

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
 Data File : VX029465.D
 Acq On : 14 Jun 2022 15:20
 Operator : JC/MD
 Sample : N3248-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EXYY2

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
 Title : VOC Analysis

Signal : TIC: VX029465.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.367	40	46	56	rVB3	224278	279246	13.15%	1.562%
2	1.685	91	98	108	rVB2	404056	642761	30.26%	3.596%
3	2.306	194	200	209	rBV	308089	506953	23.87%	2.836%
4	2.386	209	213	216	rVV	17643	29848	1.41%	0.167%
5	2.422	216	219	226	rVB	20677	33185	1.56%	0.186%
6	2.568	238	243	244	rBV4	507	670	0.03%	0.004%
7	2.654	251	257	263	rBV3	1231	2466	0.12%	0.014%
8	2.709	263	266	267	rBV	716	599	0.03%	0.003%
9	2.928	300	302	303	rBV2	956	711	0.03%	0.004%
10	2.971	305	309	320	rVV4	4671	12611	0.59%	0.071%
11	3.093	324	329	341	rVB5	7963	18432	0.87%	0.103%
12	3.331	363	368	377	rVB8	2487	6121	0.29%	0.034%
13	3.611	403	414	430	rBV	444638	1028364	48.41%	5.753%
14	4.093	491	493	496	rBV2	347	508	0.02%	0.003%
15	4.367	533	538	543	rVB4	1777	3396	0.16%	0.019%
16	4.489	545	558	586	rBV2	709741	2124148	100.00%	11.883%
17	5.013	633	644	648	rBV	213613	625439	29.44%	3.499%
18	5.391	693	706	725	rVV	165953	513492	24.17%	2.873%
19	5.556	729	733	742	rVB4	2473	6536	0.31%	0.037%
20	5.970	789	801	818	rBV2	446770	1328932	62.56%	7.434%
21	6.367	860	866	868	rBV3	820	1456	0.07%	0.008%
22	6.562	896	898	905	rVB4	354	535	0.03%	0.003%
23	6.763	920	931	947	rBV	348477	838844	39.49%	4.693%
24	7.129	982	991	1004	rBV3	65661	150974	7.11%	0.845%
25	7.312	1012	1021	1031	rBV	266491	580856	27.35%	3.249%
26	7.482	1045	1049	1053	rBV3	619	1066	0.05%	0.006%
27	7.811	1100	1103	1105	rBV2	524	573	0.03%	0.003%
28	7.915	1115	1120	1121	rBV3	856	1400	0.07%	0.008%
29	8.330	1181	1188	1198	rVV	175634	288709	13.59%	1.615%
30	8.415	1198	1202	1210	rVB2	8425	14990	0.71%	0.084%
31	8.598	1226	1232	1234	rBV3	2175	3562	0.17%	0.020%
32	8.653	1234	1241	1248	rVV	710194	1104768	52.01%	6.180%
33	8.720	1248	1252	1264	rVB	50300	83774	3.94%	0.469%
34	8.951	1279	1290	1302	rBV	121577	186830	8.80%	1.045%
35	9.104	1310	1315	1322	rVB4	1402	2510	0.12%	0.014%
36	9.275	1337	1343	1353	rVB	229954	328077	15.45%	1.835%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
 Title : VOC Analysis

37	9.384	1356	1361	1379	rBV	454873	676843	31.86%	3.786%
38	9.622	1396	1400	1403	rBV4	1853	2640	0.12%	0.015%
39	9.842	1431	1436	1440	rVB4	704	1170	0.06%	0.007%
40	9.890	1440	1444	1447	rBV3	416	644	0.03%	0.004%
41	10.055	1465	1471	1481	rBV	718394	975230	45.91%	5.456%
42	10.195	1489	1494	1500	rBV	69359	91277	4.30%	0.511%
43	10.305	1506	1512	1518	rVB	22008	33144	1.56%	0.185%
44	10.354	1518	1520	1523	rVB2	440	556	0.03%	0.003%
45	10.457	1533	1537	1542	rVB5	1532	2410	0.11%	0.013%
46	10.604	1553	1561	1562	rBV4	2336	4031	0.19%	0.023%
47	10.646	1563	1568	1573	rVB	117200	153775	7.24%	0.860%
48	10.780	1583	1590	1594	rVV3	7982	12306	0.58%	0.069%
49	10.866	1600	1604	1608	rVV6	2966	4110	0.19%	0.023%
50	10.921	1610	1613	1615	rBV2	570	626	0.03%	0.004%
51	10.963	1616	1620	1626	rVB	21054	27668	1.30%	0.155%
52	11.097	1638	1642	1643	rBV3	3162	4362	0.21%	0.024%
53	11.128	1644	1647	1652	rVB2	10251	13632	0.64%	0.076%
54	11.195	1652	1658	1666	rBV	491251	639538	30.11%	3.578%
55	11.305	1672	1676	1683	rVB	53292	68169	3.21%	0.381%
56	11.378	1683	1688	1696	rBV	88306	132321	6.23%	0.740%
57	11.451	1696	1700	1712	rVB	117108	156342	7.36%	0.875%
58	11.549	1712	1716	1718	rBV4	974	1332	0.06%	0.007%
59	11.616	1722	1727	1733	rBV	186791	234611	11.04%	1.312%
60	11.664	1733	1735	1740	rVB6	1669	2107	0.10%	0.012%
61	11.713	1740	1743	1745	rBV2	3550	4402	0.21%	0.025%
62	11.756	1745	1750	1760	rVB	598715	723816	34.08%	4.049%
63	11.866	1762	1768	1769	rBV4	6459	9281	0.44%	0.052%
64	11.890	1769	1772	1776	rVB	18937	24857	1.17%	0.139%
65	11.933	1776	1779	1781	rBV2	2928	2932	0.14%	0.016%
66	11.969	1782	1785	1789	rVB	20689	23840	1.12%	0.133%
67	12.024	1789	1794	1800	rBV	830674	989520	46.58%	5.536%
68	12.091	1800	1805	1812	rBV	308181	388826	18.31%	2.175%
69	12.177	1817	1819	1822	rVB	3899	4042	0.19%	0.023%
70	12.219	1822	1826	1828	rBV	91476	125947	5.93%	0.705%
71	12.244	1828	1830	1838	rVV3	126381	182392	8.59%	1.020%
72	12.323	1838	1843	1850	rVB	760933	990129	46.61%	5.539%
73	12.463	1863	1866	1872	rVB	17226	23194	1.09%	0.130%
74	12.530	1873	1877	1879	rBV	20267	25613	1.21%	0.143%
75	12.615	1887	1891	1894	rBV	31358	35423	1.67%	0.198%
76	12.707	1901	1906	1913	rVB3	29933	58909	2.77%	0.330%

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77	12.853	1926	1930	1935	rVB2	18249	24948	1.17%	0.140%
78	12.920	1935	1941	1945	rBV	16093	22999	1.08%	0.129%
79	12.963	1945	1948	1955	rVB	29169	35134	1.65%	0.197%
80	13.030	1955	1959	1961	rBV4	2932	4292	0.20%	0.024%
81	13.140	1973	1977	1979	rBV3	3813	4097	0.19%	0.023%
82	13.170	1979	1982	1985	rVV2	4839	4963	0.23%	0.028%
83	13.256	1994	1996	1997	rBV2	1593	993	0.05%	0.006%
84	13.292	1997	2002	2008	rVB2	40438	57174	2.69%	0.320%
85	13.426	2019	2024	2028	rVB	16858	22937	1.08%	0.128%
86	13.512	2034	2038	2040	rBV3	1520	2163	0.10%	0.012%
87	13.542	2040	2043	2046	rBV3	1805	2218	0.10%	0.012%
88	13.585	2046	2050	2055	rBV5	3390	6159	0.29%	0.034%
89	13.780	2076	2082	2089	rVB	35312	48042	2.26%	0.269%
90	13.853	2089	2094	2097	rBV5	1198	2007	0.09%	0.011%
91	13.926	2104	2106	2108	rBV2	1053	930	0.04%	0.005%
92	14.073	2128	2130	2133	rVB3	666	734	0.03%	0.004%
93	14.103	2133	2135	2137	rBV2	609	487	0.02%	0.003%
94	14.127	2137	2139	2142	rVB2	864	920	0.04%	0.005%
95	14.225	2151	2155	2159	rBV5	1725	3040	0.14%	0.017%
96	14.481	2192	2197	2199	rBV5	1398	2004	0.09%	0.011%
97	14.640	2219	2223	2231	rVB	5106	6931	0.33%	0.039%
98	14.780	2240	2246	2250	rBV4	4609	6934	0.33%	0.039%
99	15.048	2289	2290	2293	rBV2	473	474	0.02%	0.003%
100	15.975	2436	2442	2448	rBV8	5316	9780	0.46%	0.055%

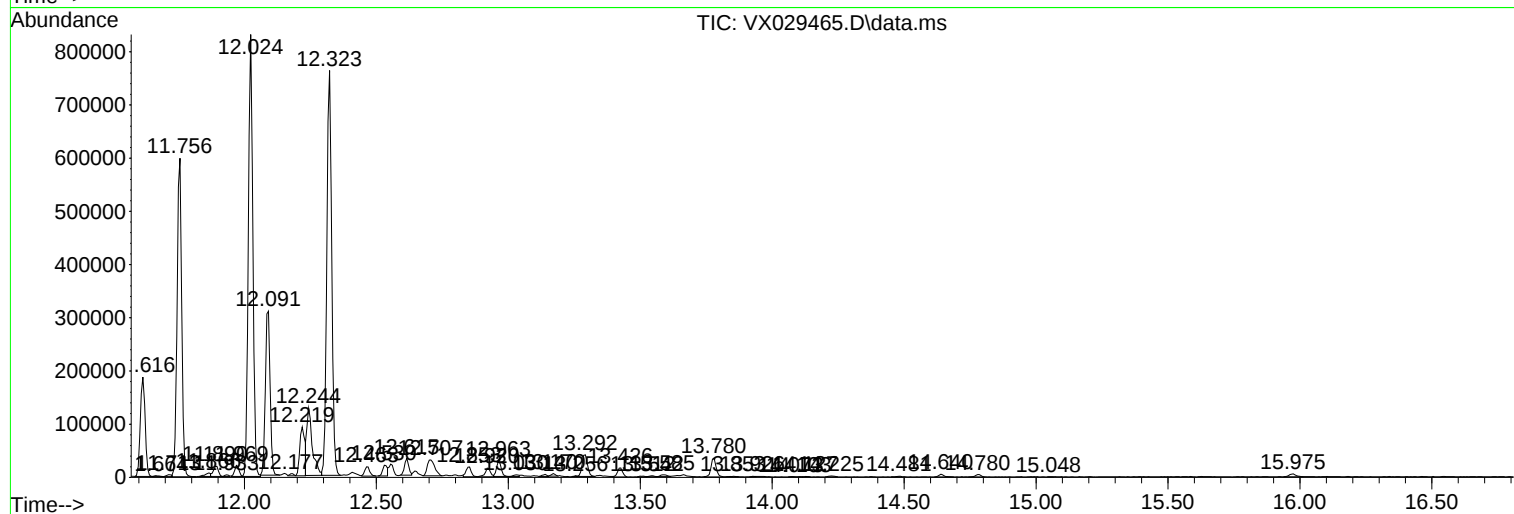
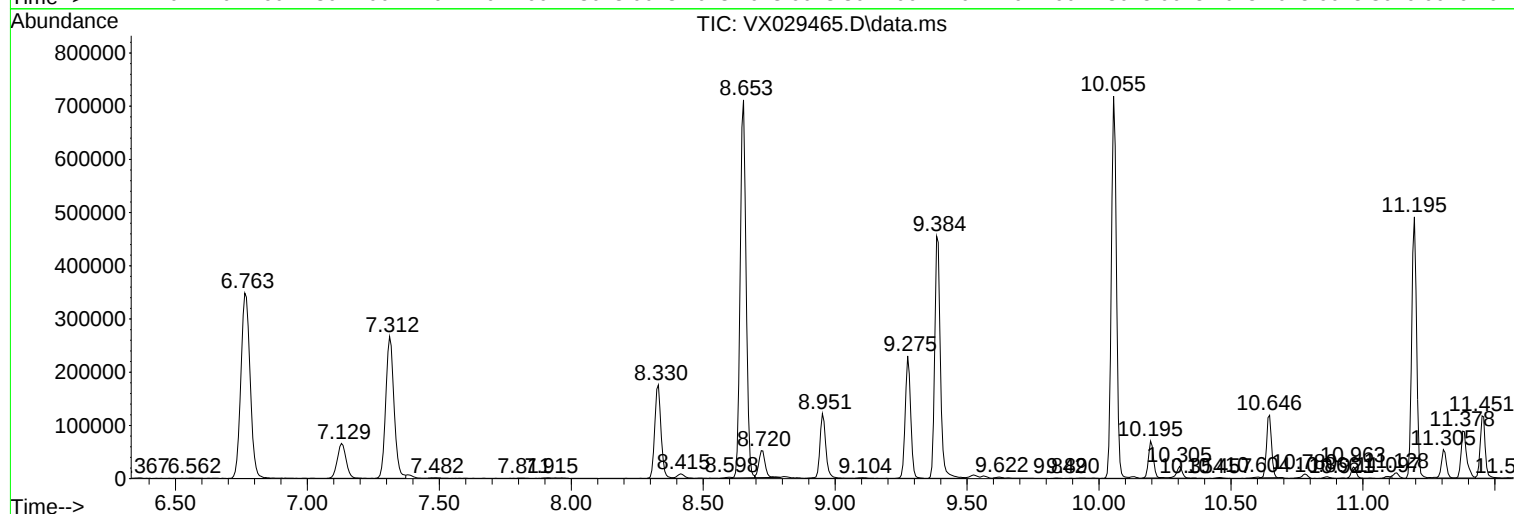
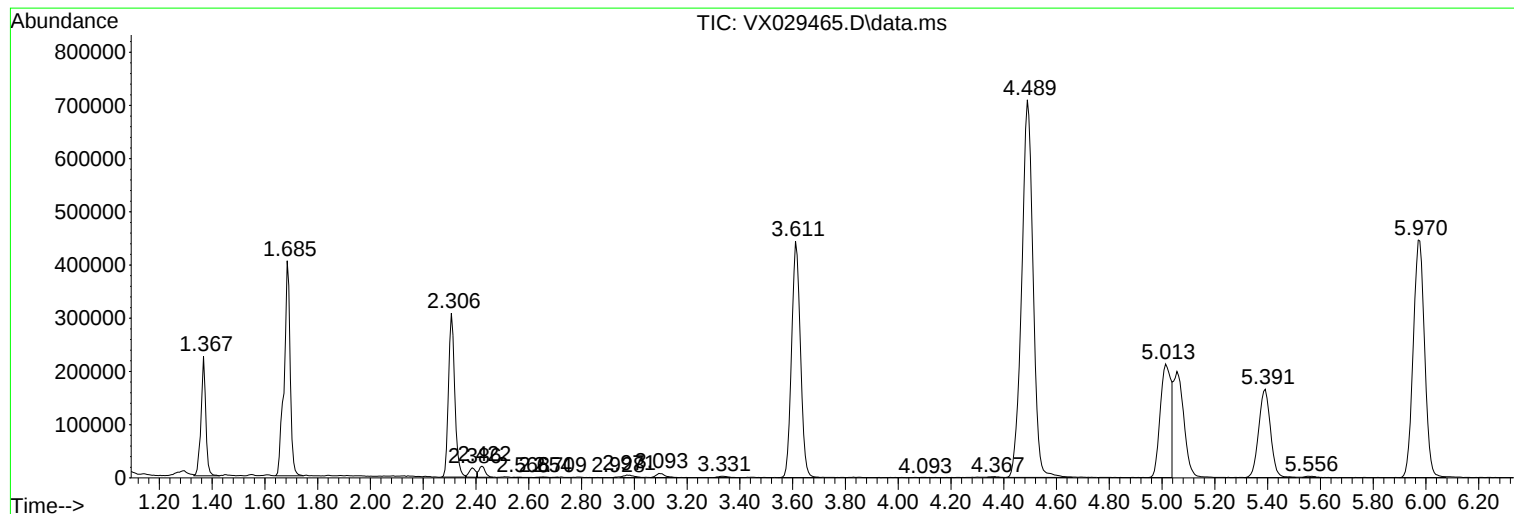
Sum of corrected areas: 17875669

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ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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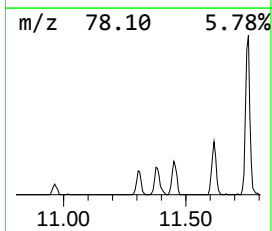
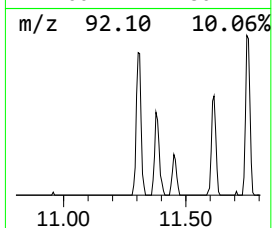
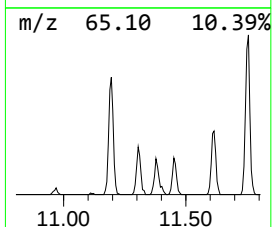
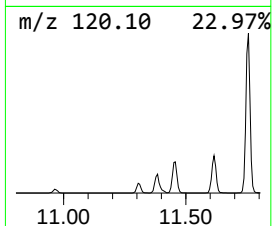
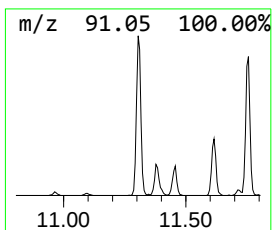
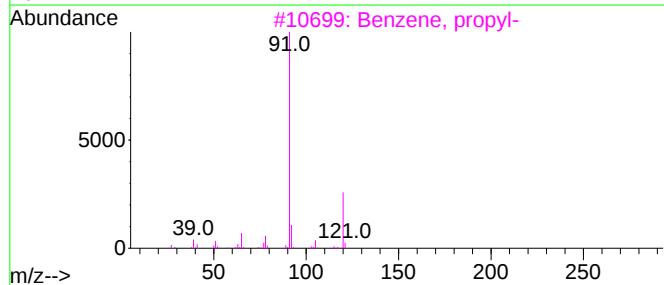
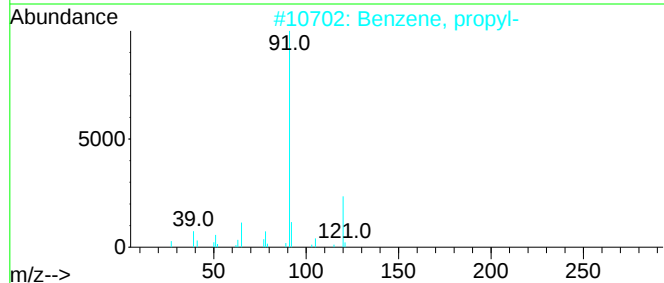
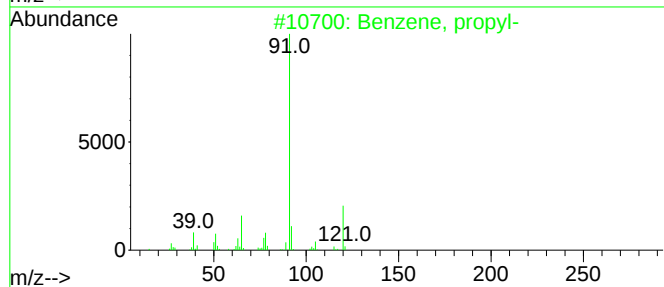
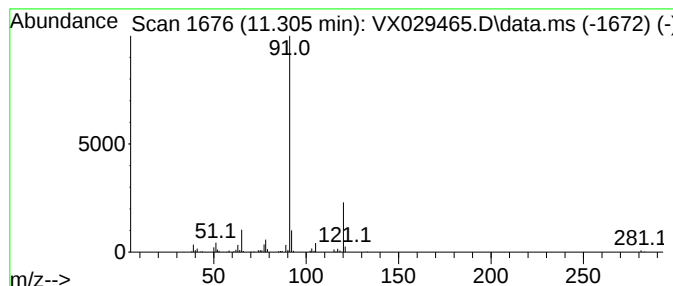
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.305	3.44 ug/L	68169	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	90
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			Benzene, propyl-	120	C9H12	000103-65-1	87
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78



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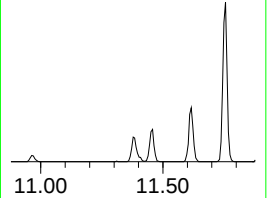
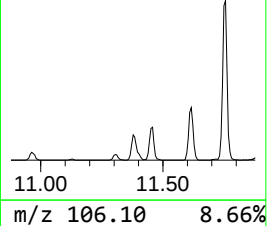
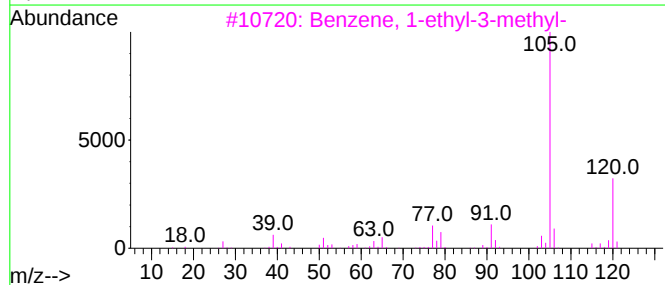
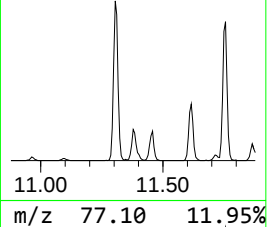
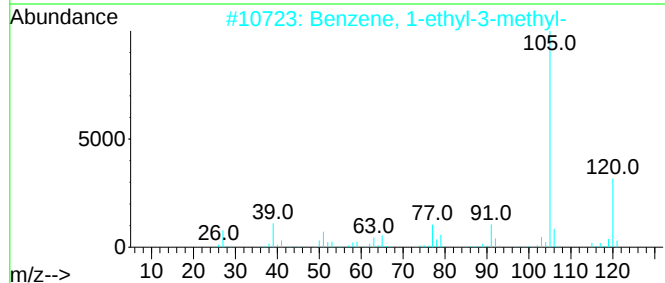
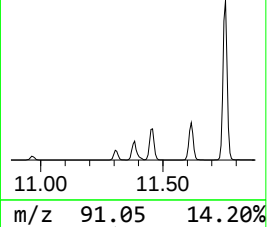
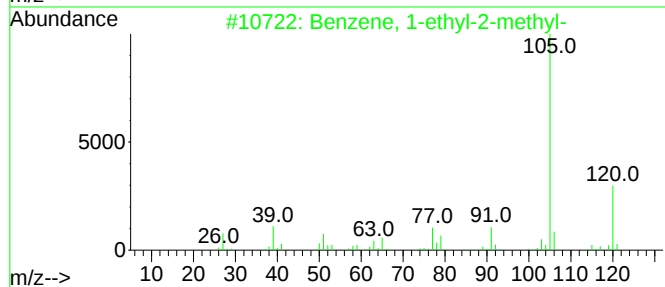
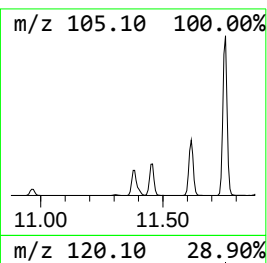
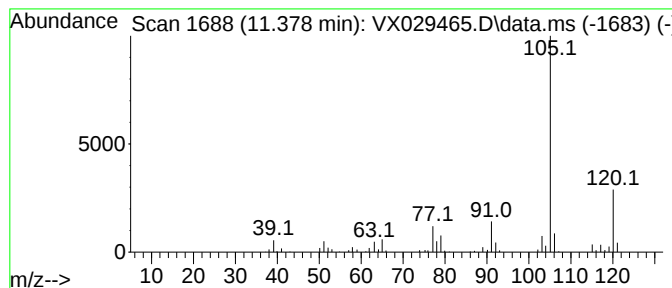
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	6.69 ug/L	132321	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



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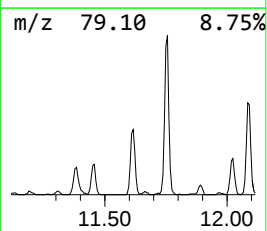
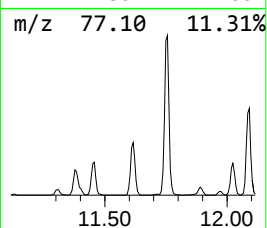
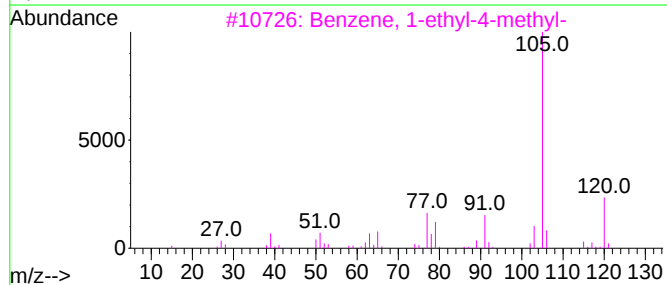
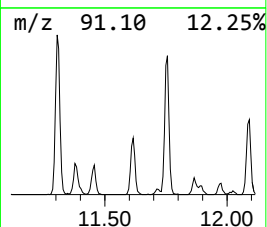
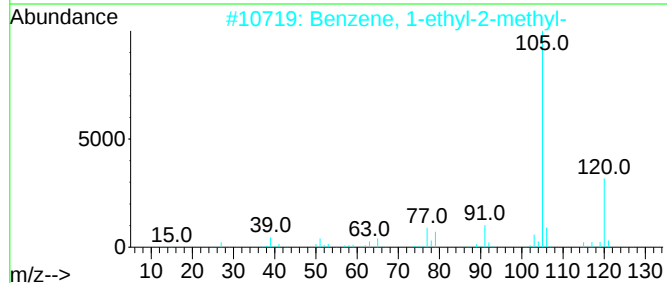
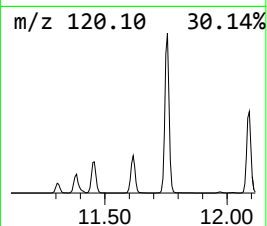
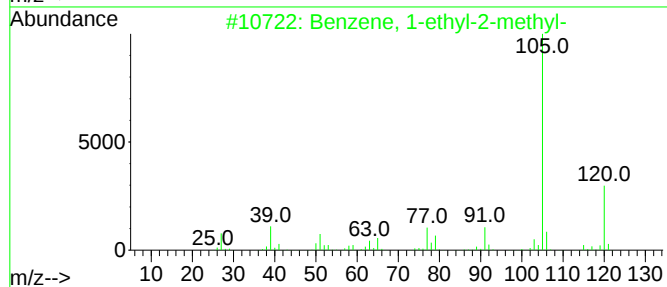
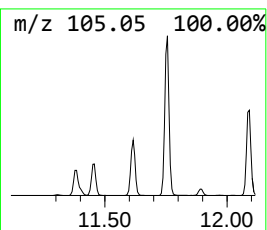
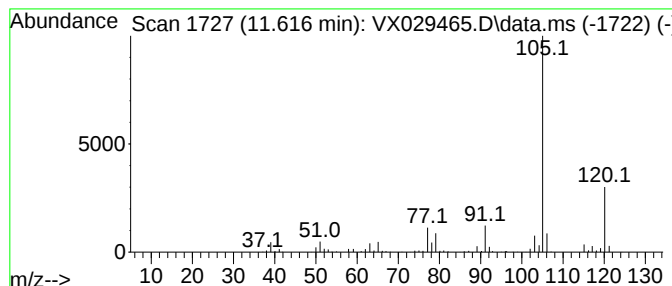
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	11.85 ug/L	234611	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
Data File : VX029465.D
Acq On : 14 Jun 2022 15:20
Operator : JC/MD
Sample : N3248-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

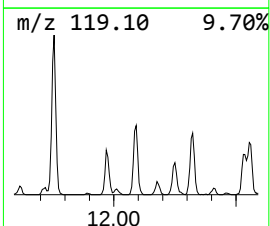
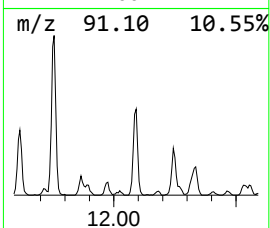
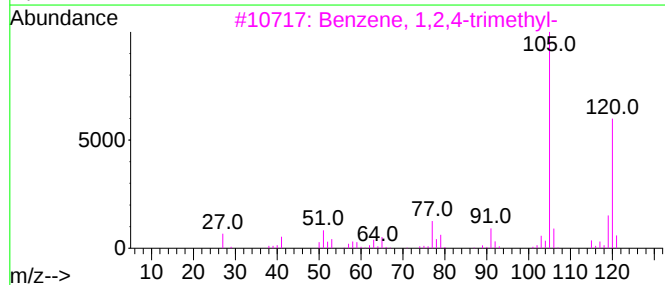
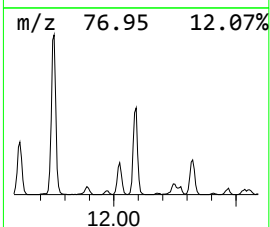
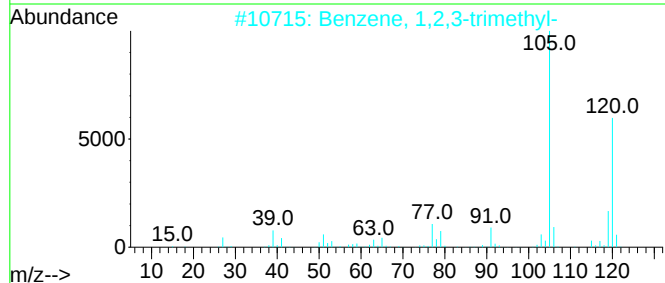
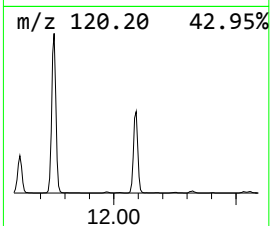
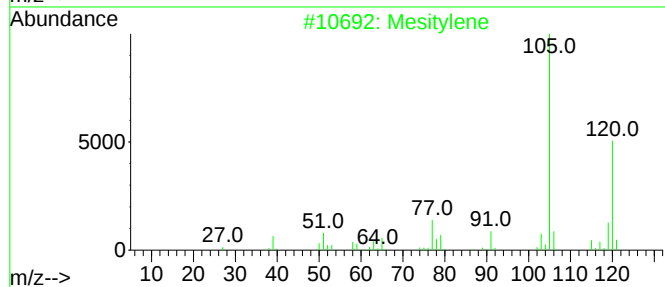
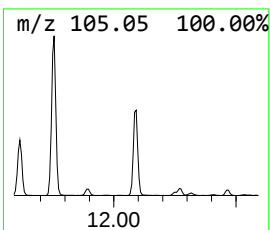
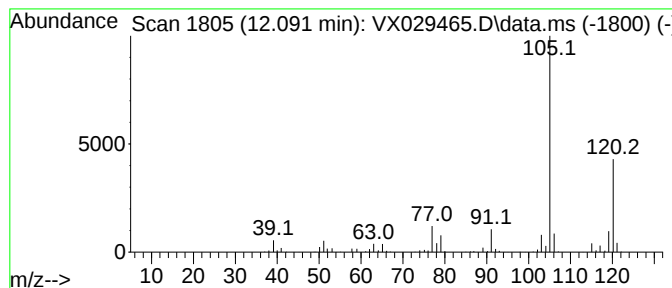
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.091	19.65 ug/L	388826	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	95
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
Data File : VX029465.D
Acq On : 14 Jun 2022 15:20
Operator : JC/MD
Sample : N3248-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

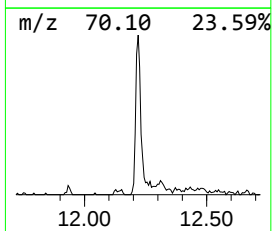
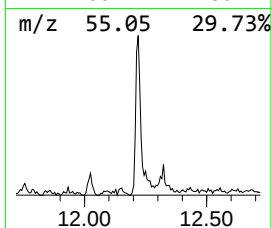
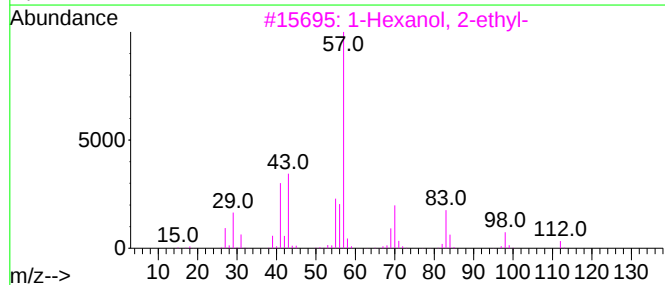
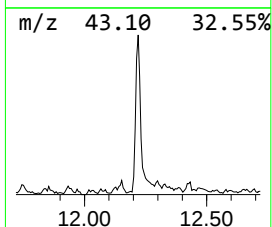
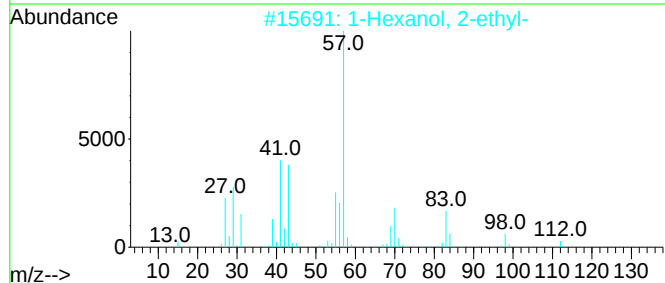
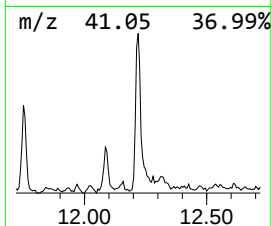
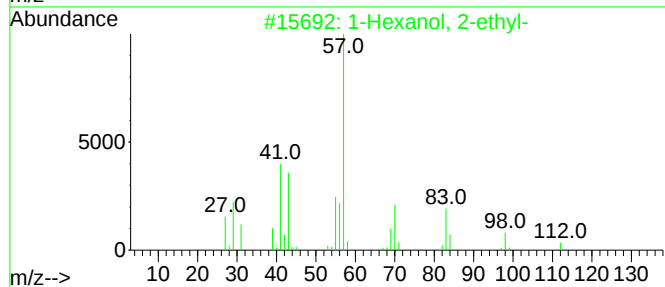
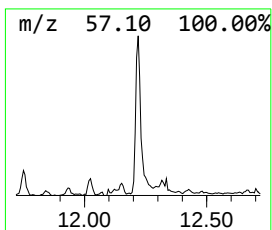
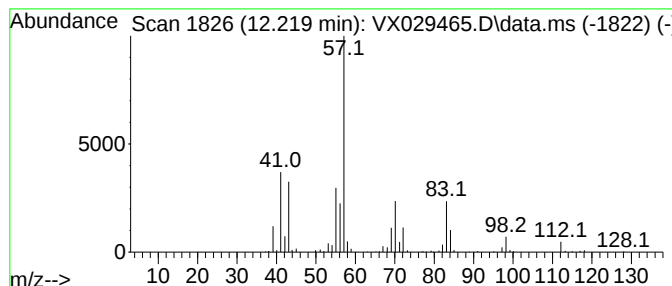
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.219	6.36 ug/L	125947	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	86
2		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
3		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	78
4		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
5		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
Data File : VX029465.D
Acq On : 14 Jun 2022 15:20
Operator : JC/MD
Sample : N3248-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

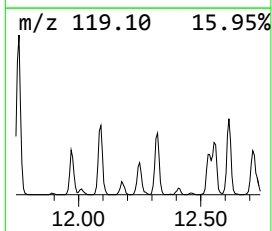
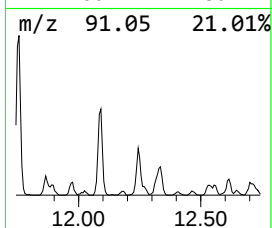
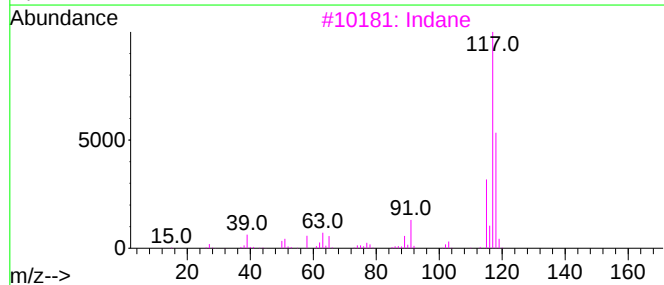
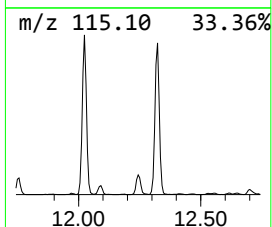
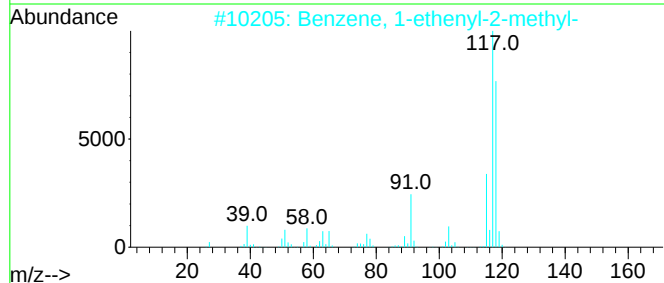
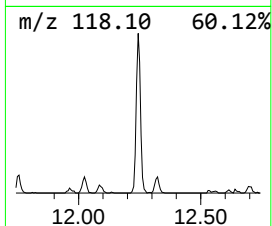
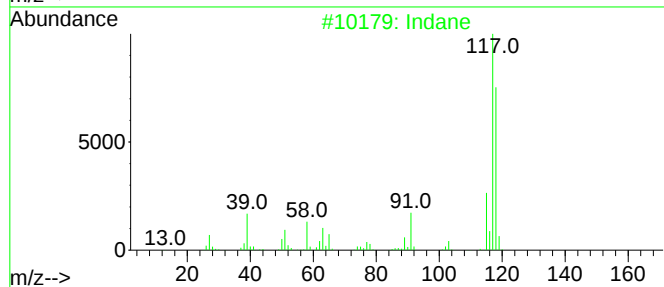
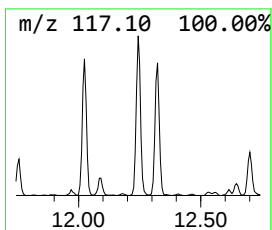
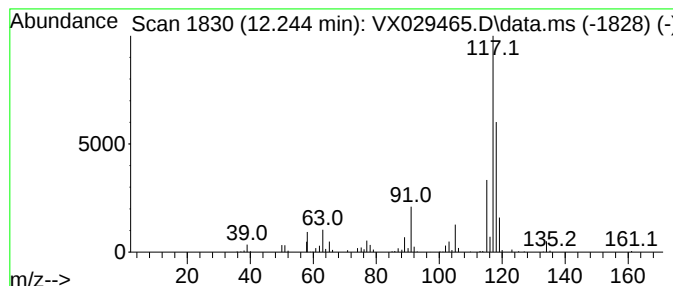
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	9.22 ug/L	182392	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	76
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70
3			Indane	118	C9H10	000496-11-7	68
4			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
Data File : VX029465.D
Acq On : 14 Jun 2022 15:20
Operator : JC/MD
Sample : N3248-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

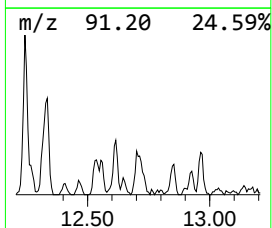
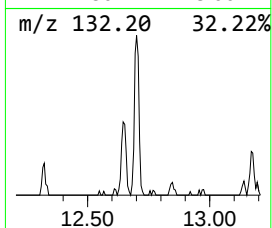
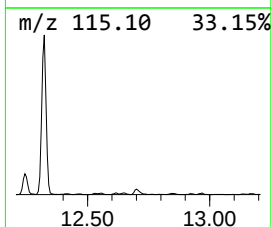
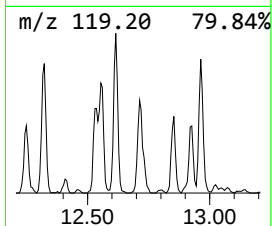
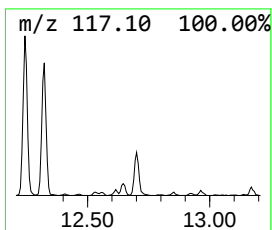
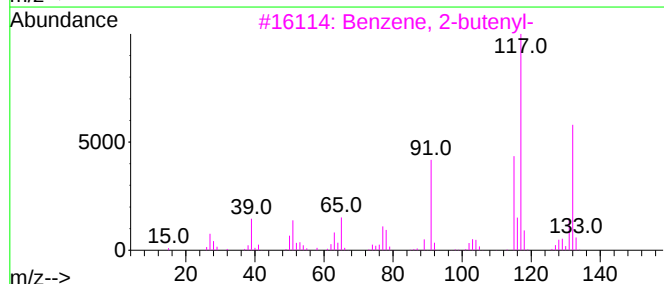
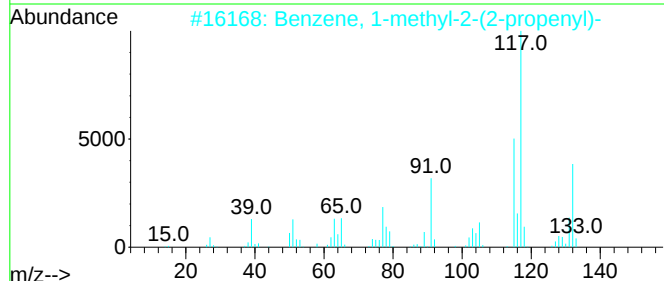
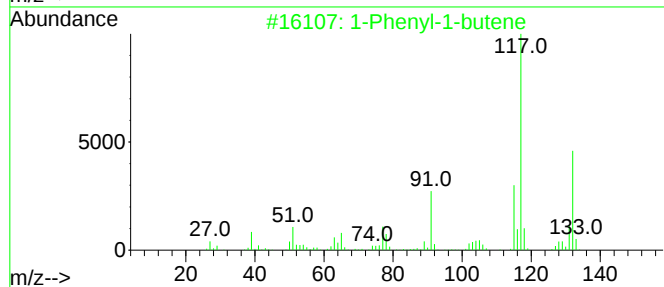
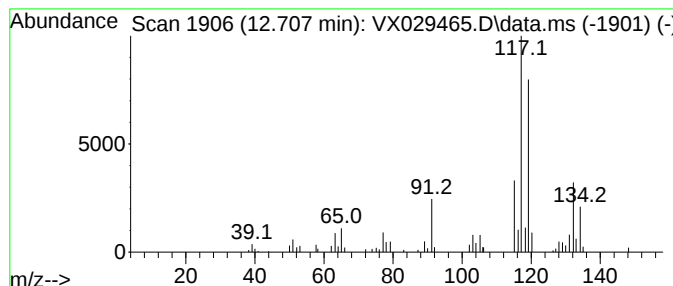
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Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.707	2.98 ug/L	58909	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	55
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
Data File : VX029465.D
Acq On : 14 Jun 2022 15:20
Operator : JC/MD
Sample : N3248-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

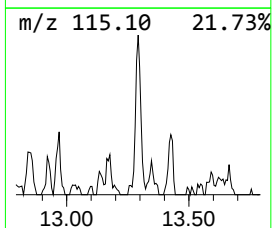
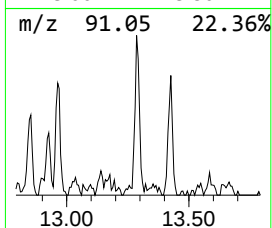
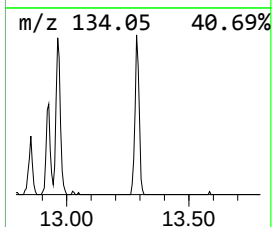
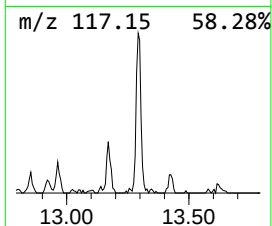
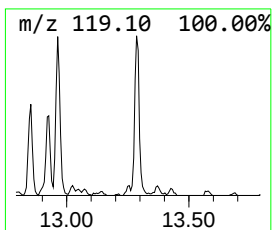
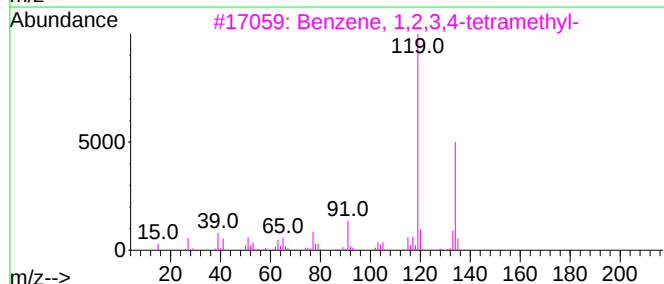
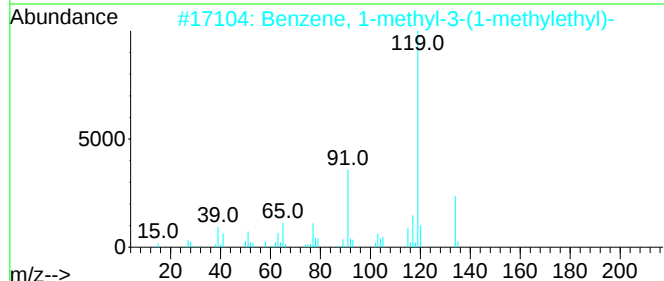
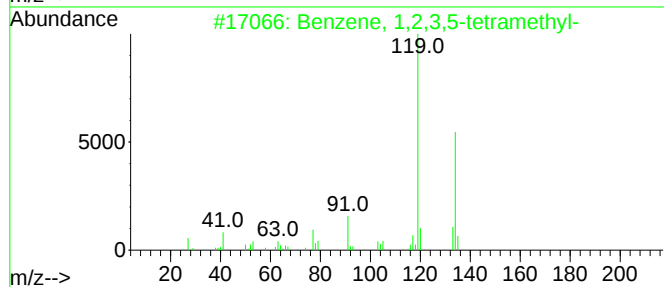
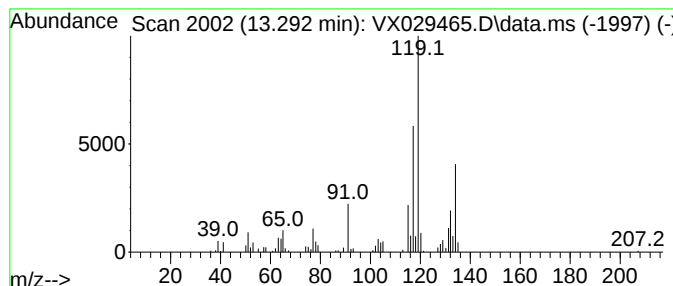
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	2.89 ug/L	57174	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	55
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	53
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	53
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	53
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX061422\
Data File : VX029465.D
Acq On : 14 Jun 2022 15:20
Operator : JC/MD
Sample : N3248-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EXYY2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXML061322WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, propyl-	11.305	3.4	ug/L	68169	3	12.024	989520	50.0
Benzene, 1-ethy...	11.378	6.7	ug/L	132321	3	12.024	989520	50.0
Benzene, 1-ethy...	11.616	11.9	ug/L	234611	3	12.024	989520	50.0
Benzene, 1-ethy...	12.091	19.6	ug/L	388826	3	12.024	989520	50.0
1-Hexanol, 2-et...	12.219	6.4	ug/L	125947	3	12.024	989520	50.0
Indane	12.244	9.2	ug/L	182392	3	12.024	989520	50.0
1-Phenyl-1-butene	12.707	3.0	ug/L	58909	3	12.024	989520	50.0
Benzene, 1,2,3,...	13.292	2.9	ug/L	57174	3	12.024	989520	50.0