

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX070319\
 Data File : VX010675.D
 Acq On : 04 Jul 2019 08:15
 Operator : JC/SP
 Sample : K3664-02
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FL-0580

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.276	16	20	25	rBV	39577	53369	1.30%	0.211%
2	1.788	100	104	110	rVB	30235	46210	1.12%	0.182%
3	2.843	269	277	285	rBV2	21818	52053	1.26%	0.205%
4	2.934	285	292	299	rVB	29262	59483	1.44%	0.235%
5	3.178	323	332	341	rVB	58541	133655	3.25%	0.527%
6	4.403	525	533	547	rVB2	56494	164694	4.00%	0.650%
7	5.501	701	713	720	rBV	135523	386270	9.38%	1.524%
8	5.580	720	726	732	rVV2	48667	127972	3.11%	0.505%
9	5.659	732	739	749	rVV	210196	561535	13.63%	2.215%
10	5.952	776	787	797	rBV3	18181	55586	1.35%	0.219%
11	6.068	797	806	814	rVV	130981	346758	8.42%	1.368%
12	6.147	814	819	828	rVV2	16001	42114	1.02%	0.166%
13	6.263	828	838	847	rVV3	32043	109781	2.67%	0.433%
14	6.360	847	854	862	rVV3	30971	87325	2.12%	0.344%
15	6.458	862	870	880	rVB2	30593	87085	2.11%	0.344%
16	6.860	925	936	948	rBV	355907	824517	20.02%	3.253%
17	7.464	1022	1035	1046	rVB2	122725	301982	7.33%	1.191%
18	7.848	1088	1098	1106	rBV3	23221	50765	1.23%	0.200%
19	8.031	1121	1128	1140	rVB2	30521	73109	1.78%	0.288%
20	8.390	1179	1187	1193	rBV3	23640	45108	1.10%	0.178%
21	8.665	1224	1232	1235	rBV2	51179	93033	2.26%	0.367%
22	8.713	1235	1240	1248	rVB	748439	1164729	28.28%	4.595%
23	8.841	1254	1261	1268	rBV	51120	90787	2.20%	0.358%
24	9.037	1288	1293	1300	rVB	102175	172154	4.18%	0.679%
25	9.165	1307	1314	1319	rBV	46242	74979	1.82%	0.296%
26	9.567	1375	1380	1385	rBV	30641	47478	1.15%	0.187%
27	9.677	1393	1398	1403	rBV	73163	112087	2.72%	0.442%
28	10.110	1463	1469	1476	rBV	737211	989757	24.03%	3.905%
29	10.183	1476	1481	1486	rVB	37949	55858	1.36%	0.220%
30	10.250	1486	1492	1502	rBV	288397	384054	9.32%	1.515%
31	10.359	1502	1510	1516	rVV	306247	427503	10.38%	1.686%
32	10.652	1547	1558	1562	rBV6	24890	67719	1.64%	0.267%
33	10.701	1562	1566	1575	rVB2	43378	67803	1.65%	0.267%
34	10.835	1584	1588	1595	rVB	30737	44100	1.07%	0.174%

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Integration Parameters: RTEINT.P

Integrator: RTE
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 Sampling : 1
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35	10.914	1595	1601	1608	rBV7	17930	41730	1.01%	0.165%
36	11.018	1613	1618	1631	rVB	187944	245130	5.95%	0.967%
37	11.134	1632	1637	1642	rBV	643117	800751	19.44%	3.159%
38	11.359	1669	1674	1683	rVB	230647	293672	7.13%	1.159%
39	11.451	1683	1689	1694	rBV	87603	119112	2.89%	0.470%
40	11.664	1715	1724	1734	rBV	154205	222236	5.40%	0.877%
41	11.804	1742	1747	1753	rVB	3494396	4118624	100.00%	16.248%
42	11.945	1767	1770	1778	rVB	68569	89058	2.16%	0.351%
43	12.079	1786	1792	1797	rBV	766938	1001783	24.32%	3.952%
44	12.140	1797	1802	1812	rVB	1667667	2009201	48.78%	7.926%
45	12.231	1812	1817	1822	rBV	63351	75038	1.82%	0.296%
46	12.298	1823	1828	1833	rBV	494931	604083	14.67%	2.383%
47	12.365	1833	1839	1849	rVB2	120392	194332	4.72%	0.767%
48	12.457	1849	1854	1859	rBV	151762	184717	4.48%	0.729%
49	12.518	1859	1864	1870	rBV	317345	384703	9.34%	1.518%
50	12.585	1870	1875	1877	rBV	347146	442941	10.75%	1.747%
51	12.609	1877	1879	1883	rVV	379965	376783	9.15%	1.486%
52	12.664	1883	1888	1892	rVV	396109	486727	11.82%	1.920%
53	12.701	1892	1894	1898	rVV	97037	110081	2.67%	0.434%
54	12.755	1898	1903	1910	rVB2	310185	494499	12.01%	1.951%
55	12.902	1922	1927	1932	rVB2	250952	333285	8.09%	1.315%
56	12.975	1932	1939	1942	rBV	280064	369900	8.98%	1.459%
57	13.018	1942	1946	1952	rVB	475452	585104	14.21%	2.308%
58	13.078	1952	1956	1962	rVB3	69994	95565	2.32%	0.377%
59	13.146	1962	1967	1971	rBV2	40382	61765	1.50%	0.244%
60	13.194	1971	1975	1977	rBV	73175	85347	2.07%	0.337%
61	13.225	1977	1980	1983	rVB	63148	64427	1.56%	0.254%
62	13.261	1983	1986	1989	rVB3	39350	49303	1.20%	0.194%
63	13.304	1989	1993	1995	rBV	74439	94496	2.29%	0.373%
64	13.341	1995	1999	2005	rVB2	857103	1163694	28.25%	4.591%
65	13.475	2015	2021	2027	rVB	390104	511063	12.41%	2.016%
66	13.560	2027	2035	2038	rBV3	32932	61673	1.50%	0.243%
67	13.597	2038	2041	2044	rVV	43710	67654	1.64%	0.267%
68	13.633	2044	2047	2053	rVV3	113163	237800	5.77%	0.938%
69	13.700	2054	2058	2059	rVV	59118	91657	2.23%	0.362%
70	13.719	2059	2061	2069	rVB2	93755	142539	3.46%	0.562%
71	13.834	2069	2080	2087	rBV	441652	687573	16.69%	2.712%

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Integration Parameters: RTEINT.P

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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72	13.908	2087	2092	2097	rVB4	45574	68122	1.65%	0.269%
73	13.981	2097	2104	2108	rBV2	81494	121559	2.95%	0.480%
74	14.286	2146	2154	2159	rVB3	82111	167933	4.08%	0.662%
75	14.426	2173	2177	2181	rVB2	33867	43874	1.07%	0.173%
76	14.536	2190	2195	2204	rBV2	102816	147979	3.59%	0.584%
77	14.694	2214	2221	2224	rBV	99690	139554	3.39%	0.551%
78	14.834	2239	2244	2249	rBV	197198	260504	6.33%	1.028%

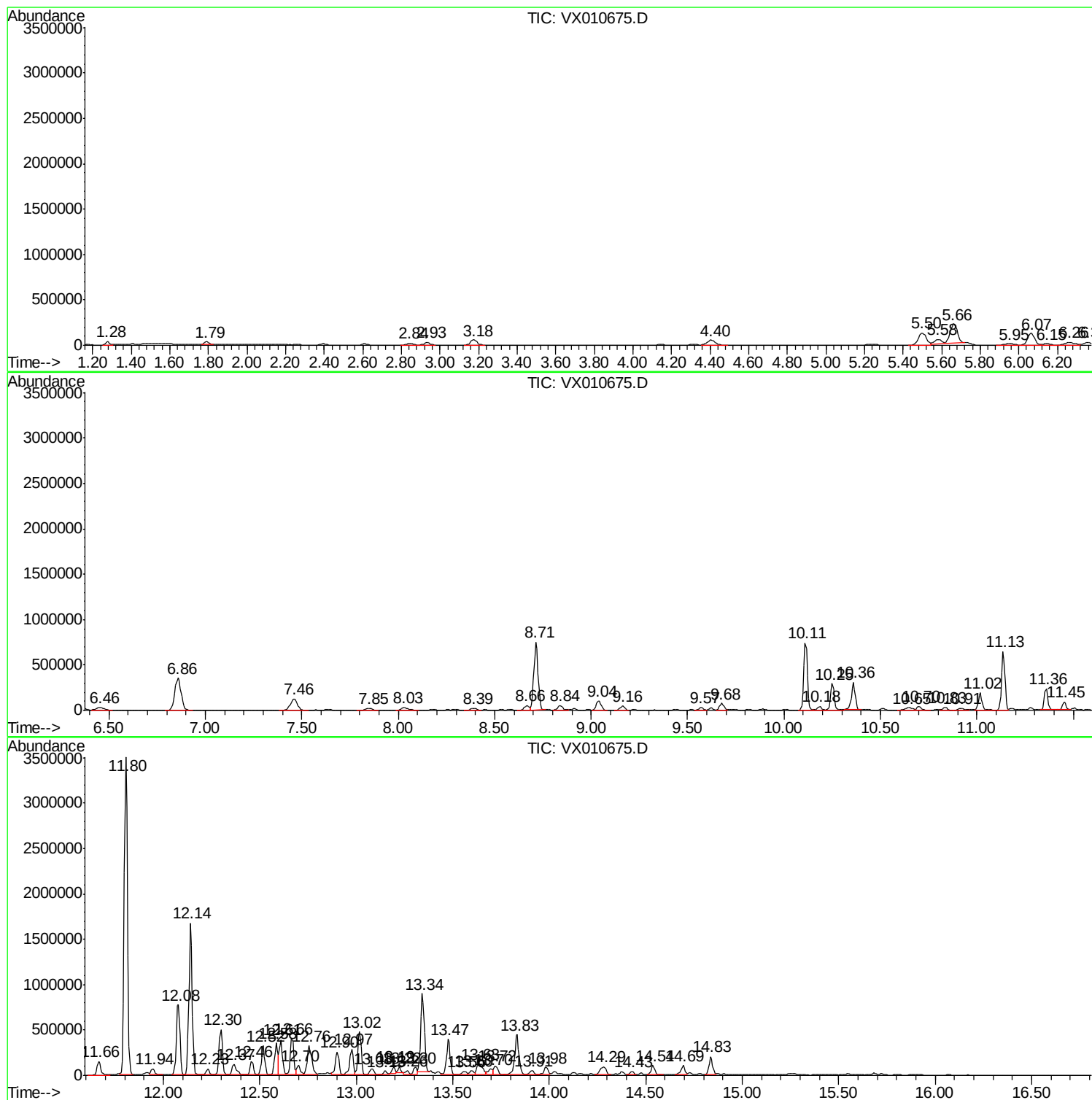
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Instrument :
MSVOA_X
ClientSampleId :
FL-0580

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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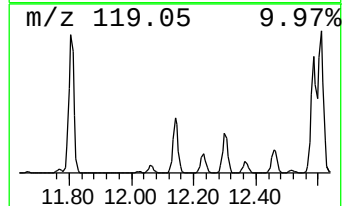
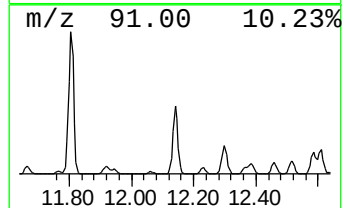
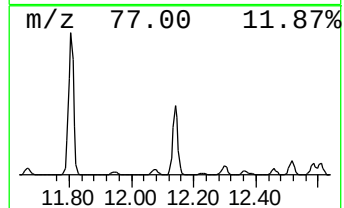
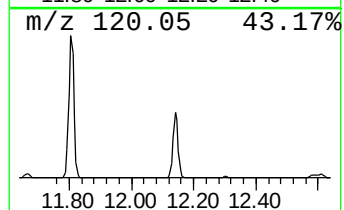
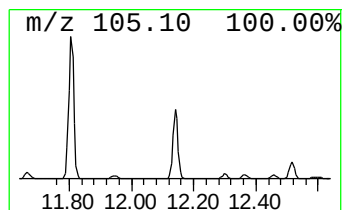
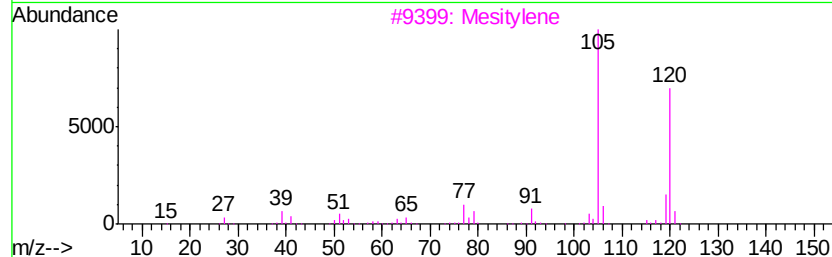
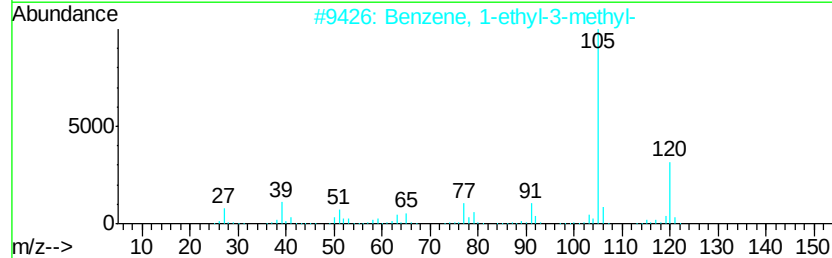
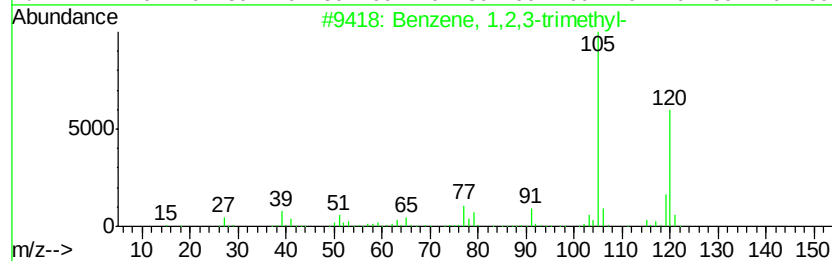
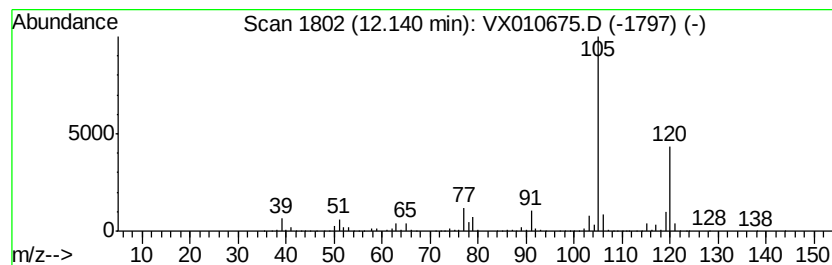
Instrument :
MSVOA_X
ClientSampled :
FL-0580

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.14	100.28 ug/l	2009200	1,4-Dichlorobenzene-d4	12.08		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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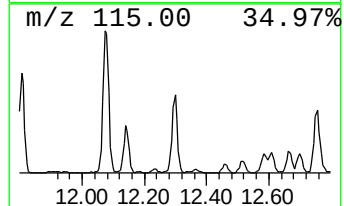
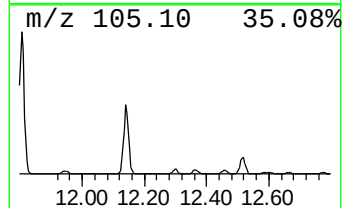
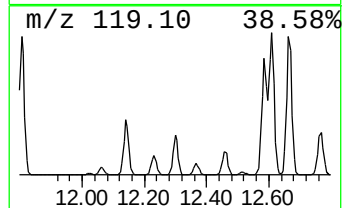
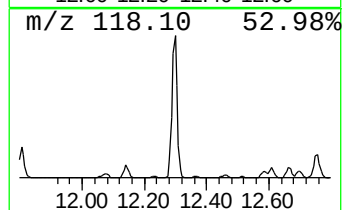
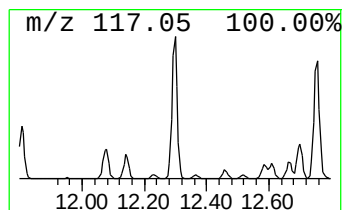
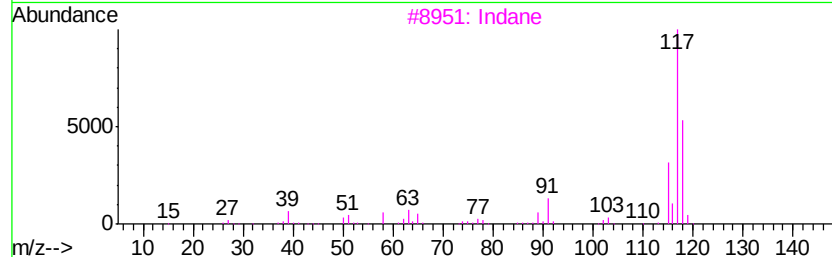
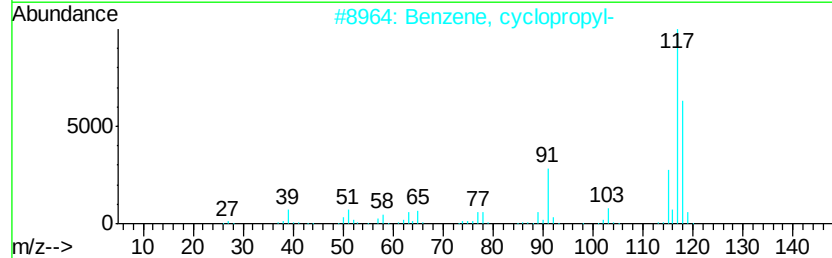
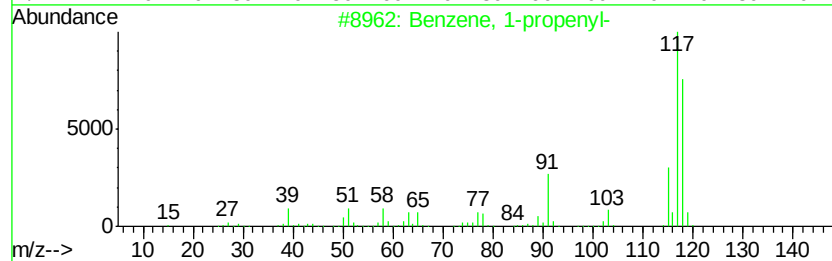
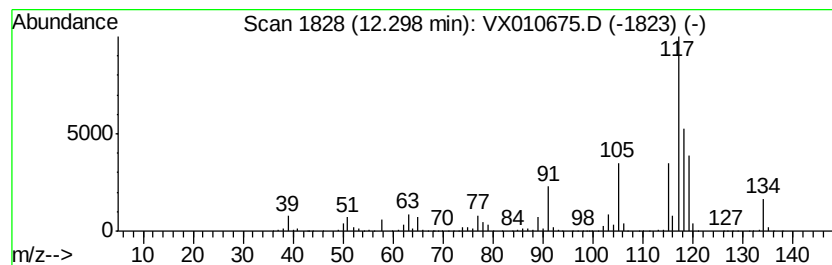
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-propenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	30.15 ug/l	604083	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-propenyl-	118 C9H10	000637-50-3 90
2		Benzene, cyclopropyl-	118 C9H10	000873-49-4 55
3		Indane	118 C9H10	000496-11-7 55
4		Benzene, 2-propenyl-	118 C9H10	000300-57-2 55
5		Benzeneethanol, .beta.-ethenyl-	148 C10H12O	006052-63-7 50



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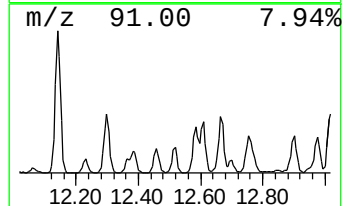
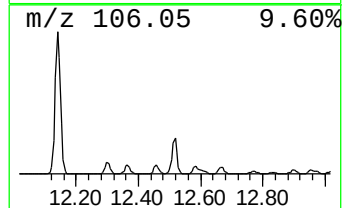
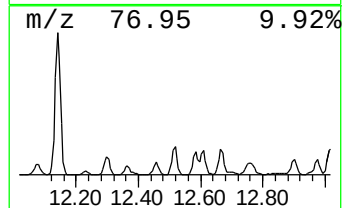
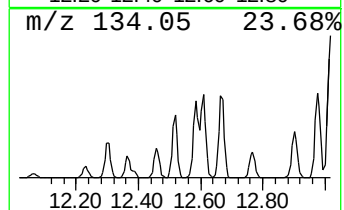
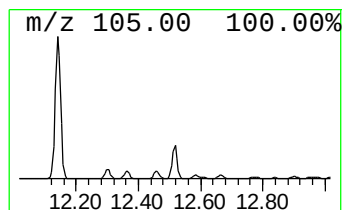
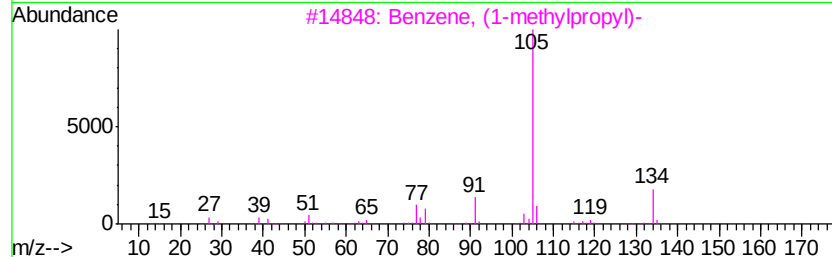
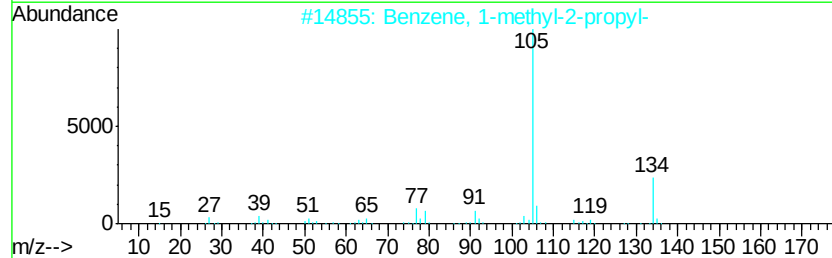
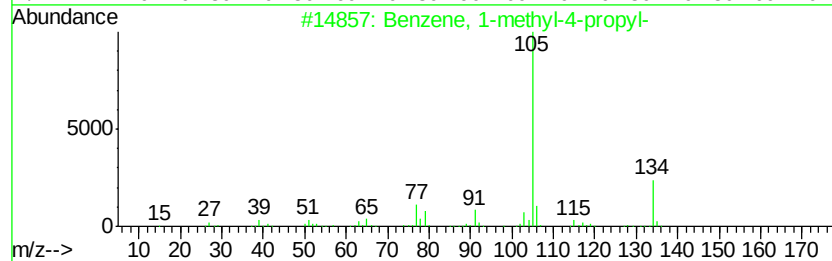
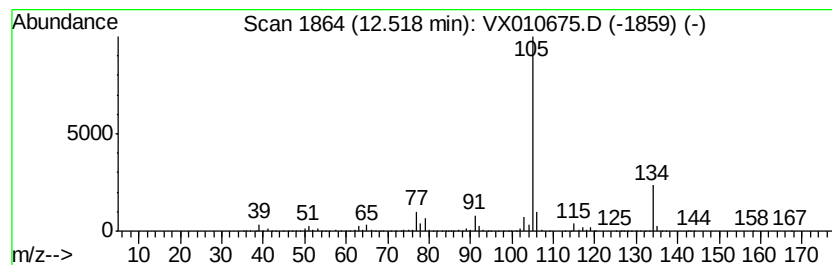
Instrument :
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-methyl-4-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	19.20 ug/l	384703	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1 94
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5 91
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8 91
4	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7 74
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8 72



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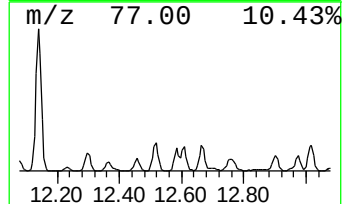
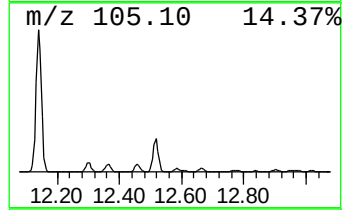
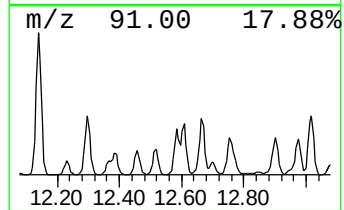
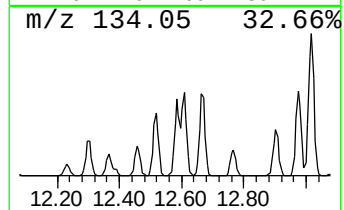
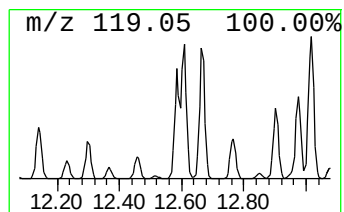
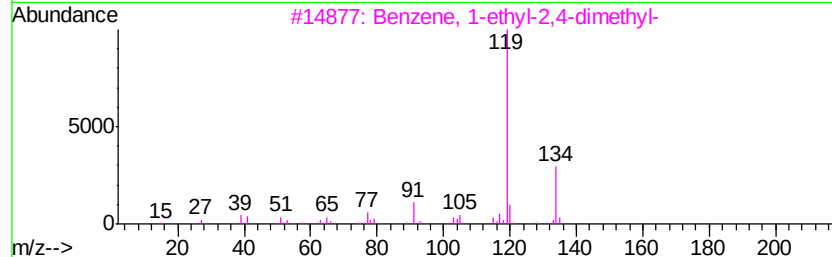
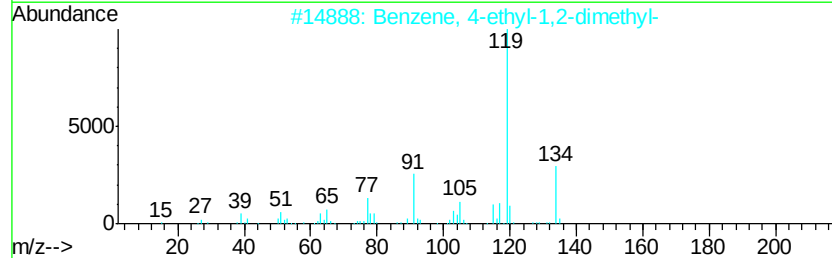
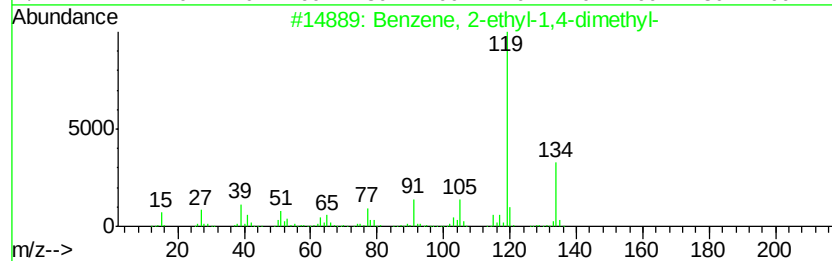
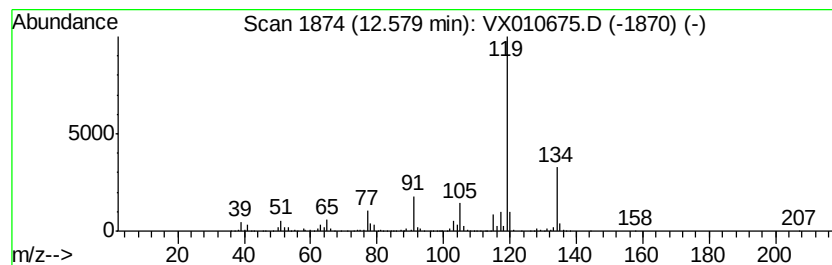
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	22.11 ug/l	442941	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010675.D
Acq On : 04 Jul 2019 08:15
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0580

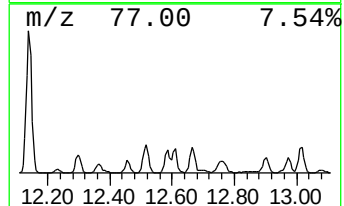
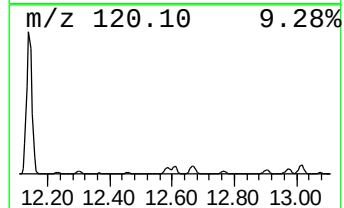
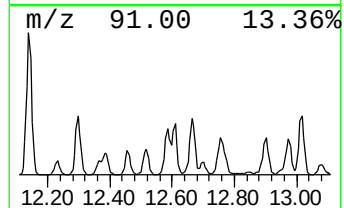
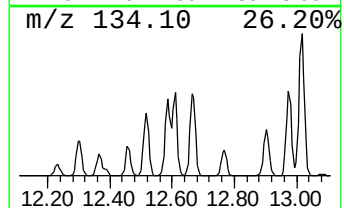
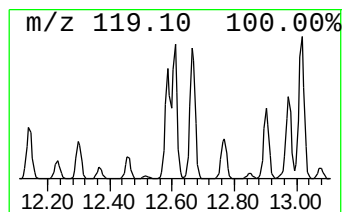
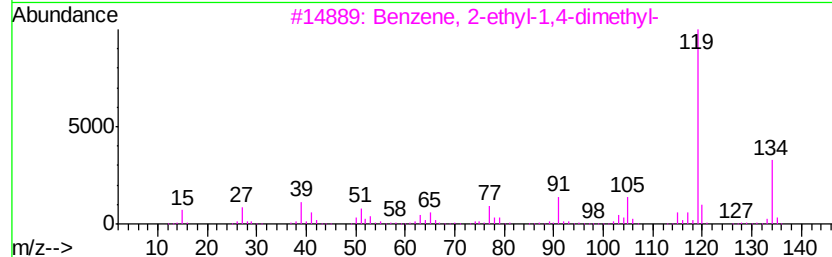
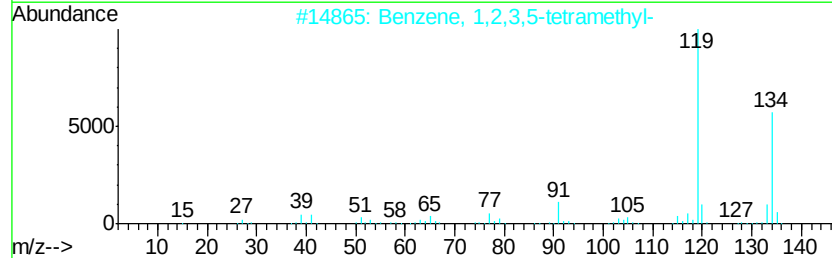
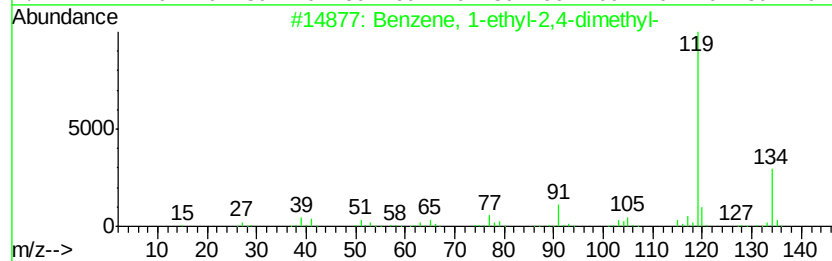
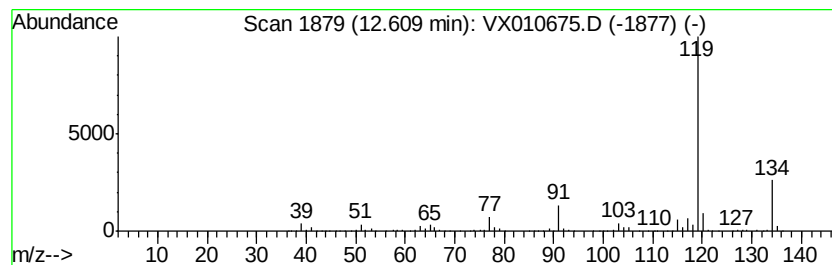
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	18.81 ug/l	376783	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010675.D
Acq On : 04 Jul 2019 08:15
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

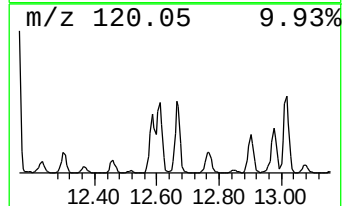
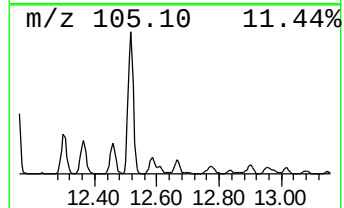
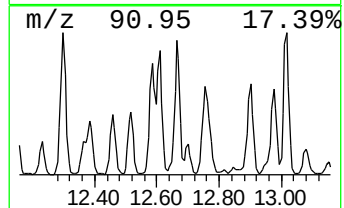
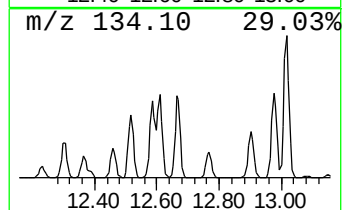
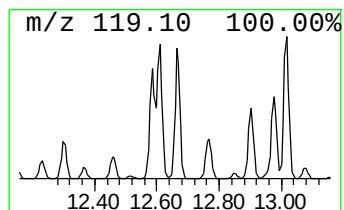
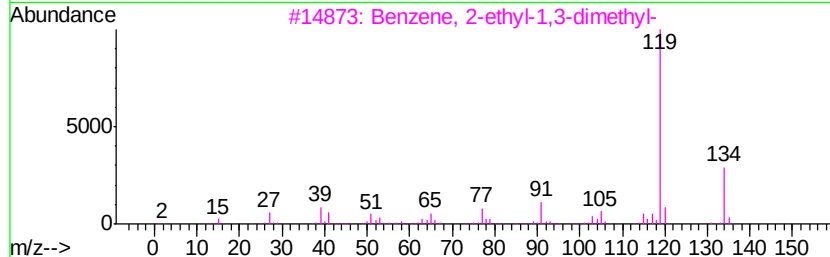
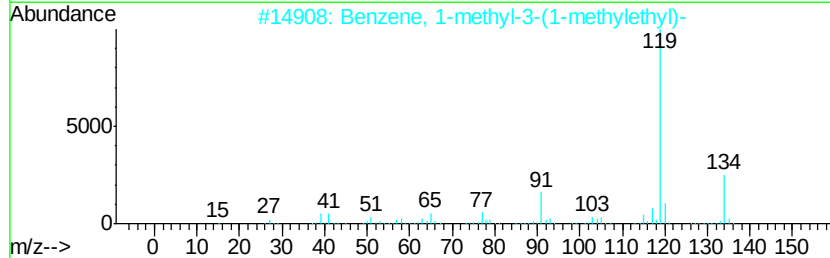
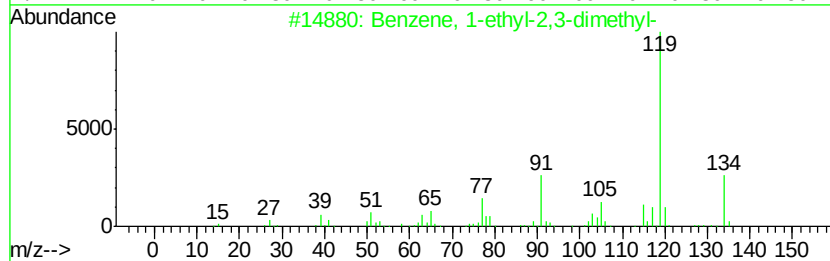
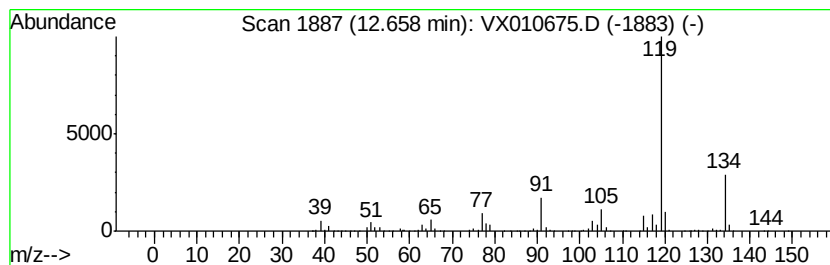
Instrument :
MSVOA_X
ClientSampleId :
FL-0580

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	24.29 ug/l	486727	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 96
2		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 95
3		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 95
4		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 95
5		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010675.D
Acq On : 04 Jul 2019 08:15
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0580

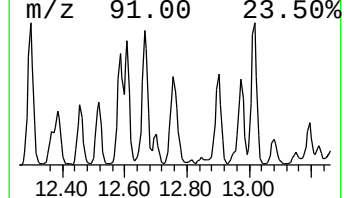
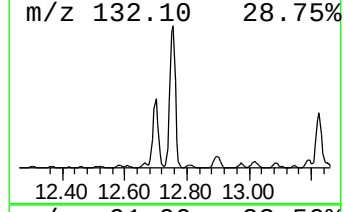
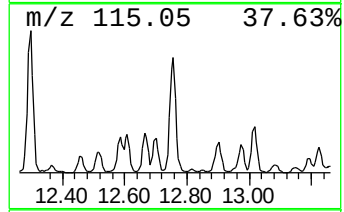
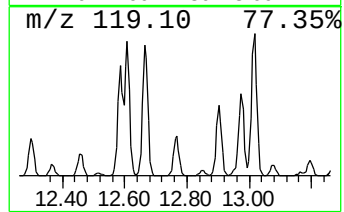
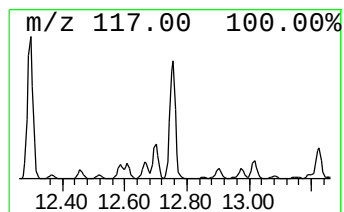
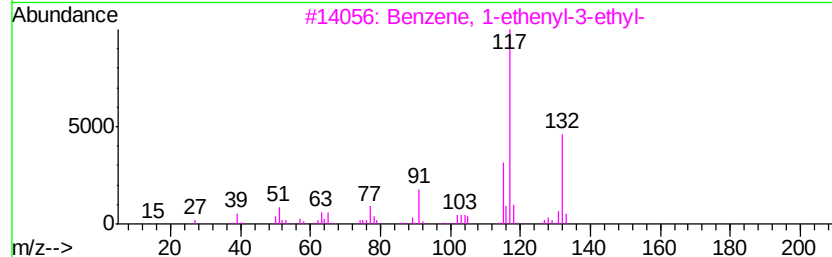
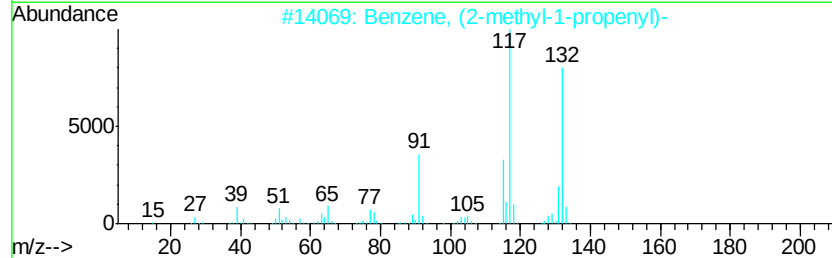
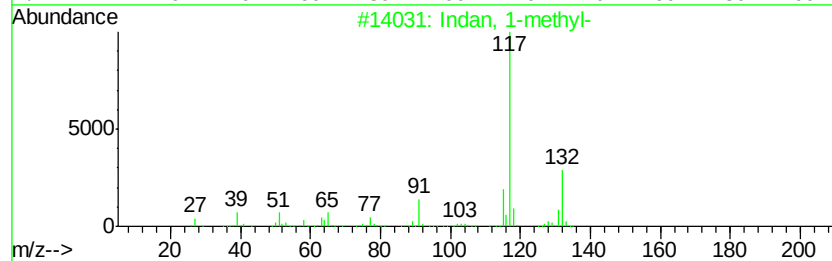
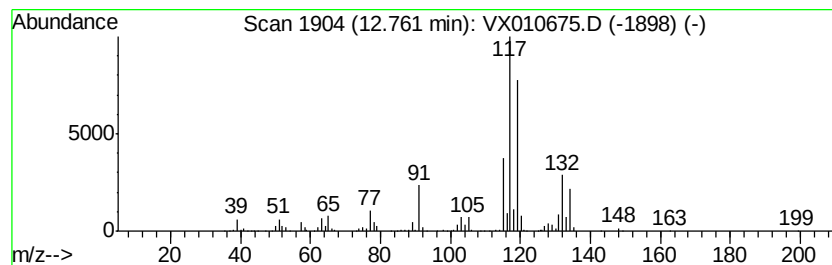
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	24.68 ug/l	494499	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	94
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4		Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010675.D
Acq On : 04 Jul 2019 08:15
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0580

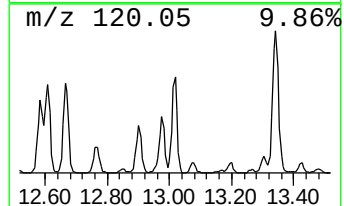
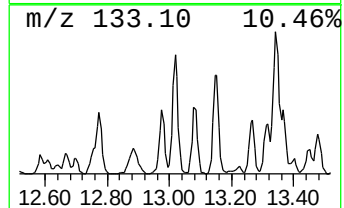
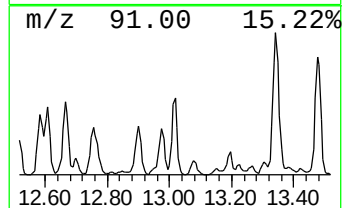
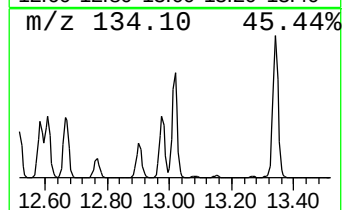
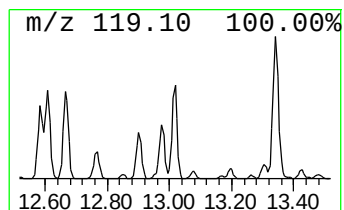
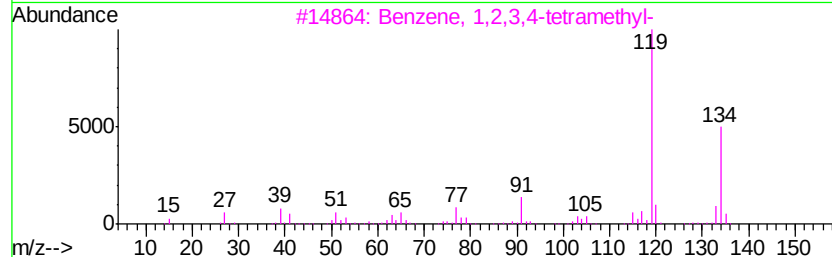
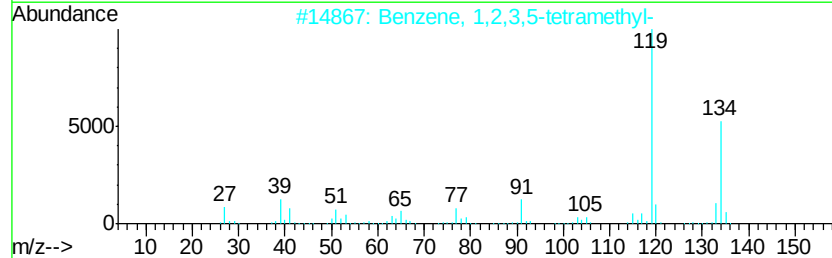
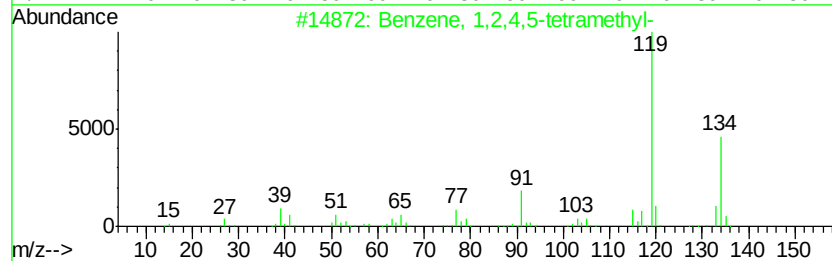
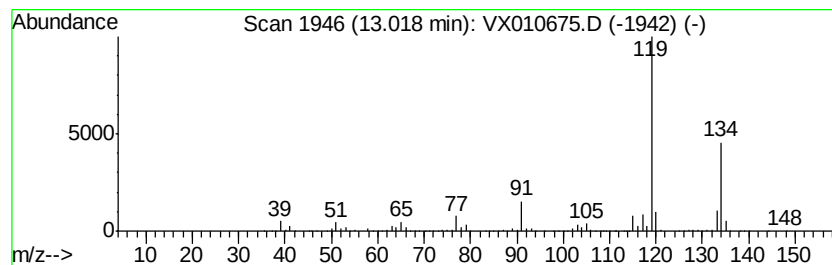
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	29.20 ug/l	585104	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX070319\
Data File : VX010675.D
Acq On : 04 Jul 2019 08:15
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
FL-0580

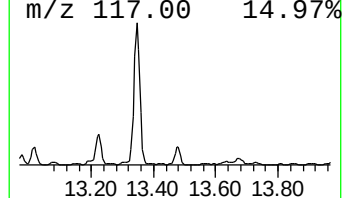
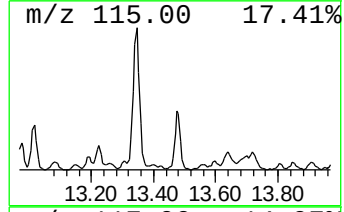
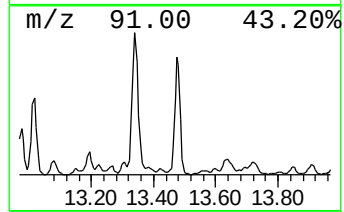
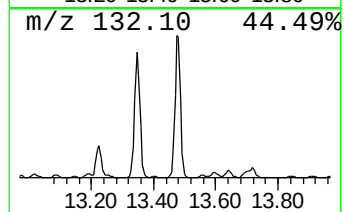
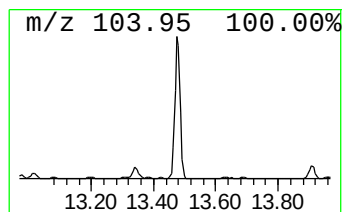
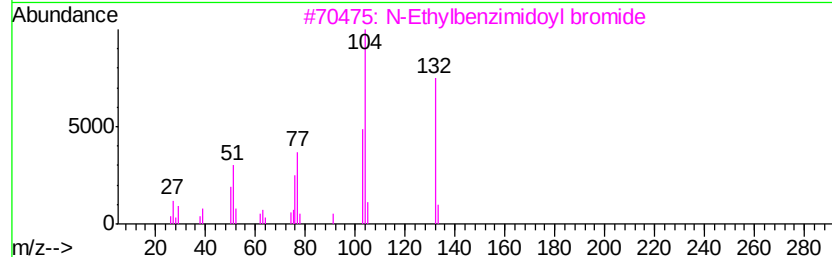
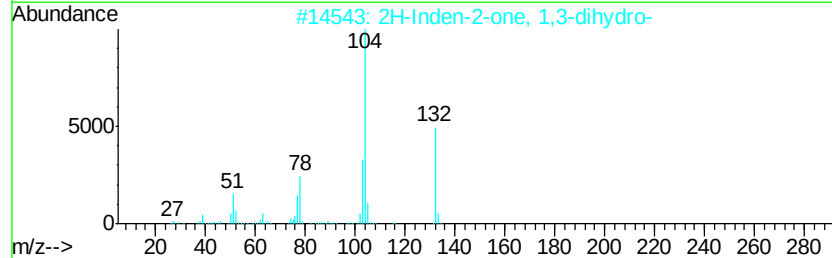
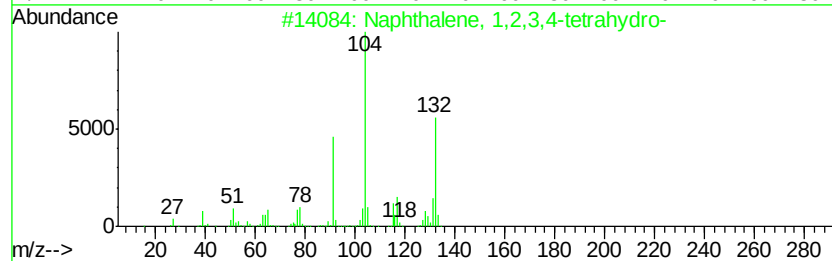
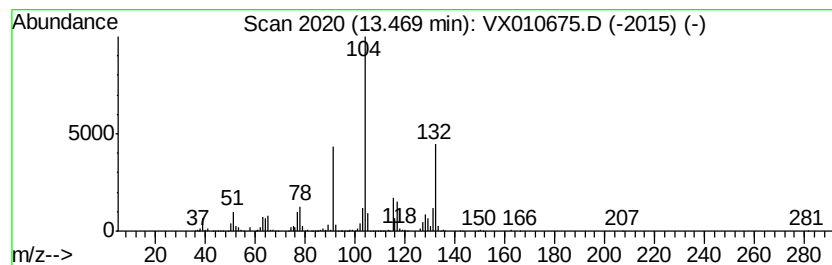
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.47	25.51 ug/l	511063	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
3		N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40
4		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	38
5		3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX070319\
Data File : VX010675.D
Acq On : 04 Jul 2019 08:15
Operator : JC/SP
Sample : K3664-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
FL-0580

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1-propenyl-	12.30	30.1	ug/l	604083	4	12.08	1001780	50.0
Benzene, 1-methyl...	12.52	19.2	ug/l	384703	4	12.08	1001780	50.0
Benzene, 2-ethyl-...	12.58	22.1	ug/l	442941	4	12.08	1001780	50.0
Benzene, 1-ethyl-...	12.61	18.8	ug/l	376783	4	12.08	1001780	50.0
Benzene, 1-ethyl-...	12.66	24.3	ug/l	486727	4	12.08	1001780	50.0
Indan, 1-methyl-	12.76	24.7	ug/l	494499	4	12.08	1001780	50.0
Benzene, 1,2,4,5-...	13.02	29.2	ug/l	585104	4	12.08	1001780	50.0
Naphthalene, 1,2,...	13.47	25.5	ug/l	511063	4	12.08	1001780	50.0