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

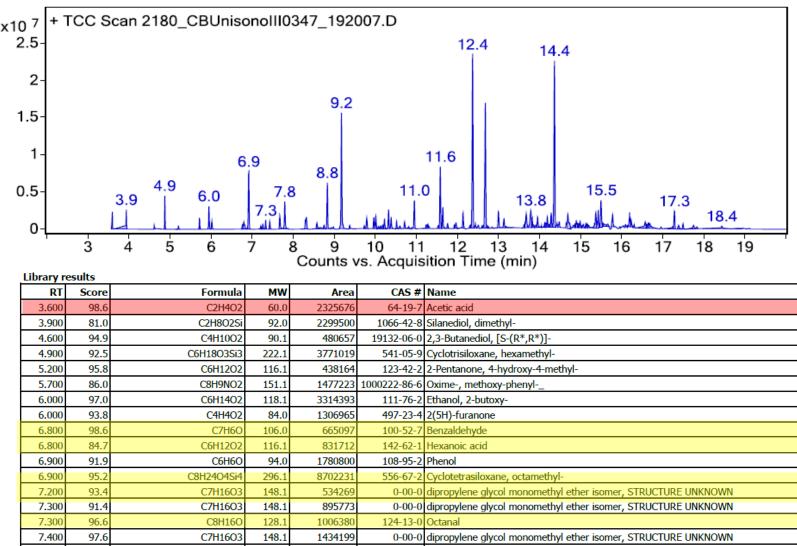
Sample: Creation Baumann Unisono III 0347 red fabric

Oddy test result: Temporary

Date collected: 05/31/2018

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



	7.300	96.6	C8H16O	128.1	1006380	124-13-0	Octanal			
	7.400	97.6	C7H16O3	148.1	1434199	0-00-0	dipropylene glycol monomethyl ether isomer, STRUCTURE UNKNOWN			
	7.700	96.9	C8H18O	130.1	2481144	104-76-7	1-Hexanol, 2-ethyl-			
	7.800	96.1	C7H8O	108.1	4496975	100-51-6	Benzyl Alcohol			
	7.800	89.8	C6H14O2	118.1	433033	6920-22-5	1,2-Hexanediol			
	8.300	97.1	C7H8O	108.1	2339927	106-44-5	Phenol, 4-methyl-			
	8.700	83.9	C9H20O	144.2	385926	999057-76-9	2,6-dimethyl-1-heptanol			
	8.800	97.2	C9H18O	142.1	7686368	124-19-6	Nonanal			
	8.900	84.5	C6H6O3	126.0	393495	37112-31-5	Levoglucosenone			
	9.000	94.9	C8H10O	122.1	359158	60-12-8	Benzeneethanol			
	9.100	85.0	C9H20O	144.2	694445	143-08-8	1-Nonanol			

						
9.200	95.4	C10H30O5Si5	370.1	20007620	541-02-6	Cyclopentasiloxane, decamethyl-
9.200	92.9	C9H20O	144.2	920489	143-08-8	1-Nonanol
9.400	96.0	C9H20O	144.2	688440	110453-78-6	(S)-(+)-6-Methyl-1-octanol
9.800	92.7	C9H20O	144.2	994078		1-Nonanol
10.000	98.1	C10H20O	156.2	2059857	15356-70-4	Cyclohexanol, 5-methyl-2-(1-methylethyl)-, (1.alpha.,2.beta.,5.alpha.)-(.+/)-
10.000	95.1	C8H18O3	162.1	1193482		Ethanol, 2-(2-butoxyethoxy)-
10.100	94.0	C10H8	128.1	490398	275-51-4	
10.200	96.9	C8H8O3	152.0	455813		Methyl salicylate
10.200	85.5	C12H26	170.2	1705860		Dodecane
10.300	91.8	C10H20O	156.2	3313790	112-31-2	
10.500	93.2	C8H10O2	138.1	1405365		Ethanol, 2-phenoxy-
10.700	80.2	C7H5NS	135.0	1351219		Benzothiazole
10.800	82.2	C10H22O3	190.2	406895		2-Propanol, 1-(2-butoxy-1-methylethoxy)-
10.900	93.9	C9H12O2	152.1	384954		1-Phenoxypropan-2-ol
11.000	97.0	C12H24O2	200.2	4932289		Hexanoic acid, 2-ethyl-, 2-methylpropyl ester
11.200	96.4	C10H22O	158.2	752649		1-Decanol
11.300	87.5	C12H25I	296.1	378243		Dodecane, 1-iodo-
11.600	95.7	C12H36O6Si6	444.1	11070502		Cyclohexasiloxane, dodecamethyl-
11.600	95.3	C13H28	184.2	4114414		Tridecane
11.800	96.8	C11H22O	170.2	888909	112-44-7	Undecanal
12.000	90.0	C16H34	226.3	502117		Nonane, 2,2,4,4,6,8,8-heptamethyl-
12.100	93.7	C9H14O6	218.1	1734626	102-76-1	
12.400	89.7	C12H24O3	216.2	19054401	74367-33-2	Propanoic acid, 2-methyl-, 2,2-dimethyl-1-(2-hydroxy-1-methylethyl)propyl ester
12.400	83.8	C10H20O4	204.1	533265		Ethanol, 2-(2-butoxyethoxy)-, acetate
12.700	93.8	C12H24O3	216.2	25017081		Propanoic acid, 2-methyl-, 3-hydroxy-2,4,4-trimethylpentyl ester
12.700	86.7	C14H42O5Si6	458.2	383748		Hexasiloxane, tetradecamethyl-
13.000	95.0	C14H30	198.2	3404034		Tetradecane
13.100	96.6	C12H240	184.2	1748337		Dodecanal
13.700	90.8	C14H28	196.2	1630790		Cyclopentane, nonyl-
13.800	88.6	C21H44	296.3	727834		Eicosane, 10-methyl-
13.900	86.4	C14H20O2	220.1	335769		2,5-Cyclohexadiene-1,4-dione, 2,6-bis(1,1-dimethylethyl)-
14.000	96.5	C12H26O	186.2	2152214		1-Dodecanol
14.100	88.9	C13H28	184.2	1043328		Undecane, 3,8-dimethyl-
14.200	95.4	C15H30	210.2	2025597	13360-61-7	1-Pentadecene
14.300	95.1	C15H32	212.3	2925980	629-62-9	pentadecane
14.400	94.6	C15H24O	220.2	33483809	128-37-0	Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl-
14.400	91.0	C13H26O	198.2	517453	10486-19-8	Tridecanal
14.700	87.4	C16H48O6Si7	532.2	569060	541-01-5	Heptasiloxane, hexadecamethyl-
14.900	86.5	C20H42O3S	362.3	1094979		Sulfurous acid, hexyl tetradecyl ester
15.000	92.4	C15H30	210.2	1262360		n-Nonylcyclohexane
15.100	84.6	C20H42	282.3	716409		Hexadecane, 2,6,10,14-tetramethyl-
15.100	92.3	C16H34	226.3	742404		Pentadecane, 3-methyl-
15.300	83.7	C12H24	168.2	367595		5-Undecene, 3-methyl-, (E)-
15.500	90.4	C16H34	226.3	2616962		Hexadecane
15.800	90.5	C16H48O8Si8	592.2	783050		Cyclooctasiloxane, hexadecamethyl-
15.800	94.0	C15H30O2	242.2	2223845		Dodecanoic acid, 1-methylethyl ester
15.900	90.8	C13H100	182.1	345456		Benzophenone
						Pentadecane, 2,6,10-trimethyl-
16.000	86.7 86.1	C18H38 C17H36	254.3	701297		
16.100			240.3			3,3-Diethyltridecane
16.200	91.6	C16H32	224.3	1471140		Cyclohexadecane
16.200	89.6	C16H34O	242.3	1229202		Octane, 1,1'-oxybis-
16.300	85.8	C15H22O2	234.2	364044		1-(4-ISOPROPYLPHENYL)-2-METHYLPROPYL ACETATE
16.600	92.0	C19H38	266.3	1431481		1-Nonadecene
16.600	89.3	C17H36	240.3	1441202		Heptadecane
17.200	81.5	C16H33I	352.2	414711		Hexadecane, 1-iodo-
17.300	86.8	C16H16	208.1	3242567	20071-09-4	Benzene, 1,1'-(1,2-cyclobutanediyl)bis-, trans-
17.400	90.6	C17H34	238.3	484941	54105-66-7	Cyclohexane, undecyl-
17.500	98.0	C14H12O2	212.1	1123998	120-51-4	Benzyl benzoate
17.800	92.9	C15H22O3	250.2	401683	118-60-5	2-Ethylhexyl salicylate
18.400	88.6	C17H36O	256.3	419521		1-Heptadecanol
19.000	81.4	C18H54O9Si9	666.2	638205		Cyclononasiloxane, octadecamethyl-
20.800	85.7	C27H42NP	411.3	338463		1,1'-bis(t-Butylimino)-1,1'-(1",1"-diethylpropyl)-3,3'-diphenyllambda(5).,
	94.4	C16H10	202.1	837264		
21.100		010110	202.1	03/201	122 00-0	