

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
 Data File : VX010391.D
 Acq On : 21 Jun 2019 17:45
 Operator : JC/SP
 Sample : K3469-20
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T11-MW-1-8.0-062019

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	1.282	16	21	27	rBV	32970	54889	2.26%	0.246%
2	2.611	233	239	247	rVB	13399	24443	1.01%	0.110%
3	3.038	301	309	317	rBV	15133	33924	1.40%	0.152%
4	3.190	324	334	345	rVB4	46749	126094	5.19%	0.566%
5	4.403	523	533	546	rVB3	18188	54732	2.25%	0.246%
6	5.500	701	713	722	rBV	161149	472407	19.44%	2.121%
7	5.580	722	726	731	rVV3	29382	77576	3.19%	0.348%
8	5.665	731	740	760	rVB	273201	803577	33.06%	3.607%
9	6.061	796	805	819	rBV	149657	394666	16.24%	1.772%
10	6.458	861	870	880	rVB7	9836	30959	1.27%	0.139%
11	6.860	925	936	946	rBV	408182	977787	40.23%	4.390%
12	7.256	993	1001	1008	rBV3	13038	29952	1.23%	0.134%
13	7.457	1025	1034	1043	rBV4	46084	118476	4.87%	0.532%
14	8.213	1145	1158	1163	rBV	37038	83313	3.43%	0.374%
15	8.573	1210	1217	1225	rVB	34719	64286	2.64%	0.289%
16	8.713	1234	1240	1249	rVB	852256	1324903	54.51%	5.948%
17	9.219	1318	1323	1333	rVB	30978	51476	2.12%	0.231%
18	10.109	1464	1469	1475	rBV	807645	1118379	46.01%	5.021%
19	10.359	1502	1510	1515	rVB2	20058	46609	1.92%	0.209%
20	11.134	1632	1637	1643	rBV	692046	896512	36.89%	4.025%
21	11.664	1721	1724	1732	rBV	23144	41011	1.69%	0.184%
22	11.768	1735	1741	1745	rBV	22027	32058	1.32%	0.144%
23	11.920	1761	1766	1768	rBV	52161	66729	2.75%	0.300%
24	11.944	1768	1770	1777	rVB	20715	25261	1.04%	0.113%
25	12.079	1786	1792	1798	rBV	877715	1101682	45.33%	4.946%
26	12.231	1813	1817	1821	rBV	60825	72297	2.97%	0.325%
27	12.298	1821	1828	1834	rVV	982582	1217522	50.09%	5.466%
28	12.365	1835	1839	1847	rVB2	65381	104506	4.30%	0.469%
29	12.457	1847	1854	1861	rBV	152134	189901	7.81%	0.853%
30	12.664	1880	1888	1892	rBV	983737	1254629	51.62%	5.632%
31	12.700	1892	1894	1898	rVV	267930	284209	11.69%	1.276%
32	12.755	1898	1903	1910	rVV2	819996	1177983	48.47%	5.288%
33	12.810	1910	1912	1918	rVB2	26363	33161	1.36%	0.149%
34	12.896	1918	1926	1931	rVB	95672	132136	5.44%	0.593%

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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

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 Title : SW846 8260

35	12.975	1931	1939	1944	rBV	798813	997104	41.02%	4.476%
36	13.024	1944	1947	1951	rVB	19006	24929	1.03%	0.112%
37	13.078	1951	1956	1963	rBV2	80953	126530	5.21%	0.568%
38	13.152	1963	1968	1971	rBV	47413	64951	2.67%	0.292%
39	13.194	1971	1975	1976	rVV2	101822	125732	5.17%	0.564%
40	13.225	1976	1980	1990	rVB	455048	611171	25.15%	2.744%
41	13.347	1990	2000	2006	rBV2	1653111	2430500	100.00%	10.911%
42	13.402	2006	2009	2010	rVV2	47696	53616	2.21%	0.241%
43	13.426	2010	2013	2018	rVV	63969	88502	3.64%	0.397%
44	13.481	2018	2022	2026	rVB	54096	76124	3.13%	0.342%
45	13.560	2028	2035	2038	rBV	72497	102264	4.21%	0.459%
46	13.597	2038	2041	2044	rVV	170483	228018	9.38%	1.024%
47	13.639	2044	2048	2052	rVV3	229745	343677	14.14%	1.543%
48	13.694	2052	2057	2059	rVV2	144286	252806	10.40%	1.135%
49	13.719	2059	2061	2070	rVB2	259091	398927	16.41%	1.791%
50	13.804	2070	2075	2078	rBV	40332	52349	2.15%	0.235%
51	13.853	2078	2083	2089	rBV	61152	80402	3.31%	0.361%
52	13.908	2089	2092	2097	rVB	102002	122330	5.03%	0.549%
53	13.981	2097	2104	2108	rBV2	185880	266469	10.96%	1.196%
54	14.023	2108	2111	2116	rVB	80767	98510	4.05%	0.442%
55	14.133	2123	2129	2132	rBV	129574	182967	7.53%	0.821%
56	14.157	2132	2133	2138	rVB2	42806	42103	1.73%	0.189%
57	14.273	2146	2152	2161	rBV3	204275	435050	17.90%	1.953%
58	14.377	2164	2169	2173	rVV	90904	124521	5.12%	0.559%
59	14.426	2173	2177	2182	rVB2	145493	200228	8.24%	0.899%
60	14.474	2182	2185	2190	rVB2	21836	30205	1.24%	0.136%
61	14.535	2190	2195	2200	rBV2	217925	295569	12.16%	1.327%
62	14.670	2211	2217	2219	rBV2	31216	54934	2.26%	0.247%
63	14.731	2223	2227	2231	rVB2	41080	54045	2.22%	0.243%
64	14.779	2231	2235	2240	rVB4	21141	32305	1.33%	0.145%
65	14.846	2240	2246	2254	rBV2	61860	135945	5.59%	0.610%
66	14.962	2259	2265	2272	rBV7	11181	27000	1.11%	0.121%
67	15.249	2306	2312	2319	rVB2	38701	65231	2.68%	0.293%
68	15.444	2332	2344	2346	rBV	51379	87330	3.59%	0.392%
69	15.474	2346	2349	2354	rVB	59497	79521	3.27%	0.357%
70	15.547	2354	2361	2366	rBV	184791	296011	12.18%	1.329%
71	15.682	2375	2383	2390	rBV	372253	582540	23.97%	2.615%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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 >
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72	15.804	2397	2403	2410	rBV2	18218	37444	1.54%	0.168%
73	15.913	2410	2421	2432	rVB2	102847	273962	11.27%	1.230%
74	16.066	2439	2446	2451	rBV	60328	103726	4.27%	0.466%
75	16.413	2496	2503	2511	rVB2	34768	71676	2.95%	0.322%

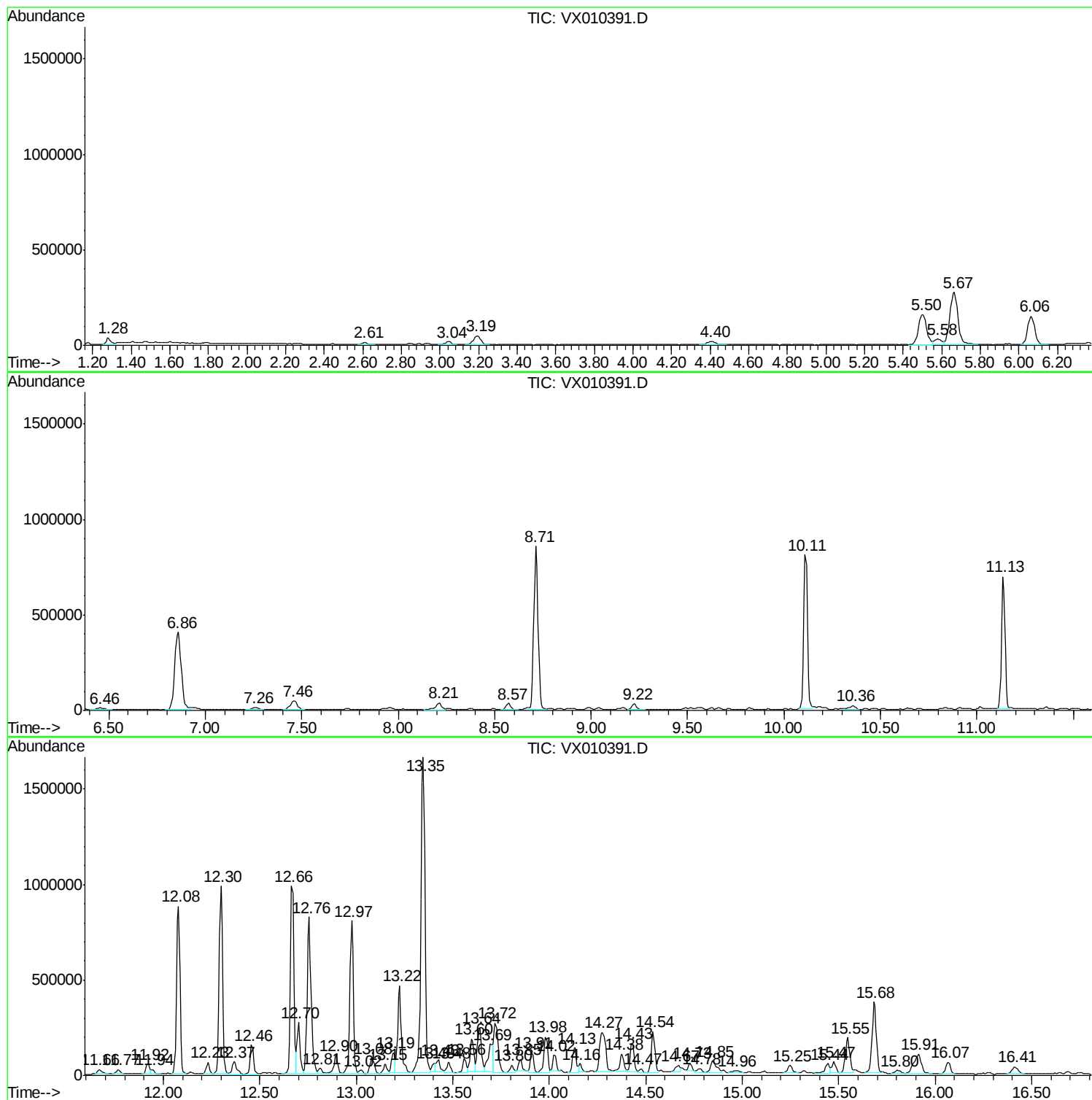
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Instrument :
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ClientSampleId :
T11-MW-1-8.0-062019

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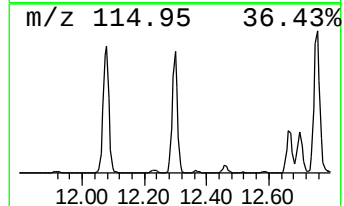
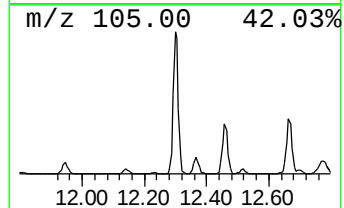
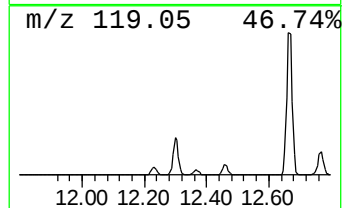
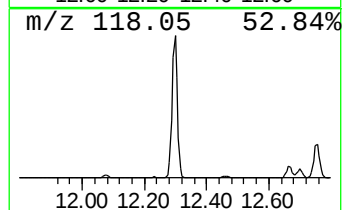
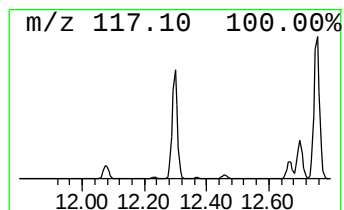
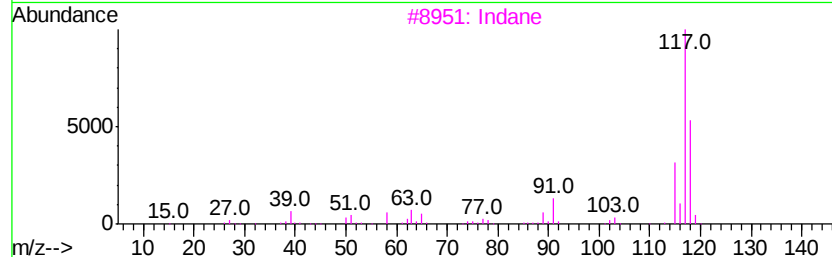
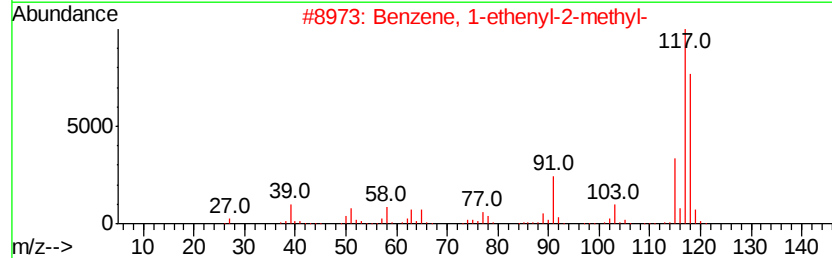
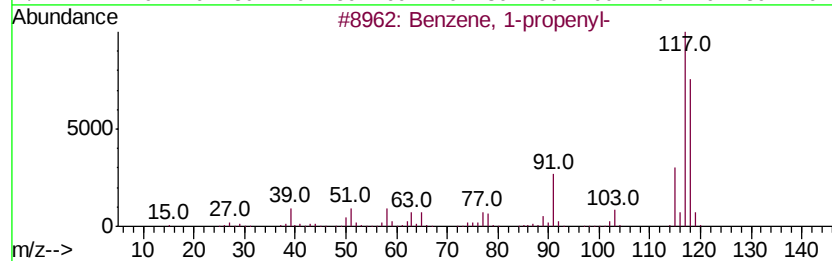
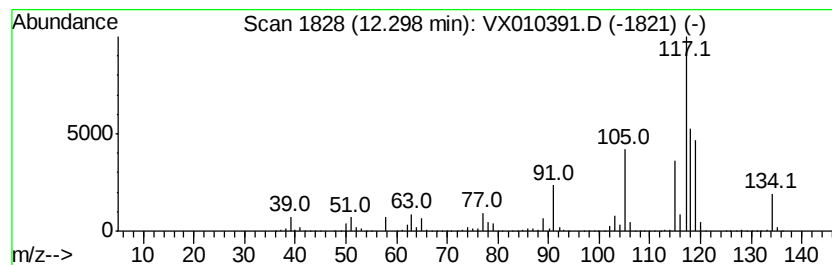
Instrument :
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ClientSampleId :
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	55.26 ug/l	1217520	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, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 55
3		Indane	118 C9H10	000496-11-7 55
4		Benzene, cyclopropyl-	118 C9H10	000873-49-4 55
5		Benzene, 2-propenyl-	118 C9H10	000300-57-2 55



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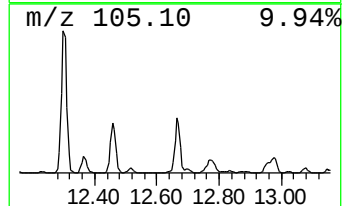
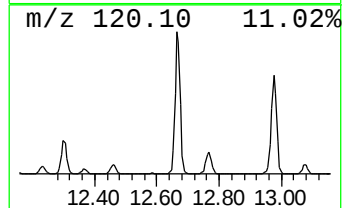
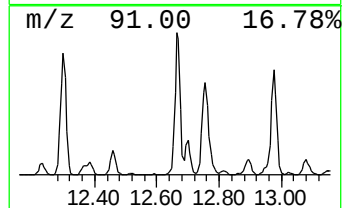
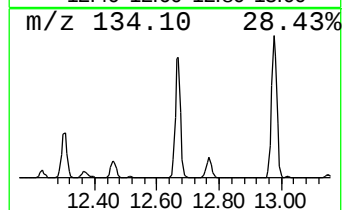
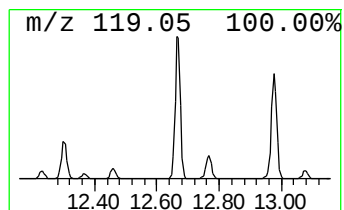
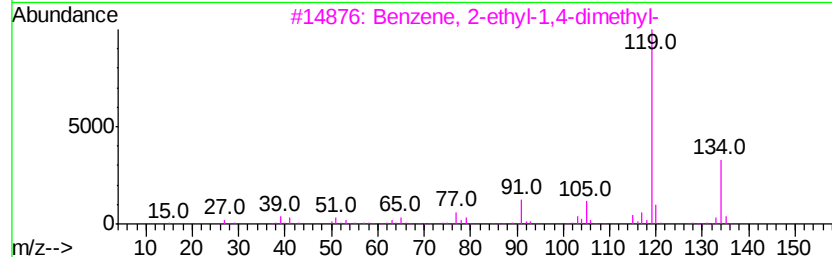
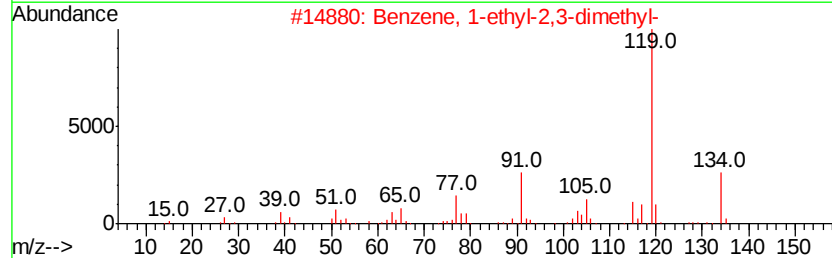
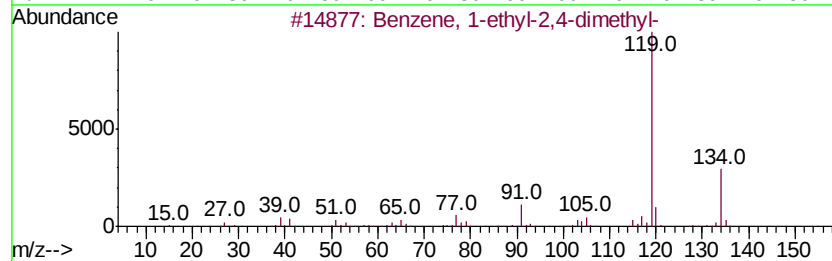
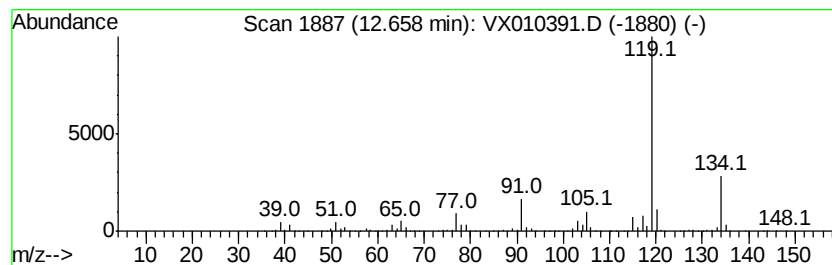
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.66	56.94 ug/l	1254630	1,4-Dichlorobenzene-d4	12.08	
Hit# of	5	Tentative ID	MW MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



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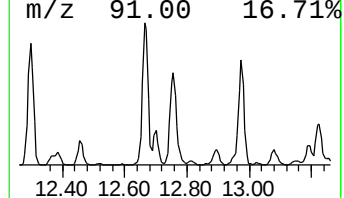
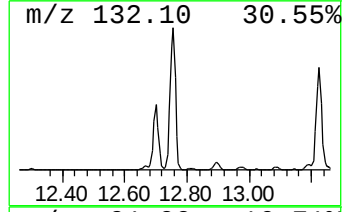
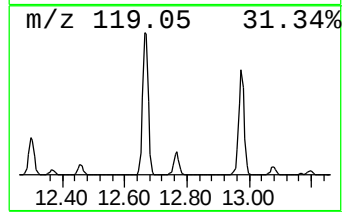
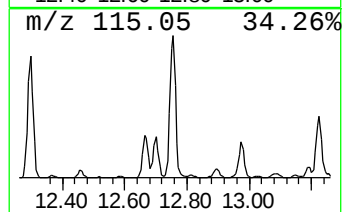
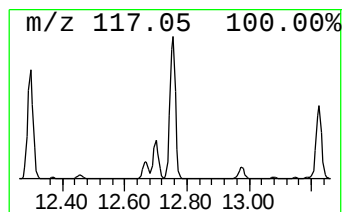
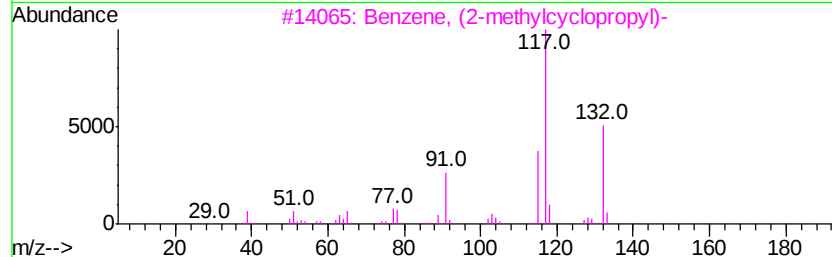
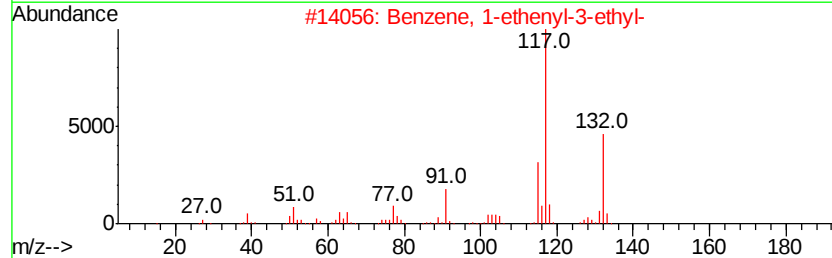
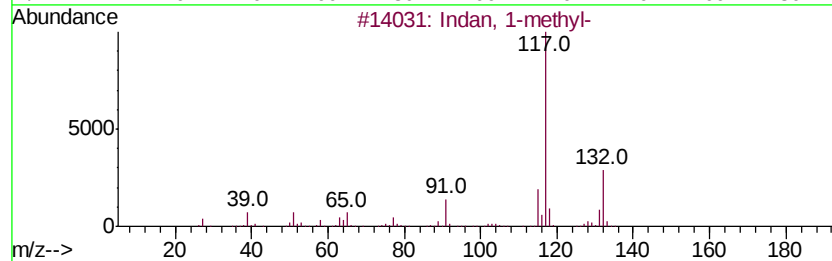
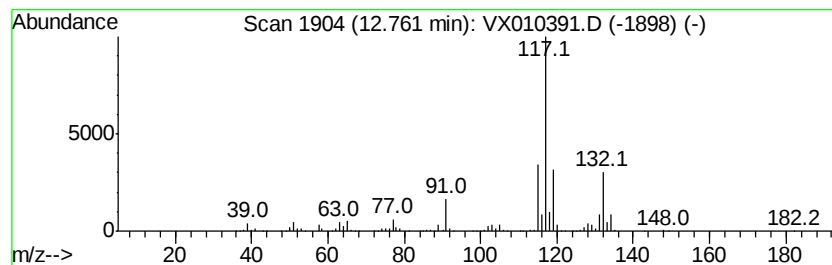
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Quant Title : SW846 8260

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

Peak Number 3 Indan, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	53.46 ug/l	1177980	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indan, 1-methyl-	132 C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4	90
3	Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0	81
4	Benzene, (2-methyl-2-propenyl)-	132 C10H12	003290-53-7	80
5	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3	74



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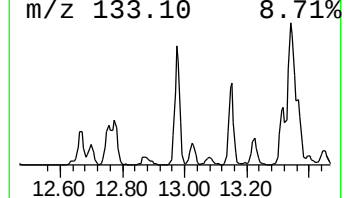
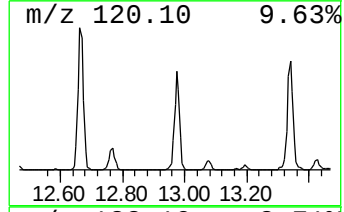
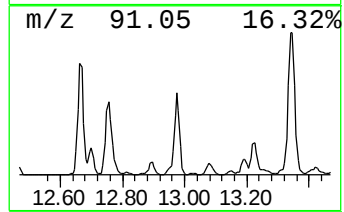
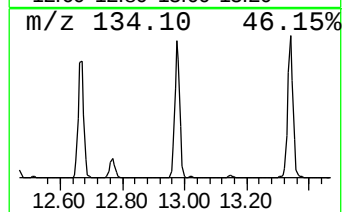
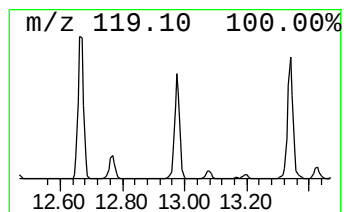
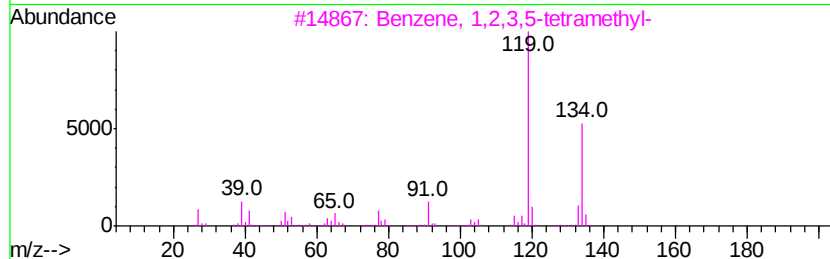
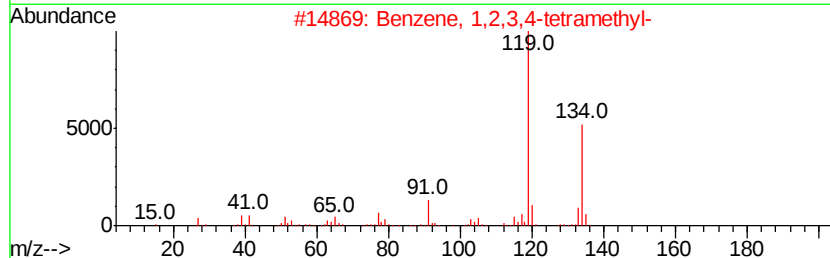
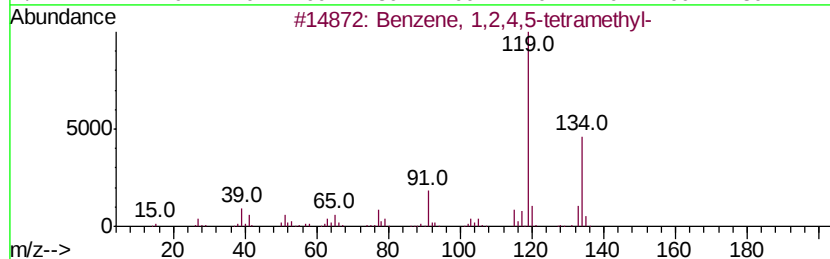
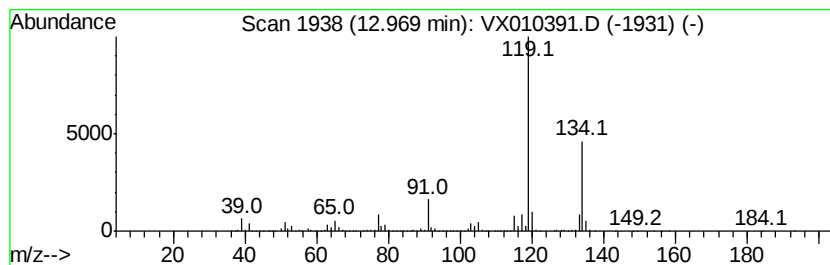
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	45.25 ug/l	997104	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, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010391.D
Acq On : 21 Jun 2019 17:45
Operator : JC/SP
Sample : K3469-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-062019

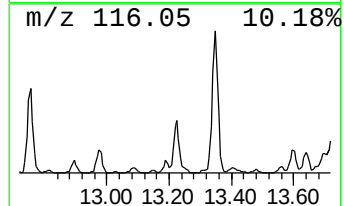
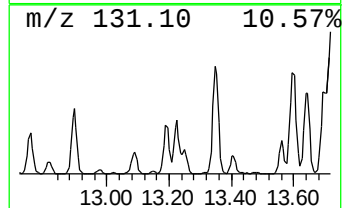
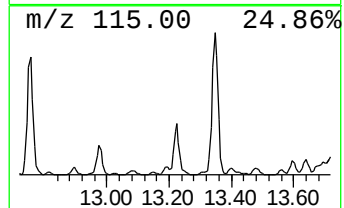
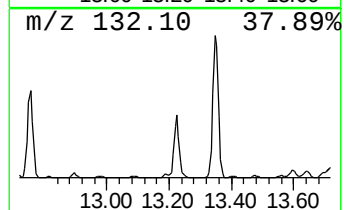
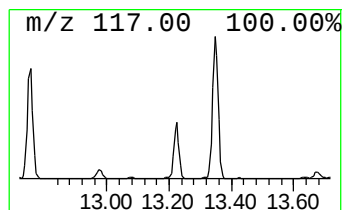
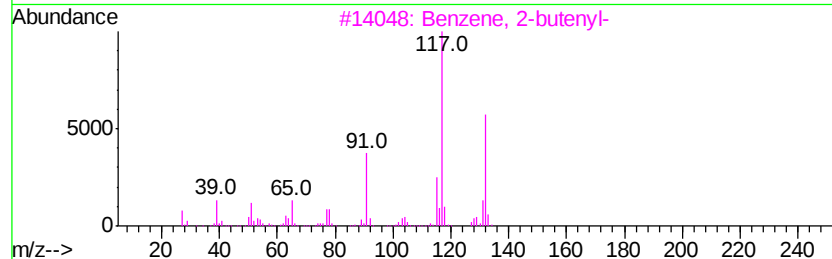
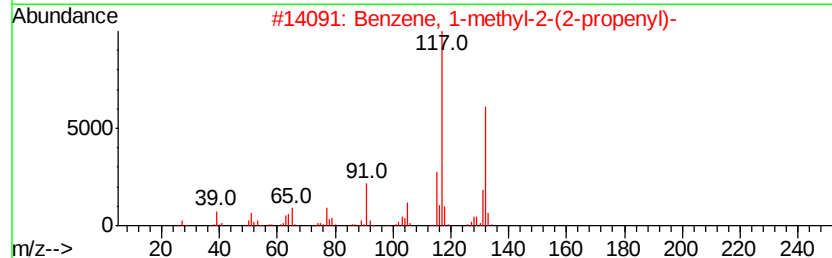
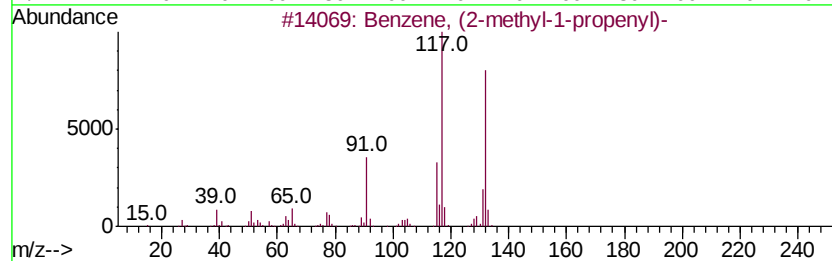
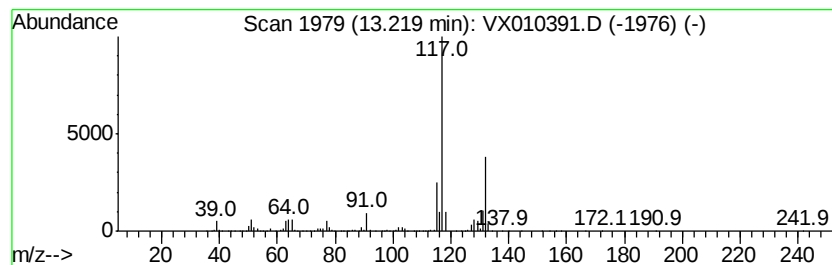
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-propen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	27.74 ug/l	611171	1,4-Dichlorobenzene-d4	12.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
4		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	91
5		Indan, 1-methyl-	132	C10H12	000767-58-8	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010391.D
Acq On : 21 Jun 2019 17:45
Operator : JC/SP
Sample : K3469-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-062019

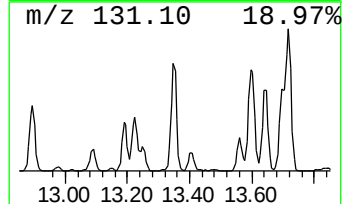
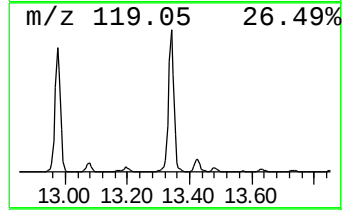
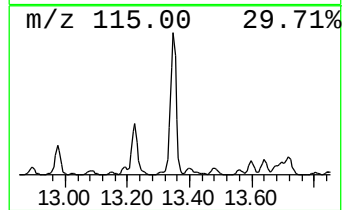
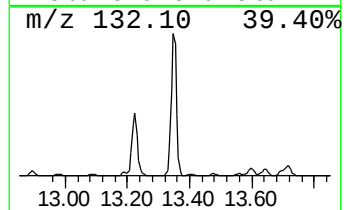
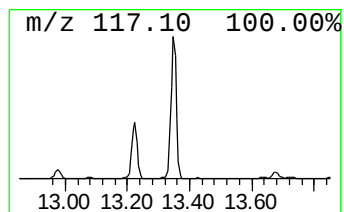
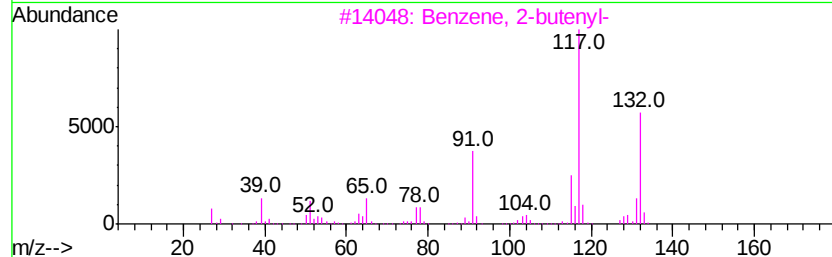
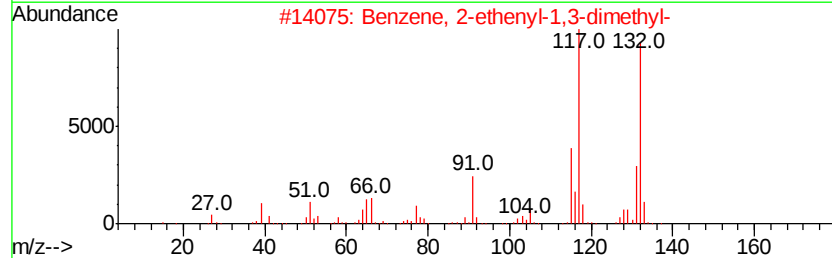
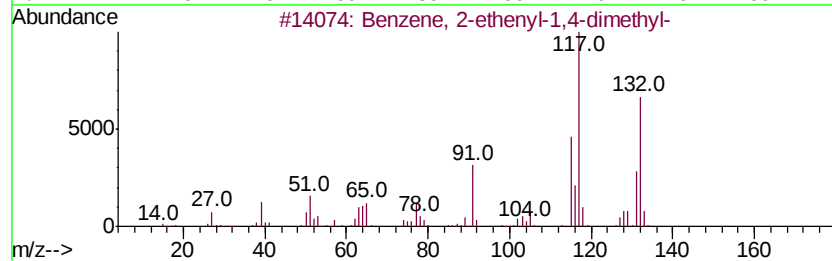
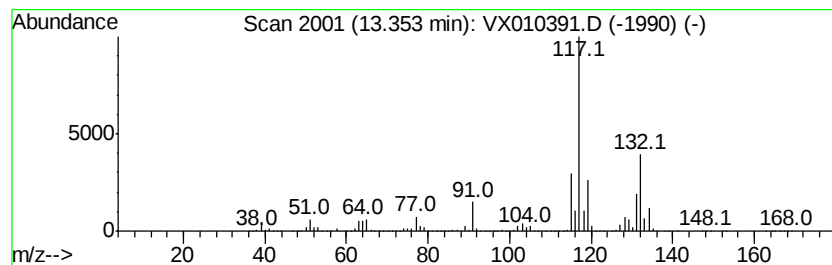
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	110.31 ug/l	2430500	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	93
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010391.D
Acq On : 21 Jun 2019 17:45
Operator : JC/SP
Sample : K3469-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

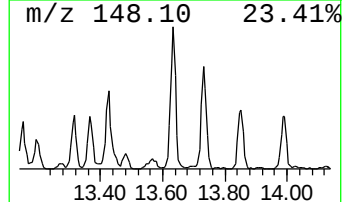
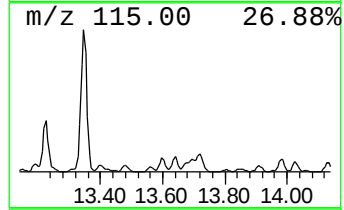
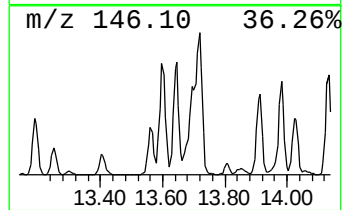
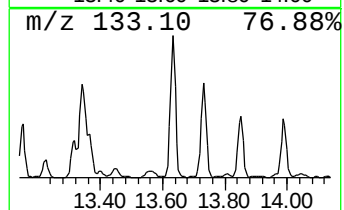
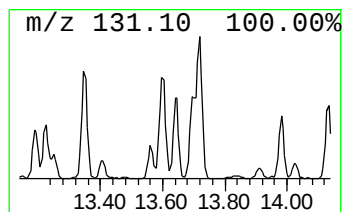
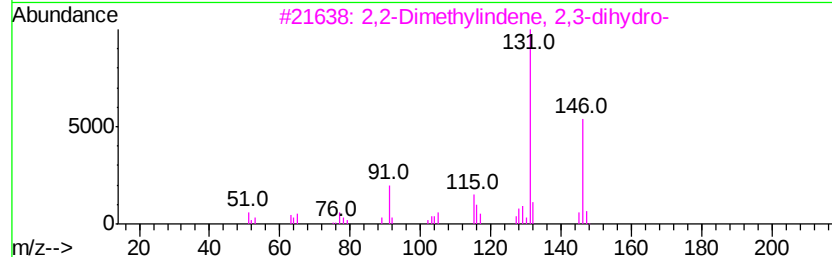
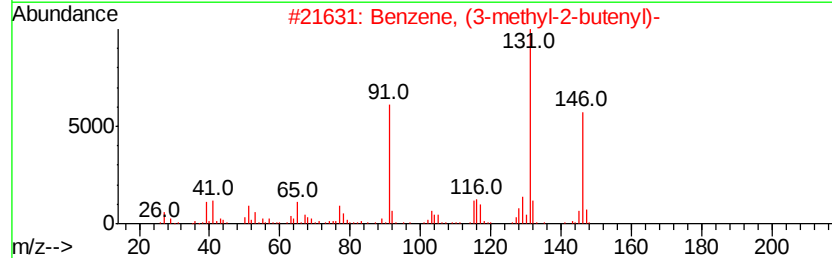
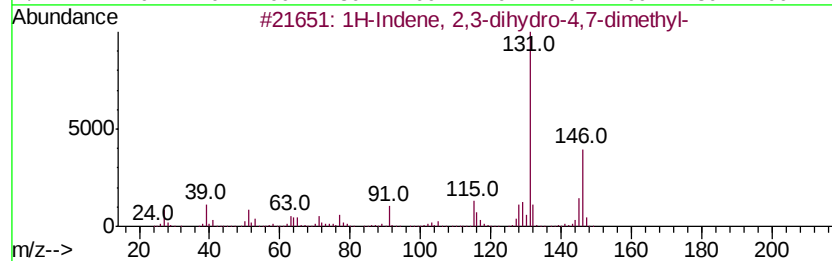
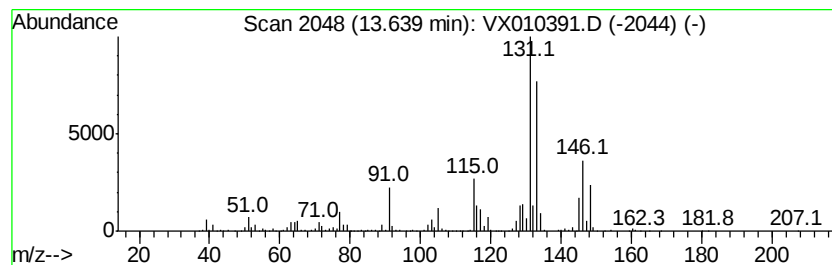
Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-062019

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

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

Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	15.60 ug/l	343677	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 90
2		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 78
3		2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7 78
4		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 78
5		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 60



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010391.D
Acq On : 21 Jun 2019 17:45
Operator : JC/SP
Sample : K3469-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-062019

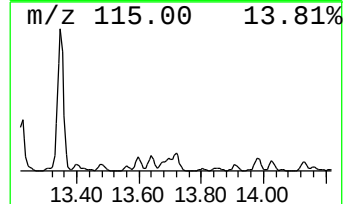
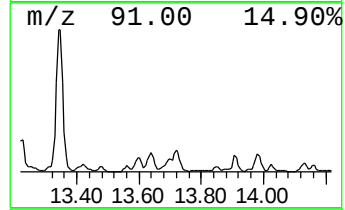
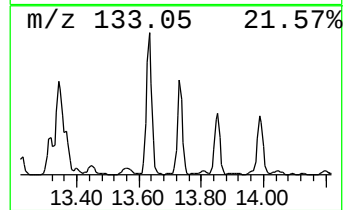
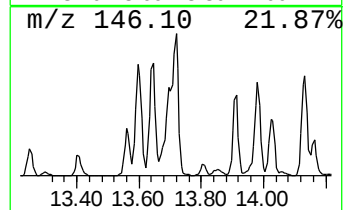
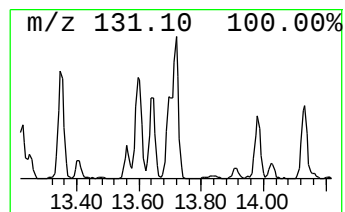
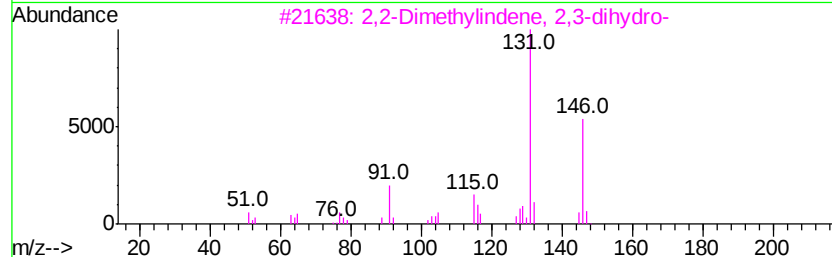
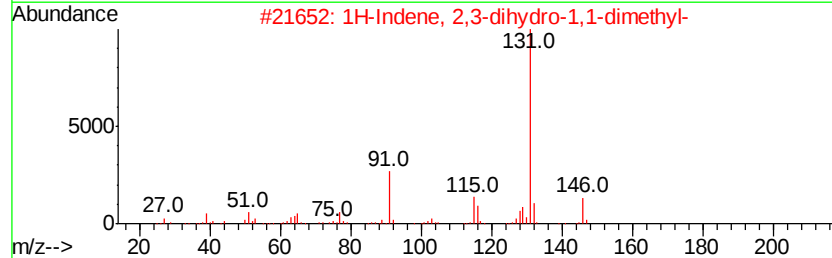
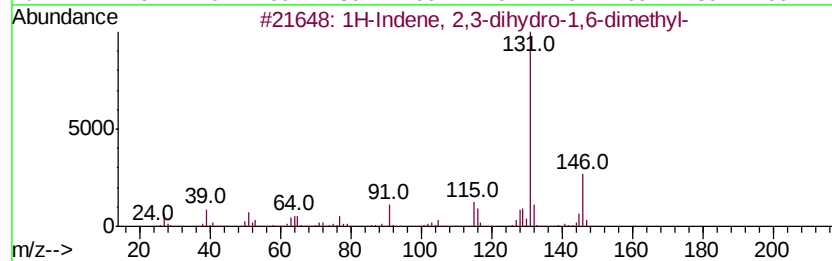
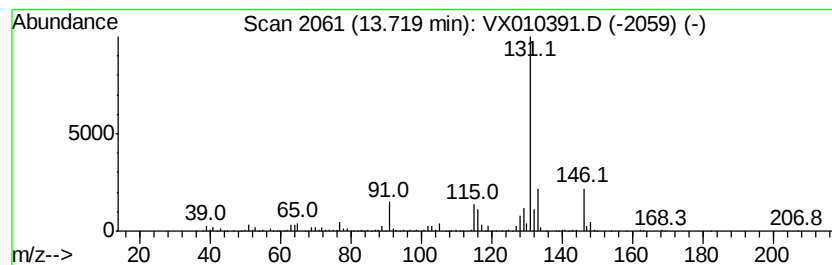
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	18.11 ug/l	398927	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	83
2		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	81
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81
4		Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	81
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	74



Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010391.D
Acq On : 21 Jun 2019 17:45
Operator : JC/SP
Sample : K3469-20
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 19 Sample Multiplier: 1

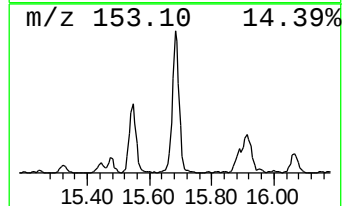
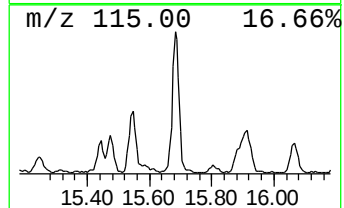
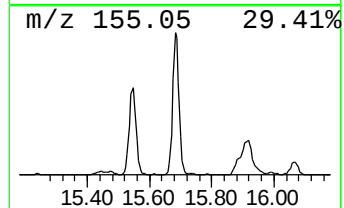
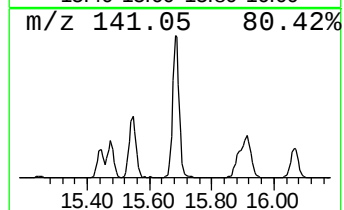
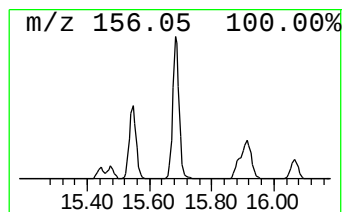
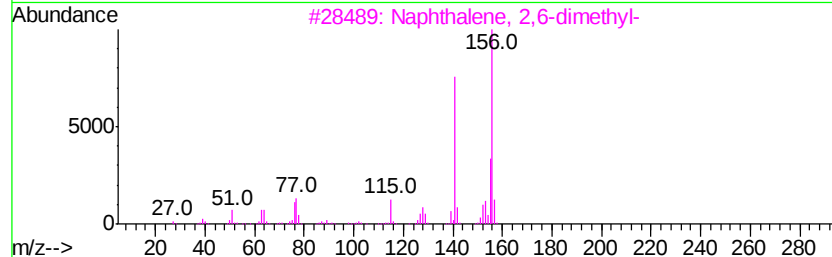
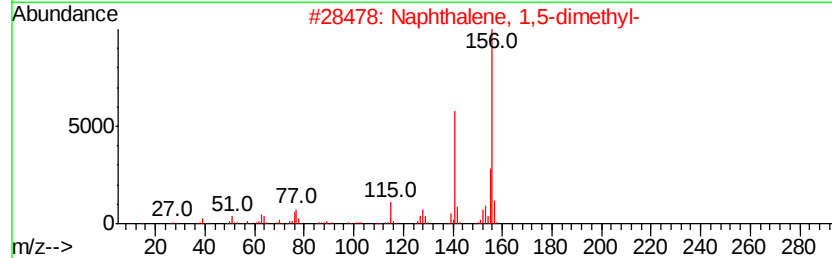
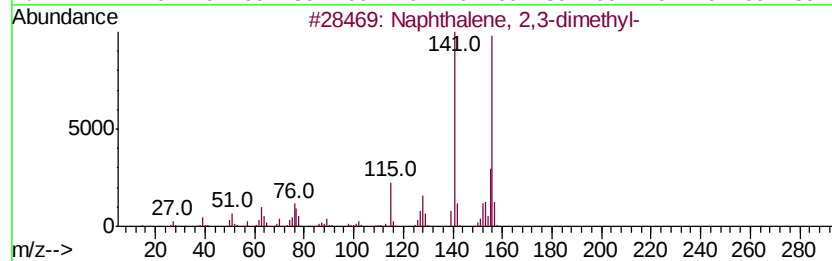
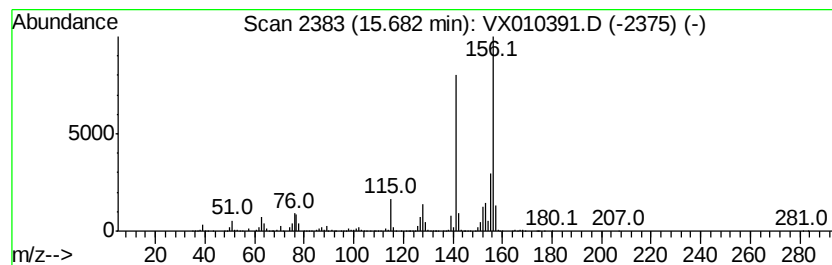
Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-062019

Quant Method : Z:\VOASRV\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, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	26.44 ug/l	582540	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,3-dimethyl-	156 C12H12	000581-40-8 98
2		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 97
3		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 97
4		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 97
5		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 97



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010391.D
Acq On : 21 Jun 2019 17:45
Operator : JC/SP
Sample : K3469-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T11-MW-1-8.0-062019

Quant Method : Z:\VOASRV\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
Benzene, 1-propenyl-	12.30	55.3	ug/l	1217520	4	12.08	1101680	50.0
Benzene, 1-ethyl-...	12.66	56.9	ug/l	1254630	4	12.08	1101680	50.0
Indan, 1-methyl-	12.76	53.5	ug/l	1177980	4	12.08	1101680	50.0
Benzene, 1,2,4,5-...	12.97	45.3	ug/l	997104	4	12.08	1101680	50.0
Benzene, (2-methy...	13.22	27.7	ug/l	611171	4	12.08	1101680	50.0
Benzene, 2-etheny...	13.35	110.3	ug/l	2430500	4	12.08	1101680	50.0
1H-Indene, 2,3-di...	13.64	15.6	ug/l	343677	4	12.08	1101680	50.0
1H-Indene, 2,3-di...	13.72	18.1	ug/l	398927	4	12.08	1101680	50.0
Naphthalene, 2,3-...	15.68	26.4	ug/l	582540	4	12.08	1101680	50.0