:A:810:LOP:H332	1.98	0.46
1:A:154:LEU:CD1	1:A:156:LEU:HD22	2.45	0.46
1:A:170:TYR:O	1:A:174:VAL:HG23	2.16	0.46
1:A:239:TYR:CZ	1:A:241:ASP:HB2	2.51	0.46
7:A:809:LOP:O6	7:A:809:LOP:C6	2.64	0.46
1:A:298:LEU:HD13	1:A:299:THR:H	1.80	0.46
1:A:292:ASP:OD2	1:A:295:ARG:HD2	2.17	0.45
1:A:54:GLU:HA	1:A:64:MET:O	2.16	0.45
1:A:217:THR:HA	1:A:220:LEU:HD13	1.96	0.45
1:A:692:MET:HE1	1:A:748:ILE:HG23	1.97	0.45
1:A:463:GLU:OE2	1:A:471:HIS:HD2	1.99	0.45
1:A:225:ARG:HB2	1:A:228:ASP:CB	2.47	0.45
1:A:81:LEU:HD21	1:A:175:PHE:CE1	2.52	0.45
1:A:111:GLN:CA	1:A:134:ILE:HD11	2.47	0.45
1:A:395:TRP:C	1:A:398:VAL:HG12	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:THR:O	1:A:103:LEU:HB2	2.16	0.45
1:A:127:THR:HG22	1:A:128:ASP:N	2.32	0.45
1:A:549:LEU:HD13	1:A:549:LEU:O	2.17	0.45
1:A:126:LEU:HB3	1:A:130:GLU:CG	2.44	0.44
1:A:679:GLU:H	1:A:679:GLU:HG3	1.52	0.44
1:A:28:PHE:CZ	1:A:32:ILE:HG13	2.52	0.44
1:A:435:ARG:NH1	1:A:438:GLN:OE1	2.51	0.44
1:A:600:ARG:HA	1:A:600:ARG:HD3	1.55	0.44
1:A:94:MET:CE	1:A:329:TYR:HB3	2.46	0.44
1:A:506:ILE:O	1:A:510:TRP:HB3	2.18	0.44
1:A:526:LEU:O	1:A:530:LEU:HD13	2.18	0.44
1:A:134:ILE:CG1	1:A:135:ARG:N	2.81	0.44
1:A:423:LEU:CD2	1:A:436:ILE:HG12	2.48	0.44
1:A:363:ILE:CG2	1:A:515:PHE:CZ	2.99	0.44
1:A:306:GLY:HA3	1:A:675:ILE:HD13	2.00	0.44
1:A:84:TYR:C	1:A:92:SER:HB2	2.37	0.44
1:A:112:ASP:O	1:A:115:ALA:HB3	2.16	0.44
1:A:522:VAL:O	1:A:526:LEU:HD13	2.18	0.44
1:A:56:ARG:HD3	1:A:62:VAL:HG12	1.99	0.44
1:A:706:LEU:HB3	1:A:707:PRO:CD	2.47	0.44
1:A:64:MET:HE1	1:A:212:ASP:CB	2.48	0.44
1:A:433:LEU:HD13	1:A:438:GLN:HG3	1.98	0.44
1:A:55:VAL:HG22	1:A:64:MET:CB	2.46	0.44
1:A:236:TRP:CG	1:A:237:PRO:HA	2.53	0.44
1:A:339:ILE:HG22	1:A:340:ASN:H	1.82	0.44
1:A:71:GLY:HA3	1:A:209:GLN:OE1	2.18	0.44
1:A:440:ILE:HD12	1:A:440:ILE:C	2.38	0.44
1:A:731:GLN:NE2	8:A:1006:HOH:O	2.51	0.44
1:A:524:GLY:O	1:A:528:VAL:HG23	2.18	0.43
1:A:56:ARG:CB	1:A:61:GLU:O	2.65	0.43
1:A:744:VAL:HB	1:A:745:PRO:HD3	1.99	0.43
1:A:590:LEU:HD23	1:A:590:LEU:C	2.38	0.43
1:A:442:VAL:HG23	1:A:443:VAL:N	2.32	0.43
1:A:646:PHE:O	1:A:649:PRO:HD2	2.17	0.43
1:A:294:ARG:NH2	1:A:678:PRO:HB2	2.33	0.43
1:A:549:LEU:HD13	1:A:549:LEU:C	2.39	0.43
1:A:131:LYS:HA	1:A:134:ILE:HG23	2.00	0.43
1:A:596:TYR:CD1	7:A:810:LOP:H81	2.54	0.43
1:A:762:LYS:O	1:A:762:LYS:HG3	2.17	0.43
1:A:293:LEU:HA	1:A:296:GLN:CD	2.39	0.43
1:A:51:ARG:CD	1:A:495:ASN:ND2	2.80	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:810:LOP:H112	7:A:810:LOP:C27	2.40	0.43
1:A:298:LEU:HD11	1:A:303:VAL:HG13	2.01	0.43
1:A:181:TRP:O	1:A:181:TRP:CD1	2.72	0.42
1:A:44:ILE:CG1	1:A:567:PRO:HD2	2.46	0.42
1:A:578:SER:O	1:A:581:GLU:HB2	2.19	0.42
1:A:64:MET:HE1	1:A:212:ASP:CA	2.49	0.42
1:A:750:LEU:O	1:A:753:VAL:CG1	2.67	0.42
1:A:144:LYS:HA	1:A:144:LYS:HZ3	1.83	0.42
1:A:433:LEU:HD13	1:A:438:GLN:HG2	2.01	0.42
1:A:47:GLU:OE2	5:A:807:BOG:H62	2.19	0.42
1:A:143:ARG:HD2	1:A:243:GLY:HA3	2.00	0.42
1:A:685:TRP:HE3	1:A:685:TRP:HA	1.85	0.42
1:A:773:GLY:C	1:A:775:GLY:H	2.21	0.42
1:A:86:THR:HG22	1:A:91:GLY:C	2.40	0.42
1:A:258:VAL:O	1:A:262:ILE:HG12	2.19	0.42
1:A:331:THR:HG21	1:A:740:LEU:CD1	2.49	0.42
1:A:646:PHE:HA	1:A:649:PRO:HD2	2.01	0.42
7:A:810:LOP:C10	7:A:810:LOP:H251	2.50	0.42
1:A:93:TYR:CE2	1:A:649:PRO:CG	3.02	0.42
1:A:336:PHE:O	1:A:337:PHE:HB2	2.20	0.42
1:A:347:PRO:CD	1:A:350:ILE:HD12	2.48	0.42
1:A:367:TRP:CH2	1:A:519:ALA:HB1	2.55	0.42
1:A:588:LEU:HD12	1:A:588:LEU:HA	1.87	0.41
1:A:691:TRP:CE2	7:A:810:LOP:H282	2.54	0.41
1:A:176:THR:O	1:A:186:GLY:N	2.53	0.41
1:A:462:ARG:HG2	1:A:462:ARG:NH1	2.35	0.41
1:A:48:MET:SD	1:A:498:MET:HB2	2.61	0.41
1:A:370:MET:CE	1:A:526:LEU:HD21	2.49	0.41
1:A:62:VAL:C	1:A:63:LEU:HD13	2.41	0.41
1:A:341:TRP:HE3	1:A:345:ILE:HG21	1.85	0.41
1:A:321:MET:HE2	1:A:750:LEU:HD21	2.01	0.41
1:A:50:PRO:O	1:A:52:PRO:HD3	2.21	0.41
1:A:56:ARG:HB3	1:A:62:VAL:CA	2.36	0.41
1:A:198:ARG:NH2	1:A:204:SER:O	2.52	0.41
1:A:400:LEU:HG	1:A:440:ILE:CD1	2.50	0.41
1:A:56:ARG:HD2	1:A:62:VAL:HG12	2.03	0.41
1:A:113:TYR:HB2	1:A:162:TYR:HD2	1.83	0.41
1:A:187:LEU:H	1:A:187:LEU:CD1	2.34	0.41
1:A:208:GLN:O	1:A:212:ASP:OD1	2.38	0.41
1:A:391:ASP:O	1:A:394:PHE:HB3	2.20	0.41
1:A:127:THR:N	1:A:130:GLU:HB3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:ASP:HA	1:A:343:TYR:CE2	2.55	0.41
1:A:56:ARG:CB	1:A:62:VAL:HA	2.35	0.41
1:A:654:ALA:O	1:A:658:GLY:HA3	2.21	0.41
1:A:750:LEU:O	1:A:753:VAL:HG13	2.20	0.41
7:A:810:LOP:O7	7:A:810:LOP:C9	2.59	0.41
1:A:675:ILE:C	1:A:768:ARG:HD3	2.42	0.41
2:A:801:HEM:HBA2	2:A:801:HEM:HHA	2.03	0.41
1:A:230:ILE:HD12	1:A:234:ASN:HA	2.03	0.40
1:A:309:PHE:HA	1:A:312:VAL:CG1	2.50	0.40
1:A:310:VAL:HG21	1:A:764:LEU:HD21	2.03	0.40
1:A:417:GLY:C	1:A:419:GLU:H	2.23	0.40
1:A:81:LEU:HD21	1:A:175:PHE:HE1	1.85	0.40
1:A:239:TYR:CE1	1:A:241:ASP:HB2	2.57	0.40
1:A:326:LEU:HD11	1:A:352:LYS:HA	2.04	0.40
1:A:393:LEU:HG	1:A:447:LEU:HD13	2.03	0.40
1:A:63:LEU:HD13	1:A:63:LEU:N	2.36	0.40
1:A:596:TYR:CD1	7:A:810:LOP:C8	3.04	0.40
1:A:761:ILE:O	1:A:764:LEU:CD2	2.69	0.40
1:A:761:ILE:C	1:A:763:ALA:H	2.25	0.40

There are no symmetry-related clashes.

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	750/800 (94%)	695 (93%)	49 (6%)	6 (1%)	19	35

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	ASP
1	A	157	THR

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Mol	Chain	Res	Type
1	A	199	ALA
1	A	204	SER
1	A	194	PRO
1	A	603	GLY

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	631/675 (94%)	549 (87%)	82 (13%)	4 7

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
1	A	37	LEU
1	A	51	ARG
1	A	53	LYS
1	A	55	VAL
1	A	63	LEU
1	A	68	THR
1	A	81	LEU
1	A	87	VAL
1	A	96	PRO
1	A	103	LEU
1	A	105	VAL
1	A	126	LEU
1	A	134	ILE
1	A	135	ARG
1	A	148	ASN
1	A	151	THR
1	A	154	LEU
1	A	156	LEU
1	A	164	LEU
1	A	166	LYS
1	A	187	LEU

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Mol	Chain	Res	Type
1	A	198	ARG
1	A	207	ILE
1	A	210	ILE
1	A	220	LEU
1	A	231	THR
1	A	251	VAL
1	A	287	LYS
1	A	293	LEU
1	A	295	ARG
1	A	298	LEU
1	A	308	TYR
1	A	361	PHE
1	A	382	GLU
1	A	391	ASP
1	A	393	LEU
1	A	397	LEU
1	A	400	LEU
1	A	433	LEU
1	A	435	ARG
1	A	436	ILE
1	A	440	ILE
1	A	441	LEU
1	A	446	LEU
1	A	447	LEU
1	A	462	ARG
1	A	469	LEU
1	A	478	ILE
1	A	483	PHE
1	A	497	THR
1	A	512	GLU
1	A	527	LEU
1	A	530	LEU
1	A	545	GLN
1	A	550	LEU
1	A	559	HIS
1	A	587	LEU
1	A	588	LEU
1	A	590	LEU
1	A	591	GLU
1	A	615	LEU
1	A	618	THR
1	A	622	ASN

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Mol	Chain	Res	Type
1	A	635	LEU
1	A	656	MET
1	A	663	PHE
1	A	672	LEU
1	A	673	ARG
1	A	679	GLU
1	A	685	TRP
1	A	706	LEU
1	A	711	LEU
1	A	738	LEU
1	A	739	LEU
1	A	740	LEU
1	A	742	ARG
1	A	750	LEU
1	A	756	LEU
1	A	757	LEU
1	A	768	ARG
1	A	777	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	148	ASN
1	A	340	ASN
1	A	381	GLN
1	A	471	HIS
1	A	529	GLN
1	A	595	GLN
1	A	622	ASN
1	A	682	ASN
1	A	731	GLN

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 ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BOG	A	805	-	20,20,20	0.92	2 (10%)	25,25,25	0.91	1 (4%)
7	LOP	A	810	-	44,44,44	0.71	0	47,49,49	1.17	3 (6%)
6	EPE	A	808	-	15,15,15	1.33	1 (6%)	18,20,20	1.55	3 (16%)
7	LOP	A	809	-	44,44,44	0.69	0	47,49,49	1.20	3 (6%)
2	HEM	A	802	1,8,4	27,50,50	1.45	3 (11%)	17,82,82	1.65	4 (23%)
5	BOG	A	807	-	20,20,20	0.89	2 (10%)	25,25,25	0.99	2 (8%)
2	HEM	A	801	1,4	27,50,50	1.33	2 (7%)	17,82,82	1.91	3 (17%)
5	BOG	A	806	-	20,20,20	1.01	2 (10%)	25,25,25	0.96	1 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BOG	A	805	-	-	0/11/31/31	0/1/1/1
7	LOP	A	810	-	-	13/48/48/48	-
6	EPE	A	808	-	-	1/9/19/19	0/1/1/1
7	LOP	A	809	-	-	13/48/48/48	-
2	HEM	A	802	1,8,4	-	0/6/54/54	-
5	BOG	A	807	-	-	0/11/31/31	0/1/1/1
2	HEM	A	801	1,4	-	2/6/54/54	-
5	BOG	A	806	-	-	0/11/31/31	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	808	EPE	C10-S	3.99	1.83	1.77
2	A	802	HEM	C3C-CAC	-3.47	1.40	1.47
2	A	802	HEM	C4B-NB	3.02	1.42	1.36
5	A	806	BOG	O1-C1	3.00	1.45	1.40
2	A	801	HEM	C4B-NB	2.87	1.42	1.36
2	A	801	HEM	C3C-CAC	-2.68	1.42	1.47
5	A	807	BOG	O1-C1	2.44	1.44	1.40
2	A	802	HEM	CAD-C3D	2.36	1.56	1.52
5	A	805	BOG	O1-C1	2.34	1.44	1.40
5	A	805	BOG	O5-C1	2.28	1.47	1.41
5	A	806	BOG	O5-C1	2.19	1.47	1.41
5	A	807	BOG	O5-C1	2.10	1.47	1.41

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	CBA-CAA-C2A	5.12	121.92	112.49
6	A	808	EPE	O1S-S-C10	4.88	112.79	106.92
2	A	802	HEM	CBD-CAD-C3D	4.16	120.14	112.48
7	A	809	LOP	C4-O5-C6	-4.08	107.76	117.79
7	A	810	LOP	O5-C6-C7	4.01	120.15	111.50
7	A	809	LOP	O5-C6-C7	3.97	120.07	111.50
5	A	807	BOG	C1'-O1-C1	-2.93	108.98	113.84
2	A	801	HEM	C4C-C3C-C2C	-2.79	104.95	106.90
5	A	806	BOG	C1'-O1-C1	-2.64	109.46	113.84
6	A	808	EPE	O3S-S-O2S	-2.37	105.49	111.27
7	A	810	LOP	O6-C24-C25	2.35	119.28	111.91
2	A	802	HEM	C1D-C2D-C3D	2.29	108.59	107.00
2	A	802	HEM	C4C-C3C-C2C	-2.26	105.32	106.90
2	A	801	HEM	CAD-CBD-CGD	2.23	116.41	112.67
7	A	809	LOP	C5-O6-C24	-2.18	109.04	117.12
6	A	808	EPE	O3S-S-O1S	-2.17	105.98	111.27
5	A	805	BOG	C1'-O1-C1	-2.16	110.27	113.84
2	A	802	HEM	CMB-C2B-C3B	2.14	128.69	124.68
5	A	807	BOG	C6-C5-C4	-2.04	108.22	113.00
7	A	810	LOP	O6-C5-C4	-2.03	102.52	108.43

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	810	LOP	N1-C1-C2-O1
7	A	810	LOP	C3-O2-P1-O3
7	A	810	LOP	C4-C3-O2-P1
7	A	809	LOP	C1-C2-O1-P1
7	A	809	LOP	C2-O1-P1-O4
7	A	809	LOP	C4-C3-O2-P1
2	A	801	HEM	C1A-C2A-CAA-CBA
2	A	801	HEM	C3A-C2A-CAA-CBA
6	A	808	EPE	N4-C7-C8-O8
7	A	809	LOP	C24-C25-C26-C27
7	A	810	LOP	C3-O2-P1-O1
7	A	809	LOP	C2-O1-P1-O2
7	A	810	LOP	C29-C30-C31-C32
7	A	810	LOP	C15-C16-C17-C18
7	A	809	LOP	C30-C31-C32-C33
7	A	810	LOP	O2-C3-C4-C5
7	A	810	LOP	C24-C25-C26-C27
7	A	809	LOP	C32-C33-C34-C35
7	A	810	LOP	O2-C3-C4-O5
7	A	810	LOP	C3-O2-P1-O4
7	A	809	LOP	C2-O1-P1-O3
7	A	810	LOP	C2-O1-P1-O2
7	A	809	LOP	C3-O2-P1-O1
7	A	810	LOP	C28-C29-C30-C31
7	A	809	LOP	C27-C28-C29-C30
7	A	809	LOP	C12-C13-C14-C15
7	A	810	LOP	C12-C13-C14-C15
7	A	809	LOP	C6-C7-C8-C9
7	A	809	LOP	C3-O2-P1-O4

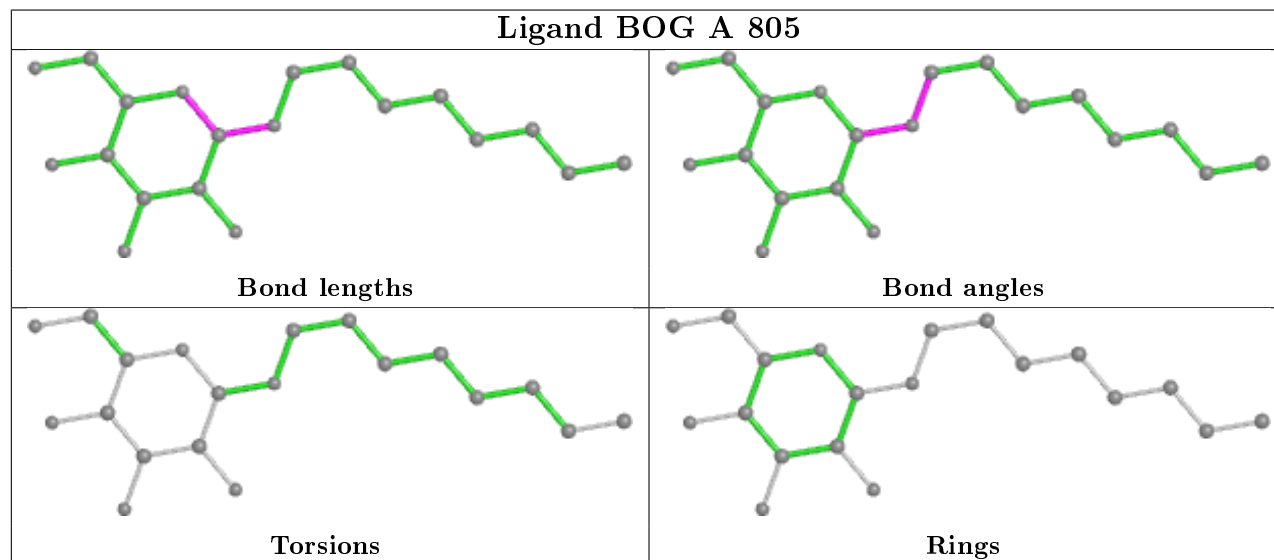
There are no ring outliers.

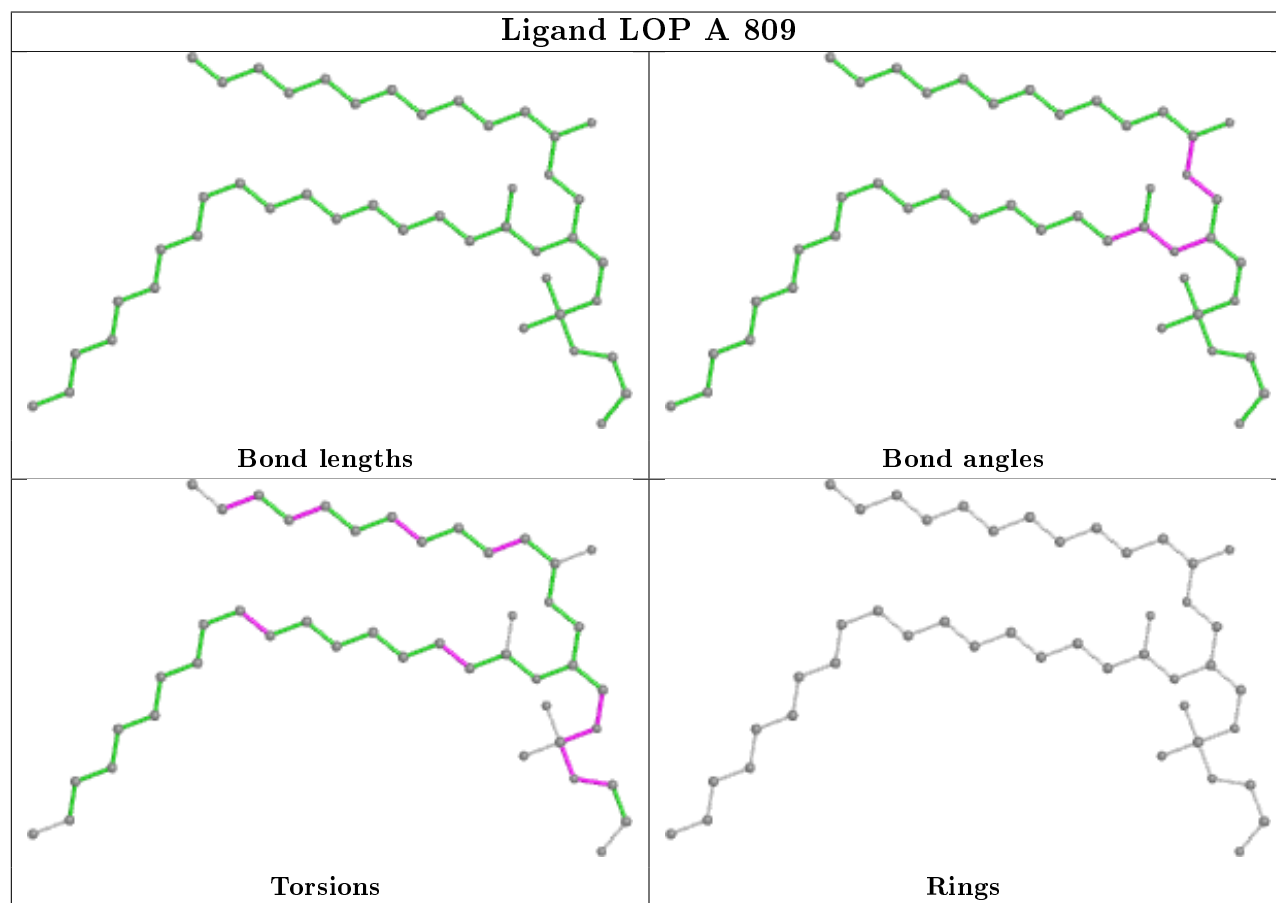
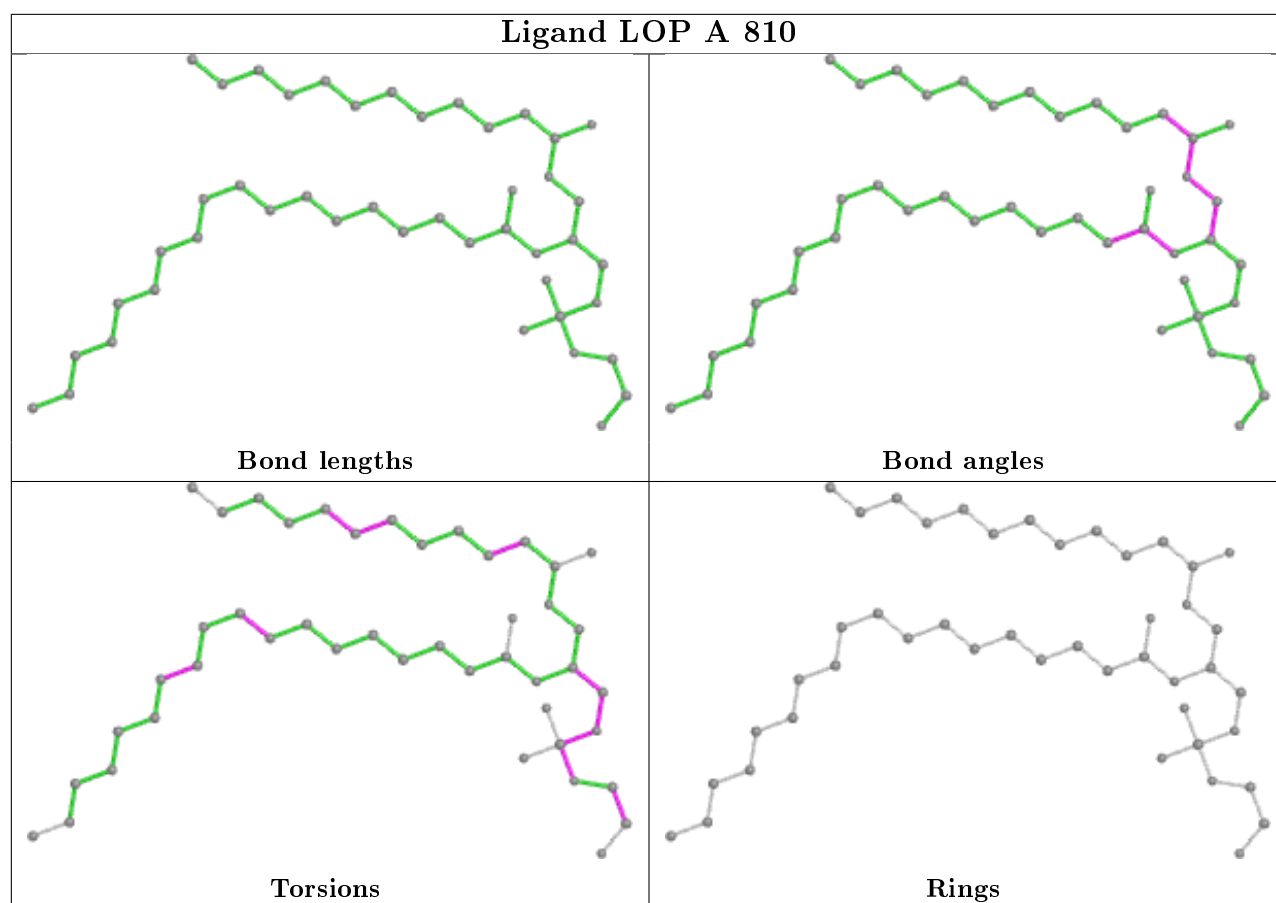
6 monomers are involved in 38 short contacts:

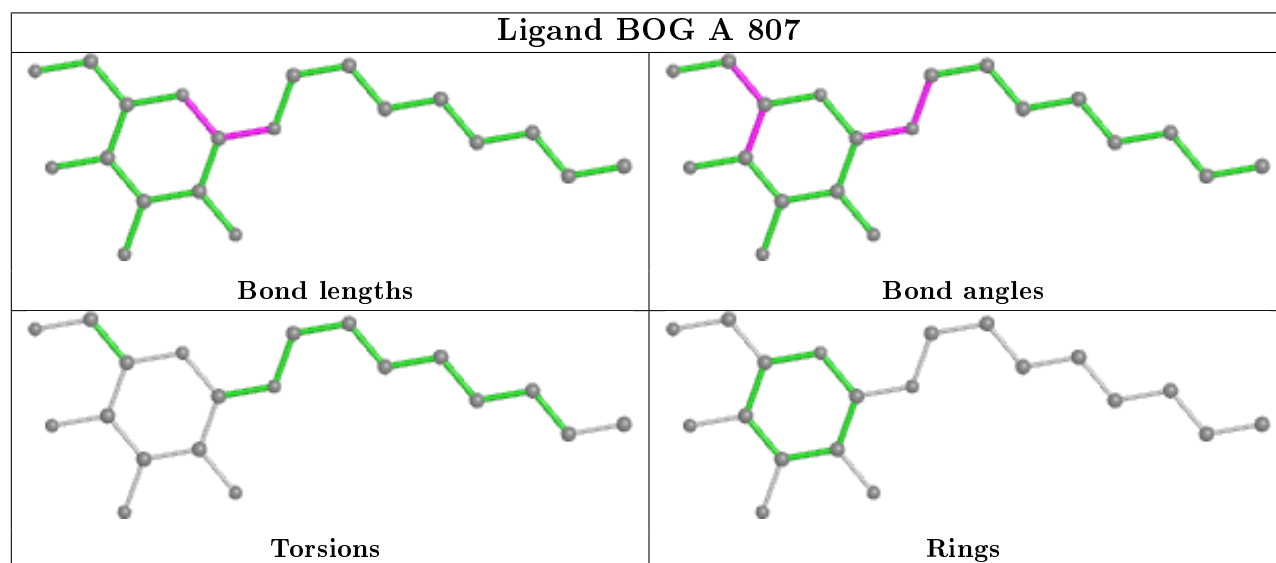
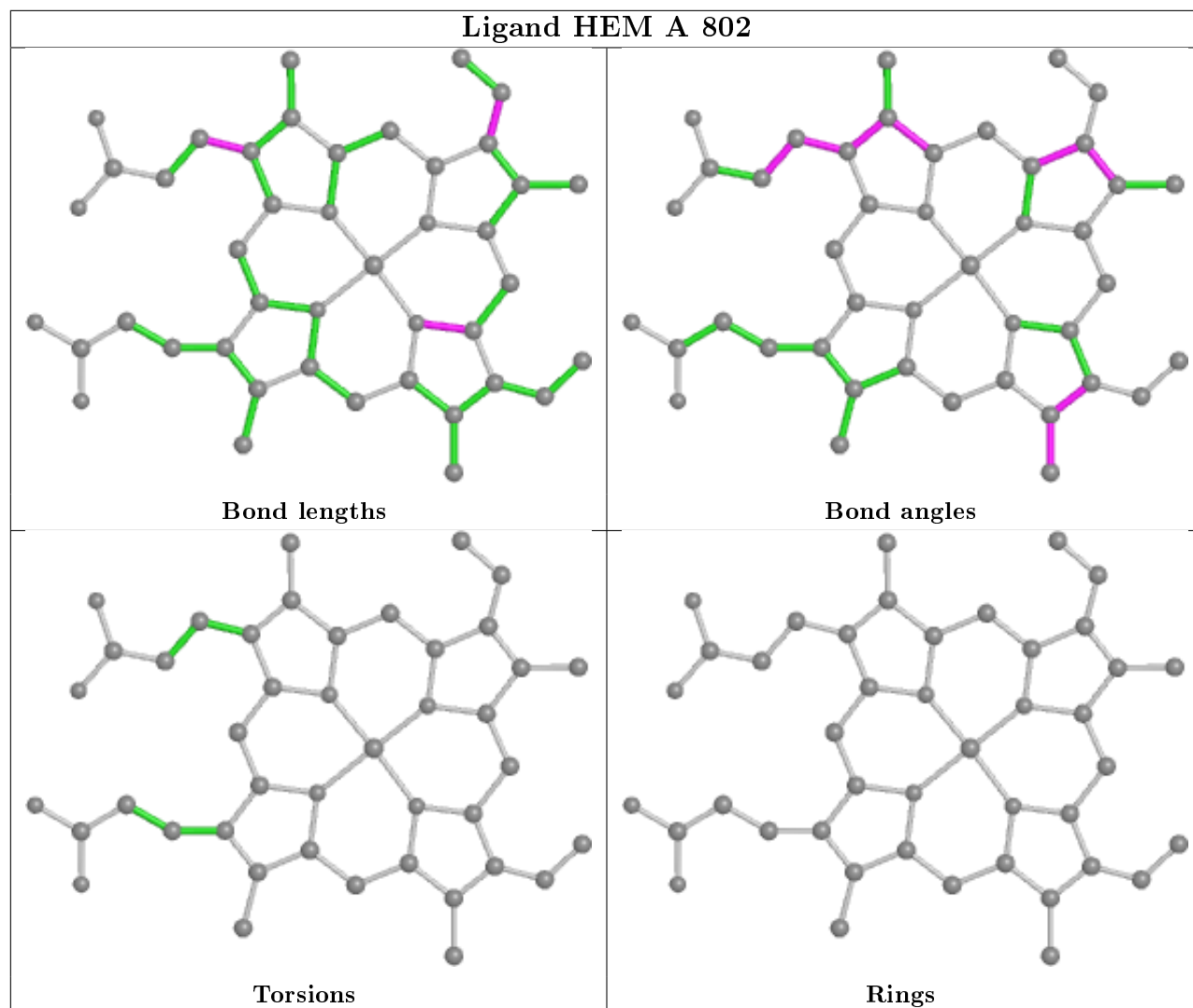
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	810	LOP	20	0
7	A	809	LOP	11	0
2	A	802	HEM	2	0
5	A	807	BOG	3	0
2	A	801	HEM	1	0
5	A	806	BOG	1	0

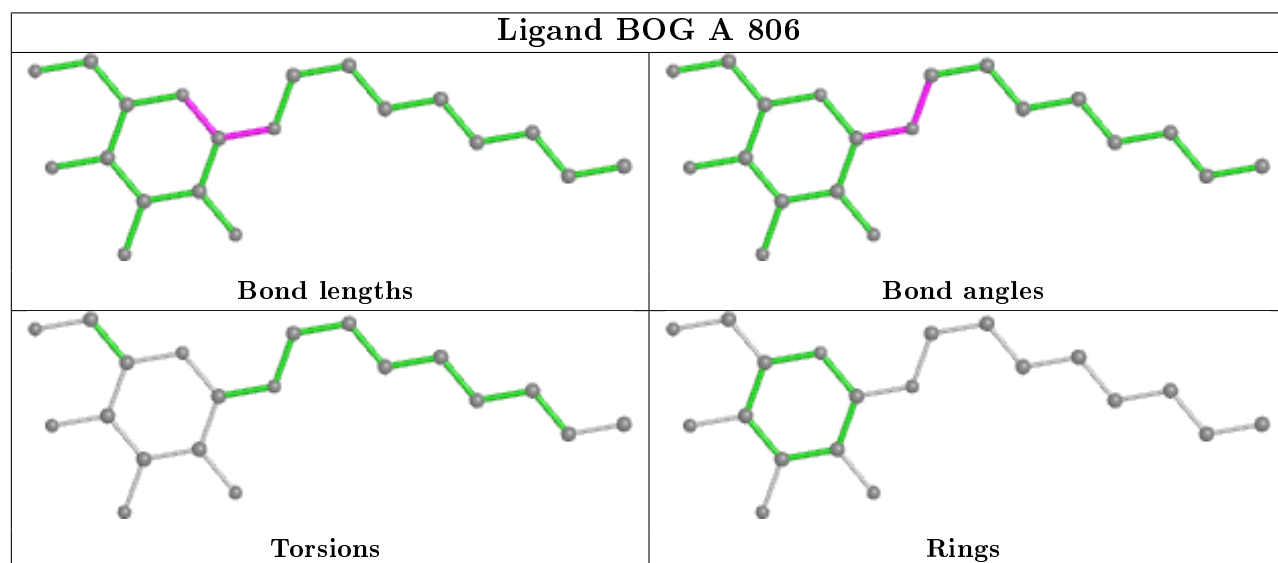
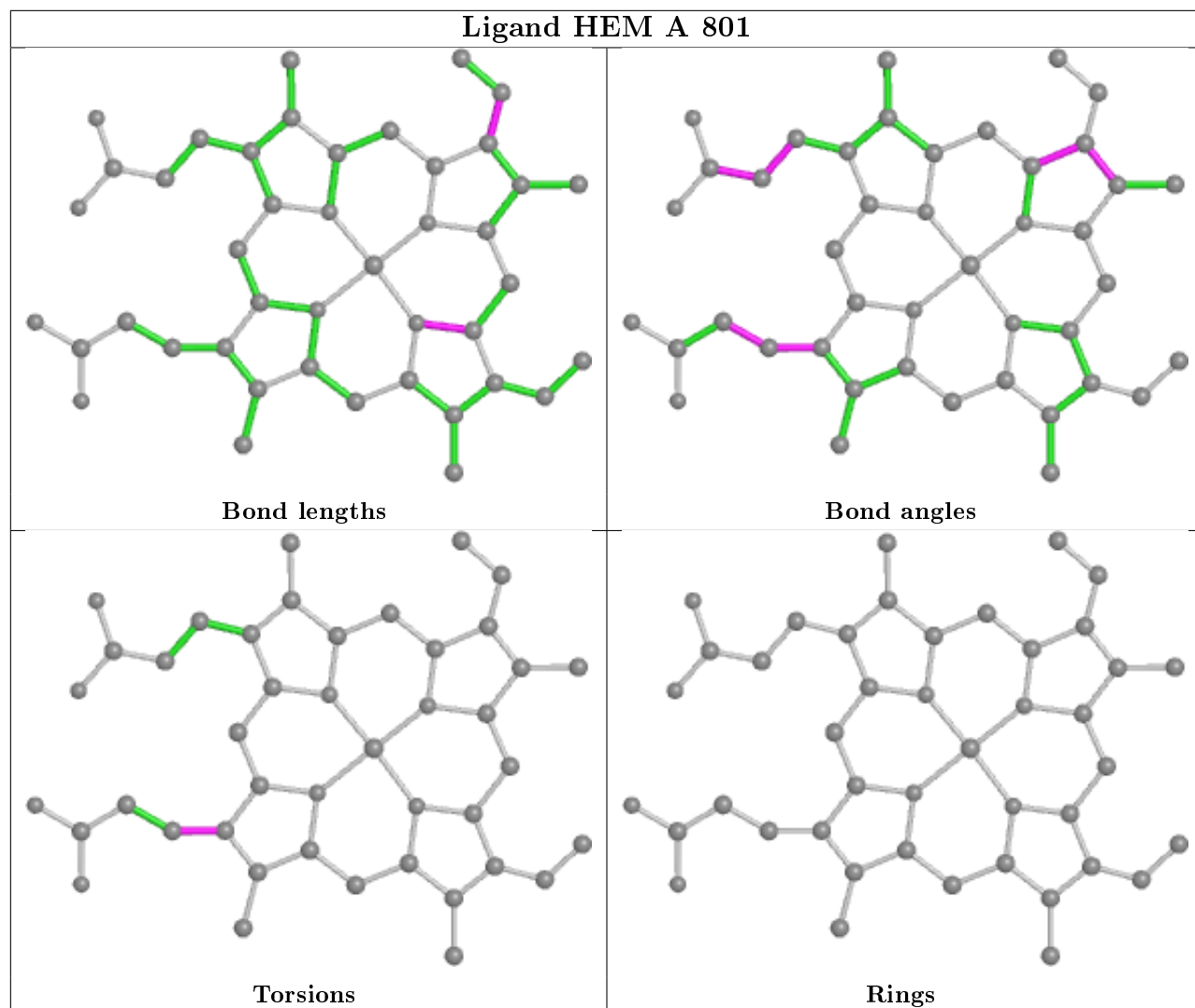
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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.









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	754/800 (94%)	0.52	78 (10%) 6 6	34, 62, 93, 104	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	154	LEU	7.3
1	A	113	TYR	6.4
1	A	394	PHE	5.8
1	A	134	ILE	5.6
1	A	130	GLU	4.9
1	A	416	LEU	4.7
1	A	158	ASP	4.4
1	A	414	GLY	4.3
1	A	155	VAL	4.2
1	A	513	GLY	4.0
1	A	514	ILE	4.0
1	A	163	GLY	4.0
1	A	139	ILE	3.9
1	A	199	ALA	3.8
1	A	776	GLU	3.7
1	A	126	LEU	3.6
1	A	115	ALA	3.6
1	A	59	SER	3.6
1	A	150	VAL	3.6
1	A	300	PRO	3.5
1	A	64	MET	3.4
1	A	141	GLU	3.4
1	A	162	TYR	3.3
1	A	127	THR	3.3
1	A	159	ALA	3.2
1	A	192	ASP	3.2
1	A	161	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	777	GLU	3.2
1	A	743	ALA	3.2
1	A	773	GLY	3.1
1	A	137	GLN	3.1
1	A	198	ARG	3.1
1	A	273	TYR	3.0
1	A	112	ASP	2.9
1	A	114	LYS	2.9
1	A	21	PHE	2.8
1	A	303	VAL	2.8
1	A	421	PHE	2.8
1	A	17	THR	2.7
1	A	109	GLY	2.7
1	A	392	LEU	2.6
1	A	423	LEU	2.6
1	A	194	PRO	2.6
1	A	418	ASN	2.6
1	A	160	GLN	2.6
1	A	135	ARG	2.6
1	A	52	PRO	2.5
1	A	660	TYR	2.5
1	A	153	VAL	2.5
1	A	193	MET	2.5
1	A	60	GLY	2.4
1	A	31	LEU	2.4
1	A	771	THR	2.3
1	A	536	LYS	2.3
1	A	388	LEU	2.3
1	A	691	TRP	2.3
1	A	169	ASP	2.3
1	A	39	VAL	2.3
1	A	395	TRP	2.3
1	A	415	TYR	2.2
1	A	390	VAL	2.2
1	A	290	VAL	2.2
1	A	391	ASP	2.2
1	A	63	LEU	2.2
1	A	195	LYS	2.2
1	A	143	ARG	2.2
1	A	632	LEU	2.1
1	A	772	HIS	2.1
1	A	458	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	125	ASP	2.1
1	A	308	TYR	2.1
1	A	147	TYR	2.1
1	A	633	ILE	2.0
1	A	630	GLY	2.0
1	A	53	LYS	2.0
1	A	258	VAL	2.0
1	A	129	ASP	2.0
1	A	703	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [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.

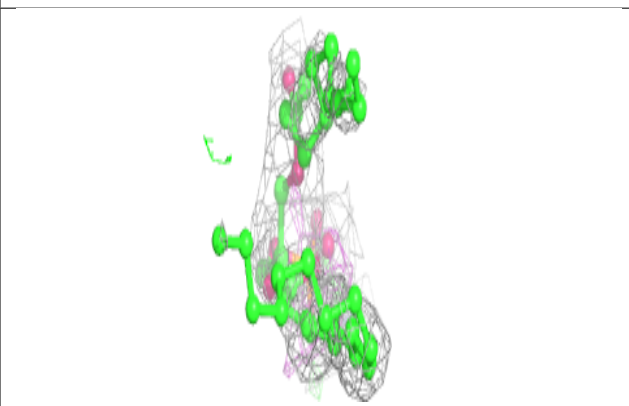
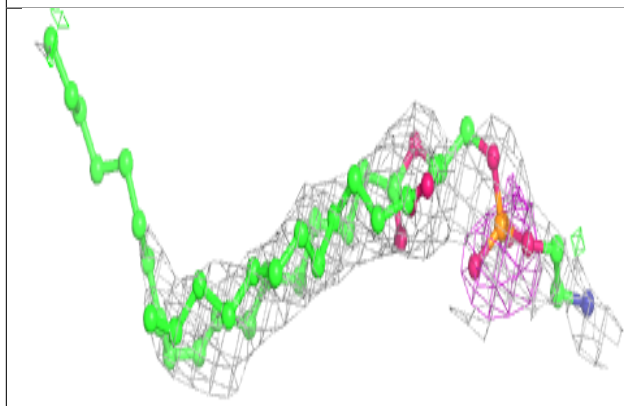
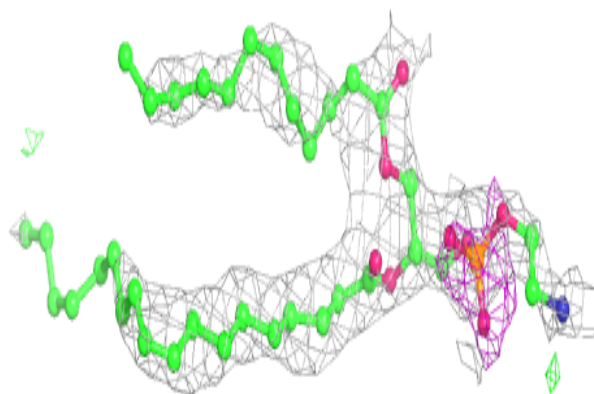
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	EPE	A	808	15/15	0.52	0.46	88,93,109,109	0
7	LOP	A	809	45/45	0.72	0.38	73,85,101,104	0
5	BOG	A	805	20/20	0.73	0.24	75,89,93,95	0
7	LOP	A	810	45/45	0.74	0.35	59,70,102,104	0
5	BOG	A	807	20/20	0.77	0.26	73,85,88,90	0
5	BOG	A	806	20/20	0.84	0.37	76,87,90,91	0
2	HEM	A	801	43/43	0.95	0.23	38,46,52,54	0
2	HEM	A	802	43/43	0.97	0.21	29,38,50,54	0
4	CA	A	804	1/1	0.98	0.31	49,49,49,49	0
3	ZN	A	803	1/1	1.00	0.16	55,55,55,55	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.

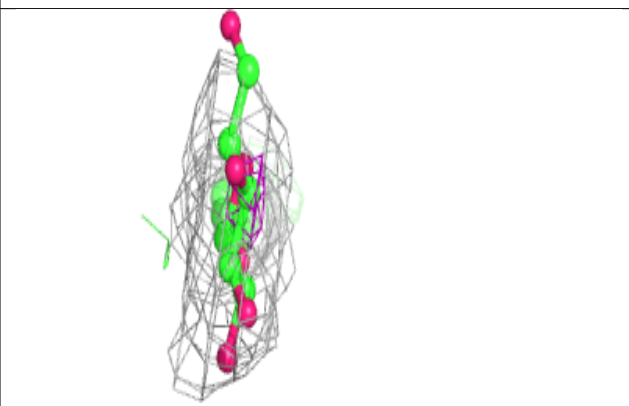
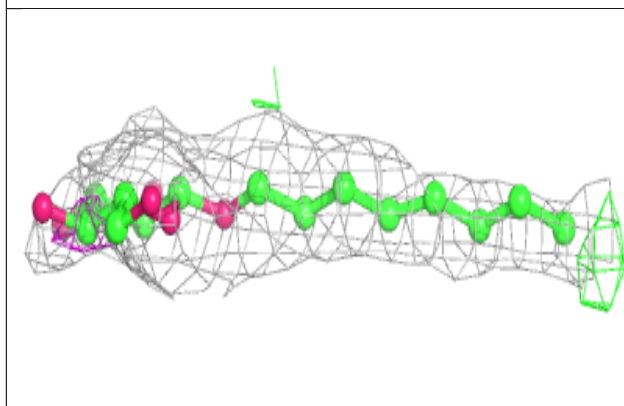
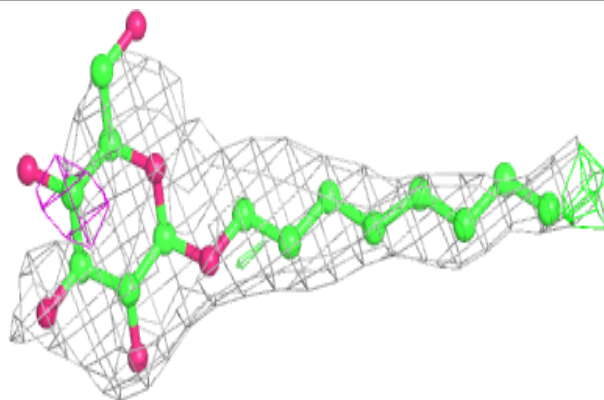
Electron density around LOP A 809:

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



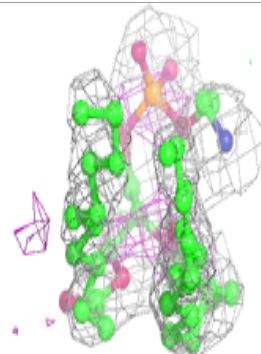
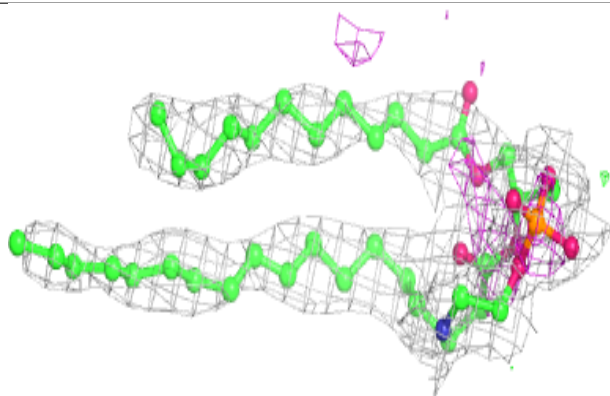
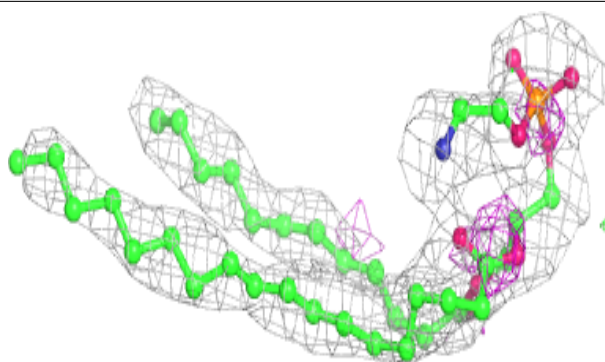
Electron density around BOG A 805:

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

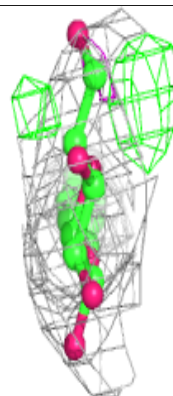
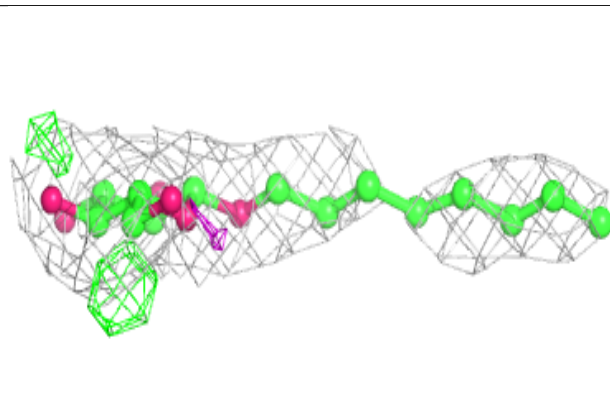
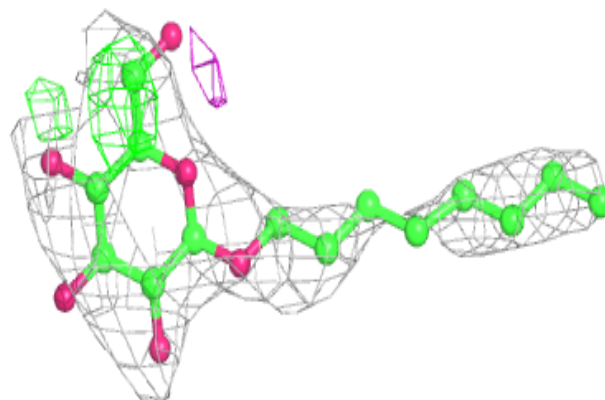


Electron density around LOP A 810:

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

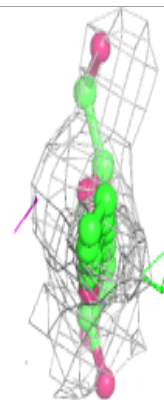
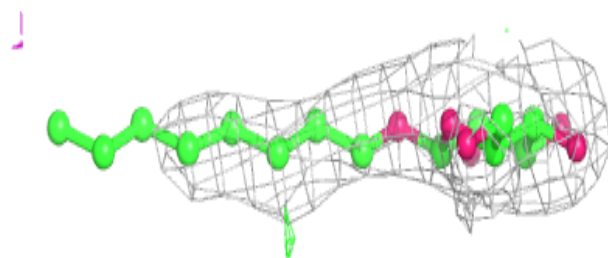
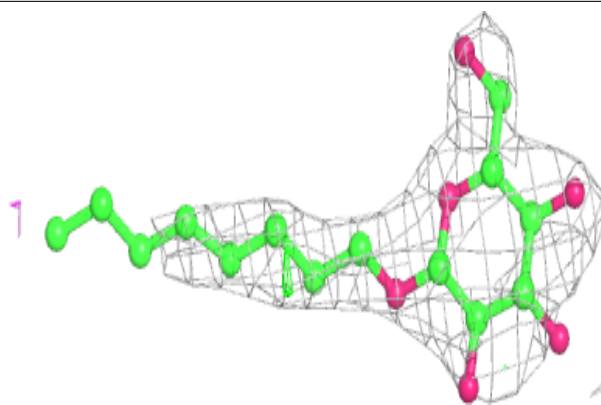
**Electron density around BOG A 807:**

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



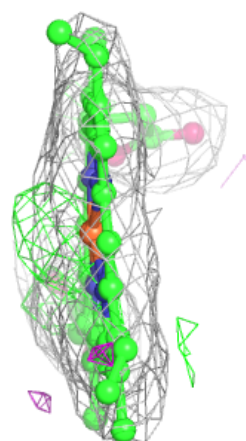
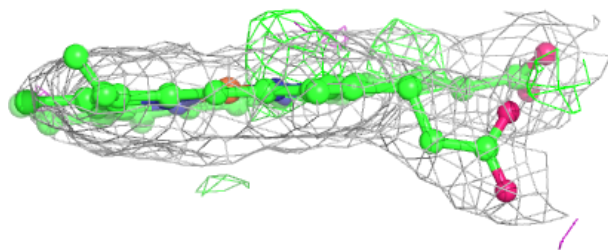
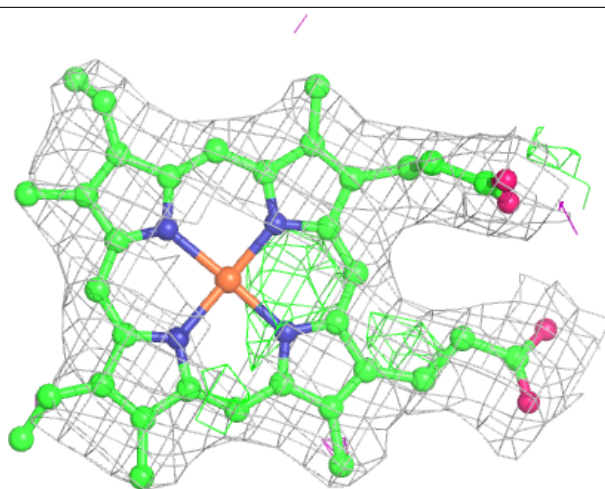
Electron density around BOG A 806:

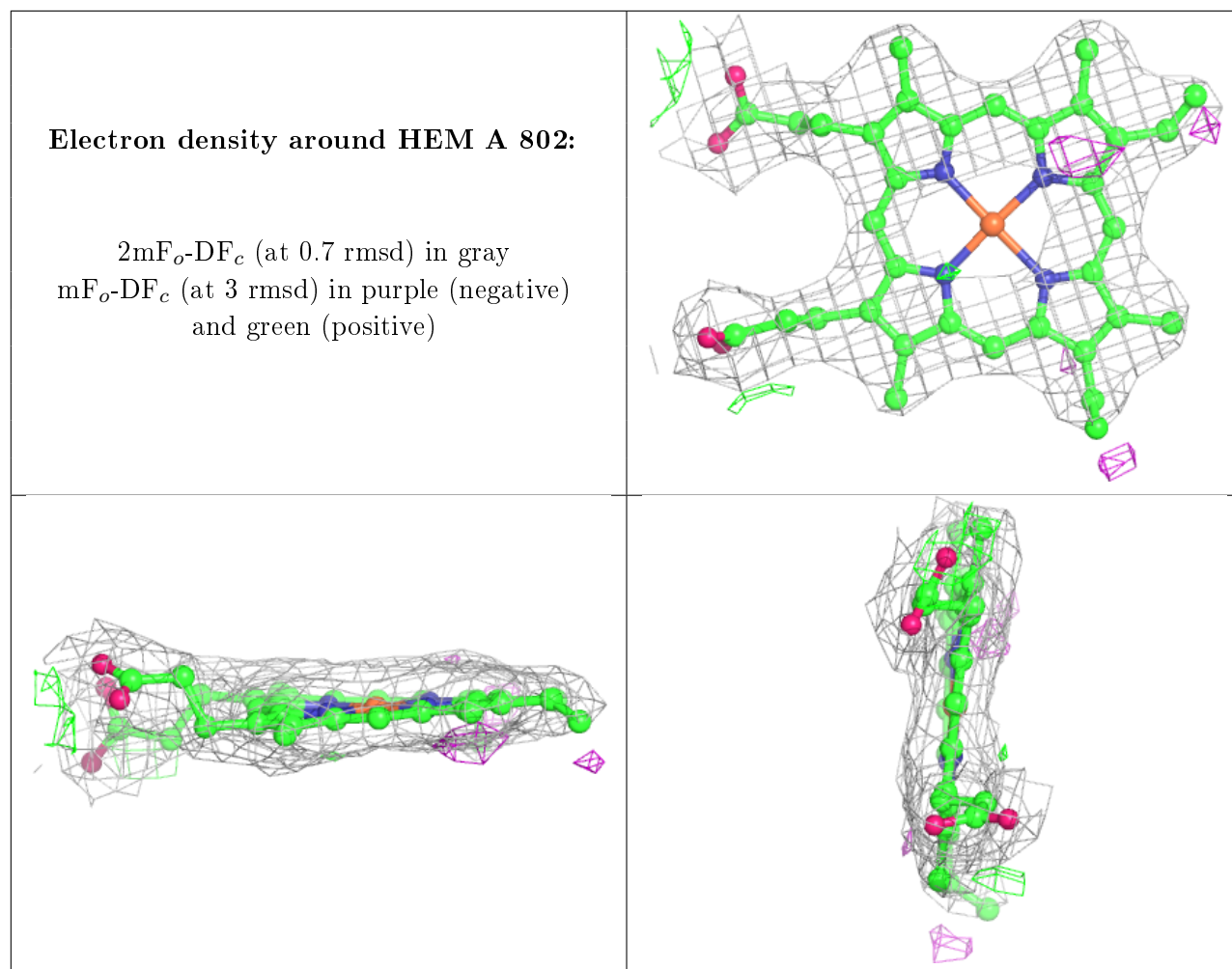
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 801:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.