## Metropolitan Museum of Art Gas Chromatography- Mass Spectrometry (GC-MS) Results from Material Analysis

This document includes (1) a mass spectrum and (2) the volatile organic compounds (VOCs) emitted from samples using GC-MS analysis. The data is not interpreted; however, several classes of chemicals are highlighted because they are potential risks for artwork in an enclosed environment. A basic key, provided below, indicates those classes. The amount of each chemical identified has not been determined; similarly, it is not known how much of each chemical is necessary to do damage to art. Finally, peaks may be present that are the result of the sample adsorbing chemicals from the air and reemitting them during testing rather than being inherent to the sample. Research is ongoing to determine specifically which chemicals and amounts are required to negatively affect artifacts.

## Highlighted data:

Pink – chemicals currently known to be hazardous to art

Green – amines; can raise the pH, are suspected to react with acids and may form crystals in an enclosed environment

Yellow – chemicals of the following type, which may be hazardous to art:

Acids – lower the pH, corrosive to metals, degrade organic materials

Aldehydes – can convert to acids with heat or exposure to UV light

Esters – can hydrolyze into acids with heat and humidity

Sulfur-containing compounds – known to tarnish and corrode some metals

Halogenated compounds – can become reactive with exposure to heat and UV light

Nitrogen-containing, not amine – can react with other off-gassed chemicals

Alkynes – can become reactive when exposed to heat or UV light

Sample: Quality Media and Laminating Solutions (QMLS): Instabond C double-sided Adhesive

Date collected: 10/24/2018

8.990

9.330

9.510

9.600

9.820

10.230

11.120

11.170

11.470

11.570

97.5

95.9

91.4

94.1

93.3

95.9

93.2

93.5

95.8

93.7

91.2

C25H23NO10

C10H12O2

C10H12O2

C11H20O2

C10H12O2

C11H14O2

C12H24O2

C9H14O6

C12H24O3

C12H36O6Si6

C10H20O

497.1

164.1

164.1

156.2

184.1

164.1

178.1

200.2

444.1

218.1

216.2

9972588

3528688

1796701

1655202

4359926

6278494

3789488

3194516

3706628

16891217

71127-22-5

TETRACARBOX..

103-11-7 2-Ethylhexyl acrylate

102-76-1 Triacetin

939-48-0 Benzoic acid, 1-methylethyl ester

103-45-7 Acetic acid, 2-phenylethyl ester

25415-84-3 n-Butyric acid 2-ethylhexyl ester

540-97-6 Cyclohexasiloxane, dodecamethyl-

21862-63-5 Cyclohexanol, 4-(1,1-dimethylethyl)-, trans-

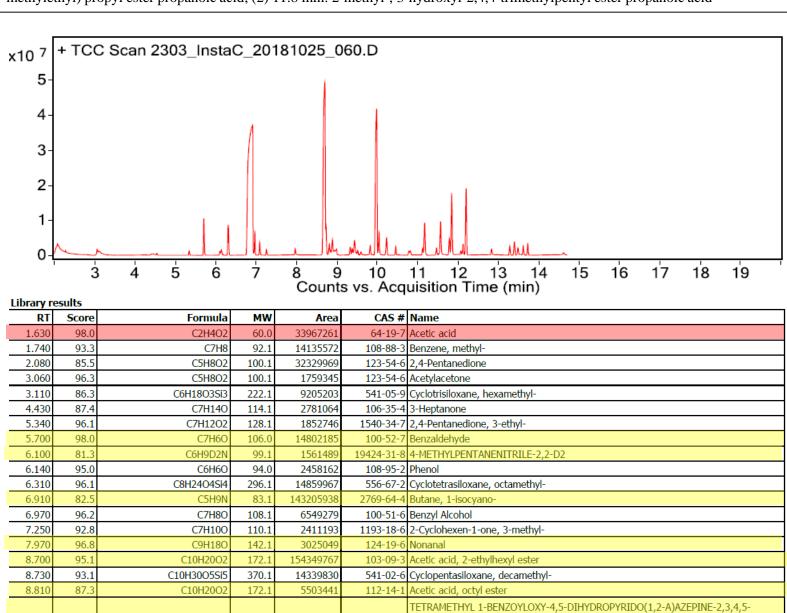
Benzenemethanol, .alpha.-methyl-, acetate

(.+/-.)-1-Phenyl-propan-2-ol, N-(acetyl)

17619057 74367-33-2 Propanoic acid, 2-methyl-, 2,2-dimethyl-1-(2-hydroxy-1-methylethyl)propyl ester

Technique used: SPME with a PDMS/DVB fiber; Agilent 7890B GC and 5977B MS fitted with a GL Sciences OPTIC-4 multimode inlet and LEAP PAL RTC auto-sampler; Pre-heated at 60°C for 20 minutes; fiber exposure at 60°C for 20 minutes; sample injected into 220°C inlet and cryotrapped for 2 min at -15°C; GC ramped from 40°C to 225 °C at 10°C/min. Data analyzed using the Masshunter Qualitative program. Samples > 80% match with a NIST 17.0 library are reported.

VOCs not highlighted are because they were also observed in blanks: (1) 11.6 min: 2-methyl-, 2,2-dimethyl-1-(2-hydroxyl-1-methylethyl) propyl ester propanoic acid; (2) 11.8 min: 2-methyl-, 3-hydroxyl-2,4,4-trimethylpentyl ester propanoic acid



11.790	96.5	C11H14O2	178.1	7971597	122-72-5	3-Phenyl-1-propanol, acetate
11.850	93.4	C12H24O3	216.2	27526913	74367-34-3	Propanoic acid, 2-methyl-, 3-hydroxy-2,4,4-trimethylpentyl ester
12.130	95.0	C14H30	198.2	2590830	629-59-4	Tetradecane
12.200	95.4	C13H12	168.1	32099344	643-58-3	1,1'-Biphenyl, 2-methyl-
12.830	84.6	C11H12O4	208.1	3207306	999200-85-3	ETHYLIDENE MONOACETATE MONOBENZOATE
13.280	94.7	C13H12	168.1	4238519	644-08-6	1,1'-Biphenyl, 4-methyl-
13.490	92.9	C16H34O	242.3	3331106	10143-60-9	bis(2-Ethylhexyl) ether
13.620	96.5	C15H24O	220.2	4236497	128-37-0	Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl-
13.730	94.1	C14H14	182.1	5513438	103-29-7	Benzene, 1,1'-(1,2-ethanediyl)bis-
14.610	90.6	C16H30O4	286.2	1755296	6846-50-0	PENTAN-1,3-DIOLDIISOBUTYRATE, 2,2,4-TRIMETHYL-