

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY030520\
 Data File : VY001875.D
 Acq On : 05 Mar 2020 12:44
 Operator : SY/MD
 Sample : L1761-01
 Misc : 7.15G/5ML/MSVOA Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SU-04-030420

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_Y\METHODS\82Y030420S.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	3.962	356	365	375	rBV	5856	14900	1.35%	0.084%
2	4.724	482	490	509	rVB4	7589	26938	2.44%	0.153%
3	7.730	974	983	989	rBV	114280	269995	24.45%	1.529%
4	7.803	989	995	1005	rVB	192646	429678	38.91%	2.433%
5	8.151	1039	1052	1063	rBV	125667	273761	24.79%	1.550%
6	8.699	1132	1142	1153	rBV	322562	641237	58.07%	3.631%
7	10.181	1377	1385	1393	rBV	519120	897432	81.27%	5.082%
8	11.492	1590	1600	1610	rVB	478774	758814	68.72%	4.297%
9	12.485	1756	1763	1770	rBV	338774	528243	47.84%	2.991%
10	12.674	1786	1794	1798	rVB	6488	13739	1.24%	0.078%
11	12.735	1798	1804	1807	rBV	37009	59852	5.42%	0.339%
12	12.766	1807	1809	1812	rVV	24080	35141	3.18%	0.199%
13	12.814	1812	1817	1822	rVV	55463	93448	8.46%	0.529%
14	12.973	1836	1843	1850	rBV2	46881	82738	7.49%	0.468%
15	13.040	1850	1854	1860	rVB3	18566	30760	2.79%	0.174%
16	13.119	1860	1867	1873	rBV	113034	177207	16.05%	1.003%
17	13.247	1882	1888	1893	rVB5	16516	29691	2.69%	0.168%
18	13.314	1893	1899	1904	rBV2	53317	97171	8.80%	0.550%
19	13.369	1904	1908	1911	rBV2	18954	25874	2.34%	0.147%
20	13.424	1911	1917	1921	rBV	421853	699092	63.31%	3.959%
21	13.558	1933	1939	1941	rVV5	7142	14562	1.32%	0.082%
22	13.595	1941	1945	1946	rVV2	50686	59652	5.40%	0.338%
23	13.625	1946	1950	1954	rVV2	247318	391287	35.43%	2.216%
24	13.668	1954	1957	1966	rVB3	190768	350281	31.72%	1.983%
25	13.753	1966	1971	1976	rBV3	98178	159135	14.41%	0.901%
26	13.826	1976	1983	1989	rBV	91781	153816	13.93%	0.871%
27	13.887	1989	1993	1995	rBV	95148	130021	11.77%	0.736%
28	13.918	1995	1998	2003	rVB	120729	164430	14.89%	0.931%
29	13.973	2003	2007	2012	rVB	198433	285140	25.82%	1.615%
30	14.034	2012	2017	2020	rBV	46006	66599	6.03%	0.377%
31	14.083	2020	2025	2030	rVB	247775	381604	34.56%	2.161%
32	14.143	2030	2035	2036	rBV3	28209	45180	4.09%	0.256%
33	14.217	2042	2047	2050	rBV2	64065	101281	9.17%	0.573%
34	14.259	2050	2054	2057	rVV3	67764	107108	9.70%	0.606%

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35	14.302	2057	2061	2064	rVV	115537	182814	16.55%	1.035%
36	14.339	2064	2067	2071	rVB	186665	240284	21.76%	1.361%
37	14.387	2071	2075	2078	rBV	107961	147755	13.38%	0.837%
38	14.424	2078	2081	2085	rVV2	92879	120097	10.88%	0.680%
39	14.485	2087	2091	2093	rVV2	33663	45140	4.09%	0.256%
40	14.528	2093	2098	2101	rVV2	124458	189355	17.15%	1.072%
41	14.570	2101	2105	2110	rVV2	243685	439162	39.77%	2.487%
42	14.619	2110	2113	2117	rVB	151620	206893	18.74%	1.172%
43	14.686	2117	2124	2130	rBV3	509040	1104284	100.00%	6.253%
44	14.741	2130	2133	2139	rVV	98618	157920	14.30%	0.894%
45	14.790	2139	2141	2144	rVV4	11239	16679	1.51%	0.094%
46	14.838	2144	2149	2157	rVV	423862	655314	59.34%	3.711%
47	14.912	2157	2161	2162	rVV2	68853	88337	8.00%	0.500%
48	14.948	2162	2167	2170	rVV3	209866	406703	36.83%	2.303%
49	14.985	2170	2173	2177	rVV	142602	241543	21.87%	1.368%
50	15.064	2179	2186	2192	rVB3	218329	495768	44.89%	2.807%
51	15.216	2204	2211	2213	rBV4	60338	121906	11.04%	0.690%
52	15.296	2219	2224	2229	rBV	243867	371786	33.67%	2.105%
53	15.357	2229	2234	2240	rBV4	151130	421037	38.13%	2.384%
54	15.442	2245	2248	2256	rVB4	94749	161562	14.63%	0.915%
55	15.533	2256	2263	2270	rBV3	195307	417116	37.77%	2.362%
56	15.607	2271	2275	2280	rVB	34442	59032	5.35%	0.334%
57	15.698	2280	2290	2291	rBV3	149750	300735	27.23%	1.703%
58	15.735	2291	2296	2304	rVB	521393	943434	85.43%	5.342%
59	15.820	2306	2310	2316	rBV3	58059	105700	9.57%	0.599%
60	15.899	2316	2323	2329	rBV2	117721	269512	24.41%	1.526%
61	16.046	2339	2347	2353	rBV	285544	598374	54.19%	3.388%
62	16.241	2369	2379	2384	rBV3	271728	609856	55.23%	3.453%
63	16.302	2384	2389	2395	rVV2	122300	218797	19.81%	1.239%
64	16.503	2415	2422	2427	rBV6	100189	270928	24.53%	1.534%
65	16.558	2428	2431	2437	rVV3	63849	115351	10.45%	0.653%
66	16.637	2438	2444	2453	rVV2	84344	210275	19.04%	1.191%
67	16.710	2454	2456	2461	rVB5	12278	19426	1.76%	0.110%
68	16.838	2470	2477	2488	rVB6	31856	111530	10.10%	0.632%

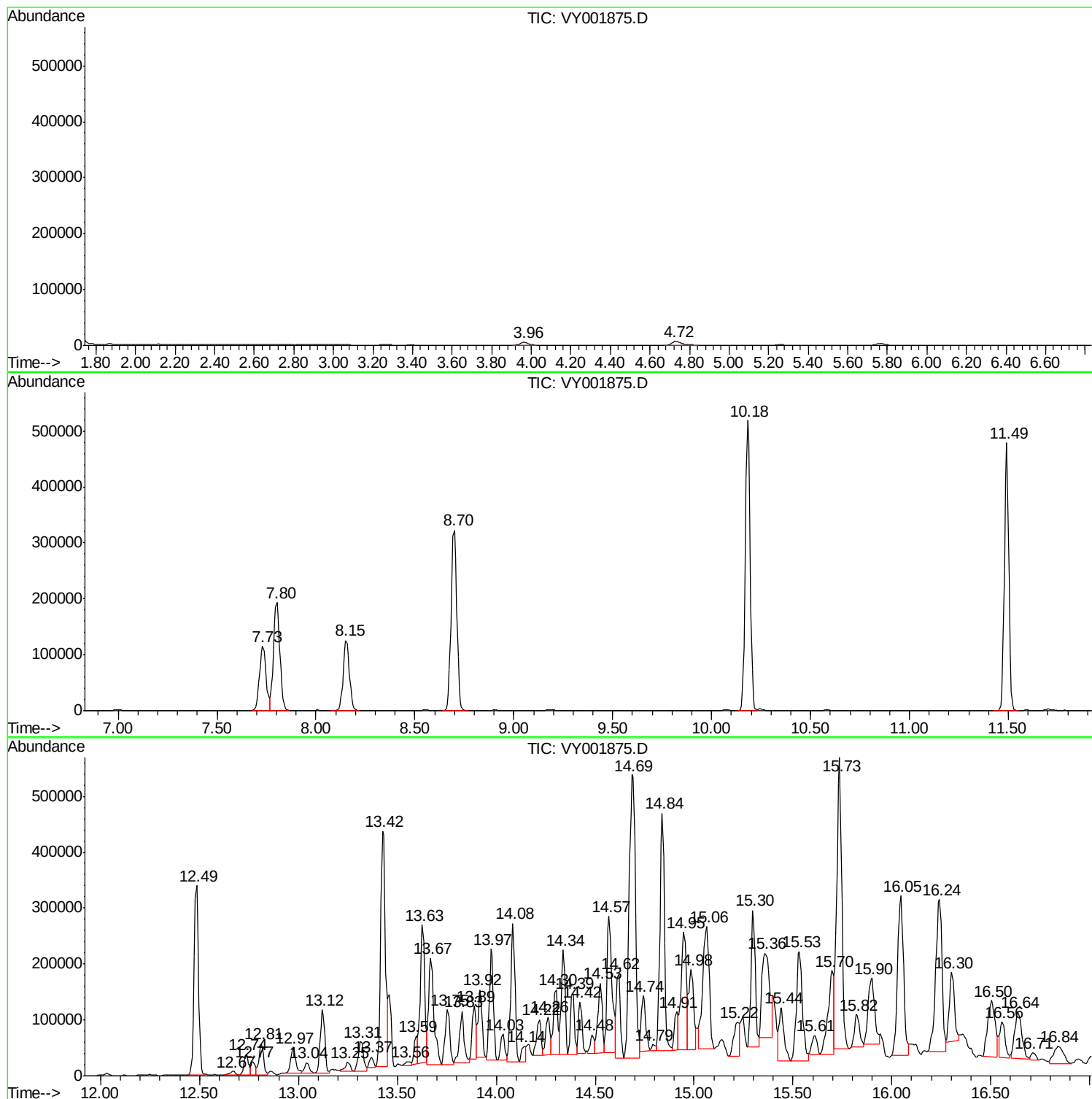
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.M
Quant Title : SW846 8260

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



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ClientSampleId :
SU-04-030420

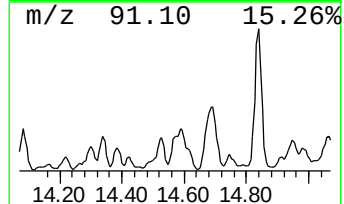
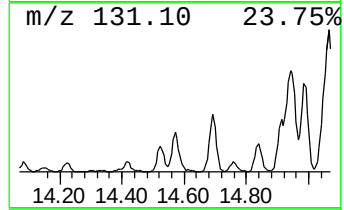
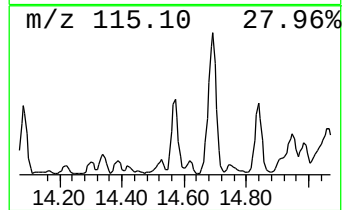
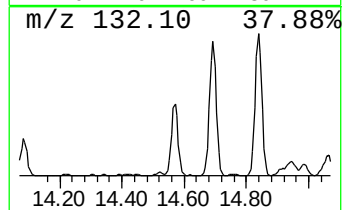
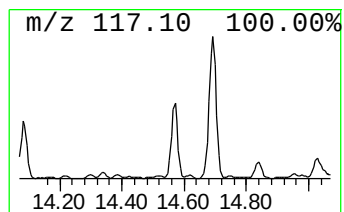
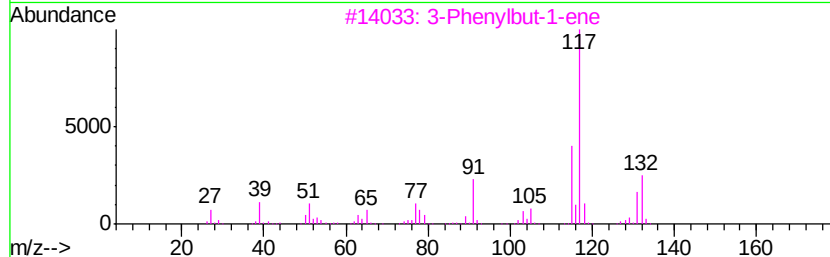
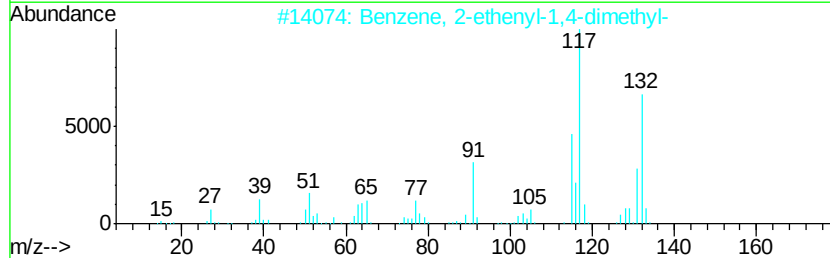
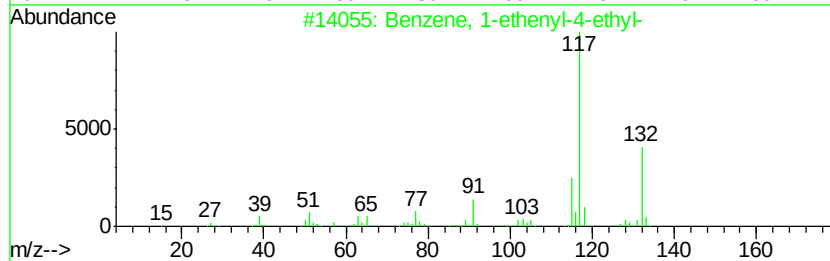
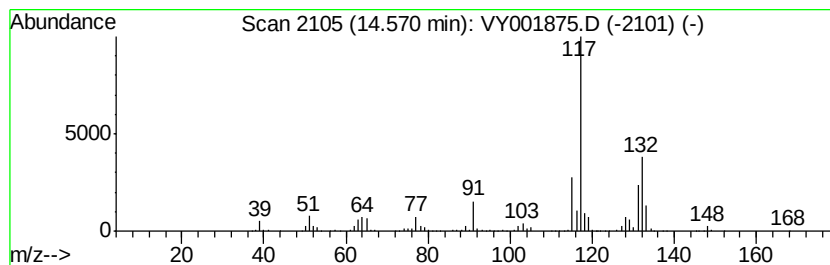
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	31.41 ug/l	439162	1,4-Dichlorobenzene-d4	13.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	81



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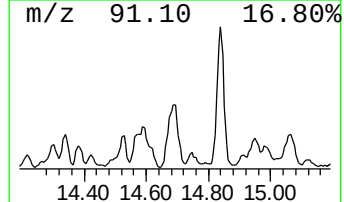
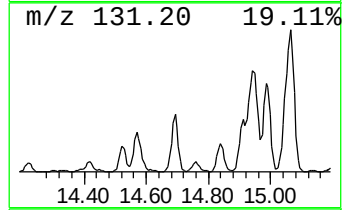
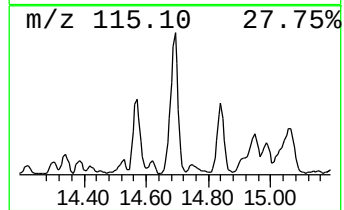
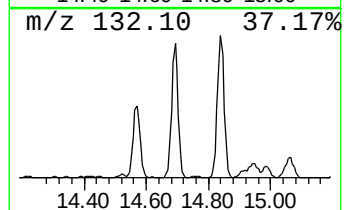
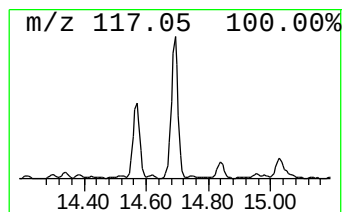
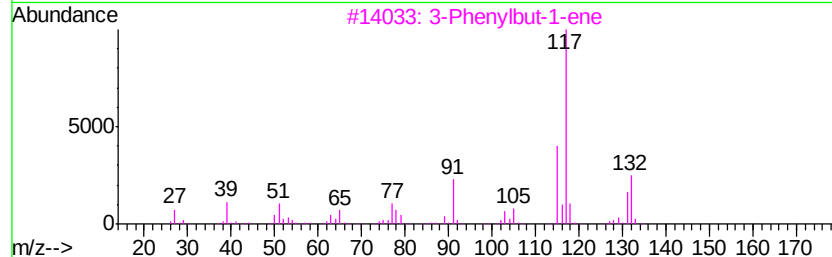
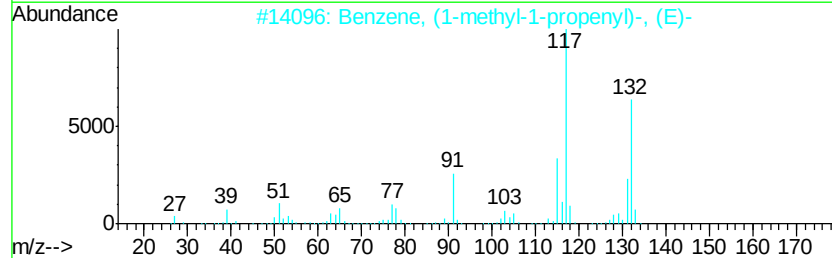
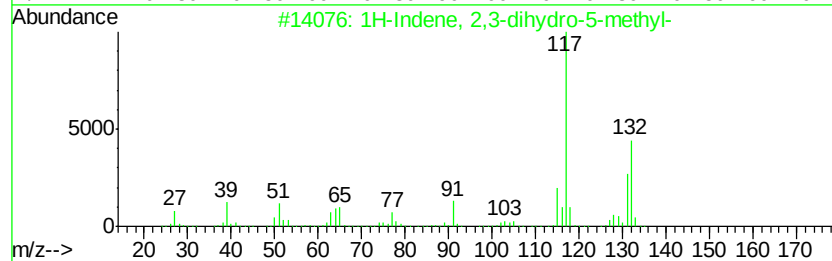
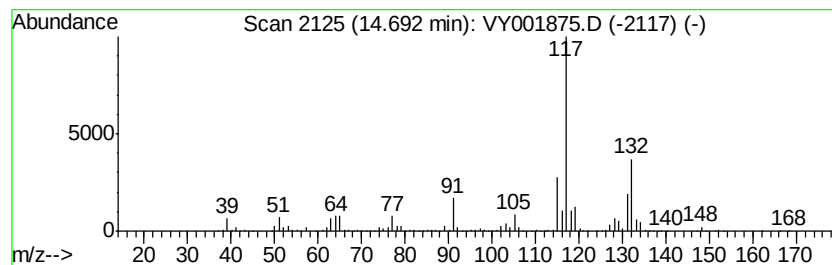
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	78.98 ug/l	1104280	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 92
2		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3 81
3		3-Phenylbut-1-ene	132 C10H12	000934-10-1 76
4		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 76
5		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 76



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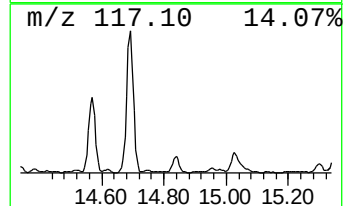
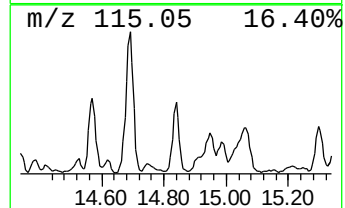
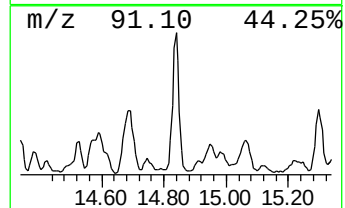
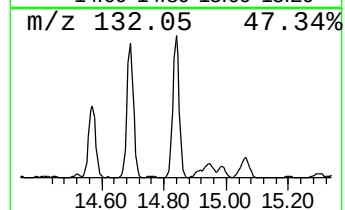
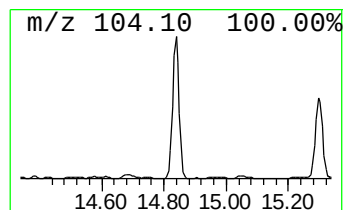
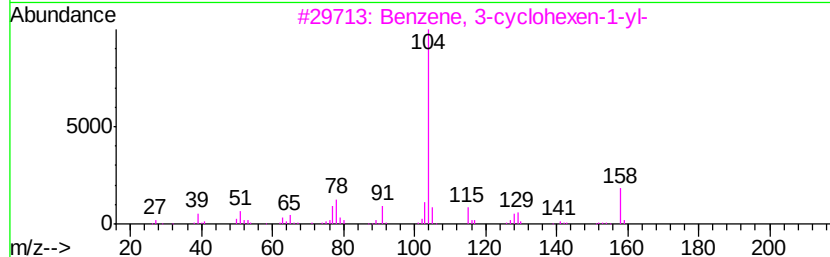
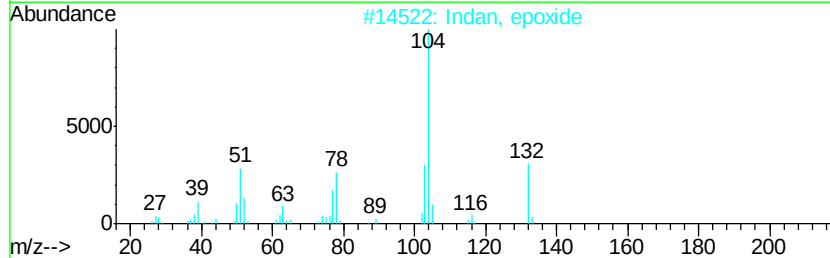
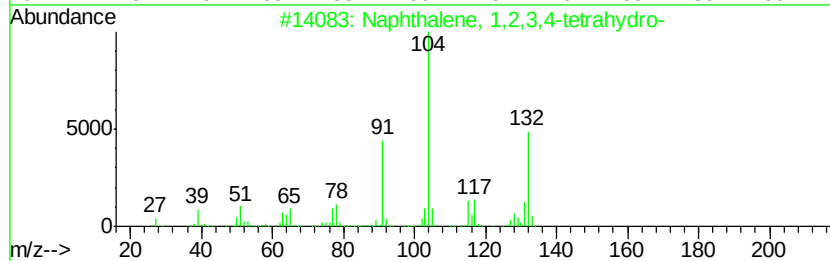
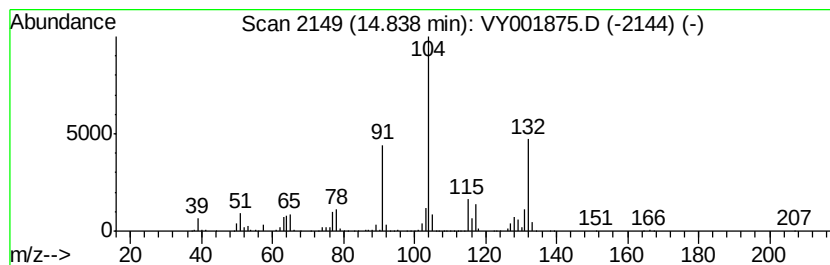
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Peak Number 3 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	46.87 ug/l	655314	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 97
2		Indan, epoxide	132 C9H8O	000768-22-9 58
3		Benzene, 3-cyclohexen-1-yl-	158 C12H14	004994-16-5 47
4		Isoquinoline, 1,2,3,4-tetrahydro-	133 C9H11N	000091-21-4 43
5		Acetic acid, trifluoro-, 2-phenyl...	218 C10H9F3O2	055419-66-4 38



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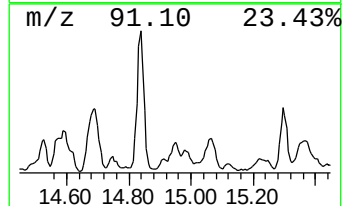
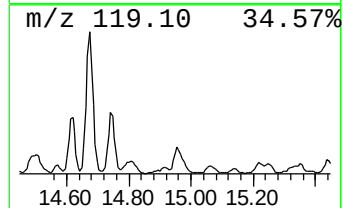
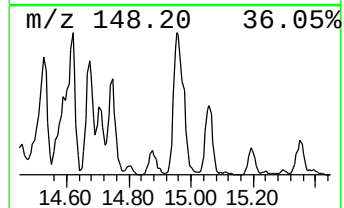
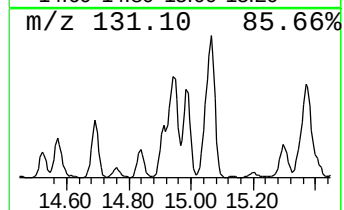
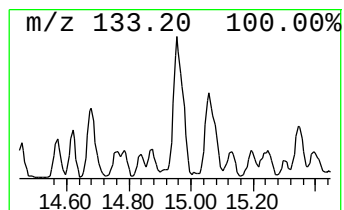
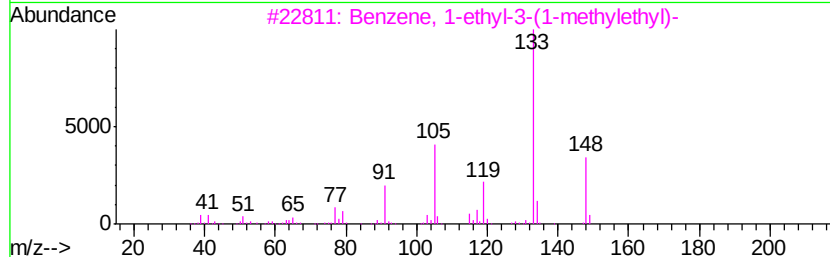
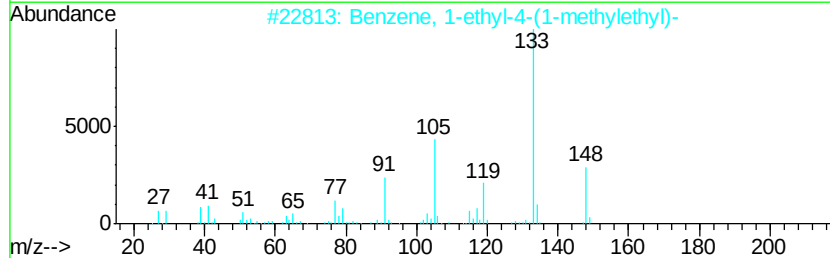
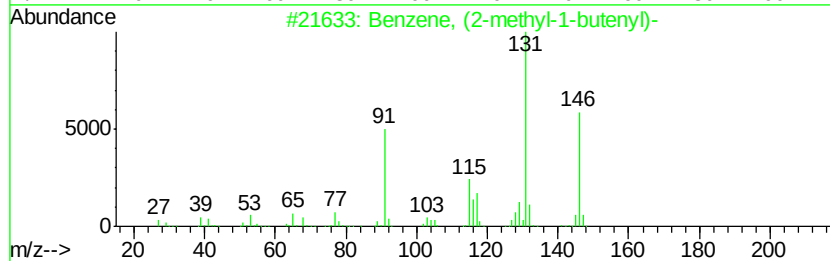
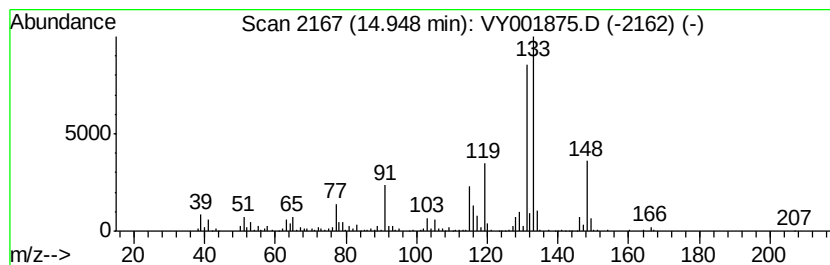
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Peak Number 4 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.95	29.09 ug/l	406703	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 64
2		Benzene, 1-ethyl-4-(1-methylethyl)-	148 C11H16	004218-48-8 49
3		Benzene, 1-ethyl-3-(1-methylethyl)-	148 C11H16	004920-99-4 46
4		Benzene, 1-(1,1-dimethylethyl)-3...	148 C11H16	001075-38-3 43
5		Benzene, pentamethyl-	148 C11H16	000700-12-9 43



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Data File : VY001875.D
Acq On : 05 Mar 2020 12:44
Operator : SY/MD
Sample : L1761-01
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SU-04-030420

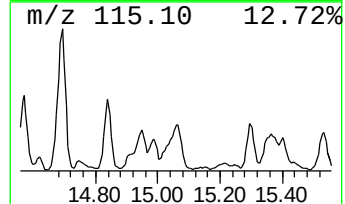
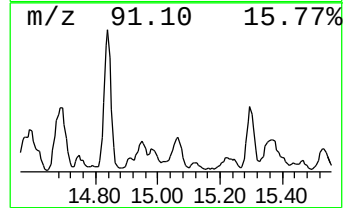
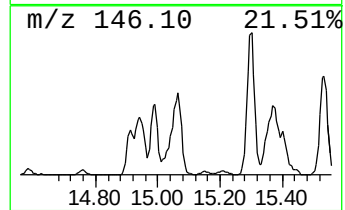
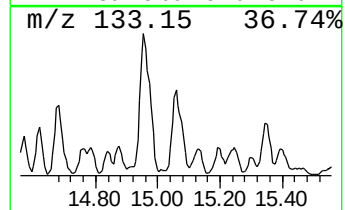
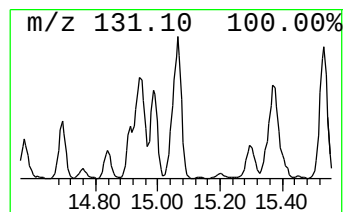
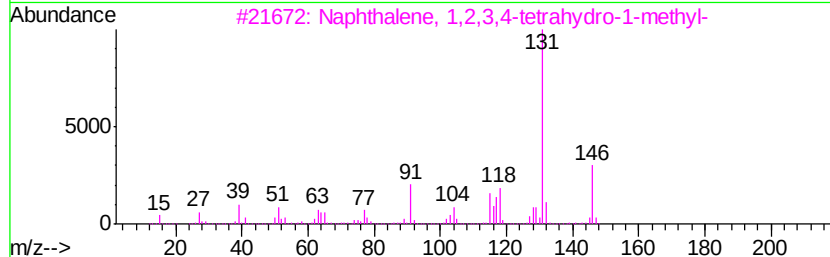
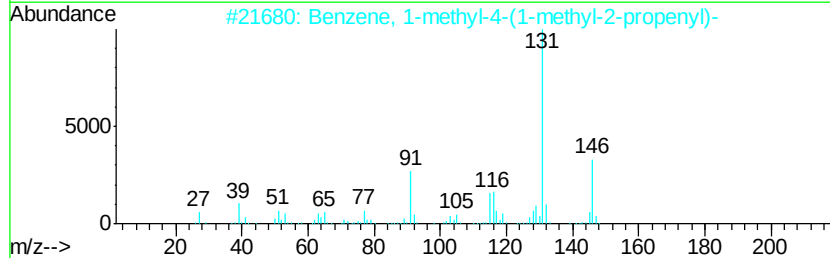
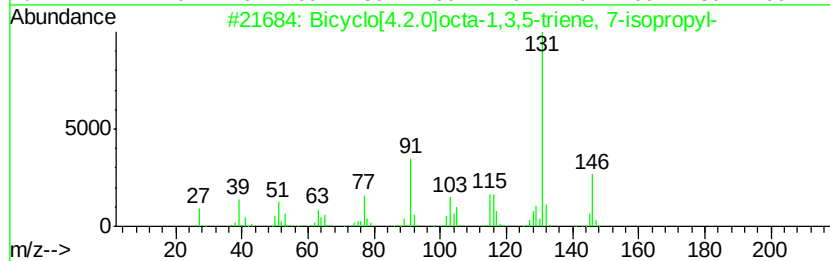
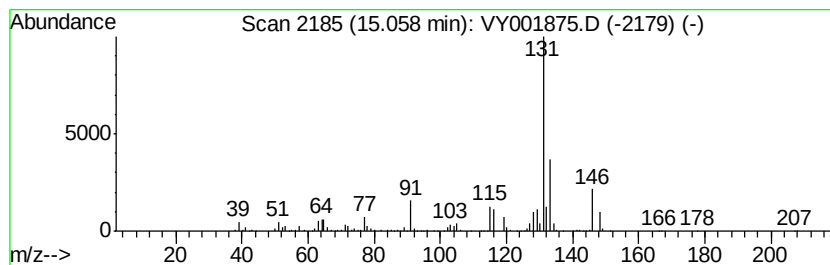
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.M
Quant Title : SW846 8260

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

Peak Number 5 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	35.46 ug/l	495768	1,4-Dichlorobenzene-d4	13.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	93
2		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	87
3		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	81
4		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	81
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY030520\
Data File : VY001875.D
Acq On : 05 Mar 2020 12:44
Operator : SY/MD
Sample : L1761-01
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SU-04-030420

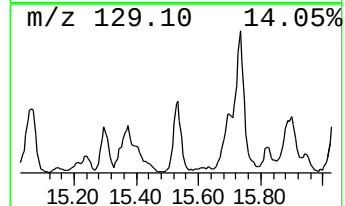
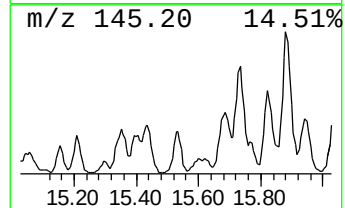
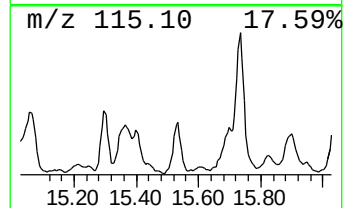
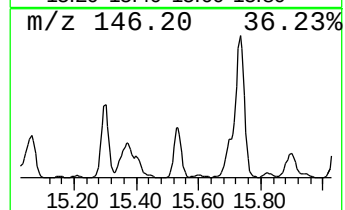
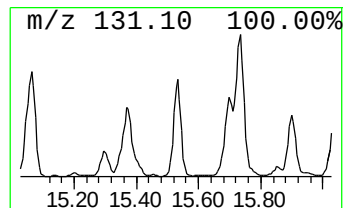
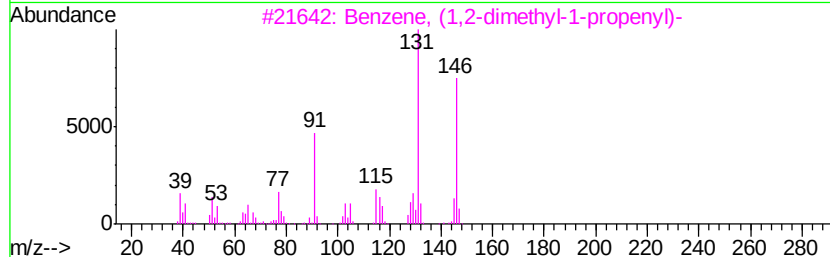
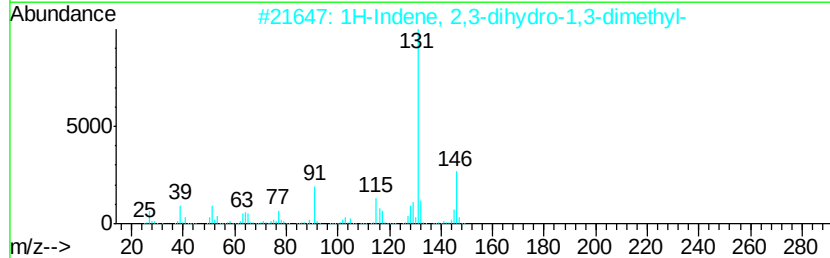
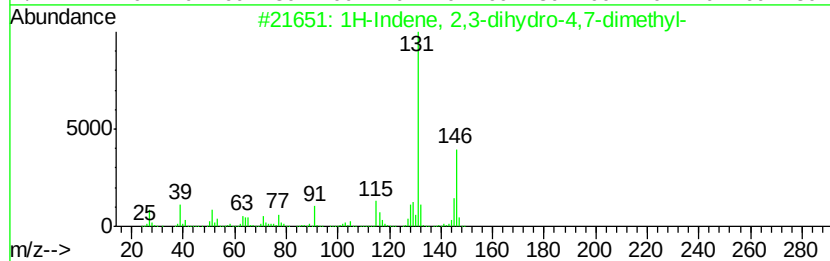
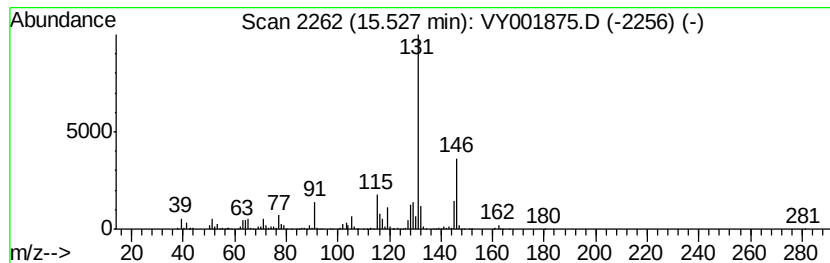
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	29.83 ug/l	417116	1,4-Dichlorobenzene-d4	13.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94
3		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
4		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY030520\
Data File : VY001875.D
Acq On : 05 Mar 2020 12:44
Operator : SY/MD
Sample : L1761-01
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

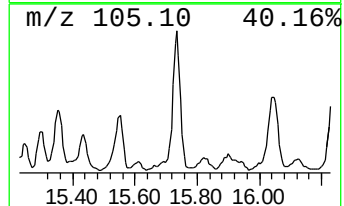
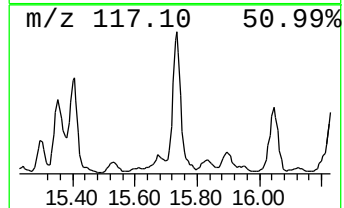
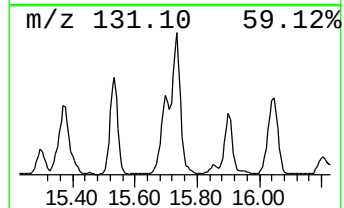
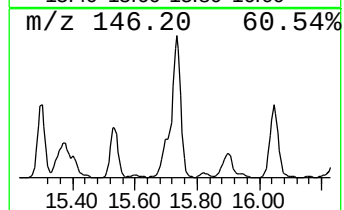
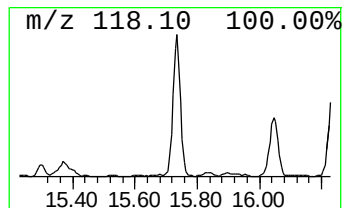
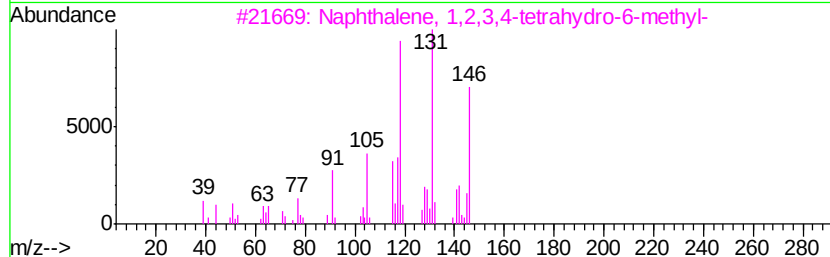
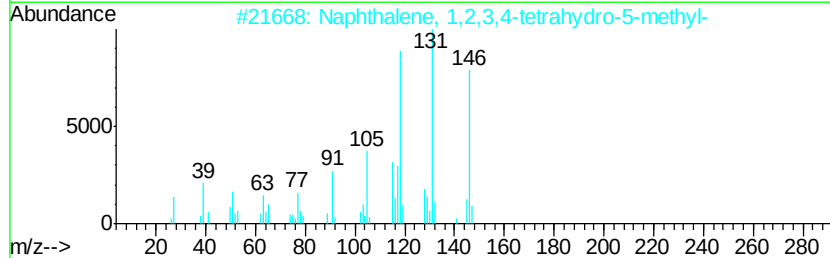
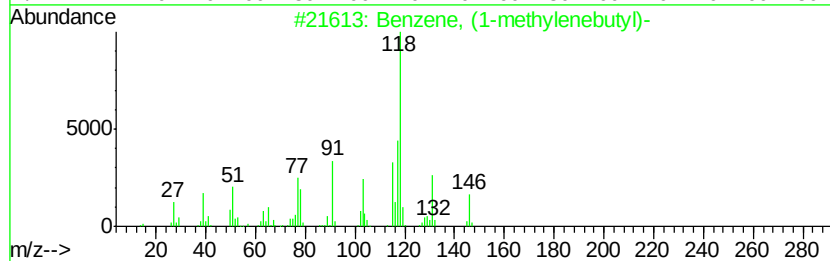
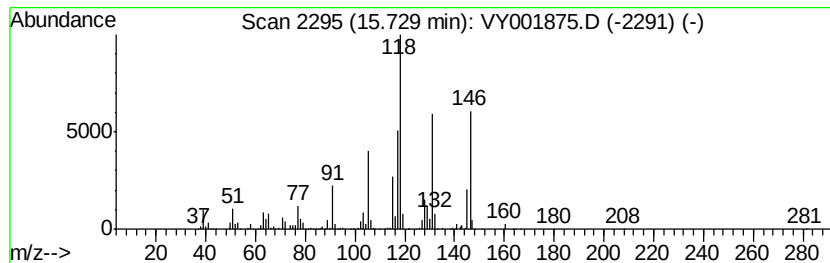
Instrument :
MSVOA_Y
ClientSampleId :
SU-04-030420

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methylenebutyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	67.48 ug/l	943434	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (1-methylenebutyl)-	146 C11H14	005676-32-4 86
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 64
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 62
4		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 49
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY030520\
Data File : VY001875.D
Acq On : 05 Mar 2020 12:44
Operator : SY/MD
Sample : L1761-01
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

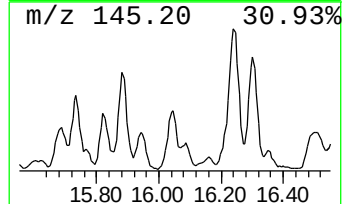
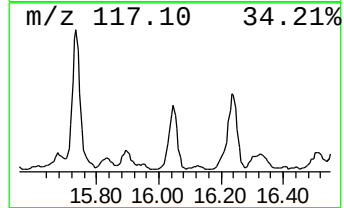
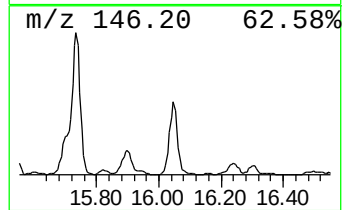
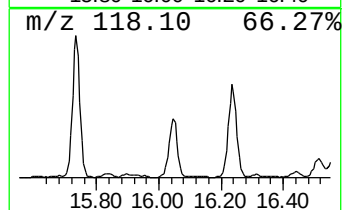
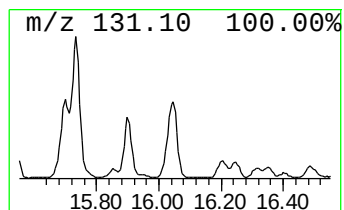
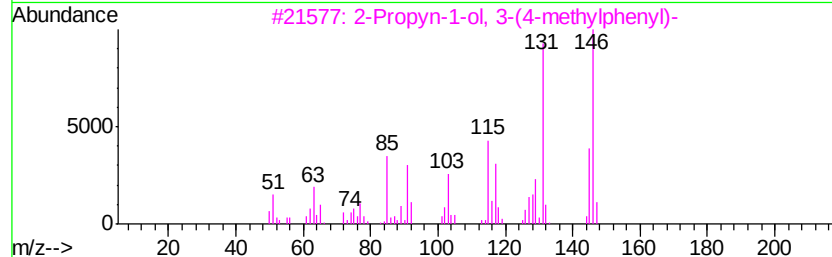
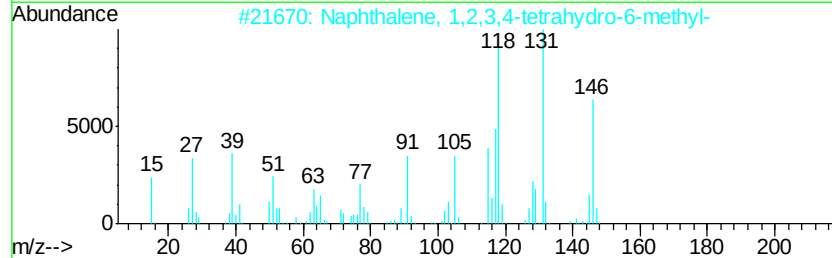
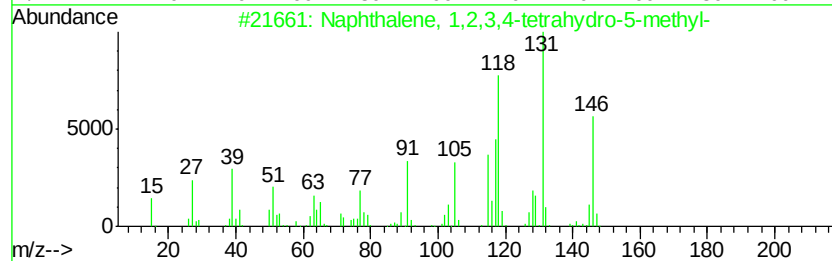
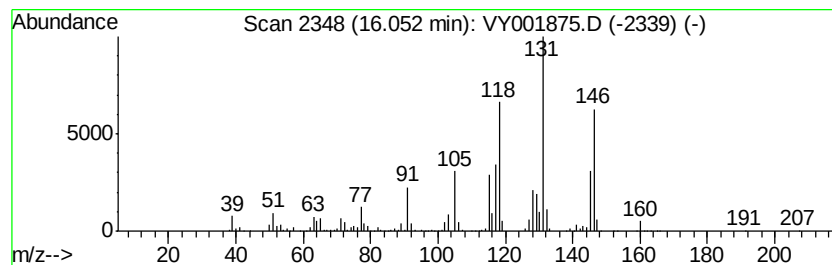
Instrument :
MSVOA_Y
ClientSampleId :
SU-04-030420

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	42.80 ug/l	598374	1,4-Dichlorobenzene-d4	13.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 95
3		2-Propyn-1-ol, 3-(4-methylphenyl)-	146 C10H10O	016017-24-6 72
4		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 68
5		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY030520\
Data File : VY001875.D
Acq On : 05 Mar 2020 12:44
Operator : SY/MD
Sample : L1761-01
Misc : 7.15G/5ML/MSVOA Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SU-04-030420

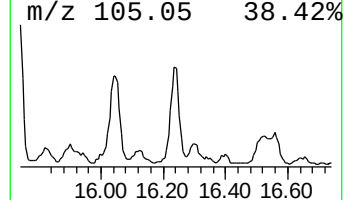
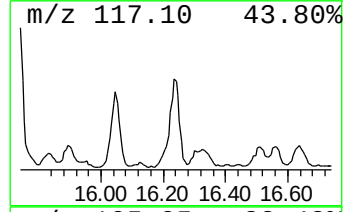
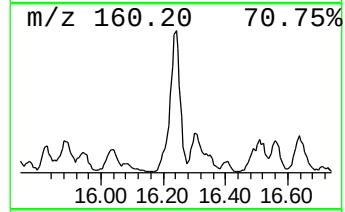
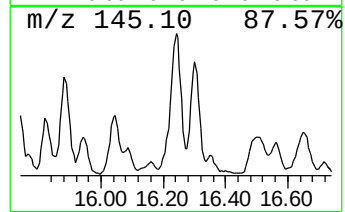
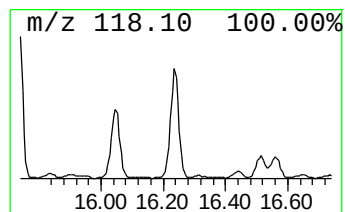
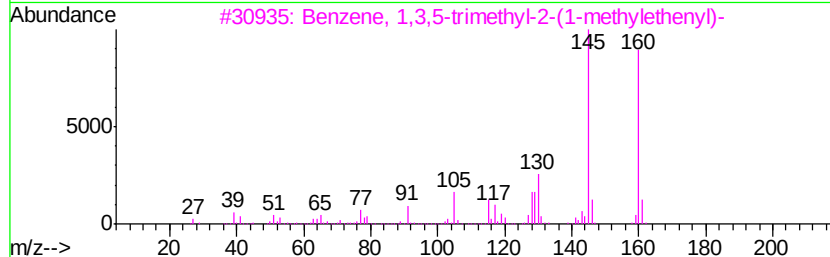
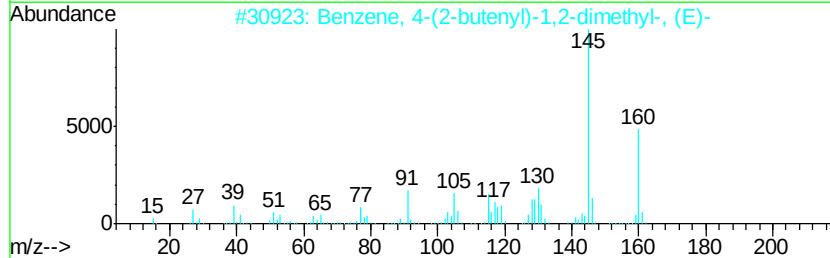
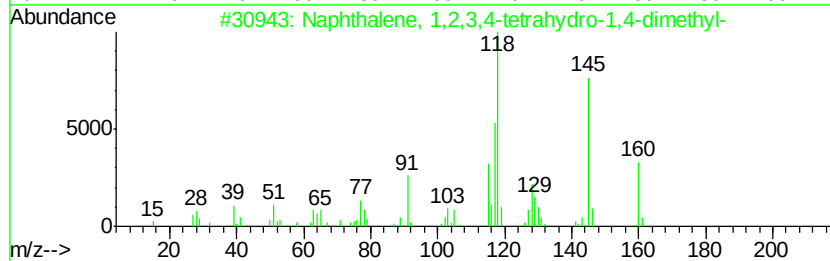
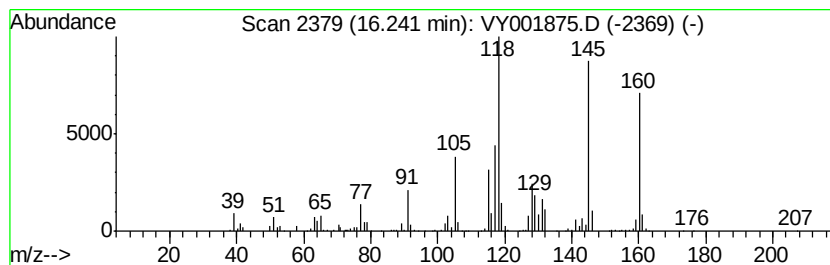
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.24	43.62 ug/l	609856	1,4-Dichlorobenzene-d4	13.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2		Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	78
3		Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	78
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	76
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_Y\DATA\VY030520\
Data File : VY001875.D
Acq On : 05 Mar 2020 12:44
Operator : SY/MD
Sample : L1761-01
Misc : 7.15G/5ML/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SU-04-030420

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y030420S.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-etheny...	14.57	31.4	ug/l	439162	4	13.43	699092	50.0
1H-Indene, 2,3-di...	14.69	79.0	ug/l	1104280	4	13.43	699092	50.0
Naphthalene, 1,2,...	14.84	46.9	ug/l	655314	4	13.43	699092	50.0
Benzene, (2-methy...	14.95	29.1	ug/l	406703	4	13.43	699092	50.0
Bicyclo[4.2.0]oct...	15.06	35.5	ug/l	495768	4	13.43	699092	50.0
1H-Indene, 2,3-di...	15.53	29.8	ug/l	417116	4	13.43	699092	50.0
Benzene, (1-methy...	15.73	67.5	ug/l	943434	4	13.43	699092	50.0
Naphthalene, 1,2,...	16.05	42.8	ug/l	598374	4	13.43	699092	50.0
Naphthalene, 1,2,...	16.24	43.6	ug/l	609856	4	13.43	699092	50.0