

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW012320\
 Data File : VW014733.D
 Acq On : 23 Jan 2020 19:14
 Operator : SY/VA
 Sample : L1243-10
 Misc : 7.25G/5ML/MSVOA W/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 CTY-SB-03-5-6

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_W\METHOD\82W010320S.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	1.959	4	9	28	rVB4	34285	84316	3.50%	0.531%
2	2.178	35	45	49	rBV2	21473	48057	1.99%	0.303%
3	2.349	63	73	81	rBV2	53844	121534	5.04%	0.765%
4	3.020	177	183	196	rVB	53327	131052	5.43%	0.825%
5	3.385	235	243	254	rVB2	37011	96320	3.99%	0.607%
6	4.117	349	363	374	rBV3	18811	62780	2.60%	0.395%
7	4.373	397	405	414	rBV2	12297	35278	1.46%	0.222%
8	4.916	481	494	497	rBV3	34694	112341	4.66%	0.708%
9	5.434	568	579	593	rBV3	19806	70609	2.93%	0.445%
10	5.934	651	661	671	rBV4	16782	49478	2.05%	0.312%
11	6.982	824	833	844	rVB4	31775	89995	3.73%	0.567%
12	7.878	968	980	985	rBV	230572	547558	22.71%	3.448%
13	7.946	985	991	1000	rVV	509350	1169890	48.52%	7.368%
14	8.031	1000	1005	1016	rVV3	17604	46331	1.92%	0.292%
15	8.153	1016	1025	1030	rVV2	19968	44676	1.85%	0.281%
16	8.220	1030	1036	1042	rVV4	10792	27190	1.13%	0.171%
17	8.305	1042	1050	1061	rVV2	322533	826481	34.28%	5.205%
18	8.433	1061	1071	1076	rVV7	16748	52417	2.17%	0.330%
19	8.506	1077	1083	1087	rVV6	13305	36211	1.50%	0.228%
20	8.567	1087	1093	1105	rVB2	25733	63591	2.64%	0.400%
21	8.689	1108	1113	1120	rBV6	13663	28012	1.16%	0.176%
22	8.836	1127	1137	1148	rBV	648938	1312610	54.44%	8.267%
23	9.329	1205	1218	1226	rBV	153507	358200	14.86%	2.256%
24	9.604	1253	1263	1269	rVB6	11106	27619	1.15%	0.174%
25	9.750	1280	1287	1294	rVB3	16251	34146	1.42%	0.215%
26	10.091	1332	1343	1353	rBV5	19825	62696	2.60%	0.395%
27	10.323	1372	1381	1388	rBV	809678	1481077	61.42%	9.328%
28	10.384	1389	1391	1397	rVB	31510	44161	1.83%	0.278%
29	10.494	1402	1409	1418	rVB5	28491	75561	3.13%	0.476%
30	10.665	1431	1437	1441	rBV4	27993	49860	2.07%	0.314%
31	10.756	1447	1452	1463	rVB5	17572	33458	1.39%	0.211%
32	11.177	1512	1521	1525	rBV3	21837	44223	1.83%	0.279%
33	11.225	1525	1529	1535	rVB2	36148	58892	2.44%	0.371%
34	11.433	1551	1563	1572	rVB9	12572	42981	1.78%	0.271%

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Integrator: RTE
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35	11.628	1587	1595	1605	rBV	633246	1081393	44.85%	6.810%
36	11.725	1606	1611	1615	rBV3	19908	33500	1.39%	0.211%
37	11.835	1621	1629	1635	rBV	58329	139628	5.79%	0.879%
38	12.164	1671	1683	1691	rBV2	45049	125845	5.22%	0.793%
39	12.341	1707	1712	1718	rBV7	15541	37420	1.55%	0.236%
40	12.463	1726	1732	1738	rVB2	37207	70089	2.91%	0.441%
41	12.615	1751	1757	1765	rBV	339314	561627	23.29%	3.537%
42	12.780	1779	1784	1791	rVB5	17402	38667	1.60%	0.244%
43	12.865	1792	1798	1801	rBV	29823	53601	2.22%	0.338%
44	12.939	1806	1810	1817	rVB	59544	104120	4.32%	0.656%
45	13.103	1829	1837	1844	rBV	25122	52746	2.19%	0.332%
46	13.249	1855	1861	1866	rBV	69221	120525	5.00%	0.759%
47	13.438	1887	1892	1896	rBV6	17981	34167	1.42%	0.215%
48	13.554	1905	1911	1921	rBV2	327717	612329	25.39%	3.856%
49	13.719	1931	1938	1939	rBV4	12092	25195	1.04%	0.159%
50	13.755	1939	1944	1957	rVB2	301302	556159	23.06%	3.503%
51	13.871	1959	1963	1969	rVB5	36407	62704	2.60%	0.395%
52	13.944	1969	1975	1979	rBV2	23713	42785	1.77%	0.269%
53	14.012	1982	1986	1988	rBV2	16937	24614	1.02%	0.155%
54	14.091	1997	1999	2005	rVB	47150	63405	2.63%	0.399%
55	14.201	2012	2017	2023	rBV3	62711	105280	4.37%	0.663%
56	14.335	2034	2039	2043	rBV2	29027	48999	2.03%	0.309%
57	14.383	2043	2047	2049	rVV5	23592	39340	1.63%	0.248%
58	14.414	2049	2052	2055	rVV2	36464	57715	2.39%	0.363%
59	14.457	2055	2059	2062	rVV	41498	64135	2.66%	0.404%
60	14.499	2062	2066	2069	rVV4	24560	36526	1.51%	0.230%
61	14.542	2069	2073	2079	rVB7	27585	56686	2.35%	0.357%
62	14.688	2092	2097	2100	rBV4	33569	53315	2.21%	0.336%
63	14.804	2108	2116	2121	rBV3	87785	217933	9.04%	1.373%
64	14.865	2122	2126	2131	rVB5	29903	52565	2.18%	0.331%
65	14.956	2133	2141	2147	rVV3	82709	176101	7.30%	1.109%
66	15.066	2153	2159	2161	rBV6	33484	63274	2.62%	0.398%
67	15.170	2171	2176	2183	rVB5	29905	72096	2.99%	0.454%
68	15.365	2197	2208	2220	rBV	1366632	2411295	100.00%	15.186%
69	15.469	2220	2225	2238	rVB7	30632	117376	4.87%	0.739%
70	15.651	2249	2255	2261	rBV5	19968	50613	2.10%	0.319%
71	16.176	2335	2341	2350	rVB9	17591	44172	1.83%	0.278%

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Sampling : 1 Min Area: 3 % of largest Peak
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72	16.438	2377	2384	2399	rVB	253720	560890	23.26%	3.532%
73	16.639	2410	2417	2429	rVB	194453	440061	18.25%	2.771%

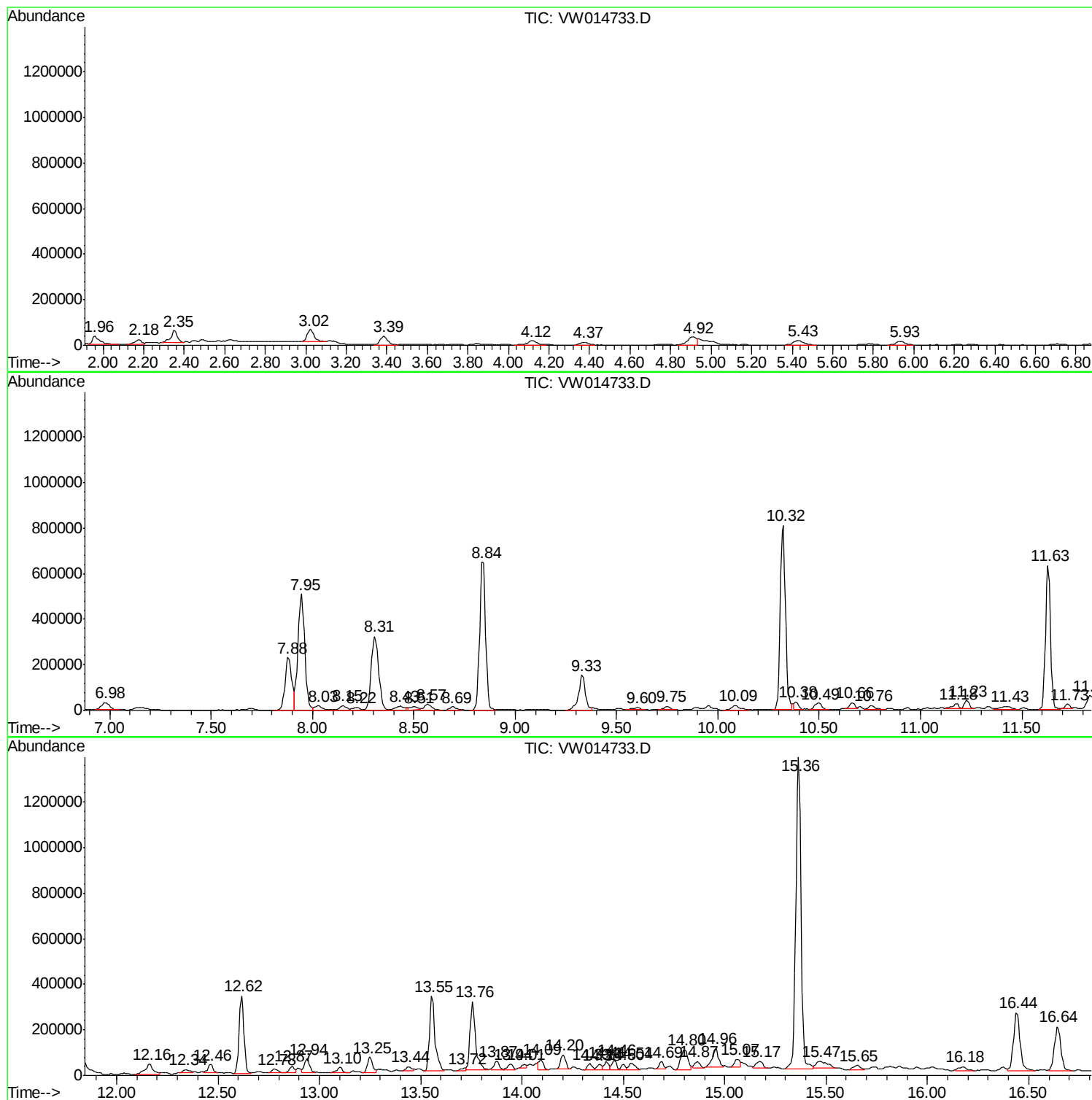
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
Quant Title : SW846 8260

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



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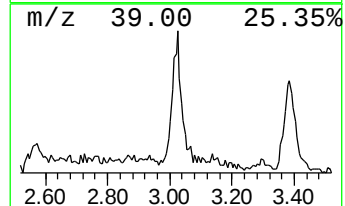
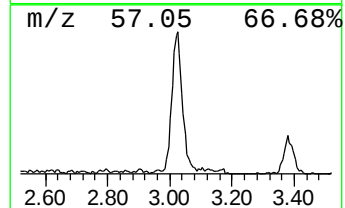
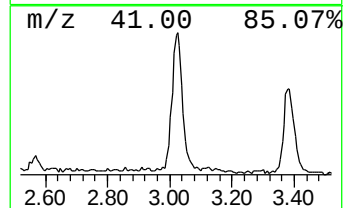
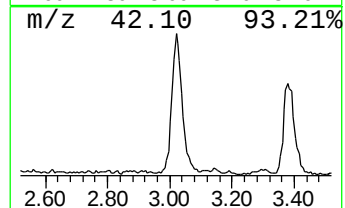
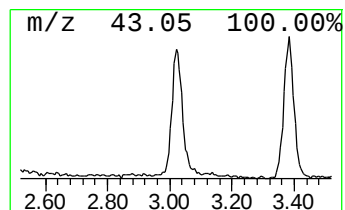
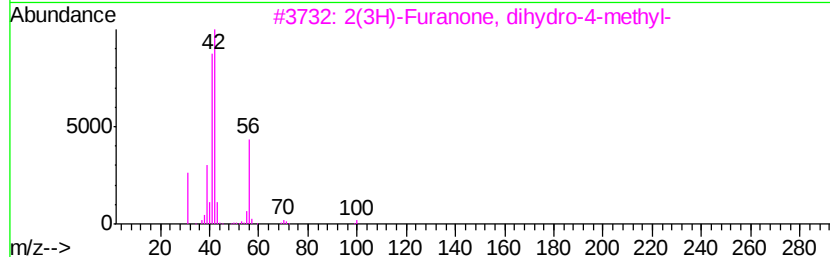
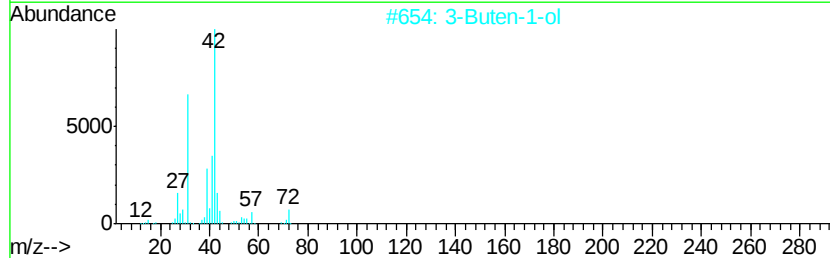
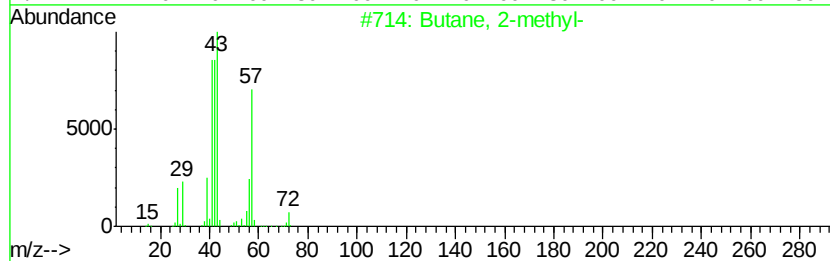
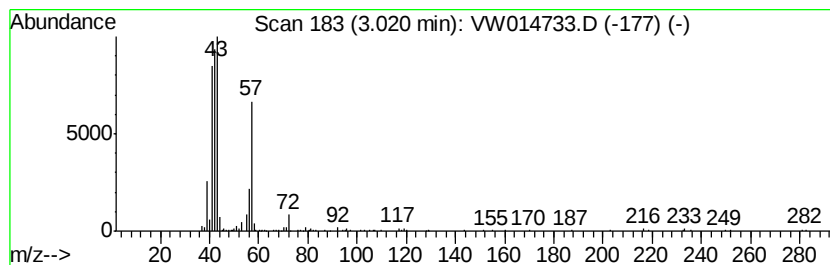
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.02	5.60 ug/l	131052	Pentafluorobenzene	7.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2		3-Buten-1-ol	72	C4H8O	000627-27-0	38
3		2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	25
4		Butane, 2-nitro-	103	C4H9NO2	000600-24-8	10
5		Propane, 1-chloro-2-methyl-	92	C4H9Cl	000513-36-0	9



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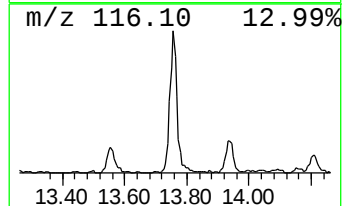
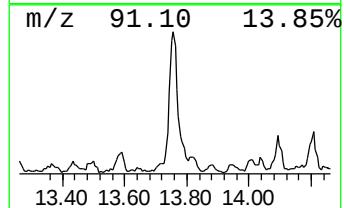
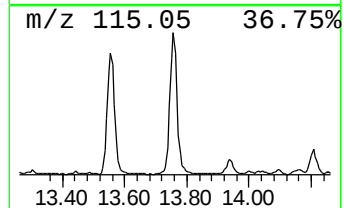
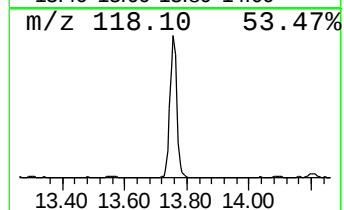
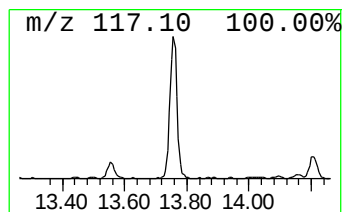
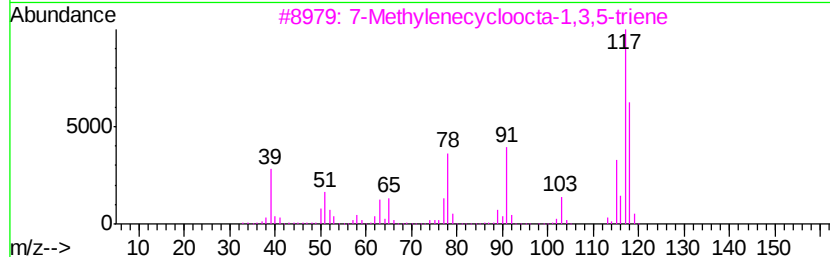
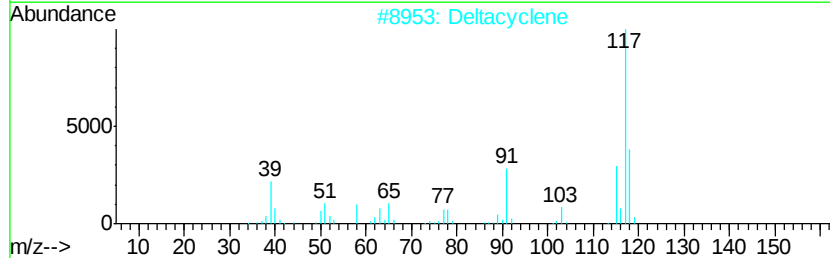
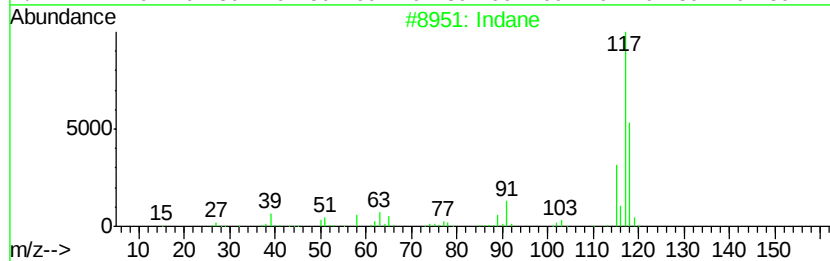
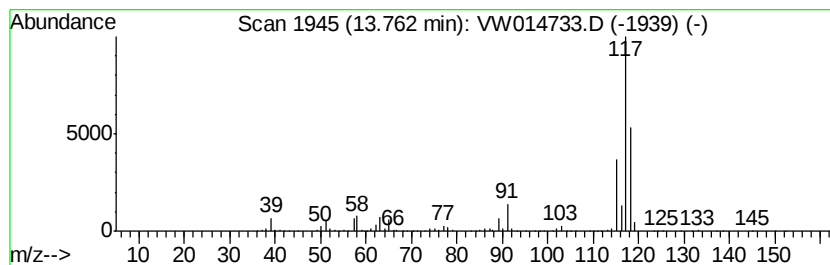
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
Quant Title : SW846 8260

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

Peak Number 2 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	45.41 ug/l	556159	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	90
2		Deltacyclene	118	C9H10	007785-10-6	64
3		7-Methylenecycloocta-1,3,5-triene	118	C9H10	002570-13-0	53
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	52
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	52



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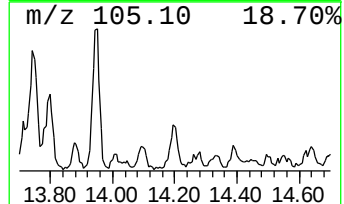
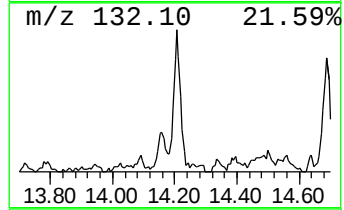
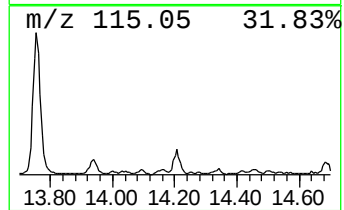
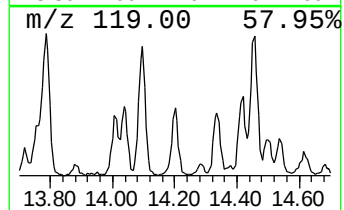
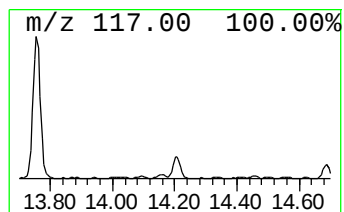
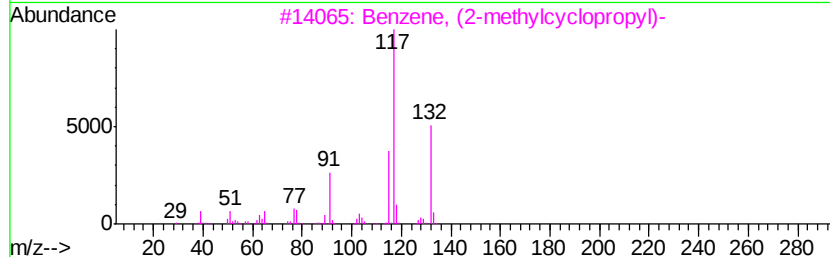
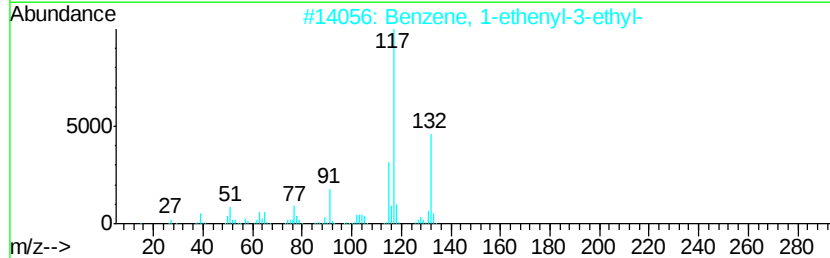
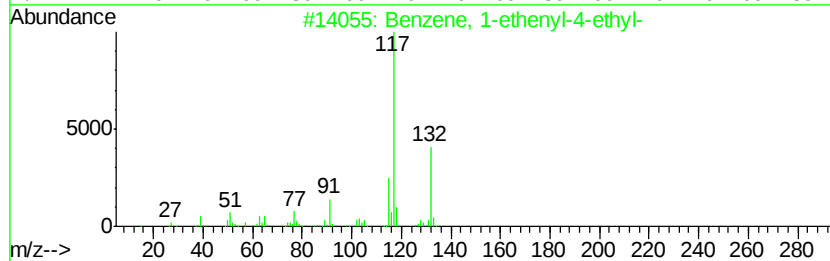
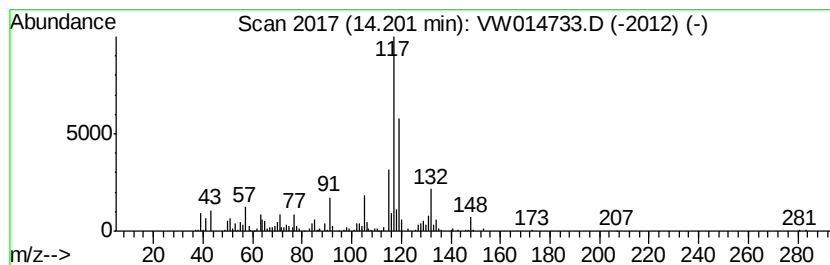
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	8.60 ug/l	105280	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	64
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	62
5		Indan, 1-methyl-	132	C10H12	000767-58-8	62



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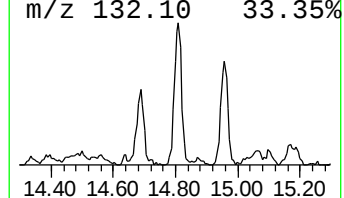
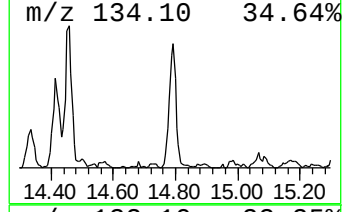
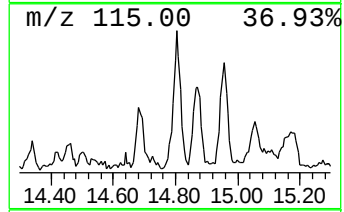
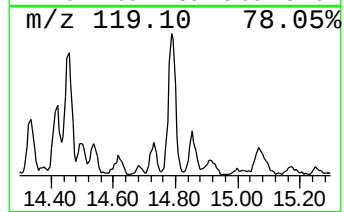
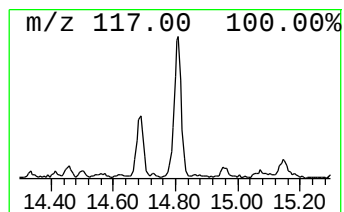
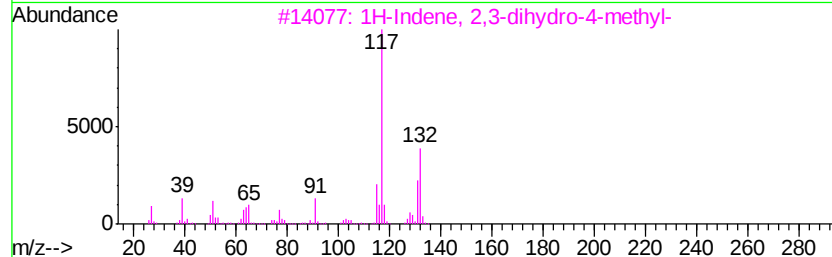
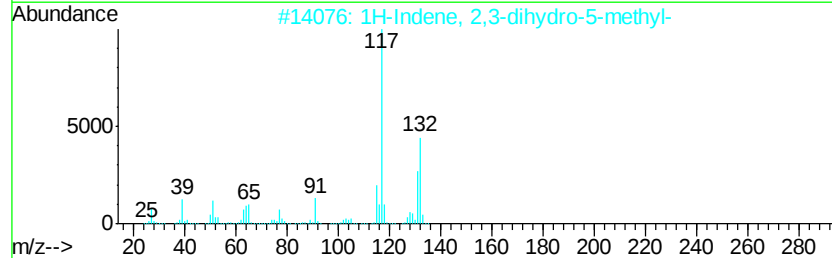
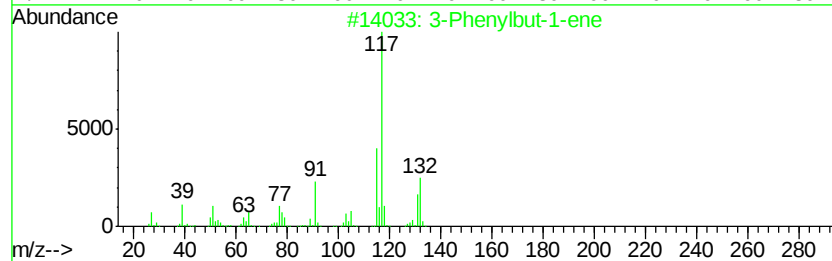
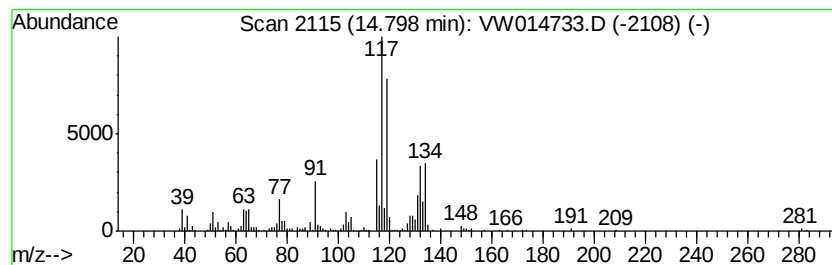
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Quant Title : SW846 8260

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

Peak Number 5 3-Phenylbut-1-ene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	17.80 ug/l	217933	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64
4		Indan, 1-methyl-	132	C10H12	000767-58-8	62
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012320\
Data File : VW014733.D
Acq On : 23 Jan 2020 19:14
Operator : SY/VA
Sample : L1243-10
Misc : 7.25G/5ML/MSVOA W/SOIL
ALS Vial : 17 Sample Multiplier: 1

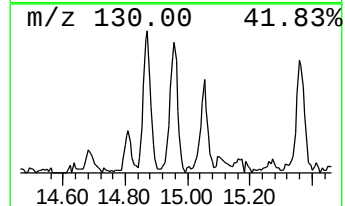
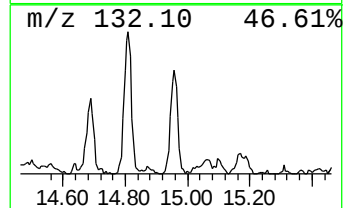
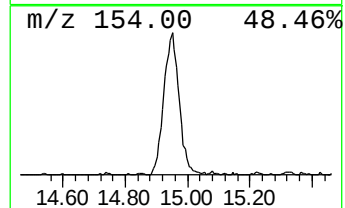
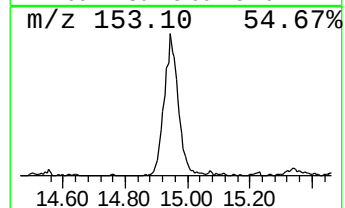
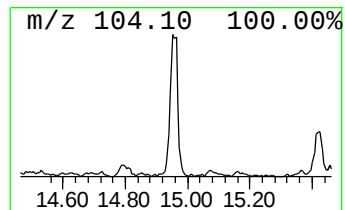
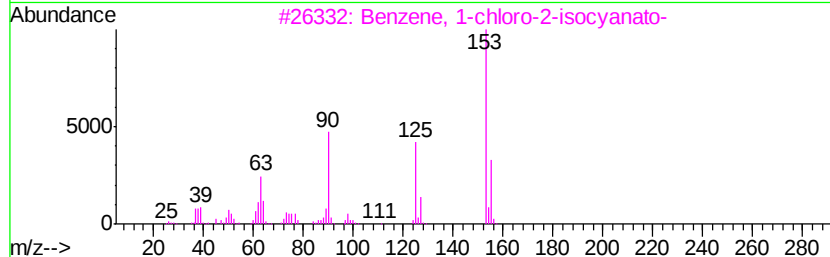
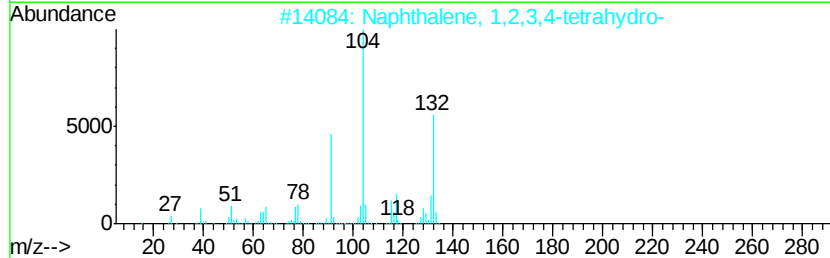
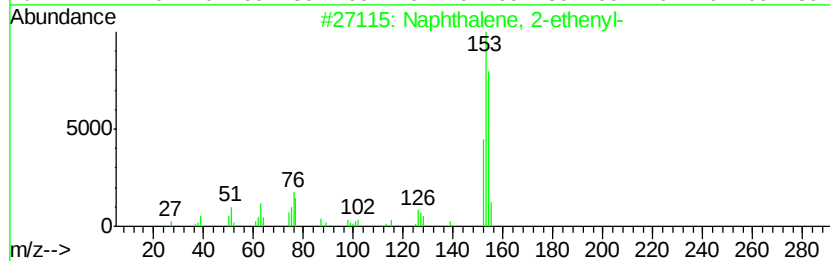
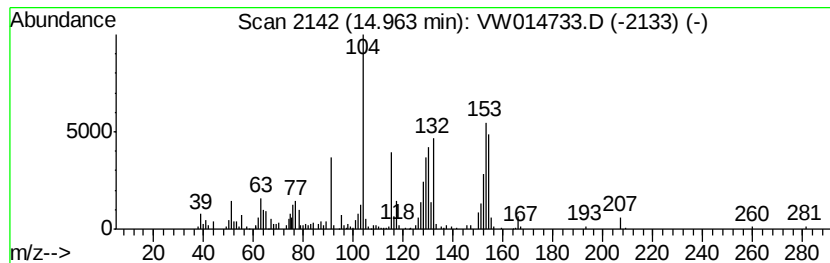
Instrument :
MSVOA_W
ClientSampleId :
CTY-SB-03-5-6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 2-ethenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.96	14.38 ug/l	176101	1,4-Dichlorobenzene-d4	13.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-ethenvl-	154	C12H10	000827-54-3	58
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	38
3	Benzene, 1-chloro-2-isocvanato-	153	C7H4ClNO	003320-83-0	25
4	2a,4a,6a,6b-Tetrahydrocyclopenta...	130	C10H10	006053-74-3	25
5	Benzene, 1-butynyl-	130	C10H10	000622-76-4	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012320\
Data File : VW014733.D
Acq On : 23 Jan 2020 19:14
Operator : SY/VA
Sample : L1243-10
Misc : 7.25G/5ML/MSVOA W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CTY-SB-03-5-6

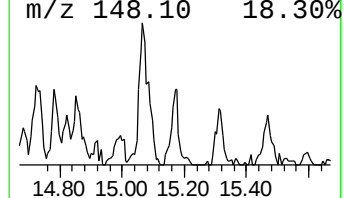
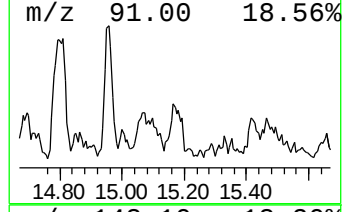
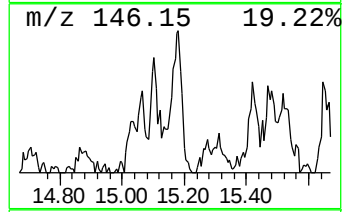
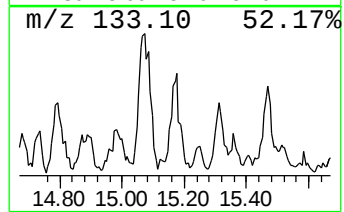
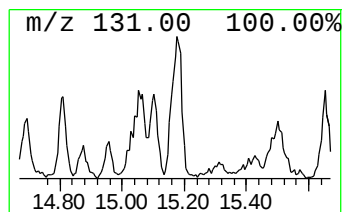
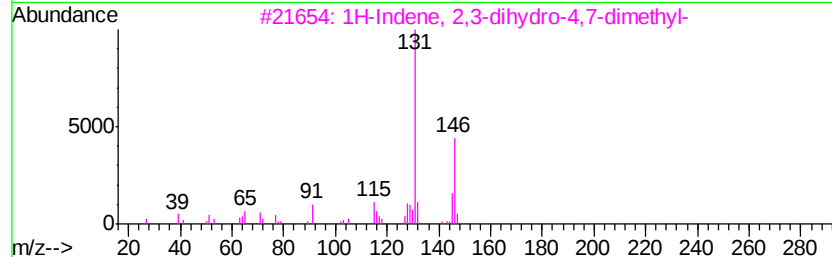
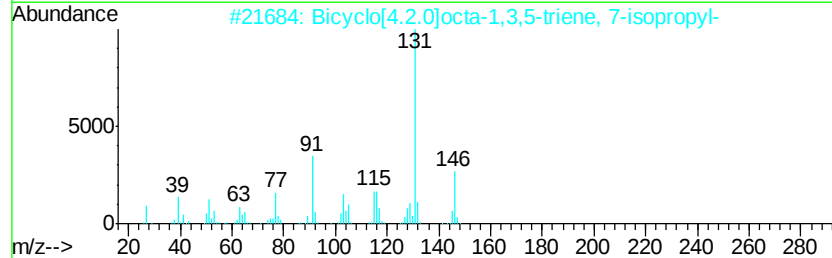
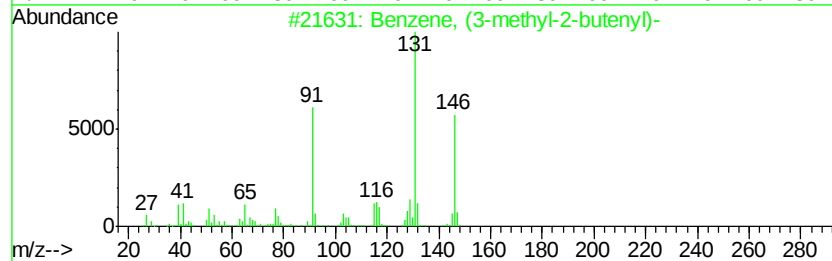
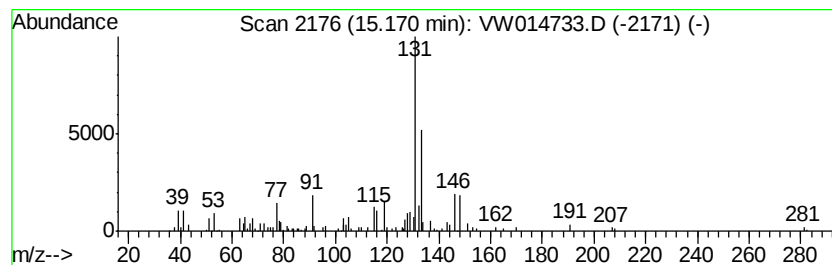
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (3-methyl-2-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	5.89 ug/l	72096	1,4-Dichlorobenzene-d4	13.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	70
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
4		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	58
5		Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012320\
Data File : VW014733.D
Acq On : 23 Jan 2020 19:14
Operator : SY/VA
Sample : L1243-10
Misc : 7.25G/5ML/MSVOA W/SOIL
ALS Vial : 17 Sample Multiplier: 1

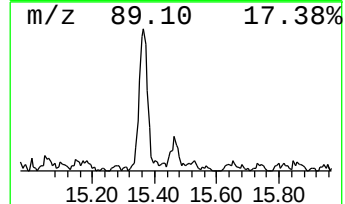
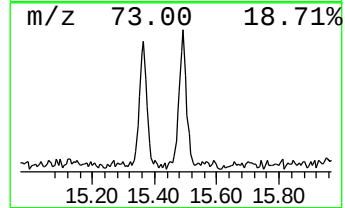
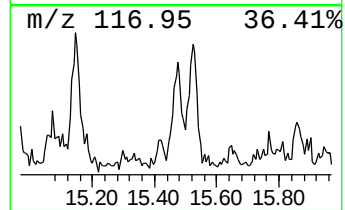
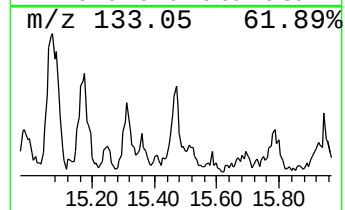
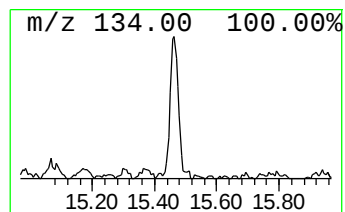
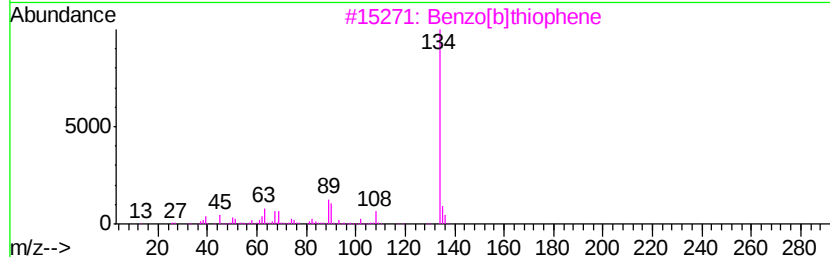
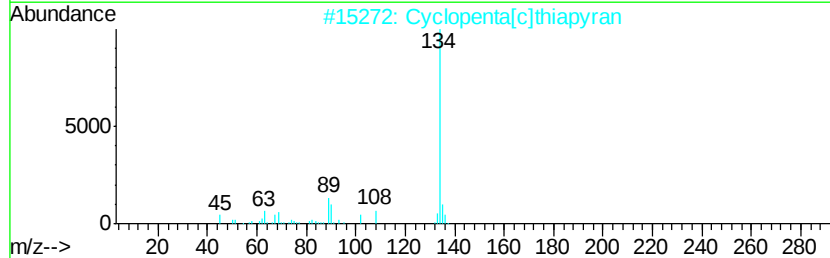
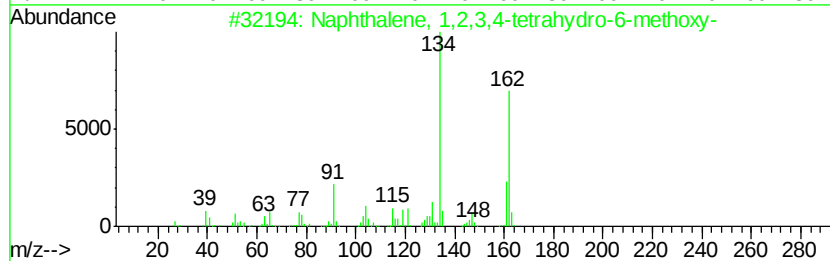
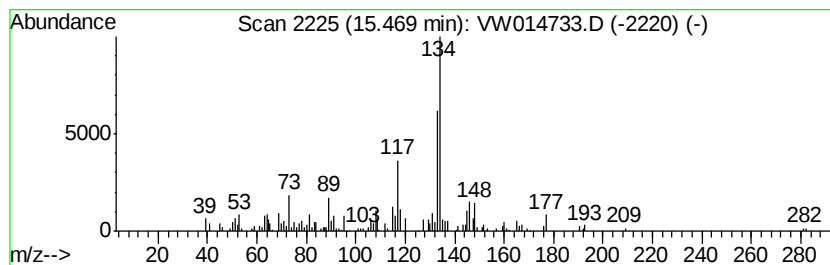
Instrument :
MSVOA_W
ClientSampleId :
CTY-SB-03-5-6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
Quant Title : SW846 8260

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

Peak Number 8 unknown15.47 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.47	9.58 ug/l	117376	1,4-Dichlorobenzene-d4	13.55	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	162	C11H14O	001730-48-9	38
2	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	38
3	Benzo[b]thiophene	134	C8H6S	000095-15-8	35
4	Pyrimidine, 2,4,6-trifluoro-	134	C4HF3N2	000696-82-2	27
5	1H-Purine, 6-methyl-	134	C6H6N4	002004-03-7	27



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW012320\
Data File : VW014733.D
Acq On : 23 Jan 2020 19:14
Operator : SY/VA
Sample : L1243-10
Misc : 7.25G/5ML/MSVOA W/SOIL
ALS Vial : 17 Sample Multiplier: 1

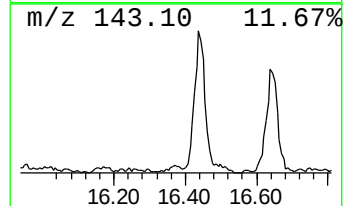
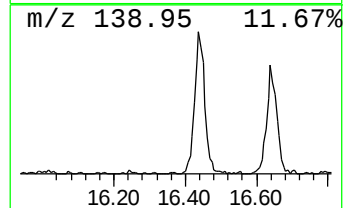
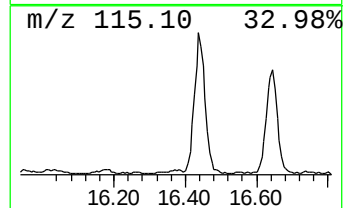
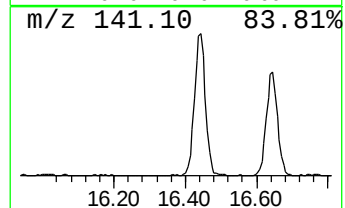
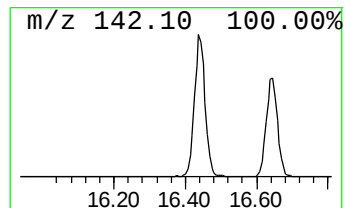
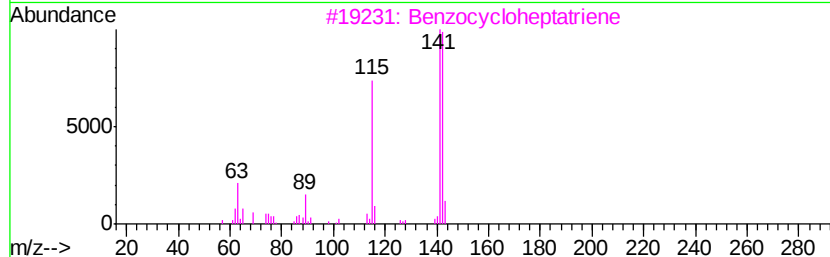
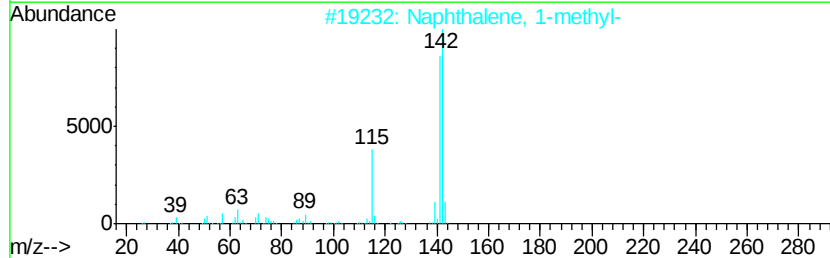
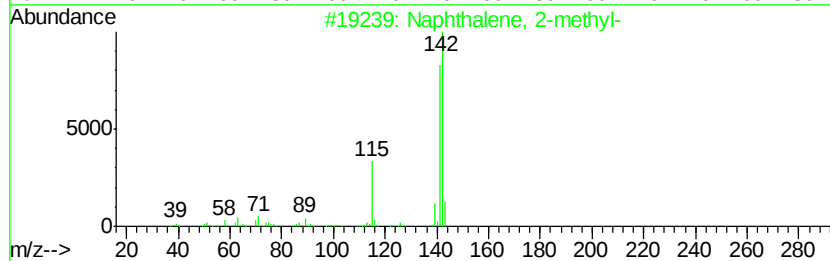
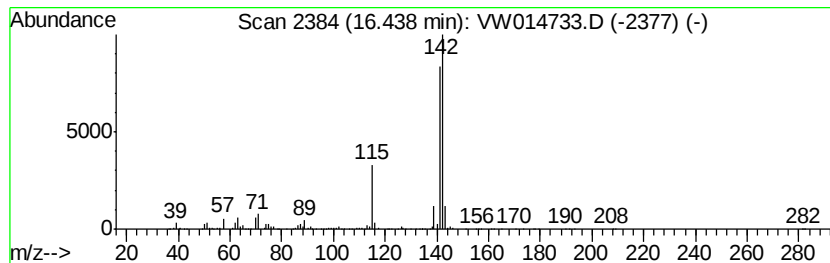
Instrument :
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ClientSampleId :
CTY-SB-03-5-6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	45.80 ug/l	560890	1,4-Dichlorobenzene-d4	13.55
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
3		Benzocycloheptatriene	142 C11H10	000264-09-5 91
4		1,4-Methanonaphthalene, 1,4-dihy...	142 C11H10	004453-90-1 91
5		1H-Indene, 1-ethylidene-	142 C11H10	002471-83-2 59



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW012320\
Data File : VW014733.D
Acq On : 23 Jan 2020 19:14
Operator : SY/VA
Sample : L1243-10
Misc : 7.25G/5ML/MSVOA_W/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
CTY-SB-03-5-6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010320S.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
Butane, 2-methyl-	3.02	5.6	ug/l	131052	1	7.95	1169890	50.0
Indane	13.76	45.4	ug/l	556159	4	13.55	612329	50.0
Benzene, 1-etheny...	14.20	8.6	ug/l	105280	4	13.55	612329	50.0
3-Phenylbut-1-ene	14.80	17.8	ug/l	217933	4	13.55	612329	50.0
Naphthalene, 2-et...	14.96	14.4	ug/l	176101	4	13.55	612329	50.0
Benzene, (3-methy...	15.17	5.9	ug/l	72096	4	13.55	612329	50.0
unknown15.47	15.47	9.6	ug/l	117376	4	13.55	612329	50.0
Naphthalene, 1-me...	16.44	45.8	ug/l	560890	4	13.55	612329	50.0