



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2023 – 11:58 PM JST

PDB ID : 7C73  
Title : Crystal structure of yak lactoperoxidase using data obtained from crystals soaked in MgCl<sub>2</sub> at 2.70 Å resolution  
Authors : Singh, P.K.; Pandey, S.N.; Rani, C.; Ahmad, N.; Viswanathan, V.; Sharma, P.; Kaur, P.; Sharma, S.; Singh, T.P.  
Deposited on : 2020-05-22  
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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

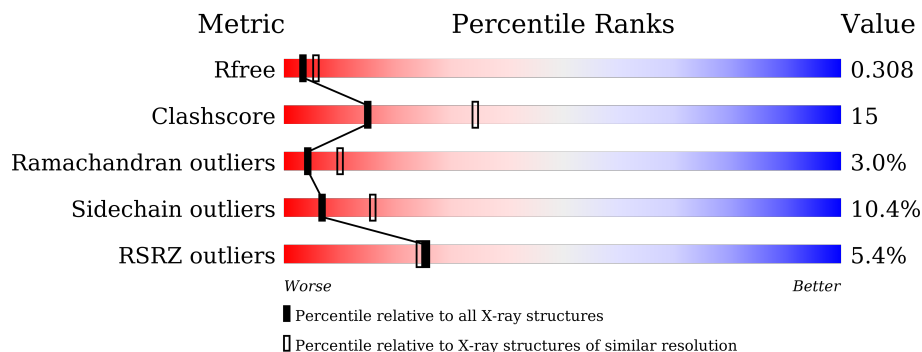
# 1 Overall quality at a glance

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))
$R_{free}$	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (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  $\geq 3$ , 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	595	

## 2 Entry composition [i](#)

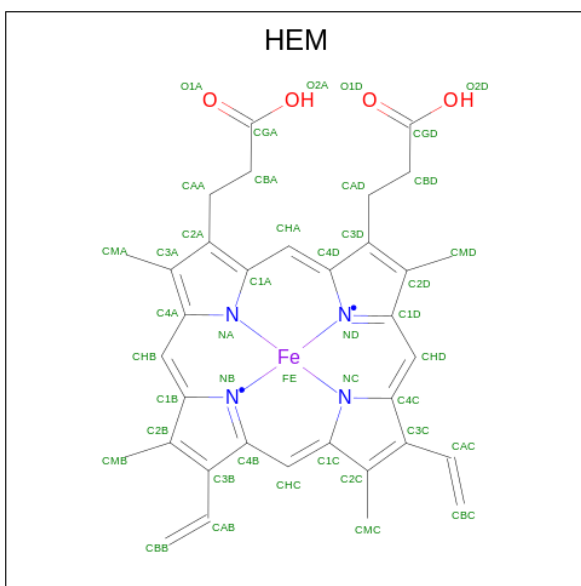
There are 5 unique types of molecules in this entry. The entry contains 5103 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lactoperoxidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	595	4774	3037	847	863	1	26	0	0	0

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ) (labeled as "Ligand of Interest" by depositor).

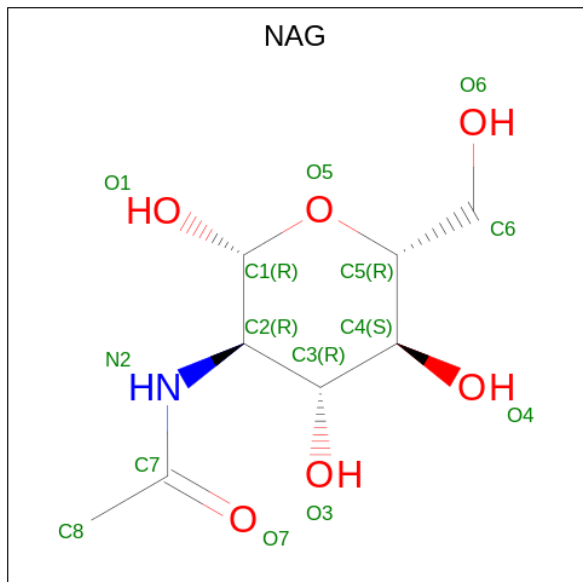


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Ca	0	0
			2	2		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	14	8	1	5	0	0
4	A	1	14	8	1	5	0	0
4	A	1	14	8	1	5	0	0

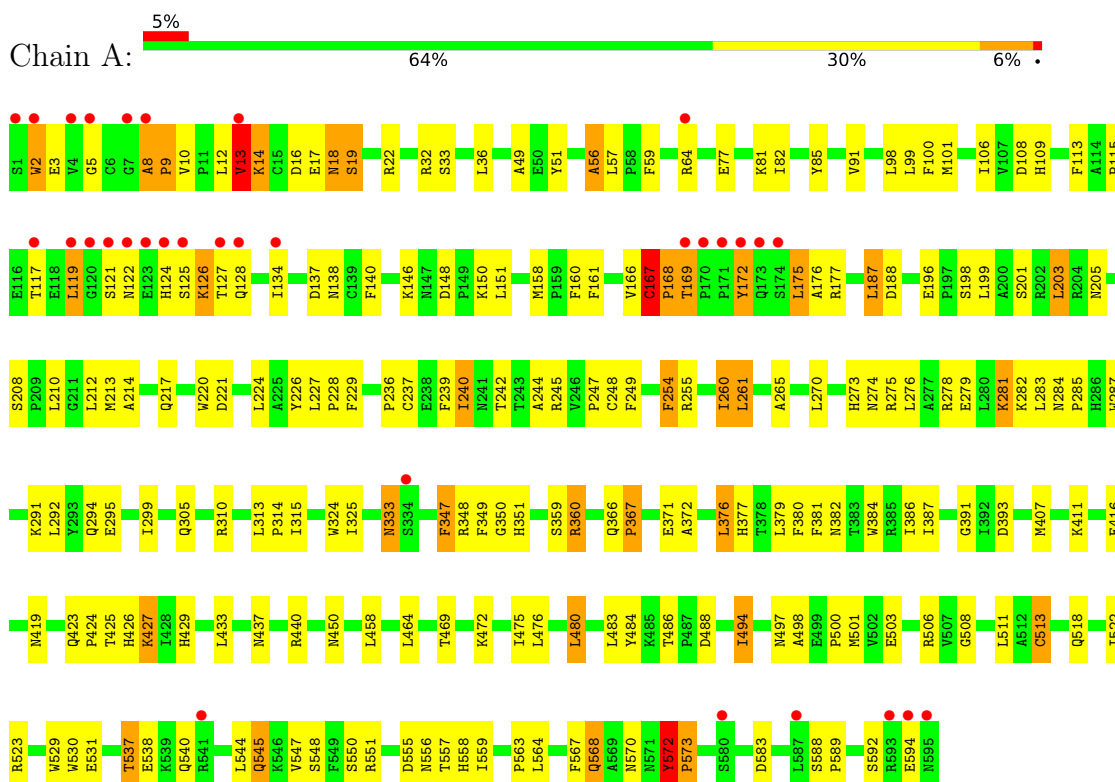
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	242	Total	O	0	0
			242	242		

### 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: Lactoperoxidase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.88Å 79.93Å 75.05Å 90.00° 101.35° 90.00°	Depositor
Resolution (Å)	47.58 – 2.70 47.58 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.6 (47.58-2.70) 97.6 (47.58-2.70)	Depositor EDS
$R_{merge}$	0.20	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.62 (at 2.69Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.218 , 0.306 0.220 , 0.308	Depositor DCC
$R_{free}$ test set	843 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.1	Xtrriage
Anisotropy	0.131	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 40.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5103	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: SEP, HEM, NAG, CA

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	0.66	0/4891	0.88	2/6634 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	CYS	CB-CA-C	5.60	121.60	110.40
1	A	254	PHE	CB-CA-C	5.45	121.30	110.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	VAL	Peptide
1	A	172	TYR	Peptide
1	A	2	TRP	Peptide
1	A	572	TYR	Peptide

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4774	0	4688	139	0
2	A	43	0	30	6	0
3	A	2	0	0	0	0
4	A	42	0	39	4	0
5	A	242	0	0	16	0
All	All	5103	0	4757	145	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 15.

All (145) 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:196:GLU:HB3	1:A:198:SEP:O1P	1.67	0.94
1:A:294:GLN:NE2	5:A:701:HOH:O	2.13	0.82
1:A:8:ALA:O	1:A:10:VAL:N	2.20	0.74
1:A:260:ILE:HD11	1:A:386:ILE:HG13	1.71	0.71
1:A:196:GLU:CB	1:A:198:SEP:O1P	2.41	0.69
1:A:8:ALA:HB3	1:A:9:PRO:HD3	1.75	0.69
4:A:604:NAG:H3	5:A:712:HOH:O	1.91	0.68
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.30	0.66
1:A:8:ALA:CB	1:A:9:PRO:HD3	2.27	0.65
1:A:205:ASN:HD22	1:A:212:LEU:HD12	1.60	0.65
1:A:59:PHE:HD1	5:A:829:HOH:O	1.80	0.65
1:A:167:CYS:HB2	1:A:168:PRO:CD	2.26	0.65
1:A:229:PHE:CD1	1:A:247:PRO:HG2	2.30	0.65
4:A:604:NAG:C3	5:A:712:HOH:O	2.44	0.65
1:A:572:TYR:CD2	1:A:573:PRO:HD3	2.32	0.65
1:A:167:CYS:HB2	1:A:168:PRO:HD3	1.79	0.64
1:A:137:ASP:O	5:A:702:HOH:O	2.15	0.64
1:A:167:CYS:CB	1:A:168:PRO:HD3	2.27	0.64
1:A:325:ILE:HG22	5:A:704:HOH:O	1.97	0.64
1:A:140:PHE:O	1:A:160:PHE:HB3	1.98	0.63
1:A:199:LEU:HD22	1:A:254:PHE:CZ	2.33	0.62
1:A:227:LEU:HD22	1:A:270:LEU:HD22	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:ASN:O	1:A:19:SER:C	2.39	0.61
1:A:220:TRP:HD1	5:A:870:HOH:O	1.81	0.61
1:A:450:ASN:ND2	1:A:488:ASP:OD1	2.26	0.60
1:A:124:HIS:O	1:A:127:THR:N	2.35	0.60
1:A:226:TYR:OH	1:A:391:GLY:HA2	2.02	0.59
1:A:550:SER:OG	1:A:563:PRO:O	2.13	0.59
1:A:429:HIS:NE2	5:A:707:HOH:O	2.32	0.59
1:A:9:PRO:HG2	1:A:167:CYS:O	2.04	0.58
1:A:351:HIS:CE1	1:A:433:LEU:HD21	2.39	0.58
1:A:393:ASP:OD1	5:A:703:HOH:O	2.17	0.58
1:A:8:ALA:HB3	1:A:9:PRO:CD	2.34	0.58
1:A:281:LYS:HD3	1:A:285:PRO:HA	1.85	0.57
1:A:426:HIS:O	1:A:427:LYS:HG2	2.04	0.57
1:A:3:GLU:HG2	1:A:175:LEU:HD13	1.86	0.57
1:A:544:LEU:O	1:A:547:VAL:HG22	2.04	0.57
2:A:601:HEM:HMC2	2:A:601:HEM:HBC2	1.87	0.57
1:A:518:GLN:O	1:A:522:ILE:HG23	2.05	0.56
1:A:217:GLN:HB2	5:A:897:HOH:O	2.05	0.55
1:A:237:CYS:HA	1:A:381:PHE:O	2.06	0.55
1:A:476:LEU:HD21	1:A:498:ALA:HB1	1.88	0.55
1:A:523:ARG:HG3	1:A:529:TRP:CE2	2.41	0.55
1:A:9:PRO:CD	1:A:167:CYS:O	2.55	0.55
2:A:601:HEM:HBB2	2:A:601:HEM:HMB2	1.88	0.55
1:A:423:GLN:HE21	1:A:423:GLN:HA	1.72	0.55
1:A:260:ILE:HG23	1:A:261:LEU:HD13	1.88	0.54
1:A:167:CYS:CB	1:A:168:PRO:CD	2.86	0.54
1:A:545:GLN:HE21	1:A:545:GLN:H	1.55	0.54
1:A:260:ILE:HG13	1:A:386:ILE:HD11	1.90	0.54
1:A:12:LEU:O	1:A:13:VAL:O	2.26	0.54
1:A:227:LEU:HB2	1:A:249:PHE:CD1	2.43	0.54
1:A:8:ALA:CB	1:A:9:PRO:CD	2.86	0.53
1:A:113:PHE:CE2	1:A:115:PRO:HG3	2.42	0.53
1:A:551:ARG:HD3	1:A:583:ASP:O	2.08	0.53
1:A:333:ASN:O	1:A:333:ASN:ND2	2.39	0.53
1:A:175:LEU:HG	1:A:176:ALA:N	2.24	0.53
1:A:295:GLU:O	1:A:299:ILE:HG13	2.08	0.53
1:A:548:SER:OG	1:A:551:ARG:HB2	2.09	0.52
1:A:82:ILE:HD11	1:A:483:LEU:HD12	1.92	0.52
1:A:472:LYS:HD2	1:A:500:PRO:HG2	1.90	0.52
1:A:124:HIS:O	1:A:126:LYS:N	2.42	0.52
1:A:494:ILE:C	1:A:494:ILE:HD13	2.30	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:ARG:HH11	1:A:437:ASN:ND2	2.08	0.52
1:A:572:TYR:HD2	1:A:573:PRO:HD3	1.73	0.52
1:A:530:TRP:CE2	1:A:531:GLU:HG3	2.45	0.51
1:A:9:PRO:CG	1:A:167:CYS:O	2.59	0.51
1:A:494:ILE:HD13	1:A:494:ILE:O	2.10	0.51
1:A:221:ASP:HB2	1:A:226:TYR:CE1	2.47	0.50
1:A:384:TRP:CZ2	4:A:605:NAG:O7	2.64	0.50
1:A:287:TRP:HA	5:A:802:HOH:O	2.11	0.50
1:A:242:THR:O	1:A:245:ARG:HG2	2.12	0.49
1:A:588:SER:N	1:A:589:PRO:HD2	2.28	0.49
1:A:376:LEU:HG	5:A:864:HOH:O	2.11	0.49
1:A:333:ASN:ND2	1:A:333:ASN:C	2.65	0.49
1:A:187:LEU:O	1:A:305:GLN:NE2	2.45	0.49
1:A:349:PHE:CB	1:A:497:ASN:HD21	2.25	0.49
1:A:360:ARG:NH1	1:A:371:GLU:O	2.46	0.49
1:A:150:LYS:NZ	1:A:419:ASN:O	2.46	0.48
1:A:161:PHE:CD1	1:A:161:PHE:N	2.81	0.48
1:A:376:LEU:HB3	1:A:416:GLU:HB3	1.96	0.48
1:A:168:PRO:HB2	1:A:169:THR:H	1.51	0.48
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.49	0.48
1:A:99:LEU:HG	1:A:567:PHE:CE1	2.49	0.47
1:A:508:GLY:HA3	5:A:777:HOH:O	2.14	0.47
1:A:359:SER:HA	1:A:372:ALA:O	2.13	0.47
1:A:108:ASP:HA	1:A:347:PHE:CE2	2.50	0.47
1:A:82:ILE:HD13	1:A:480:LEU:HD13	1.96	0.47
1:A:275:ARG:HD2	1:A:555:ASP:HB3	1.97	0.47
1:A:522:ILE:HD12	1:A:522:ILE:C	2.35	0.47
1:A:221:ASP:HB2	1:A:226:TYR:CZ	2.51	0.46
2:A:601:HEM:HBC2	2:A:601:HEM:CMC	2.45	0.46
1:A:224:LEU:HD11	1:A:558:HIS:CE1	2.49	0.46
1:A:325:ILE:O	5:A:704:HOH:O	2.20	0.46
1:A:537:THR:HG23	1:A:540:GLN:OE1	2.16	0.46
1:A:122:ASN:O	1:A:122:ASN:ND2	2.48	0.46
1:A:273:HIS:HD2	1:A:274:ASN:OD1	1.99	0.46
1:A:568:GLN:O	1:A:570:ASN:ND2	2.48	0.45
1:A:203:LEU:HB3	1:A:213:MET:CE	2.47	0.45
1:A:324:TRP:CE2	1:A:513:CYS:HA	2.52	0.45
1:A:188:ASP:N	1:A:188:ASP:OD1	2.49	0.45
1:A:239:PHE:CZ	1:A:427:LYS:HB2	2.52	0.44
1:A:347:PHE:HB3	2:A:601:HEM:HMD3	2.00	0.44
1:A:450:ASN:HB2	5:A:830:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:601:HEM:HBB2	2:A:601:HEM:CMB	2.47	0.44
1:A:160:PHE:CZ	1:A:440:ARG:NH1	2.85	0.44
1:A:187:LEU:HD23	1:A:187:LEU:HA	1.81	0.43
1:A:260:ILE:CD1	1:A:386:ILE:HG13	2.44	0.43
1:A:366:GLN:O	1:A:367:PRO:C	2.56	0.43
1:A:106:ILE:HD11	1:A:265:ALA:HB3	1.99	0.43
1:A:117:THR:OG1	1:A:138:ASN:ND2	2.52	0.43
1:A:214:ALA:O	1:A:228:PRO:CG	2.66	0.43
1:A:106:ILE:HD11	1:A:265:ALA:CB	2.48	0.43
1:A:501:MET:SD	1:A:506:ARG:HA	2.59	0.43
1:A:237:CYS:O	1:A:240:ILE:HD13	2.19	0.43
1:A:484:TYR:C	1:A:486:THR:H	2.22	0.42
1:A:347:PHE:HB3	2:A:601:HEM:CMD	2.49	0.42
1:A:458:LEU:HD23	5:A:815:HOH:O	2.19	0.42
1:A:556:ASN:O	1:A:557:THR:HG23	2.19	0.42
1:A:407:MET:HB3	1:A:501:MET:CE	2.49	0.42
1:A:244:ALA:HB2	4:A:605:NAG:O5	2.20	0.42
1:A:313:LEU:N	1:A:314:PRO:CD	2.82	0.42
1:A:236:PRO:HD2	1:A:248:CYS:SG	2.59	0.42
1:A:32:ARG:HD2	1:A:333:ASN:HD21	1.83	0.42
1:A:109:HIS:HA	1:A:255:ARG:NH2	2.35	0.41
1:A:119:LEU:HD23	1:A:119:LEU:O	2.20	0.41
1:A:100:PHE:HE1	1:A:315:ILE:HG13	1.85	0.41
1:A:379:LEU:HA	1:A:382:ASN:CB	2.50	0.41
1:A:592:SER:OG	1:A:594:GLU:OE2	2.39	0.41
1:A:229:PHE:CE1	1:A:247:PRO:HD2	2.55	0.41
1:A:148:ASP:O	1:A:151:LEU:HB2	2.21	0.41
1:A:126:LYS:H	1:A:126:LYS:HG3	1.71	0.41
1:A:77:GLU:HG2	1:A:81:LYS:HD2	2.03	0.41
1:A:199:LEU:HD22	1:A:254:PHE:HZ	1.85	0.41
1:A:284:ASN:N	1:A:285:PRO:HD3	2.35	0.41
1:A:291:LYS:O	1:A:295:GLU:HB2	2.21	0.41
1:A:348:ARG:C	1:A:350:GLY:N	2.74	0.41
1:A:49:ALA:HB1	1:A:51:TYR:CE2	2.56	0.40
1:A:56:ALA:HB1	1:A:177:ARG:HD3	2.02	0.40
1:A:138:ASN:O	1:A:161:PHE:HA	2.21	0.40
1:A:351:HIS:CD2	1:A:433:LEU:HD21	2.56	0.40
1:A:480:LEU:HD12	1:A:480:LEU:HA	1.90	0.40
1:A:85:TYR:CD1	1:A:411:LYS:HA	2.57	0.40
1:A:101:MET:C	1:A:101:MET:SD	3.00	0.40
1:A:210:LEU:O	1:A:292:LEU:HD23	2.21	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	592/595 (100%)	518 (88%)	56 (10%)	18 (3%)	<b>4</b> <b>10</b>

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	ALA
1	A	13	VAL
1	A	18	ASN
1	A	125	SER
1	A	167	CYS
1	A	168	PRO
1	A	169	THR
1	A	572	TYR
1	A	2	TRP
1	A	9	PRO
1	A	573	PRO
1	A	17	GLU
1	A	424	PRO
1	A	19	SER
1	A	56	ALA
1	A	14	LYS
1	A	367	PRO
1	A	5	GLY

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	517/517 (100%)	463 (90%)	54 (10%)	7 16

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	14	LYS
1	A	16	ASP
1	A	22	ARG
1	A	33	SER
1	A	36	LEU
1	A	57	LEU
1	A	64	ARG
1	A	91	VAL
1	A	98	LEU
1	A	119	LEU
1	A	121	SER
1	A	126	LYS
1	A	128	GLN
1	A	134	ILE
1	A	146	LYS
1	A	158	MET
1	A	172	TYR
1	A	175	LEU
1	A	187	LEU
1	A	201	SER
1	A	203	LEU
1	A	208	SER
1	A	240	ILE
1	A	260	ILE
1	A	261	LEU
1	A	276	LEU
1	A	278	ARG
1	A	279	GLU
1	A	281	LYS
1	A	282	LYS
1	A	283	LEU
1	A	310	ARG
1	A	333	ASN
1	A	347	PHE

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Mol	Chain	Res	Type
1	A	360	ARG
1	A	376	LEU
1	A	387	ILE
1	A	425	THR
1	A	427	LYS
1	A	464	LEU
1	A	469	THR
1	A	475	ILE
1	A	480	LEU
1	A	494	ILE
1	A	503	GLU
1	A	511	LEU
1	A	513	CYS
1	A	537	THR
1	A	538	GLU
1	A	545	GLN
1	A	559	ILE
1	A	564	LEU
1	A	568	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	138	ASN
1	A	154	GLN
1	A	273	HIS
1	A	322	GLN
1	A	423	GLN
1	A	497	ASN
1	A	545	GLN
1	A	574	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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
1	SEP	A	198	1	8,9,10	0.69	0	8,12,14	0.96	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
1	SEP	A	198	1	-	2/5/8/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	198	SEP	N-CA-CB-OG
1	A	198	SEP	CA-CB-OG-P

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	198	SEP	2	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	A	604	1	14,14,15	0.33	0	17,19,21	0.69	0
4	NAG	A	605	1	14,14,15	0.44	0	17,19,21	1.03	0
2	HEM	A	601	1,5	41,50,50	1.55	9 (21%)	45,82,82	2.05	13 (28%)
4	NAG	A	603	1	14,14,15	0.73	0	17,19,21	2.03	5 (29%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	604	1	-	2/6/23/26	0/1/1/1
4	NAG	A	605	1	-	2/6/23/26	0/1/1/1
2	HEM	A	601	1,5	-	4/12/54/54	-
4	NAG	A	603	1	-	1/6/23/26	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	HEM	C1B-NB	-3.98	1.33	1.40
2	A	601	HEM	C4D-ND	-3.33	1.34	1.40
2	A	601	HEM	FE-NB	3.26	2.13	1.96
2	A	601	HEM	CMB-C2B	2.28	1.55	1.50
2	A	601	HEM	C1D-ND	-2.25	1.34	1.38
2	A	601	HEM	CHB-C1B	2.24	1.40	1.35
2	A	601	HEM	C1D-C2D	2.18	1.48	1.44
2	A	601	HEM	CMD-C2D	2.04	1.55	1.50
2	A	601	HEM	FE-ND	-2.03	1.86	1.96

All (18) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	HEM	CBA-CAA-C2A	-7.10	100.50	112.62
2	A	601	HEM	CHA-C4D-ND	4.50	129.94	124.38
4	A	603	NAG	O5-C1-C2	-4.30	104.49	111.29
4	A	603	NAG	C1-C2-N2	3.98	117.28	110.49
2	A	601	HEM	CHD-C1D-ND	3.68	128.43	124.43
2	A	601	HEM	C1B-NB-C4B	3.47	108.66	105.07
2	A	601	HEM	CHA-C4D-C3D	-3.11	119.49	125.33
4	A	603	NAG	O3-C3-C2	-2.93	103.40	109.47
4	A	603	NAG	O5-C5-C6	-2.87	102.71	107.20
2	A	601	HEM	CHC-C4B-NB	2.70	127.36	124.43
2	A	601	HEM	CHB-C1B-NB	2.69	127.70	124.38
2	A	601	HEM	CBD-CAD-C3D	-2.67	105.20	112.63
2	A	601	HEM	C4D-ND-C1D	2.58	107.74	105.07
2	A	601	HEM	CHD-C1D-C2D	-2.56	120.98	124.98
2	A	601	HEM	C1D-C2D-C3D	-2.26	104.58	106.96
2	A	601	HEM	CAD-CBD-CGD	-2.21	108.85	113.60
4	A	603	NAG	C1-O5-C5	-2.12	109.31	112.19
2	A	601	HEM	CHB-C1B-C2B	-2.09	120.94	126.72

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	604	NAG	C4-C5-C6-O6
4	A	603	NAG	C8-C7-N2-C2
4	A	605	NAG	O5-C5-C6-O6
4	A	604	NAG	O5-C5-C6-O6
2	A	601	HEM	CAD-CBD-CGD-O2D
2	A	601	HEM	CAD-CBD-CGD-O1D
2	A	601	HEM	CAA-CBA-CGA-O2A
2	A	601	HEM	CAA-CBA-CGA-O1A
4	A	605	NAG	C4-C5-C6-O6

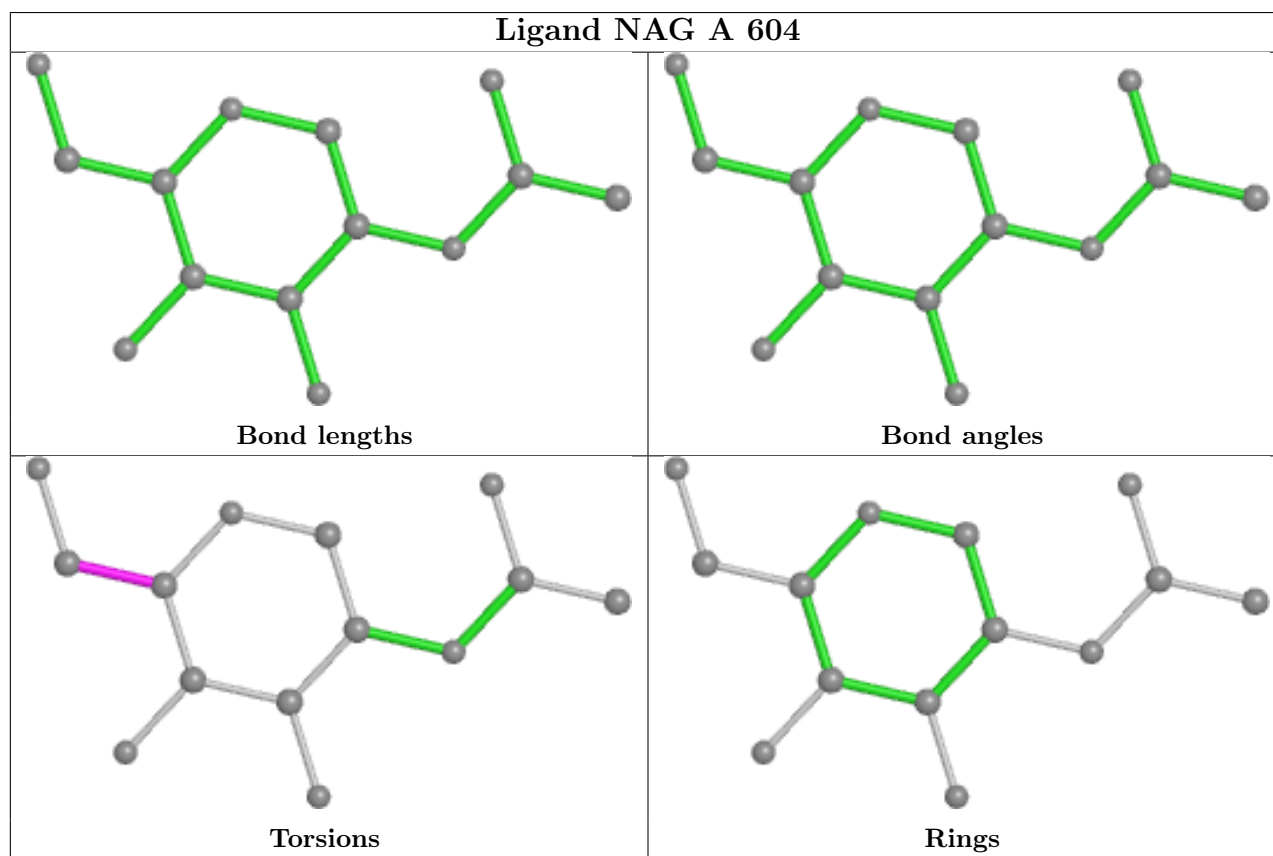
There are no ring outliers.

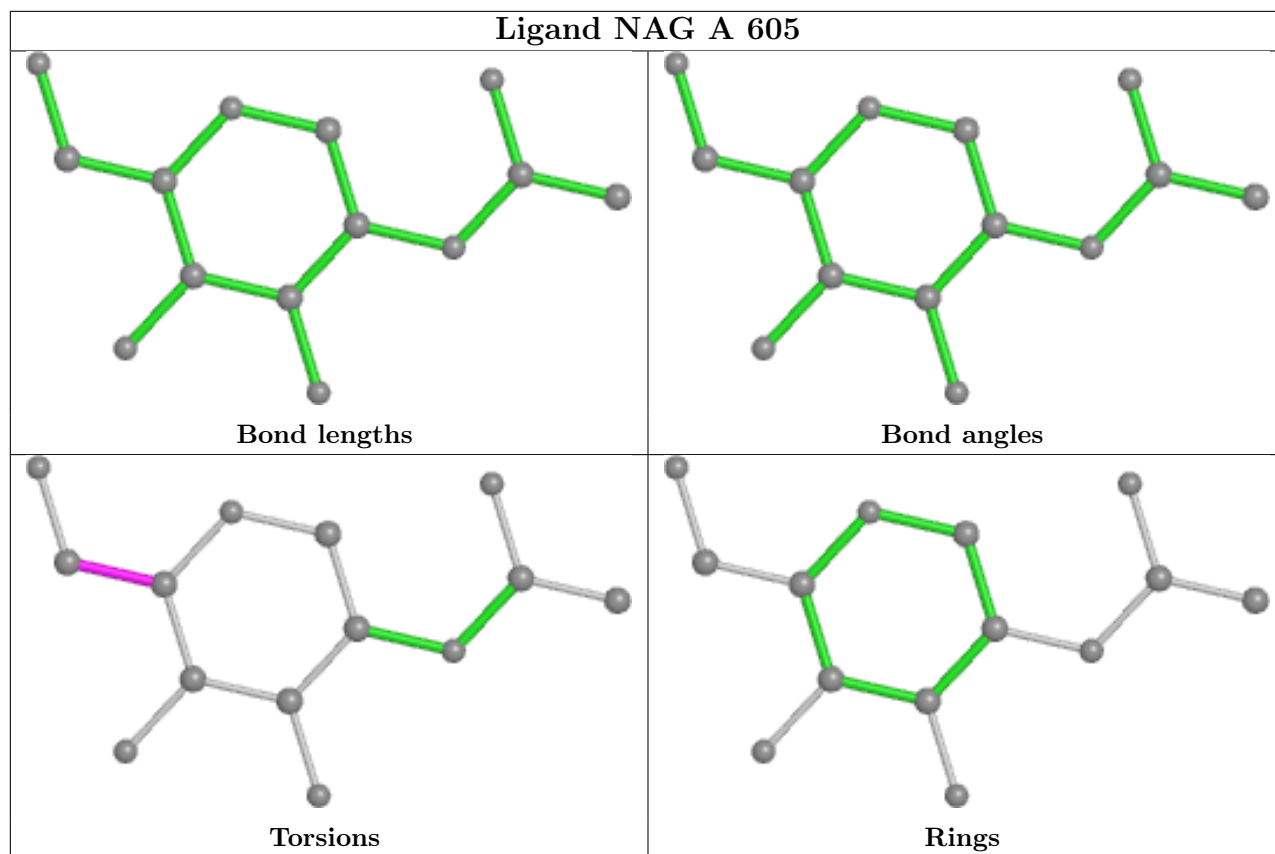
3 monomers are involved in 10 short contacts:

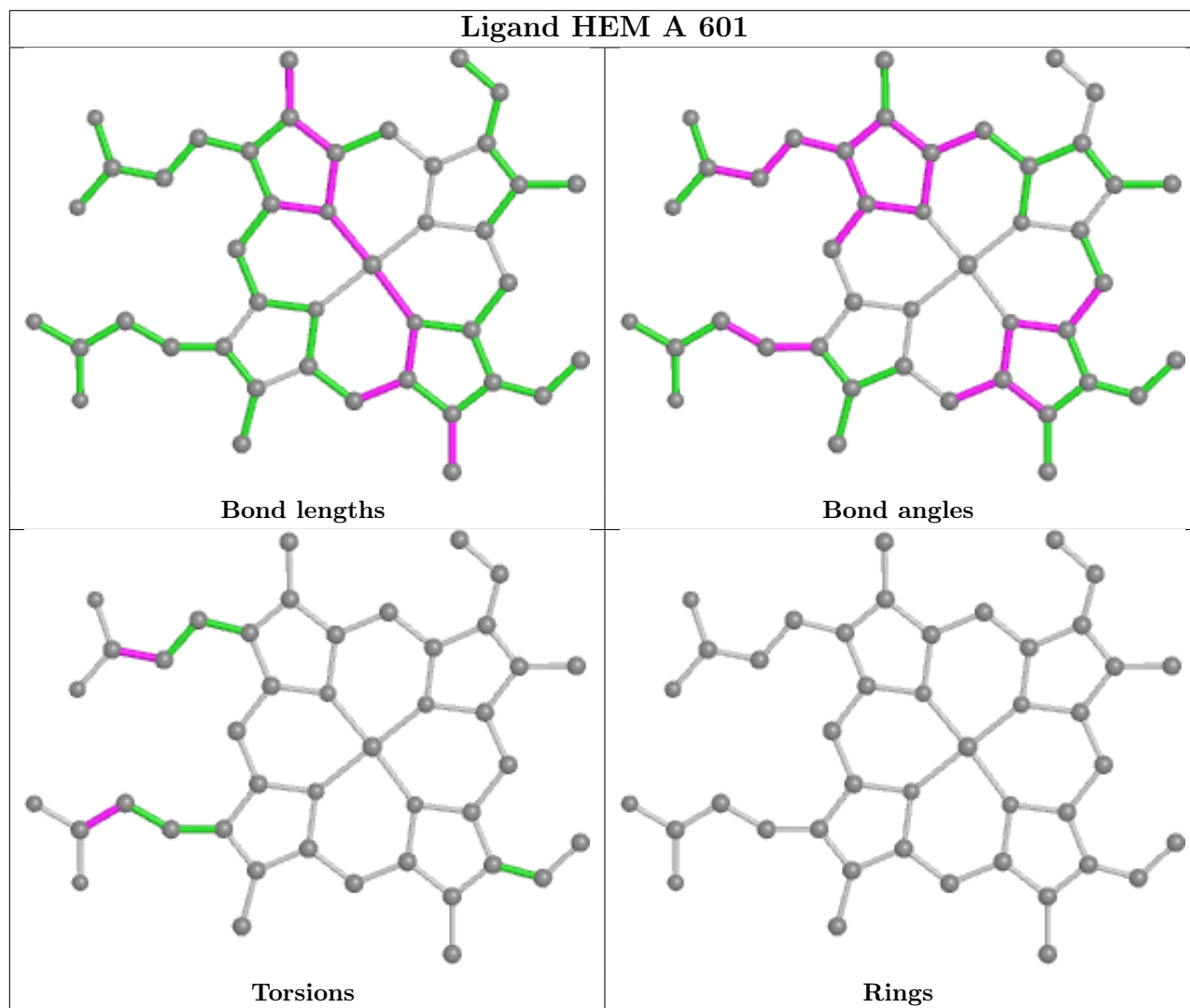
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	604	NAG	2	0
4	A	605	NAG	2	0
2	A	601	HEM	6	0

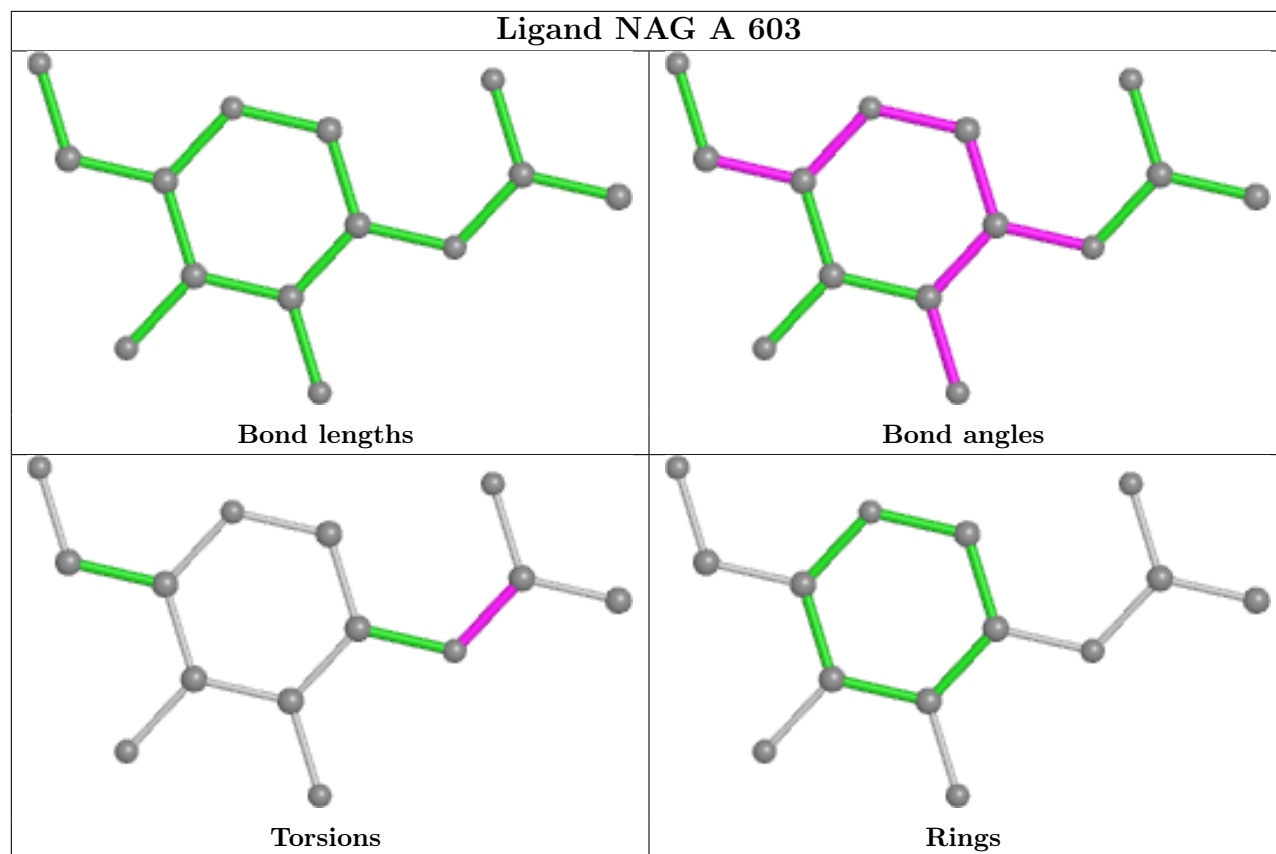
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

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.









## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	594/595 (99%)	0.06	32 (5%) <span style="border: 1px solid red; padding: 2px;">25</span> <span style="border: 1px solid red; padding: 2px;">24</span>	14, 33, 90, 138	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	SER	9.8
1	A	173	GLN	6.4
1	A	122	ASN	6.2
1	A	121	SER	5.5
1	A	4	VAL	5.4
1	A	595	ASN	5.2
1	A	8	ALA	5.1
1	A	2	TRP	4.8
1	A	170	PRO	4.8
1	A	172	TYR	4.8
1	A	123	GLU	4.7
1	A	124	HIS	4.6
1	A	13	VAL	4.6
1	A	174	SER	4.4
1	A	117	THR	3.7
1	A	171	PRO	3.6
1	A	134	ILE	3.3
1	A	593	ARG	3.1
1	A	119	LEU	3.1
1	A	120	GLY	2.8
1	A	7	GLY	2.7
1	A	127	THR	2.7
1	A	594	GLU	2.7
1	A	64	ARG	2.5
1	A	5	GLY	2.5
1	A	541	ARG	2.4
1	A	128	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	587	LEU	2.2
1	A	125	SER	2.2
1	A	580	SER	2.1
1	A	169	THR	2.1
1	A	334	SER	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	SEP	A	198	10/11	0.88	0.20	40,52,81,82	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

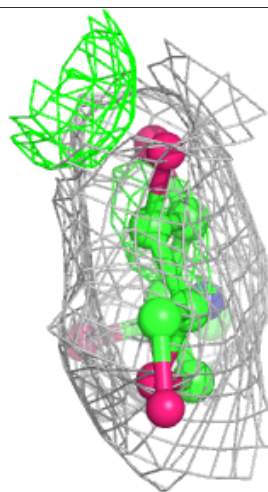
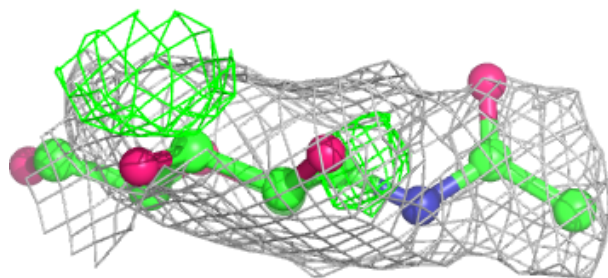
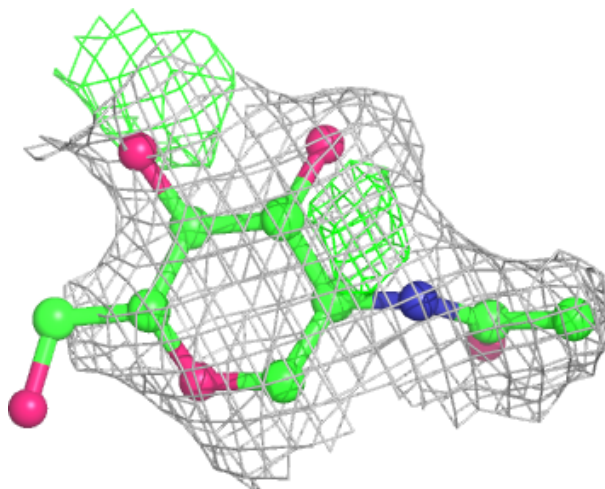
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	NAG	A	605	14/15	0.74	0.27	55,61,74,80	0
4	NAG	A	604	14/15	0.78	0.32	67,72,75,79	0
4	NAG	A	603	14/15	0.85	0.20	48,50,53,56	0
3	CA	A	602	1/1	0.96	0.07	20,20,20,20	0
2	HEM	A	601	43/43	0.96	0.12	20,25,26,27	0
3	CA	A	606	1/1	0.99	0.12	23,23,23,23	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 NAG A 605:**

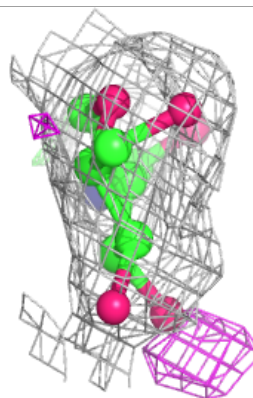
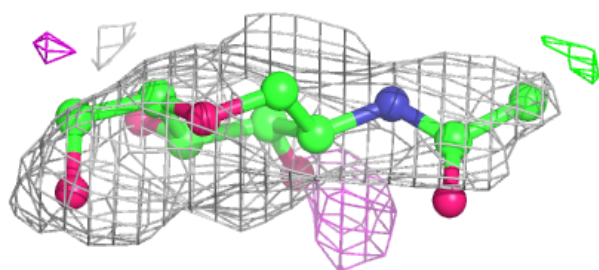
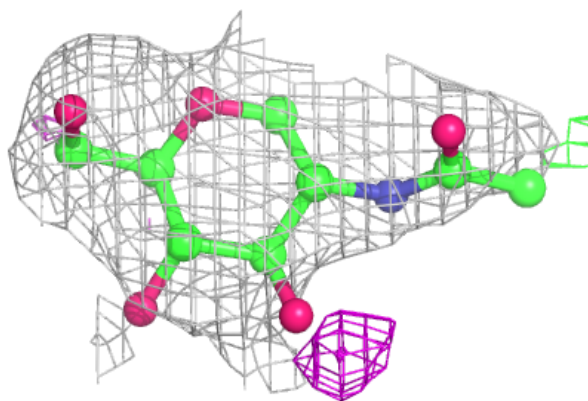
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



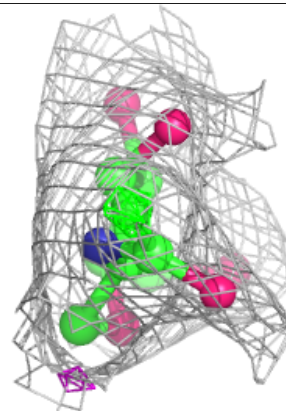
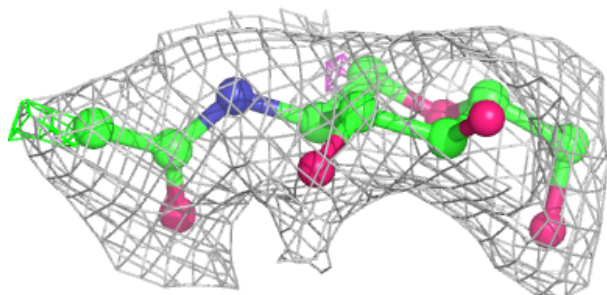
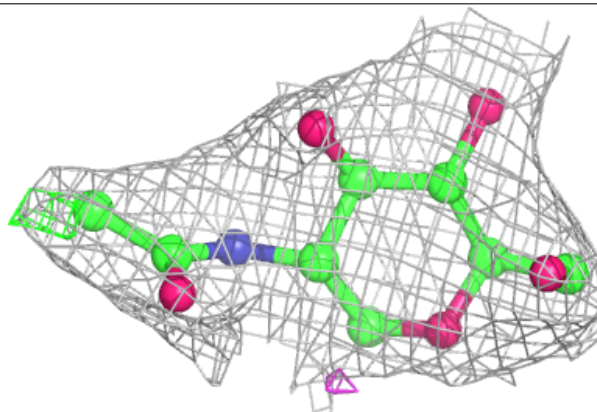


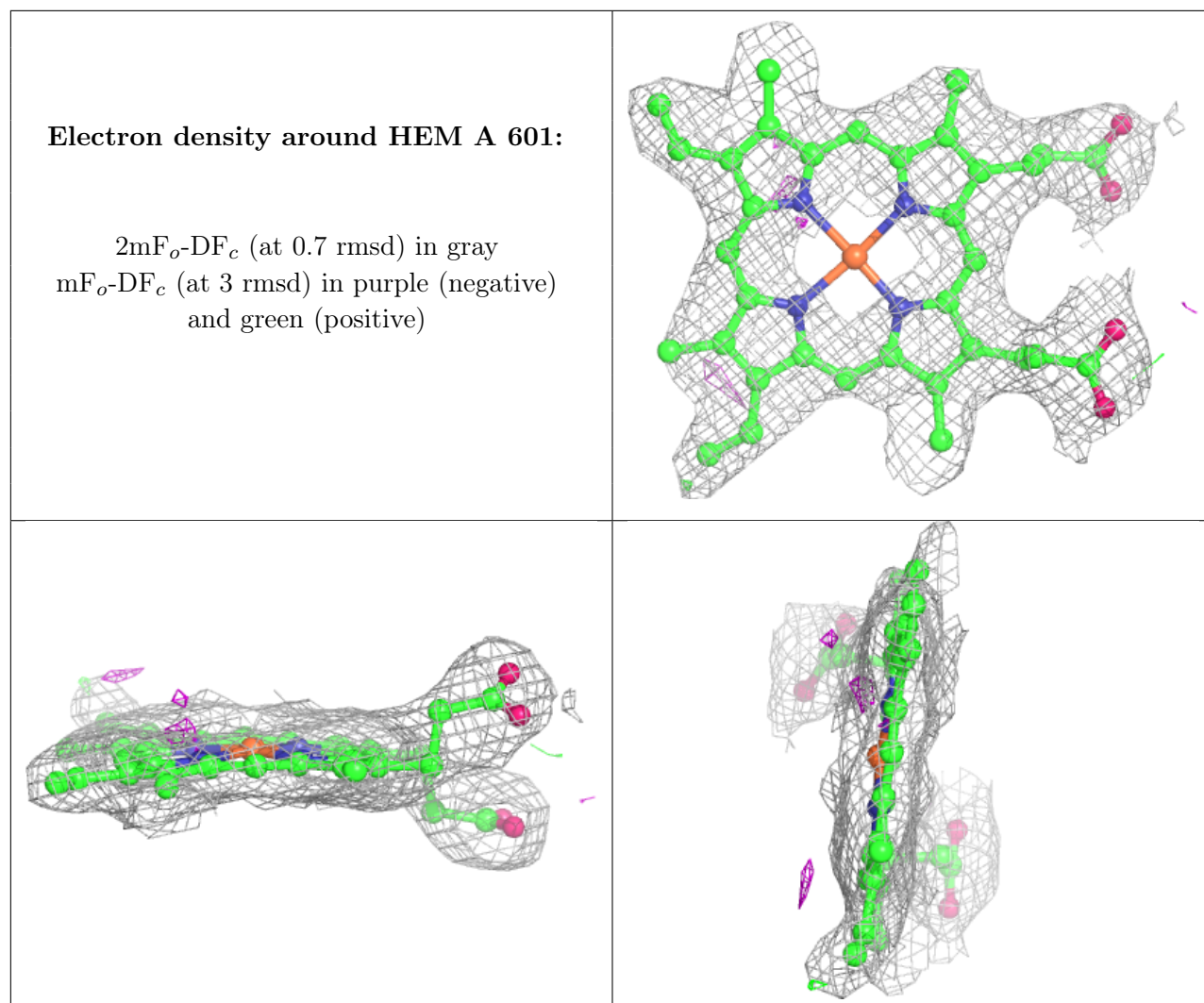
**Electron density around NAG A 604:**

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

**Electron density around NAG A 603:**

$2mF_o-DF_c$  (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.