

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052618\
 Data File : VX002018.D
 Acq On : 26 May 2018 14:48
 Operator : JC/MD
 Sample : J3064-05
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-17

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 : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.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	1.075	76	80	93	rBV	3631079	4167114	82.35%	8.264%
2	3.208	424	430	438	rVB	31311	74888	1.48%	0.149%
3	5.507	794	807	823	rBV	263029	773410	15.28%	1.534%
4	5.665	823	833	848	rVB	296573	835482	16.51%	1.657%
5	6.074	888	900	918	rBV	294374	762531	15.07%	1.512%
6	6.866	1020	1030	1044	rBV	457124	1075107	21.25%	2.132%
7	7.464	1118	1128	1140	rVB	50370	118896	2.35%	0.236%
8	8.720	1328	1334	1349	rVB	1359245	2137171	42.23%	4.238%
9	9.043	1381	1387	1394	rVB2	31132	51404	1.02%	0.102%
10	10.116	1557	1563	1571	rBV	969805	1325142	26.19%	2.628%
11	10.646	1641	1650	1657	rBV5	32039	92953	1.84%	0.184%
12	10.841	1671	1682	1687	rBV2	47567	85843	1.70%	0.170%
13	10.914	1690	1694	1698	rBV4	29991	53728	1.06%	0.107%
14	11.024	1706	1712	1716	rBV	144589	189136	3.74%	0.375%
15	11.140	1726	1731	1736	rBV	1197680	1573706	31.10%	3.121%
16	11.366	1763	1768	1778	rBV	190407	258282	5.10%	0.512%
17	11.683	1809	1820	1826	rBV5	33968	81068	1.60%	0.161%
18	11.774	1826	1835	1838	rBV	65019	100946	1.99%	0.200%
19	11.804	1838	1840	1851	rVB10	23678	52970	1.05%	0.105%
20	11.926	1851	1860	1861	rBV3	44490	76436	1.51%	0.152%
21	11.951	1861	1864	1870	rVB	109043	134204	2.65%	0.266%
22	12.085	1880	1886	1891	rBV	1203841	1500534	29.65%	2.976%
23	12.237	1905	1911	1917	rBV	156420	190129	3.76%	0.377%
24	12.304	1917	1922	1927	rVB	1516860	1820906	35.98%	3.611%
25	12.371	1929	1933	1943	rVB2	183313	313961	6.20%	0.623%
26	12.463	1943	1948	1953	rBV	433399	554624	10.96%	1.100%
27	12.670	1971	1982	1986	rBV	1225679	1607129	31.76%	3.187%
28	12.707	1986	1988	1993	rVV2	264206	339633	6.71%	0.674%
29	12.762	1993	1997	2004	rVV2	981350	1679743	33.19%	3.331%
30	12.816	2004	2006	2010	rVB2	108772	129373	2.56%	0.257%
31	12.877	2011	2016	2017	rBV	115859	143750	2.84%	0.285%
32	12.902	2017	2020	2024	rVB	193818	249678	4.93%	0.495%
33	12.981	2025	2033	2038	rBV	1146868	1587165	31.37%	3.148%
34	13.030	2038	2041	2046	rVB4	40296	79785	1.58%	0.158%

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35	13.085	2046	2050	2056	rBV3	189868	298538	5.90%	0.592%
36	13.158	2057	2062	2065	rBV	199470	261818	5.17%	0.519%
37	13.201	2065	2069	2072	rVV2	409481	521245	10.30%	1.034%
38	13.231	2072	2074	2076	rVV	259256	263512	5.21%	0.523%
39	13.255	2076	2078	2082	rVB	127771	144870	2.86%	0.287%
40	13.316	2082	2088	2089	rBV3	204132	253079	5.00%	0.502%
41	13.353	2089	2094	2101	rVV2	3456836	5060293	100.00%	10.035%
42	13.408	2101	2103	2104	rVV	57273	53049	1.05%	0.105%
43	13.432	2104	2107	2109	rVV	136285	145111	2.87%	0.288%
44	13.487	2109	2116	2121	rVB4	167342	307659	6.08%	0.610%
45	13.572	2126	2130	2132	rVV2	169858	233361	4.61%	0.463%
46	13.603	2132	2135	2138	rVV	255430	333541	6.59%	0.661%
47	13.646	2138	2142	2146	rVV3	629104	905723	17.90%	1.796%
48	13.700	2146	2151	2153	rVV3	424684	721235	14.25%	1.430%
49	13.731	2153	2156	2165	rVB3	662354	1267921	25.06%	2.514%
50	13.816	2165	2170	2173	rBV	183275	255320	5.05%	0.506%
51	13.859	2173	2177	2182	rVB2	358054	494854	9.78%	0.981%
52	13.914	2182	2186	2192	rVB	419755	569915	11.26%	1.130%
53	13.993	2192	2199	2202	rBV2	704482	1028914	20.33%	2.040%
54	14.036	2202	2206	2215	rVB4	360908	687726	13.59%	1.364%
55	14.139	2218	2223	2225	rBV2	221763	293931	5.81%	0.583%
56	14.164	2225	2227	2231	rVB2	128319	153914	3.04%	0.305%
57	14.225	2231	2237	2240	rBV6	80649	149658	2.96%	0.297%
58	14.292	2241	2248	2255	rVB2	803577	1648253	32.57%	3.269%
59	14.383	2259	2263	2267	rVB	320700	407610	8.06%	0.808%
60	14.438	2267	2272	2276	rVB	379652	517360	10.22%	1.026%
61	14.481	2276	2279	2282	rBV	110984	146279	2.89%	0.290%
62	14.548	2285	2290	2294	rBV2	794114	1045498	20.66%	2.073%
63	14.639	2302	2305	2308	rBV3	59791	83777	1.66%	0.166%
64	14.700	2312	2315	2319	rVV	438927	645567	12.76%	1.280%
65	14.737	2319	2321	2324	rVV2	158436	187770	3.71%	0.372%
66	14.792	2324	2330	2334	rVB5	130373	238569	4.71%	0.473%
67	14.847	2334	2339	2343	rBV	998351	1368130	27.04%	2.713%
68	14.914	2347	2350	2353	rVB3	72409	84678	1.67%	0.168%
69	14.975	2353	2360	2366	rBV6	146953	367865	7.27%	0.730%
70	15.054	2371	2373	2377	rVV2	108080	133705	2.64%	0.265%
71	15.121	2381	2384	2388	rVB5	97481	117513	2.32%	0.233%

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72	15.255	2400	2406	2425	rVB4	297342	972020	19.21%	1.928%
73	15.401	2425	2430	2434	rBV	140948	223906	4.42%	0.444%
74	15.450	2434	2438	2441	rBV	115310	173436	3.43%	0.344%
75	15.487	2441	2444	2447	rVB	84094	103517	2.05%	0.205%
76	15.554	2451	2455	2460	rBV	409561	615307	12.16%	1.220%
77	15.694	2472	2478	2482	rBV	678647	1074146	21.23%	2.130%
78	15.731	2482	2484	2490	rVB	328529	455269	9.00%	0.903%
79	15.816	2492	2498	2504	rBV2	77970	181924	3.60%	0.361%
80	15.895	2506	2511	2514	rVV	162387	304644	6.02%	0.604%
81	15.932	2514	2517	2527	rVB2	171406	277791	5.49%	0.551%
82	16.078	2536	2541	2553	rVB	91249	209456	4.14%	0.415%
83	16.188	2554	2559	2568	rBV6	47121	108131	2.14%	0.214%
84	16.572	2617	2622	2628	rVB3	59319	110534	2.18%	0.219%

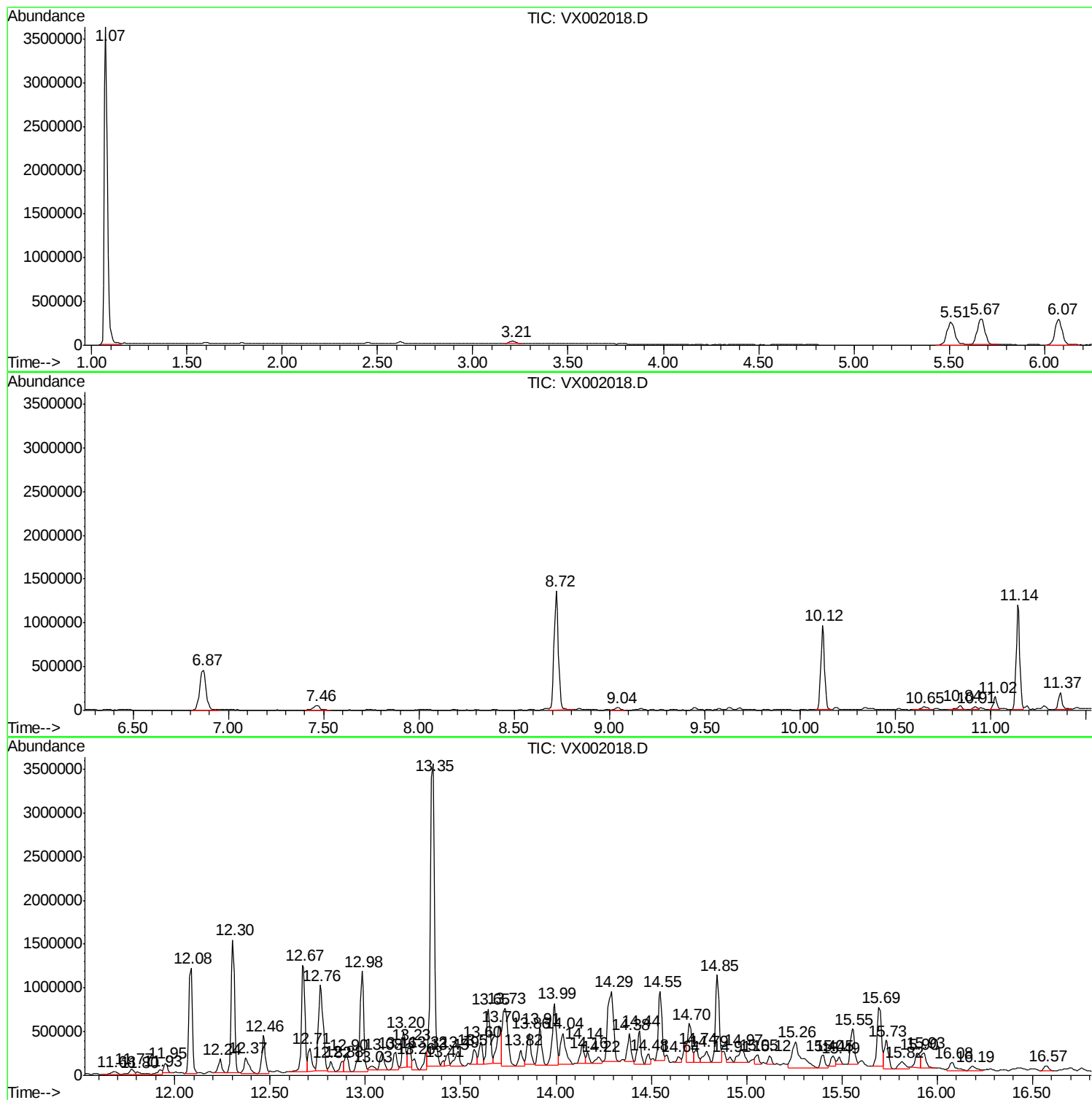
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Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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



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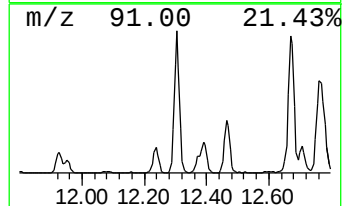
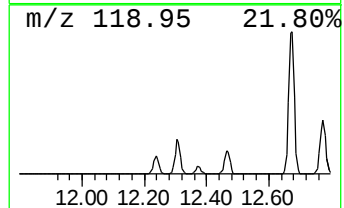
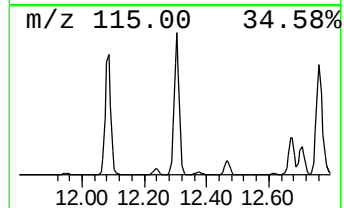
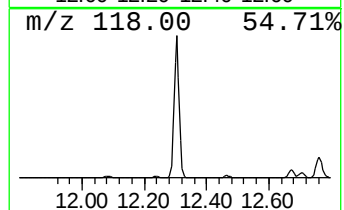
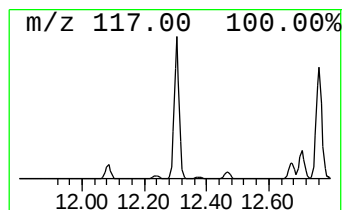
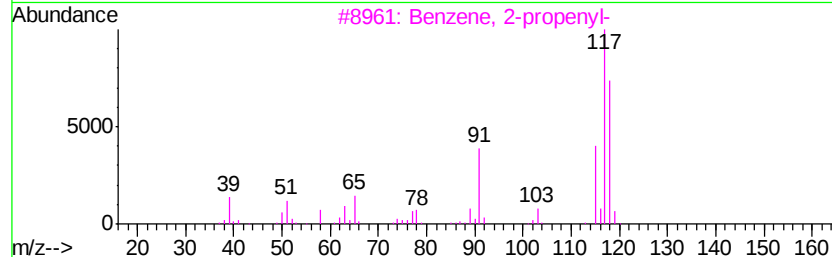
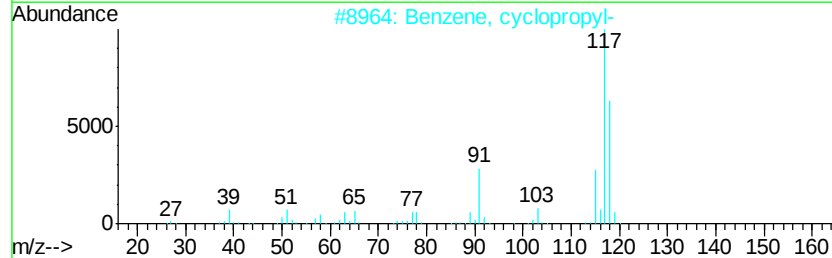
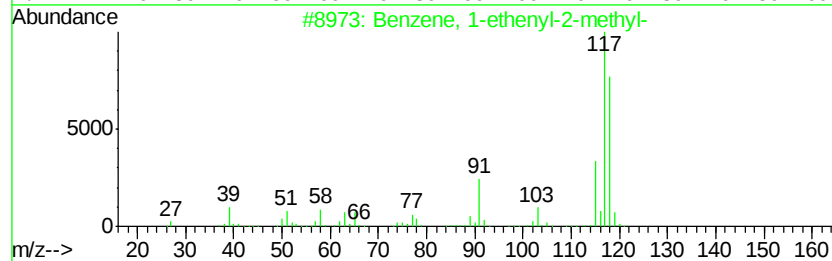
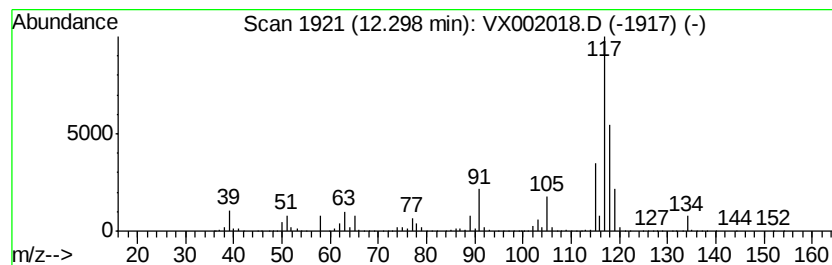
Instrument :
MSVOA_X
ClientSampled :
MW-17

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethenyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	60.68 ug/l	1820910	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 91
2		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
3		Benzene, 2-propenyl-	118 C9H10	000300-57-2 64
4		Indane	118 C9H10	000496-11-7 64
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 58



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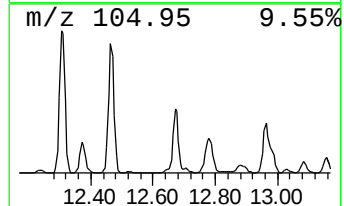
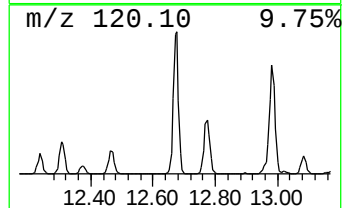
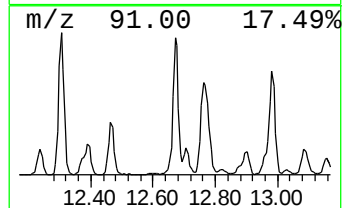
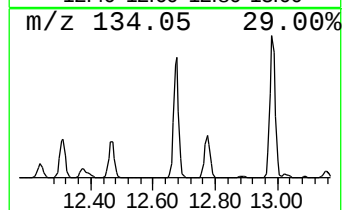
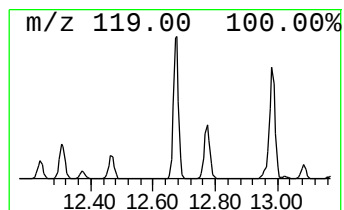
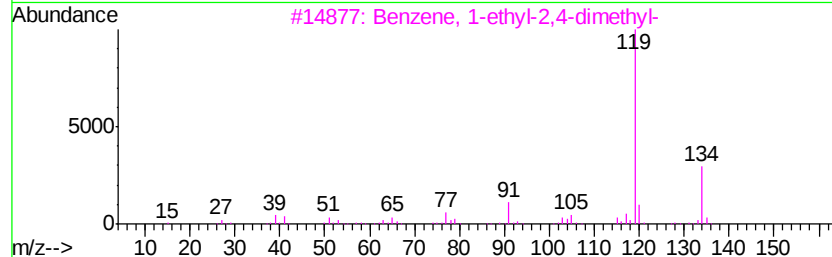
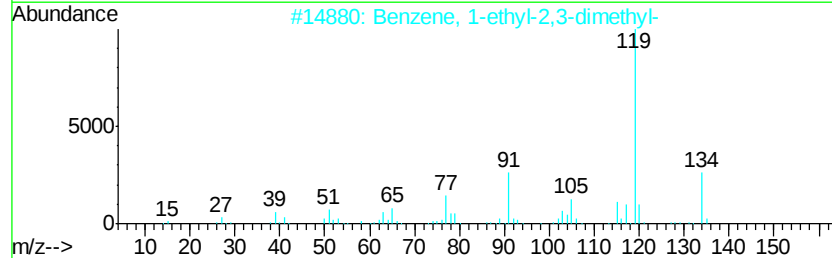
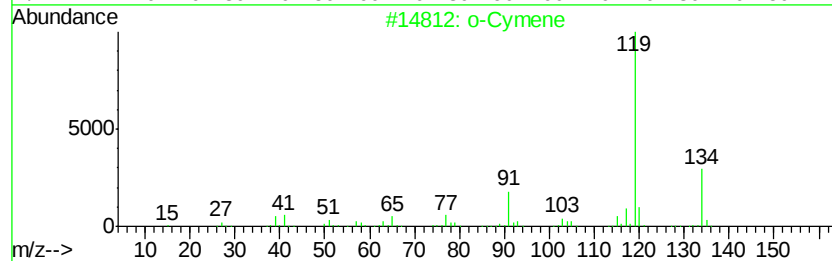
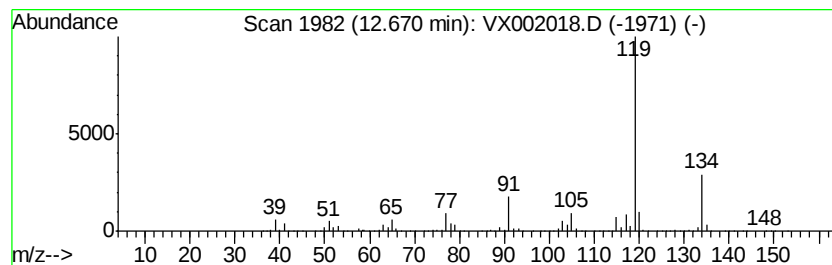
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	53.55 ug/l	1607130	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 95
2		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 95
3		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 94
4		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 94
5		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 94



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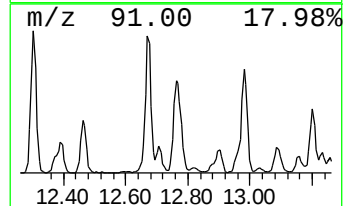
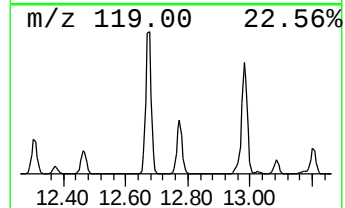
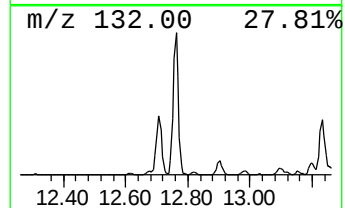
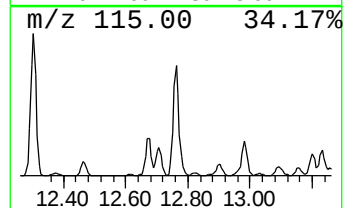
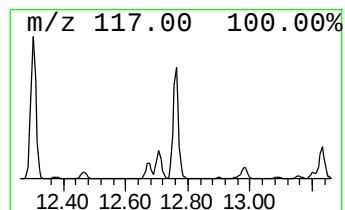
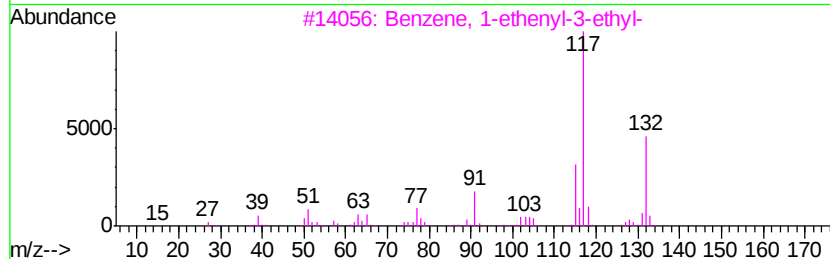
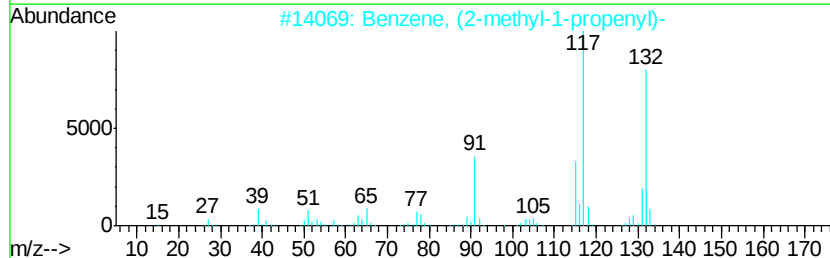
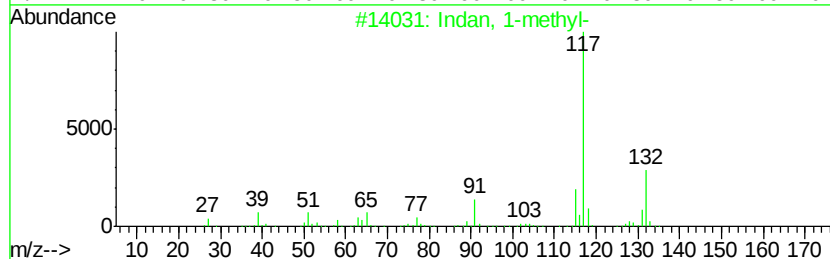
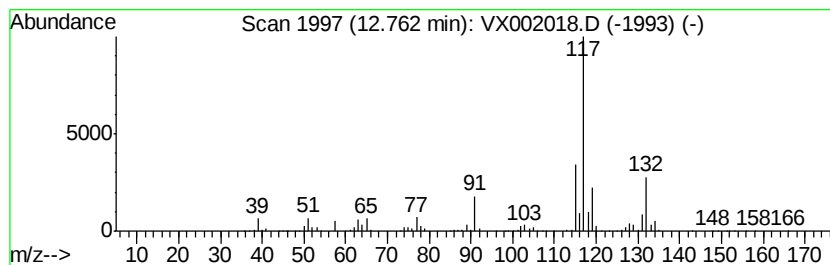
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	55.97 ug/l	1679740	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76



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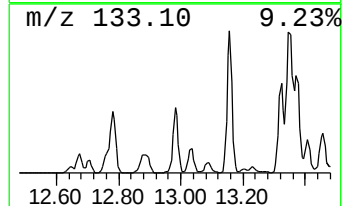
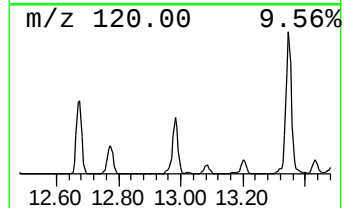
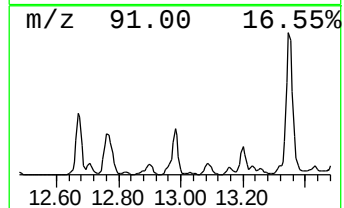
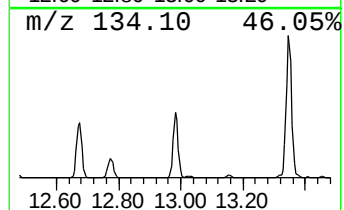
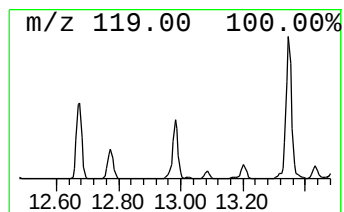
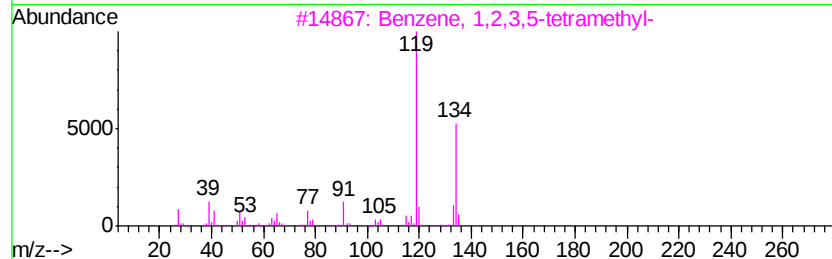
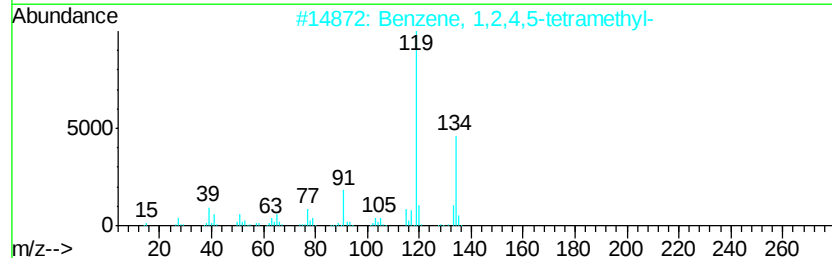
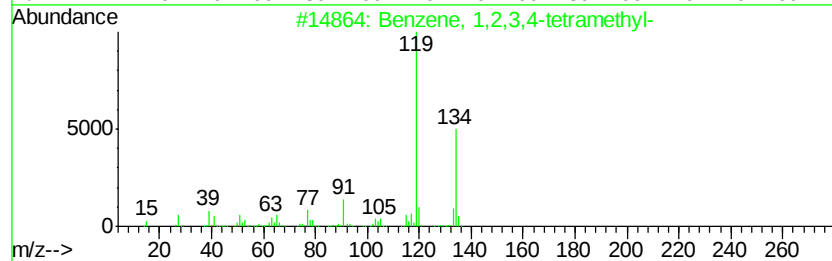
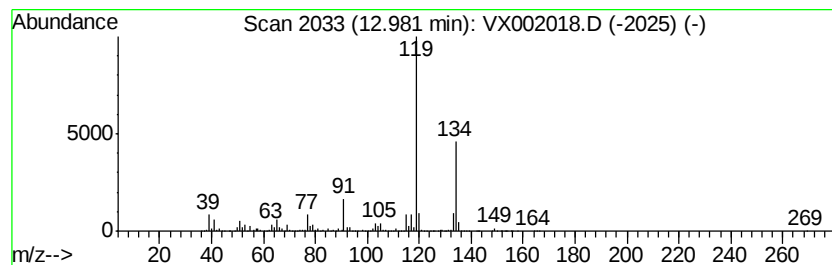
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	52.89 ug/l	1587170	1,4-Dichlorobenzene-d4	12.08

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052618\
Data File : VX002018.D
Acq On : 26 May 2018 14:48
Operator : JC/MD
Sample : J3064-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-17

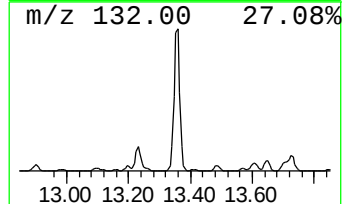
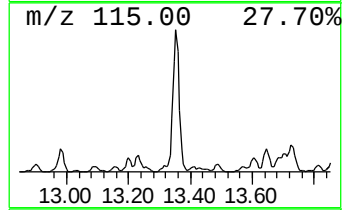
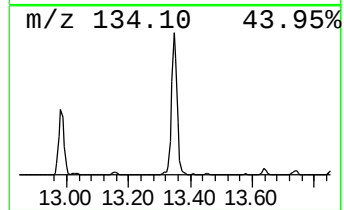
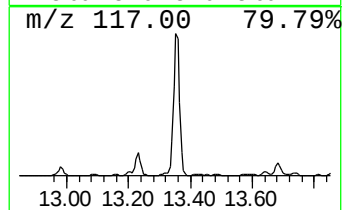
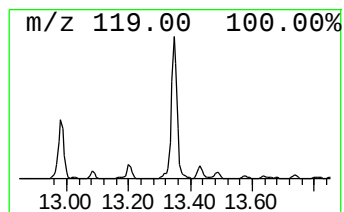
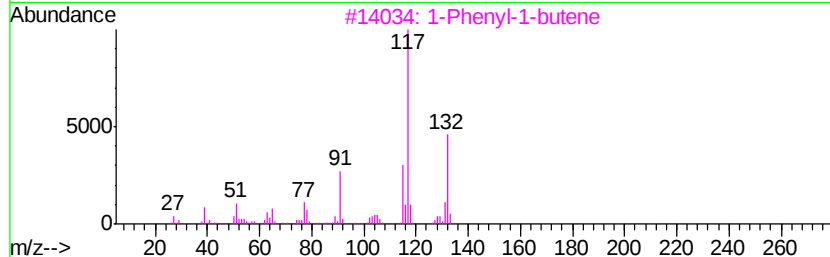
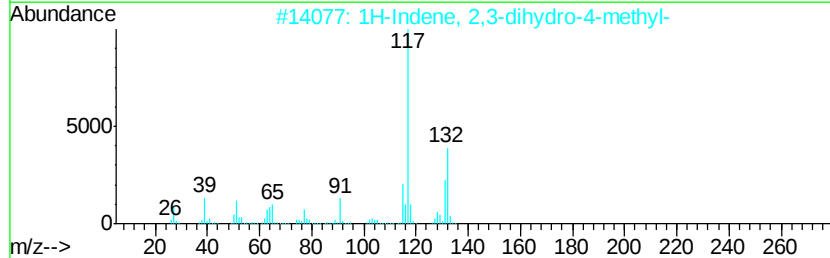
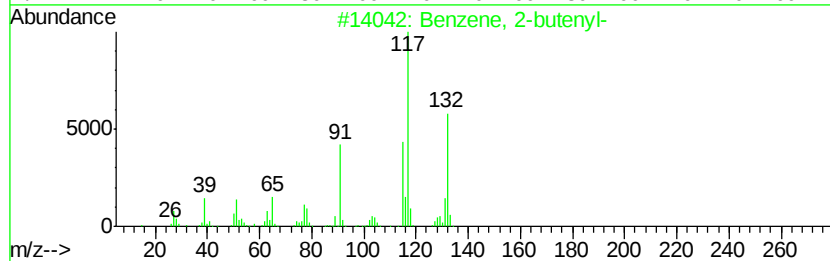
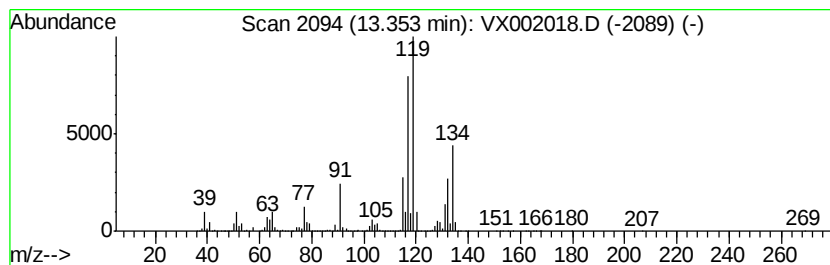
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-butenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	168.62 ug/l	5060290	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	86
4		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052618\
Data File : VX002018.D
Acq On : 26 May 2018 14:48
Operator : JC/MD
Sample : J3064-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-17

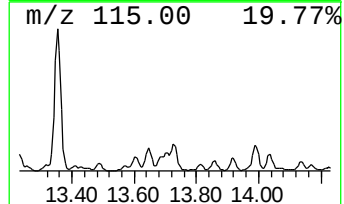
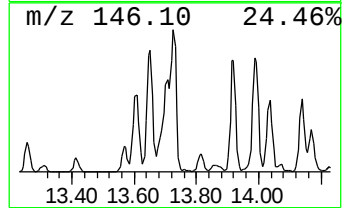
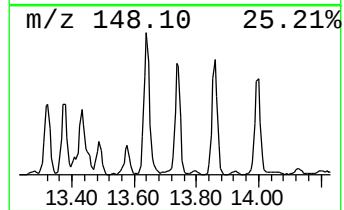
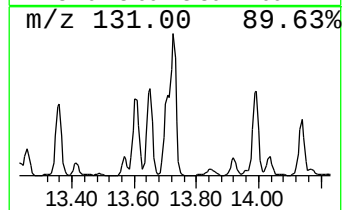
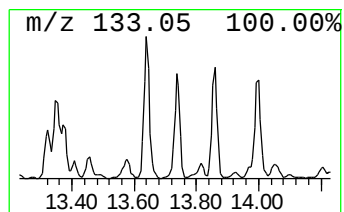
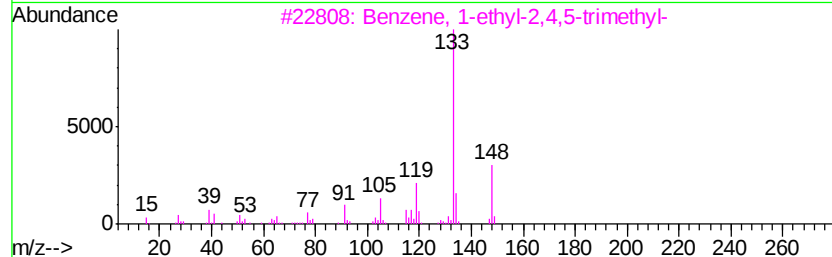
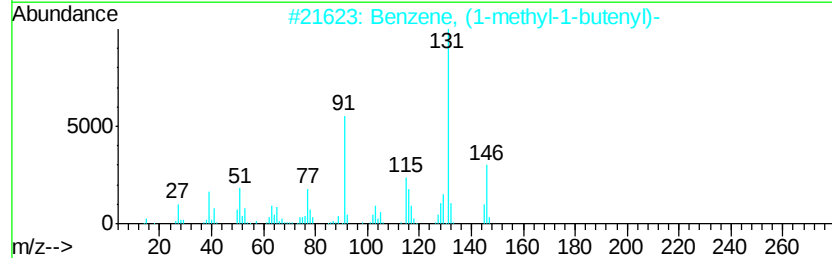
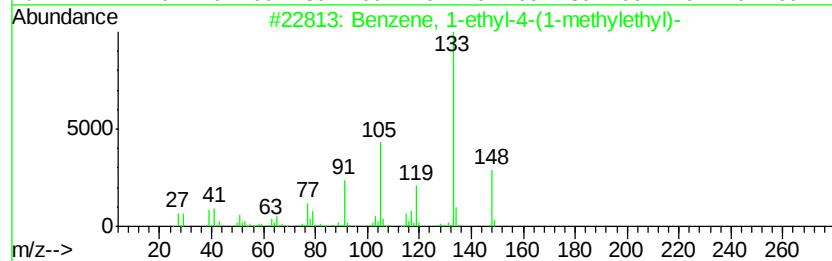
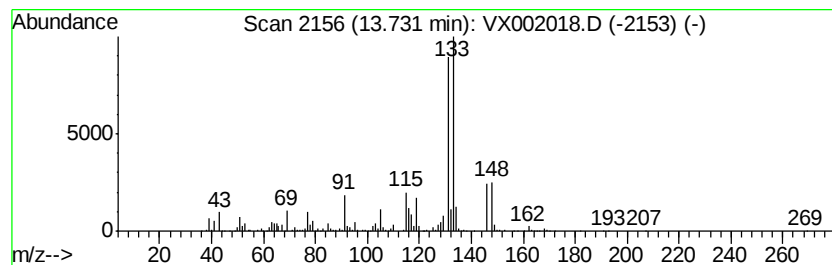
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	42.25 ug/l	1267920	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	50
2		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	45
3		Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	45
4		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	45
5		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	45



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052618\
Data File : VX002018.D
Acq On : 26 May 2018 14:48
Operator : JC/MD
Sample : J3064-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-17

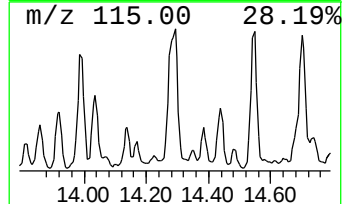
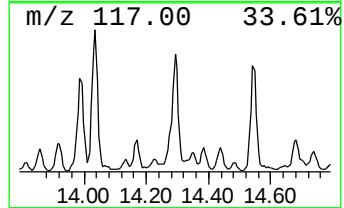
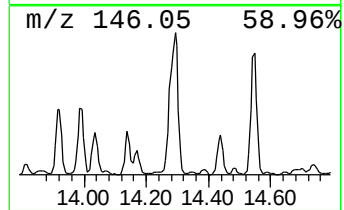
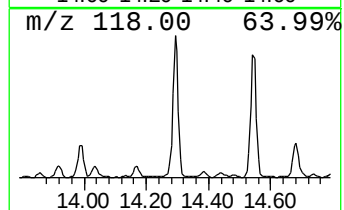
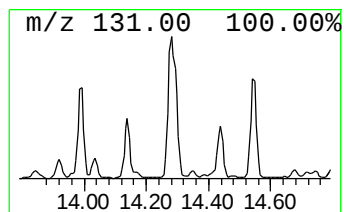
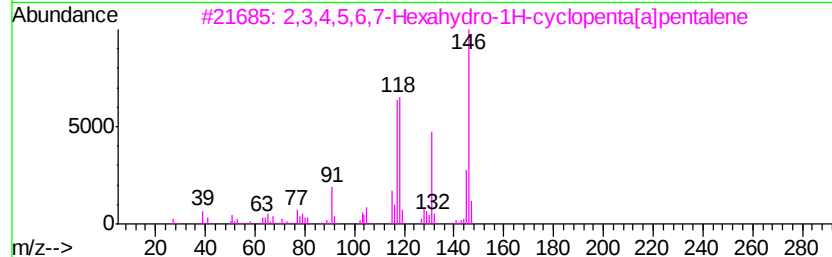
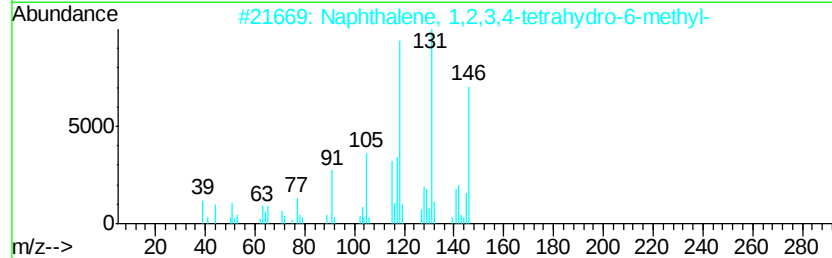
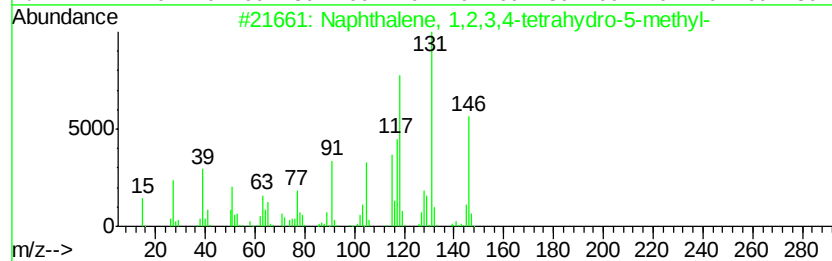
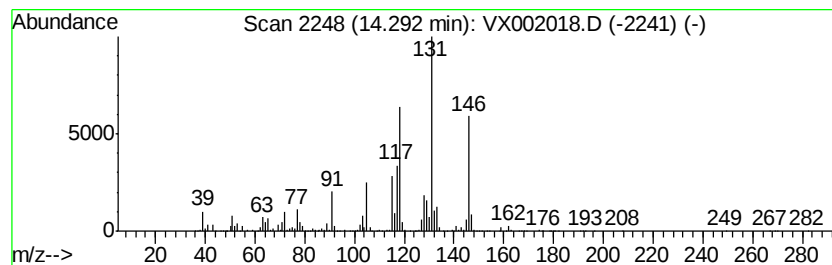
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	54.92 ug/l	1648250	1,4-Dichlorobenzene-d4	12.08

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	94
3		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	72
4		Benzene, (2-methyl-1-butenvl)-	146	C11H14	056253-64-6	70
5		Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	64



Data Path : W:\HPCHEM1\MSVOA X\DATA\VX052618\
Data File : VX002018.D
Acq On : 26 May 2018 14:48
Operator : JC/MD
Sample : J3064-05
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-17

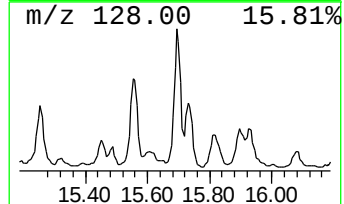
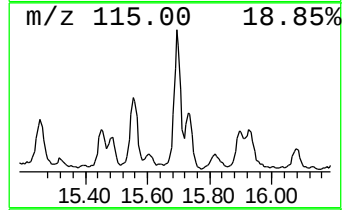
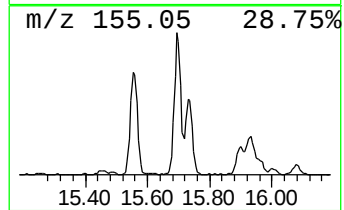
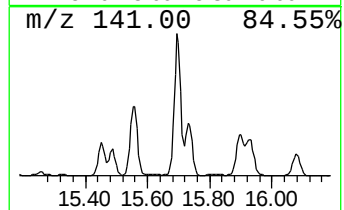
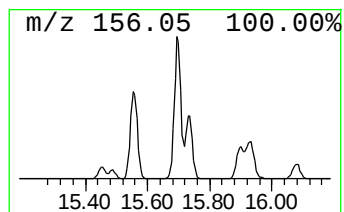
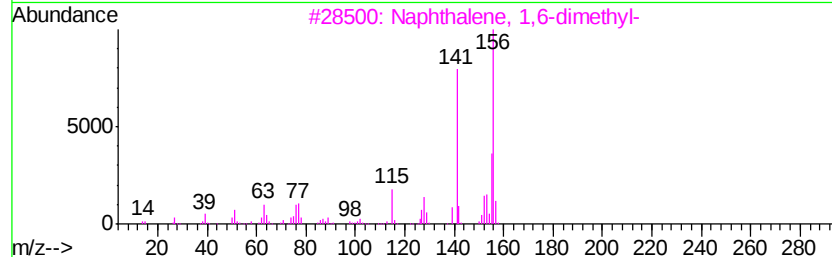
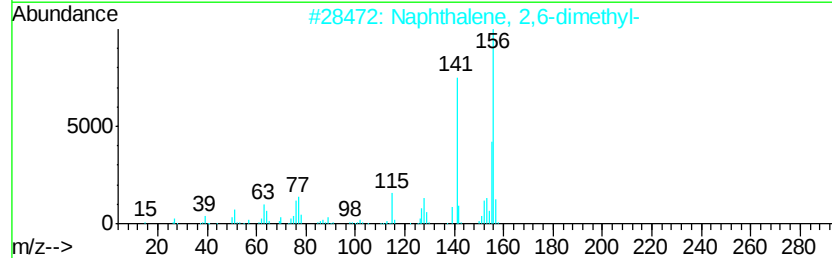
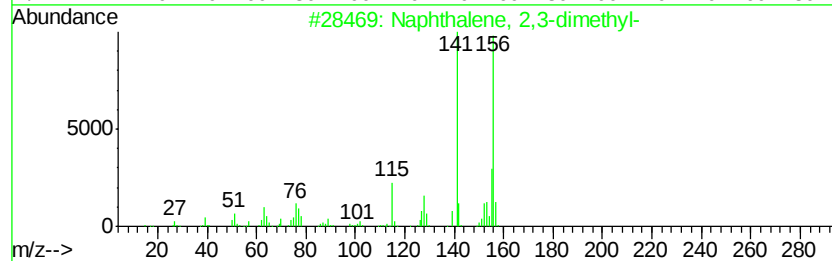
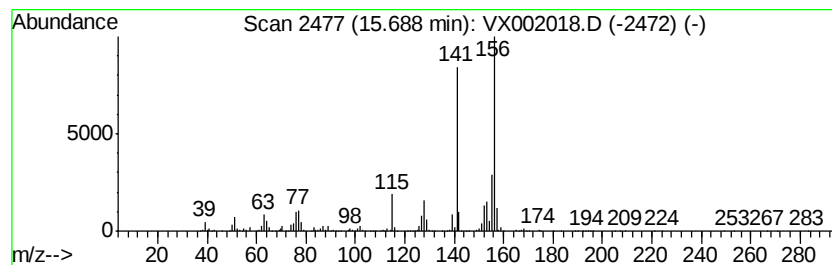
Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	35.79 ug/l	1074150	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4		Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : W:\HPCHEM1\MSVOA_X\DATA\VX052618\
Data File : VX002018.D
Acq On : 26 May 2018 14:48
Operator : JC/MD
Sample : J3064-05
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-17

Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X052418W.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...	12.30	60.7	ug/l	1820910	4	12.08	1500530	50.0
o-Cymene	12.67	53.5	ug/l	1607130	4	12.08	1500530	50.0
Indan, 1-methyl-	12.76	56.0	ug/l	1679740	4	12.08	1500530	50.0
Benzene, 1,2,3,4-...	12.98	52.9	ug/l	1587170	4	12.08	1500530	50.0
Benzene, 2-butenyl-	13.35	168.6	ug/l	5060290	4	12.08	1500530	50.0
Benzene, 1-ethyl-...	13.73	42.3	ug/l	1267920	4	12.08	1500530	50.0
Naphthalene, 1,2,...	14.29	54.9	ug/l	1648250	4	12.08	1500530	50.0
Naphthalene, 2,3-...	15.69	35.8	ug/l	1074150	4	12.08	1500530	50.0