

## 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

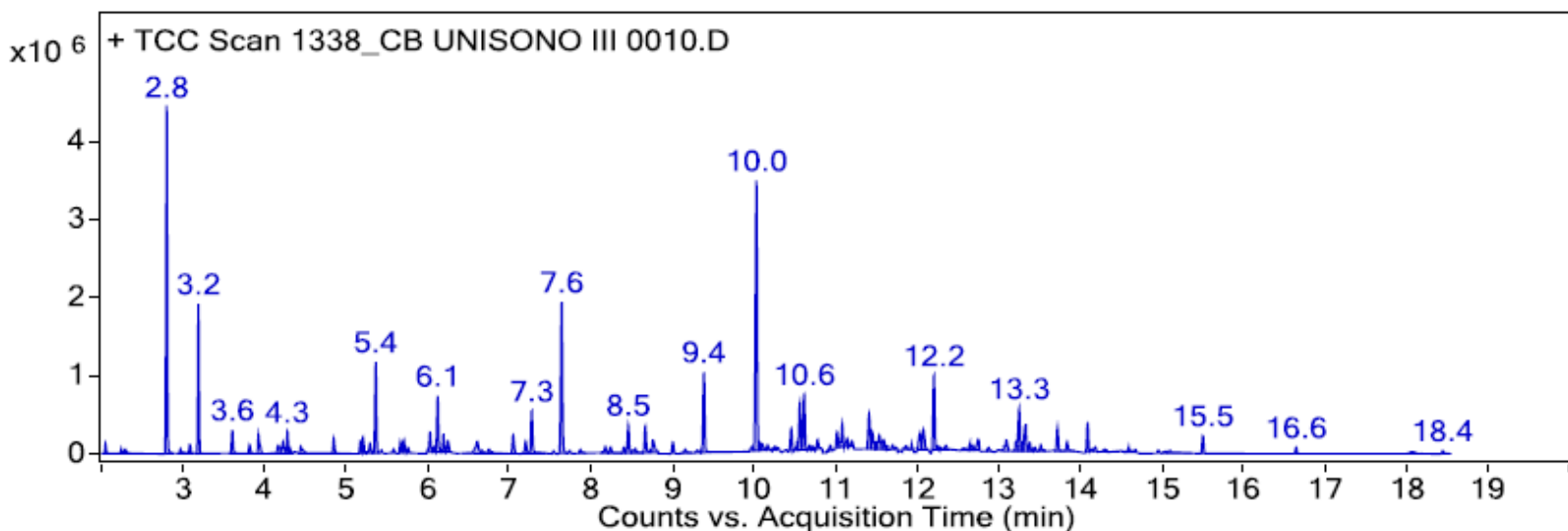
Sample: Creation Baumann Unisono III 0010 (green) cotton fabric

Oddy test result: Permanent

Date GC-MS collected: 6/13/2016

Technique used: SPME Arrow 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 sample at 60°C for 20 minutes; fiber exposure to sample at 60°C for 20 minutes; fiber 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 in Masshunter Qualitative. Samples > 80% match with a NIST library are reported.

VOCs not highlighted are because they were also observed in blanks: (1) 4.2 min: methoxy-phenyl-oxime (2) ~10.8 min: 2-methyl-, 2,2-dimethyl-1-(2-hydroxyl-1-methylethyl) propyl ester propanoic acid



#### Library results

RT	Score	Formula	MW	Area	CAS #	Name
1.400	93.4	CH4O	32.0	304142	67-56-1	Methanol
1.500	98.0	C2H6O	46.0	54750	64-17-5	Ethanol
1.500	94.1	C3H6O	58.0	284900	67-64-1	2-Propanone
1.900	93.9	C4H8O	72.1	86626	109-99-9	Furan, tetrahydro-
1.900	97.8	C2H4O2	60.0	166078	999000-92-5	Acetic acid
2.200	92.2	C7H16	100.1	51071	142-82-5	Heptane
2.800	97.4	C7H8	92.1	4605802	108-88-3	Benzene, methyl-
3.000	91.9	C8H16	112.1	63021	1632-16-2	Heptane, 3-methylene-
3.100	90.6	C8H18	114.1	152472	111-65-9	Octane
3.200	91.5	C6H18O3Si3	222.1	1870875	541-05-9	Cyclotrisiloxane, hexamethyl-
3.600	97.0	C6H12O2	116.1	370032	123-42-2	2-Pentanone, 4-hydroxy-4-methyl-
3.900	97.7	C8H10	106.1	365075	1330-20-7	XYLENE
4.200	83.9	C8H9NO2	151.1	374302	1000222-86-6	Oxime-, methoxy-phenyl-
4.200	91.1	C8H10	106.1	132625	1330-20-7	XYLENE
4.300	97.4	C9H20	128.2	160668	111-84-2	Nonane
5.200	82.2	C10H10O3	178.1	130276	67519-26-0	1,4-[13C]-4-Oxo-4-phenylbutanoic acid
5.400	94.4	C8H24O4Si4	296.1	1296992	556-67-2	Cyclotetrasiloxane, octamethyl-
5.400	92.2	C9H12	120.1	63467	0-00-0	unidentified C3-benzene
5.700	94.7	C9H12	120.1	193560	526-73-8	Benzene, 1,2,3-trimethyl-
5.700	96.5	C10H22	142.2	168043	124-18-5	Decane
5.800	91.6	C8H16O	128.1	87819	124-13-0	Octanal
6.000	88.0	C7H9N	107.1	327357	100-45-8	4-Cyanocyclohexene
6.100	96.3	C9H12	120.1	75544	0-00-0	unidentified C3-benzene
6.100	95.4	C8H18O	130.1	914767	104-76-7	1-Hexanol, 2-ethyl-

6.200	98.0	C10H16	136.1	208622	138-86-3	dl-Limonene
6.200	80.1	C7H8O	108.1	152976	100-51-6	Benzyl alcohol
6.600	86.7	C6H18O3Si3	222.1	238483	541-05-9	Cyclotrisiloxane, hexamethyl-
7.200	86.7	C11H24	156.2	182522	1120-21-4	Undecane
7.300	98.2	C9H18O	142.1	674727	124-19-6	Nonanal
7.700	91.4	C16H11NO2S	281.1	70830	70453-75-7	2-methoxy[1]benzothieno[2,3-c]quinolin-6(5H)-one
7.700	92.1	C8H24O4Si4	296.1	52209	556-67-2	Cyclotetrasiloxane, octamethyl-
7.900	83.4	C10H20O2	172.1	68624	103-09-3	Acetic acid, 2-ethylhexyl ester
8.200	83.3	C12H22O	182.2	119276	999136-31-6	3-DODECEN-1-AL
8.400	93.4	C10H20O	156.2	97772	1490-04-6	Cyclohexanol, 5-methyl-2-(1-methylethyl)-
8.500	95.8	C8H18O3	162.1	330561	112-34-5	Ethanol, 2-(2-butoxyethoxy)-
8.700	94.6	C12H26	170.2	440150	112-40-3	Dodecane
8.800	95.5	C10H20O	156.2	170219	112-31-2	Decanal
8.800	94.9	C8H24O4Si4	296.1	113928	556-67-2	Cyclotetrasiloxane, octamethyl-
9.200	80.7	C11H22O2	186.2	68614	999145-46-3	2-Ethyl-1-hexyl propionate
9.400	96.9	C12H24O2	200.2	1381946	7434-89-1	Hexanoic acid, 2-ethyl-, 2-methylpropyl ester
10.300	87.6	C16H32	224.3	90246	629-73-2	1-Hexadecene
10.500	85.0	C12H16	160.1	389340	13065-07-1	Naphthalene, 1,2,3,4-tetrahydro-2,7-dimethyl-
10.600	83.6	C12H16	160.1	816345	13065-07-1	Naphthalene, 1,2,3,4-tetrahydro-2,7-dimethyl-
10.600	86.8	C12H16	160.1	1024192	1076-69-3	5,6,7,8,9,10-Hexahydrobenzocyclooctene
10.800	84.8	C12H24O3	216.2	163851	74367-33-2	Propanoic acid, 2-methyl-, 2,2-dimethyl-1-(2-hydroxy-1-methylethyl)propyl ester
10.900	83.2	C21H44	296.3	107041	54833-23-7	Eicosane, 10-methyl-
11.000	90.8	C14H30	198.2	319541	6418-41-3	Tridecane, 3-methyl-
11.100	80.5	C17H17ClF3N	327.1	53929	999524-60-3	(R)-2-(N,N-Dibenzylamino)-3-chloro-1,1,1-trifluoropropane
11.400	96.0	C14H30	198.2	681632	629-59-4	Tetradecane
11.500	93.9	C14H22	190.2	263844	0-00-0	C8 - alkyl - benzene (structure not known)
11.700	81.9	C15H24	204.2	119353	475-20-7	Junipene
11.900	84.0	C13H18	174.1	108913	62379-92-4	1,4-Dimethyl-2-cyclopentylbenzene
12.000	86.0	C17H36	240.3	150551	6008-17-9	5,5-Dibutylnonane
12.100	92.4	C14H28	196.2	340943	2882-98-6	Cyclopentane, nonyl-
12.300	80.1	C14H20O2	220.1	89240	719-22-2	2,5-Cyclohexadiene-1,4-dione, 2,6-bis(1,1-dimethylethyl)-
12.400	82.3	C14H28	196.2	119451	1120-36-1	1-Tetradecene
13.300	88.7	C14H20	188.2	604823	30628-31-0	1,3,5-Trimethyl-2-cyclopentylbenzene
13.300	88.1	C14H20	188.2	450108	30628-31-0	1,3,5-Trimethyl-2-cyclopentylbenzene
13.500	86.7	C16H34	226.3	109199	2882-96-4	Pentadecane, 3-methyl-
13.800	90.2	C17H36	240.3	159494	1000360-41-3	5,5-Diethyltridecane
18.400	93.0	C15H16O2	228.1	82012	80-05-7	Phenol, 4,4'-(1-methylethylidene)bis-