## 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

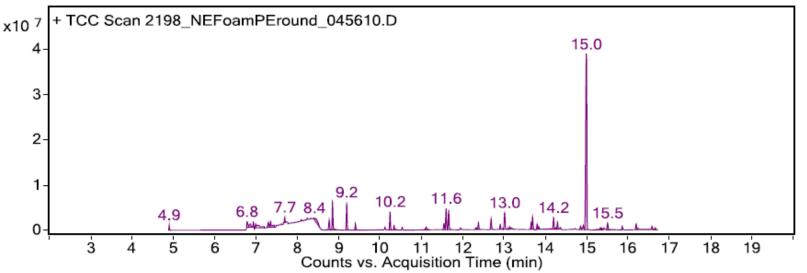
Sample: Uline Foam core, black, S-12861

Oddy test result: Temporary

Date collected: 06/26/2018

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 autosampler; Pre-heated at 60°C for 20 minutes; fiber exposure at 60°C for 20 minutes; sample injected into 220°C inlet and crotrapped for 2 min at -15°C; GC ramped from 40°C to 225 °C at 10°C/min. Data analyzed in masshunter Qualitative. Samples > 80% match with a NIST library are reported.

VOCs not highlighted are because they were also observed in blanks: : (1) 5.7 min: methoxyphenyl oxime; (2) 12.4 min: 2-methyl-, 2,2-dimethyl-1-(2-hydroxyl-1-methylethyl) propyl ester propanoic acid; (3) 12.7 min: 2-methyl-, 3-hydroxyl-2,4,4-trimethylpentyl ester propanoic acid



Library results

LIDITARY I	Library results										
RT	Score	Formula	MW	Area	CAS #	Name					
3.900	93.7	C2H8O2Si	92.0	1199947	1066-42-8	Silanediol, dimethyl-					
4.800	96.5	C6H12O	100.1	667569	66-25-1	Hexanal					
4.900	92.5	C6H18O3Si3	222.1	2853956	541-05-9	Cyclotrisiloxane, hexamethyl-					
5.500	96.3	C8H10	106.1	23705012	0-00-0	METHYLLAURATE					
5.600	97.9	C8H10	106.1	1250474	0-00-0	unidentified C2-benzene					
5.700	85.0	C8H9NO2	151.1	925682	1000222-86-6	Oxime-, methoxy-phenyl					
5.900	87.8	C6H10O	98.1	1051277	108-94-1	Cyclohexanone					
5.900	82.6	C7H14O	114.1	673709	111-71-7	Heptanal					
6.000	90.9	C6H14O2	118.1	1274294	111-76-2	Ethanol, 2-butoxy-					
6.300	96.8	C9H12	120.1	7765205	98-82-8	Benzene, (1-methylethyl)-					
6.600	95.5	C9H10	118.1	1137074	873-49-4	Benzene, cyclopropyl-					
6.700	95.4	C9H12	120.1	4621192	0-00-0	unidentified C3-benzene					
6.800	97.3	C7H6O	106.0	21597724	100-52-7	Benzaldehyde					
6.900	96.2	C9H12	120.1	764827	0-00-0	unidentified C3-benzene					

6.900	85.7	C6H6O	94.0			
6.900	94.3		296.1			Cyclotetrasiloxane, octamethyl-
7.000	80.2		120.1	1320899		Benzene, 1-ethyl-2-methyl-
7.100	96.4		118.1			Benzene, (1-methylethenyl)-
7.200	90.2		120.1	1477661		unidentified C3-benzene
7.300	94.6		142.2		124-18-5	
7.300	97.6		128.1	2530209	124-13-0	
7.500	94.8	C10H14	134.1	2591939		Benzene, (1-methylpropyl)-
7.700	97.3	C8H18O	130.1	9939992	104-76-7	7 1-Hexanol, 2-ethyl-
7.800	96.8	C10H16	136.1	2111440	138-86-3	B dl-Limonene
7.800	86.1	C7H8O	108.1	1449842		Benzyl alcohol
8.000	91.7	C8H8O	120.1	1671188		PHENYL ACETALDEHYDE
8.000	95.3	C10H14	134.1		135-01-3	o-Diethyl benzene
8.200	80.2		134.1	<del>                                     </del>		o-Diethyl benzene
8.300	94.6		120.1			Ethanone, 1-phenyl-
8.500	81.3		168.2			3-Decene, 2,2-dimethyl-, (E)-
8.700	83.5		168.2			3-UNDECENE, 4-METHYL-
8.800	88.1	C11H24	156.2			Undecane
8.800	97.3	C9H18O	142.1	30157931		5 Nonanal
8.900	84.1		168.2			3-Undecene, 9-methyl-, (E)-
9.100	87.1		138.1	<del>                                     </del>		2-Cyclohexen-1-one, 3,5,5-trimethyl-
9.200	95.4		370.1			Cyclopentasiloxane, decamethyl-
9.400	85.7	C10H20O2	172.1	890624		Acetic acid, 2-ethylhexyl ester
9.600	91.8		132.1	1291568		3 1,3,5,7-Cyclooctatetraene-1-carboxaldehyde
9.800	86.7	C20H42O	298.3			P Decane, 1,1'-oxybis-
10.000	97.6		156.2	1581712		Cyclohexanol, 5-methyl-2-(1-methylethyl)-
10.000	94.9		162.1	1043927		Ethanol, 2-(2-butoxyethoxy)-
10.100	87.0		224.3			2 1-Hexadecene
10.100	95.2		128.1			
10.200	92.9		170.2			B Dodecane
10.300	97.8		156.2		112-40-3	
10.600	85.9		184.1	650125		7 2-Propenoic acid, 2-ethylhexyl ester
11.200	93.1	C11H2002	270.3			B Decyl octyl ether
11.500	80.5		238.2			Cyclopentadecanone, 4-methyl-
11.600	95.9		444.1			Cyclohexasiloxane, dodecamethyl-
11.600	93.9		184.2			Tridecane
11.800	95.8		170.2			7 Undecanal
11.000	93.0	CITIZZO	1/0.2	050550	112 717	Ondecanal
12.400	89.8	C12H24O3	216.2	14001455	7/1267-33-2	Propanoic acid, 2-methyl-, 2,2-dimethyl-1-(2-hydroxy-1-methylethyl)propyl ester
12.700	93.9		216.2			Propanoic acid, 2-methyl-, 2,2-dimethyl-1-(2-hydroxy-1-methylethyl)propyi ester  Propanoic acid, 2-methyl-, 3-hydroxy-2,4,4-trimethylpentyl ester
13.000	93.9		198.2			Tetradecane
13.100	94.9		198.2	1046889		Dodecanal
13.100	80.2		518.1			Cycloheptasiloxane, tetradecamethyl-
$\overline{}$						
14.000	96.7 95.2	C12H260	186.2			1 1-Biphopyl 4-methyl-
14.200			168.1			1,1'-Biphenyl, 4-methyl-
14.300	94.3		212.3			pentadecane
14.400	96.6		198.2		10486-19-8	
15.500	91.5		226.3	+ + +		Hexadecane
15.800	90.7		592.2			Cyclooctasiloxane, hexadecamethyl-
15.800	86.5		242.2			Dodecanoic acid, 1-methylethyl ester
16.700	83.4		214.1	<del>                                     </del>		(2',2'-DIMETHYL-3'-PHENYLAZETIDIN-1'-YL)OXOACETONITRILE
17.300	86.9	C16H16	208.1	4708860	20071-09-4	Benzene, 1,1'-(1,2-cyclobutanediyl)bis-, trans-