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

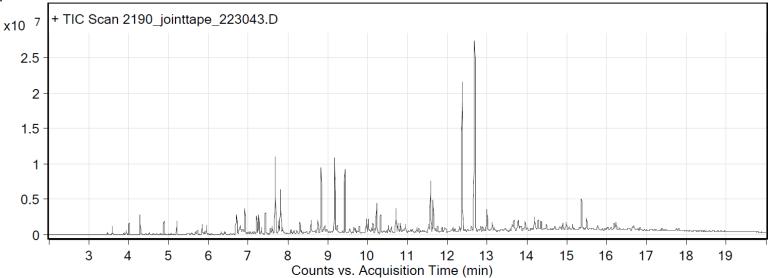
Alkynes – can become reactive when exposed to heat or UV light

Sample: Joint tape

Oddy test result: Unsuitable Date collected: 6/29/2018

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 cryo-trapped for 2 min at -15°C; GC ramped from 35°C to 250 °C at 10°C/min. Data analyzed in Masshunter Qualitative Analysis. Deconvoluted data with > 85% match with a NIST 17.0 or Wiley 9 library are reported.

VOCs not highlighted are because they were also observed in blanks: (1) 4.8 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



_		_		
Com	pound	-	ы	
COIII	ount	l a	и	ıc

und Table RT	Score (Lib)	Area	Name	Formula
3.59	97.96	856755	Acetic acid	C2H4O2
4	95.79		Triethylamine	C6H15N
4.28	94.5		1,2-Propanediol	C3H8O2
4.88	92.47	1436108	Cyclotrisiloxane, hexamethyl-	C6H18O3Si
5.21	94.56	1650028	2-Pentanone, 4-hydroxy-4-methyl-	C6H12O2
5.73	85.77	724467	Oxime-, methoxy-phenyl	C8H9NO2
5.84	94.51	1281400	2-Propenoic acid, butyl ester	C7H12O2
5.95	93.46	1273685	Ethanol, 2-butoxy-	C6H14O2
6.32	88.26	449381	Ethanol, 2-(2-methoxyethoxy)-	C5H12O3
6.7	97.03	5001453	Ethanol, 2,2'-oxybis-	C4H10O3
6.92	85.51	3732980	Cyclotetrasiloxane, octamethyl-	C8H24O4Si
7.21	85.58	2890593	dipropylene glycol monomethyl ether isomer, STRUCTURE UNKNOWN	C7H16O3
7.25	93.41	477512	unidentified C3-benzene	C9H12
7.26	94.2	3072436	dipropylene glycol monomethyl ether isomer, STRUCTURE UNKNOWN	C7H16O3
7.28	92.33	886261	Decane	C10H22
7.33	97.61	1087948	Octanal	C8H16O
7.43	97.92	3719067	dipropylene glycol monomethyl ether isomer, STRUCTURE UNKNOWN	C7H16O3
7.56	94.74	1378819	2-Propanol, 1,1'-oxybis-	C6H14O3
7.61	88.22	1229026	4-Cyanocyclohexene	C7H9N
7.68	97.28		1-Hexanol, 2-ethyl-	C8H18O
7.76	96.66		dl-Limonene	C10H16
7.86	93.48		2-Propanol, 1,1'-oxybis-	C6H14O3
8.22	90.95		Dodecane, 2,6,11-trimethyl-	C15H32
8.34	87.56		2-Octanol, 2-methyl-6-methylene-	C10H20O
8.68	85.29		dihexylsulfide	C12H26S
8.75	93.81		Undecane	C11H24
8.83	96.56	11955089		C9H18O
8.92	92.33	2472283	Hexanoic acid, 2-ethyl-	C8H16O2

9.17	95.29	12164102	Cyclopentasiloxane, decamethyl-	C10H30O5Si5
9.24	95.24		Pentanedioic acid, dimethyl ester	C7H12O4
9.42	96.47		Acetic acid, 2-ethylhexyl ester	C10H20O2
9.54	88.53		1,3-Pentanediol, 2,2,4-trimethyl-	C8H18O2
9.79	91.15		Decyl heptyl ether	C17H36O
9.96	97.67	2678193	Cyclohexanol, 5-methyl-2-(1- methylethyl)-	C10H20O
10.01	95.44	2660687	Ethanol, 2-(2-butoxyethoxy)-	C8H18O3
10.13	92.56	1630600		C10H8
10.18	96.94		Methyl salicylate	C8H8O3
10.22	92.15		Dodecane	C12H26
10.32	94.41	3441604		C10H20O
10.52	92.06		Ethanol, 2-phenoxy-	C8H10O2
10.59	90.19		2-Ethylhexyl acrylate	C11H20O2
10.71	89.2		Benzothiazole	C7H5NS
10.71	88.52		2-Propanol, 1,1'-oxybis-	C6H14O3
			2-Propanol, 1-(2-butoxy-1-methylethoxy)-	
10.82	88.1			C10H22O3
10.89	93.55	1496816	1-Phenoxypropan-2-ol Hexanoic acid, 2-ethyl-, 2-methylpropyl	C9H12O2 C12H24O2
44.00	20.42		ester	60114.000
11.09	88.42		Nonanoic acid	C9H18O2
11.24	94.33		1-Decanol	C10H22O
11.28	88.55		Hexadecane	C16H34
11.37	90.26		Benzaldehyde, 4-propyl-	C10H12O
11.53	86.29		pentadecene	C15H30
11.58	95.74		Cyclohexasiloxane, dodecamethyl-	C12H36O6Si6
11.64	94.34		Tridecane	C13H28
11.71	91		Naphthalene, 2-methyl-	C11H10
11.76	93.16	1205435	Tetradecanal	C14H28O
12.19	85.59	589092	5,6,7,8,9,10-Hexahydrobenzocyclooctene	C12H16
12.37	88.31	30629679	Propanoic acid, 2-methyl-, 2,2-dimethyl- 1-(2-hydroxy-1-methylethyl)propyl ester	C12H24O3
12.68	93.54	42365234	Propanoic acid, 2-methyl-, 3-hydroxy-	C12H24O3
			2,4,4-trimethylpentyl ester	
12.89	95.74	974887	1-Tetradecene	C14H28
13	95.63		Tetradecane	C14H30
13.13	91.6		Tetradecanal	C14H28O
13.63	85.19		5,5-Dibutylnonane	C17H36
13.67	92.1		Cyclopentane, nonyl-	C14H28
13.82	87.88		Tetracosane	C24H50
		100/431	2.5-Cyclohovadieno-1.4-diono. 2.6	
13.85	86.62	742273	2,5-Cyclohexadiene-1,4-dione, 2,6- bis(1,1-dimethylethyl)-	C14H20O2
13.95	95.61	2000250	1-Dodecanol	C12H26O
14.18	93.61		1-Pentadecene	C15H30
			pentadecane	
14.28	92.94		Phenol, 2,6-bis(1,1-dimethylethyl)-4-	C15H32
14.35	93.05	1436638	methyl-	C15H24O
14.48	89.49		Tetradecane, 2,2-dimethyl-	C16H34
14.9	87.59		Sulfurous acid, hexyl tetradecyl ester	C20H42O3S
14.99	93.66		n-Nonylcyclohexane	C15H30
15.06	88.87		Tetradecane, 3-methyl-	C15H32
15.14	86		Pentadecane, 3-methyl-	C16H34
15.49	88.55		Dodecane, 2,6,10-trimethyl-	C15H32
16.03	86.86		Tridecane, 5-propyl-	C16H34
16.11	86.01		3,3-Diethyltridecane	C17H36
46.40	90.82	1450706	Cyclopentane, undecyl-	C16H32
16.18	04.57	1500221	Methyl octyl ether	C9H20O
16.18	91.57	1009331	. rearly: eccly: earler	
	91.57 85.38		Pentadecane, 2,6,10,14-tetramethyl-	C19H40
16.23		1400541		C19H40 C12H26