

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111120\
 Data File : VD067599.D
 Acq On : 11 Nov 2020 13:46
 Operator : VA/SY
 Sample : L4702-04
 Misc : 6.22G/5.00ml/MSVOA D/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 M-99(0-10)

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_D\METHOD\82D110320S.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	7.920	1010	1020	1025	rBV	251254	610495	23.18%	1.370%
2	7.985	1025	1031	1043	rVB	449233	1000859	37.99%	2.246%
3	8.338	1081	1091	1101	rBV	198066	457424	17.36%	1.026%
4	8.867	1172	1181	1195	rBV	621462	1279874	48.59%	2.872%
5	10.344	1423	1432	1440	rBV	1003115	1806543	68.58%	4.054%
6	11.649	1647	1654	1664	rVB	968771	1725974	65.52%	3.873%
7	12.491	1792	1797	1803	rBV2	57080	96174	3.65%	0.216%
8	12.638	1815	1822	1833	rVB	832767	1419789	53.90%	3.186%
9	12.743	1835	1840	1843	rBV6	18133	29730	1.13%	0.067%
10	12.890	1859	1865	1868	rBV	68389	121126	4.60%	0.272%
11	12.967	1873	1878	1884	rBV3	101176	215059	8.16%	0.483%
12	13.126	1900	1905	1910	rVB	77262	115247	4.38%	0.259%
13	13.220	1912	1921	1924	rVV2	128735	374475	14.22%	0.840%
14	13.273	1924	1930	1939	rVB	911248	1563396	59.35%	3.508%
15	13.408	1944	1953	1957	rBV3	99269	282401	10.72%	0.634%
16	13.467	1959	1963	1968	rVV	133047	259250	9.84%	0.582%
17	13.520	1968	1972	1976	rVV2	86122	171469	6.51%	0.385%
18	13.579	1976	1982	1985	rVV	1080374	1920903	72.92%	4.311%
19	13.608	1985	1987	1997	rVB	830124	1083361	41.13%	2.431%
20	13.773	2004	2015	2019	rBV2	1055206	2278624	86.50%	5.113%
21	13.814	2019	2022	2033	rVV3	1076843	2223297	84.40%	4.989%
22	13.902	2033	2037	2043	rVV4	95315	157649	5.98%	0.354%
23	13.973	2043	2049	2054	rVV	445995	710719	26.98%	1.595%
24	14.037	2054	2060	2062	rVV	811545	1350081	51.25%	3.030%
25	14.061	2062	2064	2069	rVV	804831	1149017	43.62%	2.578%
26	14.120	2069	2074	2081	rVV	1553398	2359668	89.58%	5.295%
27	14.185	2081	2085	2086	rVV	38818	45815	1.74%	0.103%
28	14.226	2086	2092	2098	rVV2	446371	790724	30.02%	1.774%
29	14.279	2098	2101	2102	rVV	44802	54100	2.05%	0.121%
30	14.308	2102	2106	2110	rVV3	62646	110946	4.21%	0.249%
31	14.361	2110	2115	2120	rVV	439242	692684	26.30%	1.554%
32	14.443	2120	2129	2132	rVV	1122812	1802451	68.43%	4.045%
33	14.485	2132	2136	2140	rVV	1641022	2553325	96.93%	5.730%
34	14.526	2140	2143	2147	rVV	339624	500218	18.99%	1.122%

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35	14.567	2147	2150	2158	rVV2	200676	353658	13.43%	0.794%
36	14.626	2158	2160	2163	rVV2	55800	72767	2.76%	0.163%
37	14.667	2163	2167	2171	rVV	177811	263072	9.99%	0.590%
38	14.714	2171	2175	2179	rVV2	488564	801375	30.42%	1.798%
39	14.761	2179	2183	2187	rVB	322684	501689	19.05%	1.126%
40	14.826	2187	2194	2201	rBV2	1073404	2634194	100.00%	5.911%
41	14.884	2201	2204	2210	rVB	290625	439375	16.68%	0.986%
42	15.043	2217	2231	2236	rBV5	187850	739699	28.08%	1.660%
43	15.102	2236	2241	2253	rVV	459105	1162734	44.14%	2.609%
44	15.208	2254	2259	2267	rVB3	231765	451321	17.13%	1.013%
45	15.355	2280	2284	2286	rBV2	35324	52037	1.98%	0.117%
46	15.402	2286	2292	2300	rVV	778853	1399884	53.14%	3.141%
47	15.508	2304	2310	2315	rBV3	123572	215076	8.16%	0.483%
48	15.567	2315	2320	2322	rVV3	46311	77284	2.93%	0.173%
49	15.602	2322	2326	2335	rVB2	102833	189642	7.20%	0.426%
50	15.702	2336	2343	2349	rBV3	133447	297263	11.28%	0.667%
51	15.779	2349	2356	2360	rVB10	28806	64983	2.47%	0.146%
52	15.873	2367	2372	2383	rVB2	61529	157408	5.98%	0.353%
53	16.002	2388	2394	2400	rBV	85753	170015	6.45%	0.382%
54	16.073	2400	2406	2414	rVB5	62415	182205	6.92%	0.409%
55	16.149	2414	2419	2423	rBV5	36957	66201	2.51%	0.149%
56	16.202	2423	2428	2431	rBV2	47266	86791	3.29%	0.195%
57	16.284	2437	2442	2446	rVB5	54752	94453	3.59%	0.212%
58	16.343	2448	2452	2459	rVB4	61786	112137	4.26%	0.252%
59	16.496	2471	2478	2484	rBV	455071	936914	35.57%	2.102%
60	16.561	2484	2489	2502	rVV	376637	740150	28.10%	1.661%
61	16.702	2505	2513	2526	rVB	375587	988014	37.51%	2.217%

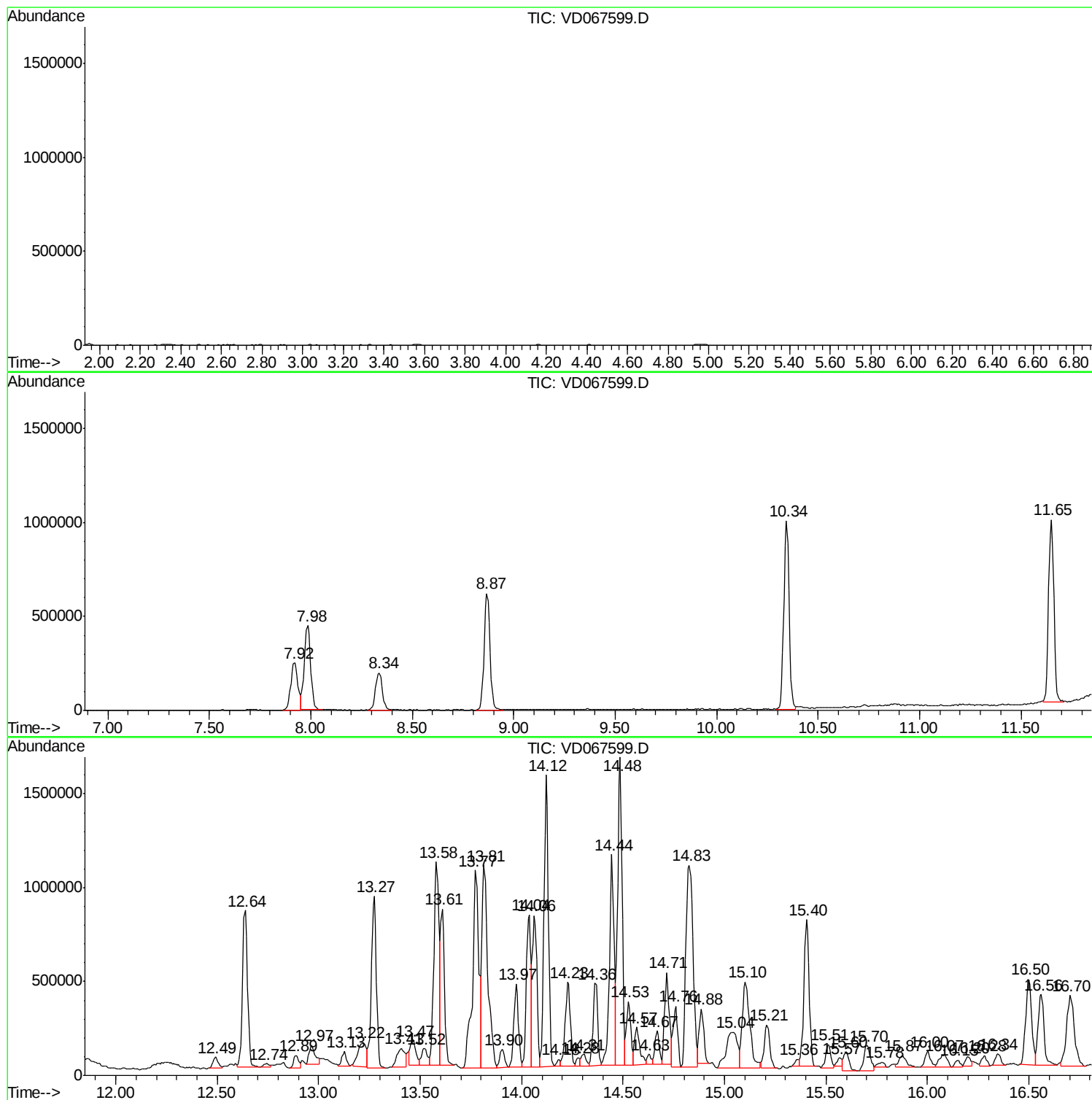
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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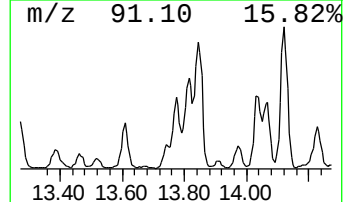
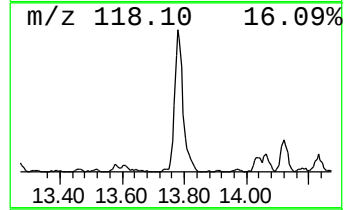
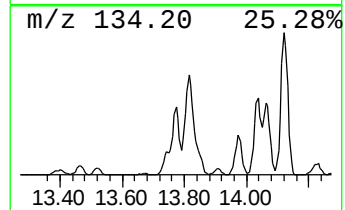
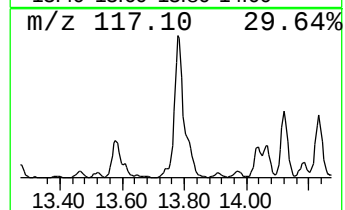
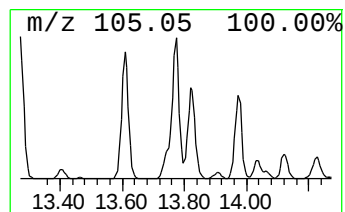
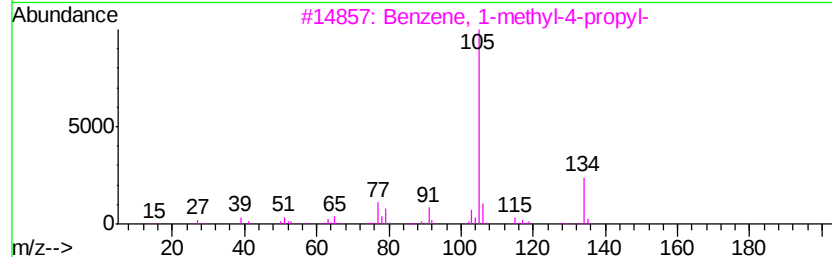
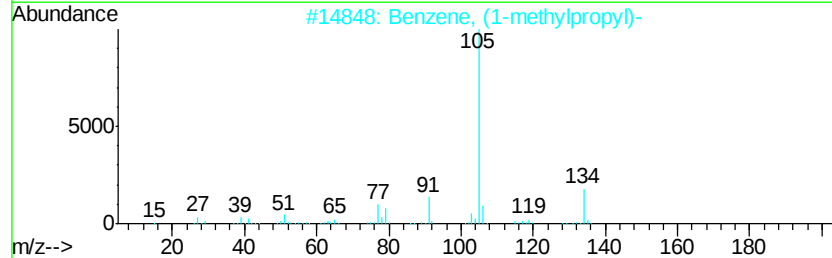
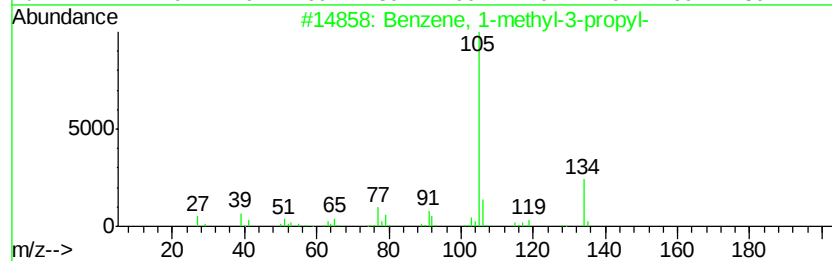
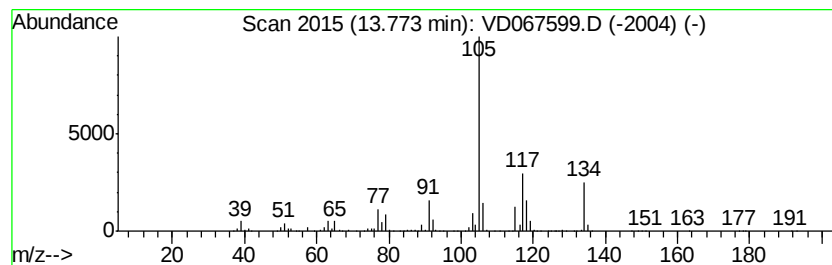
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	59.31 ug/l	2278620	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	64
2		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	64
3		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	60
4		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	55
5		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	52



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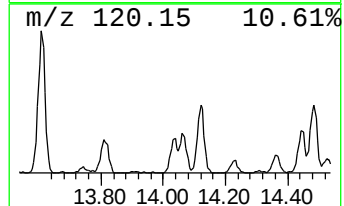
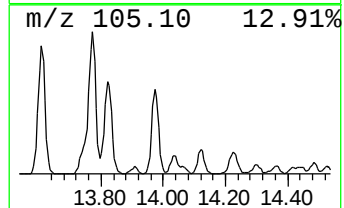
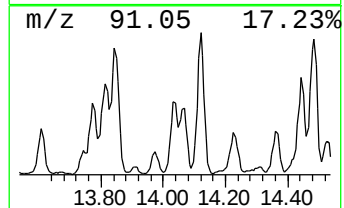
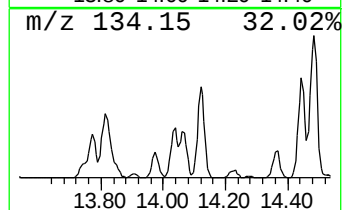
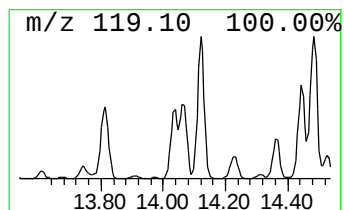
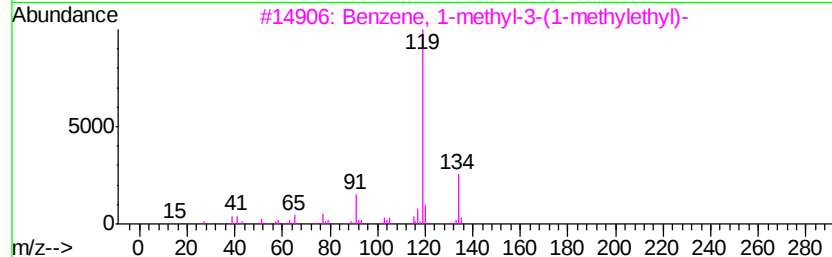
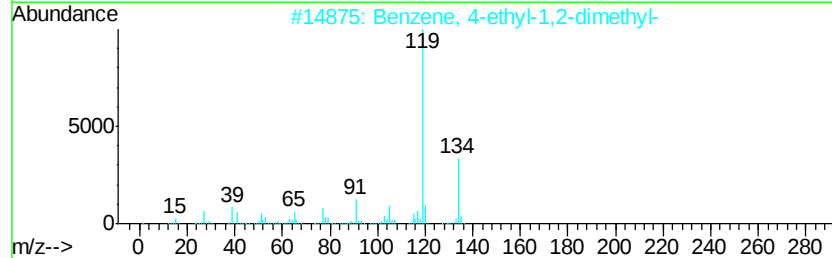
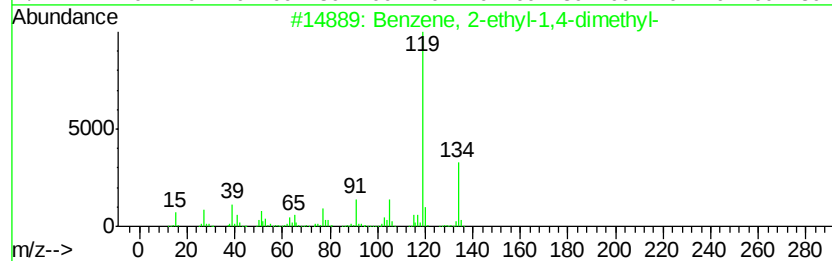
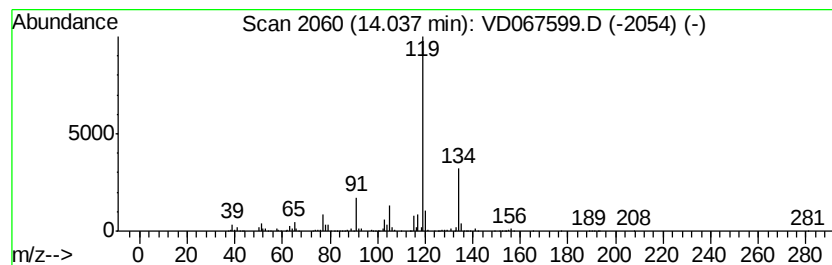
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.04	35.14 ug/l	1350080	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	94
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5		p-Cymene	134	C10H14	000099-87-6	93



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ALS Vial : 10 Sample Multiplier: 1

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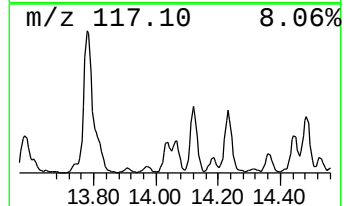
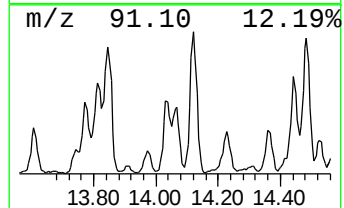
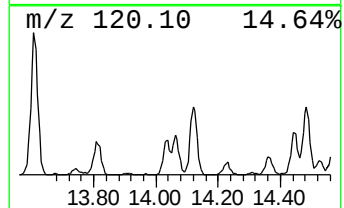
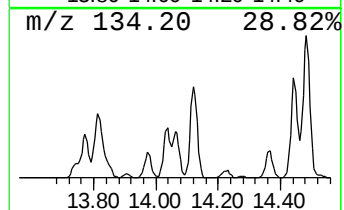
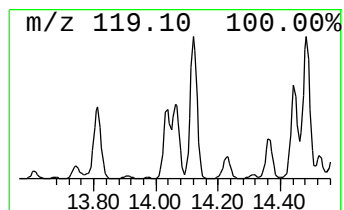
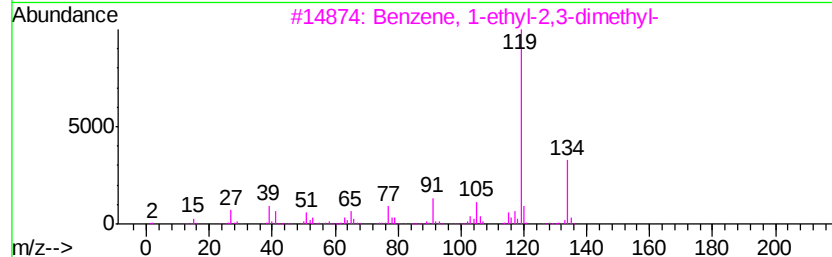
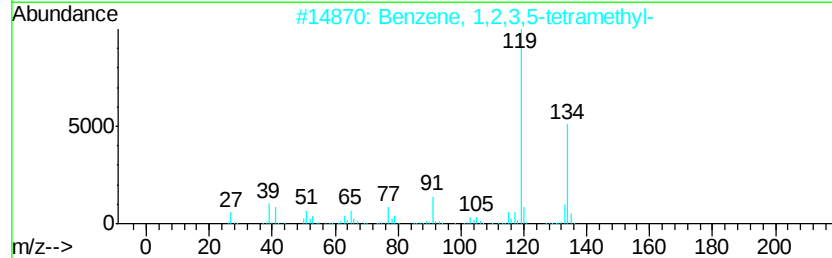
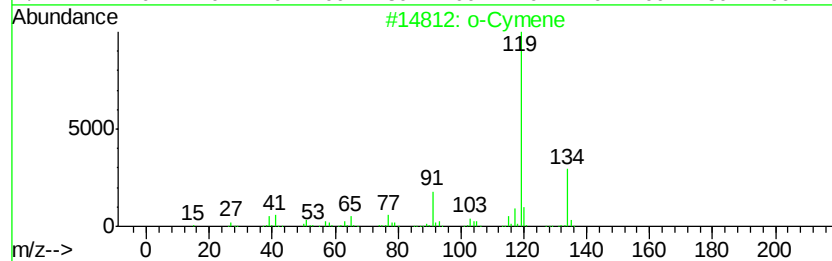
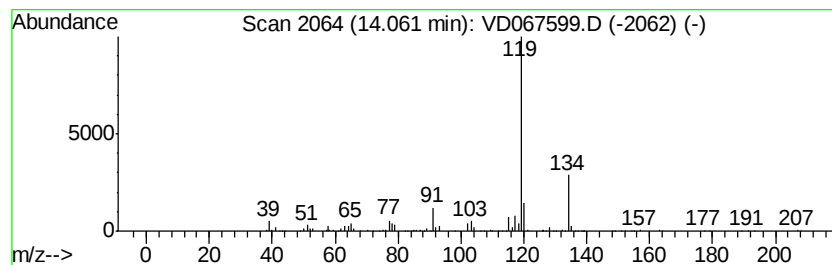
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.06	29.91 ug/l	1149020	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	93
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91



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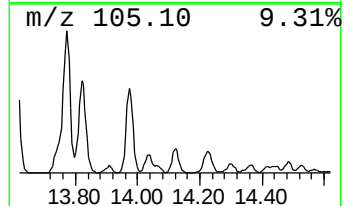
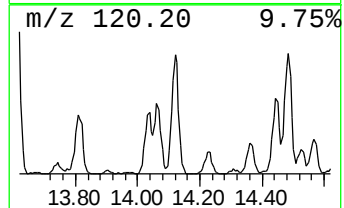
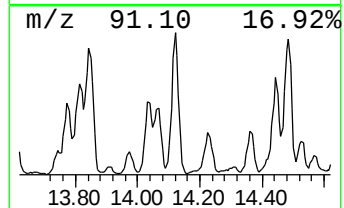
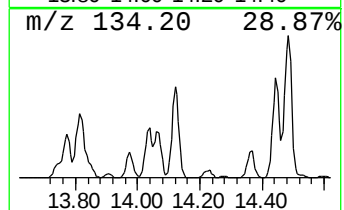
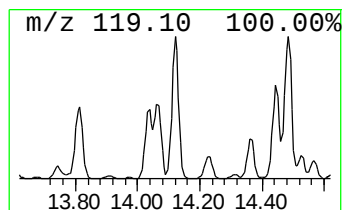
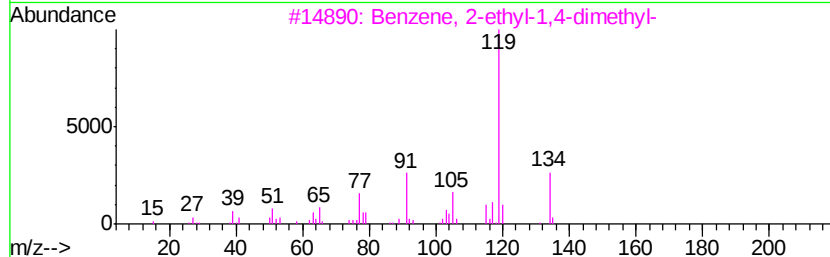
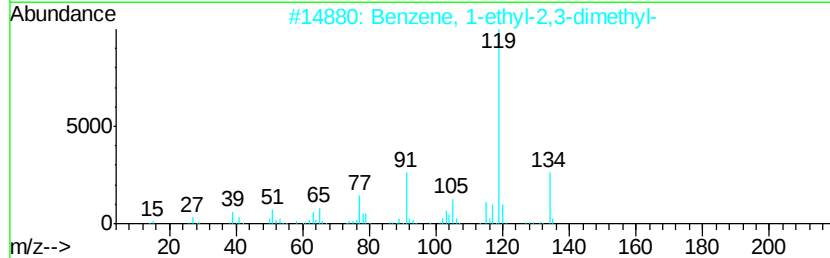
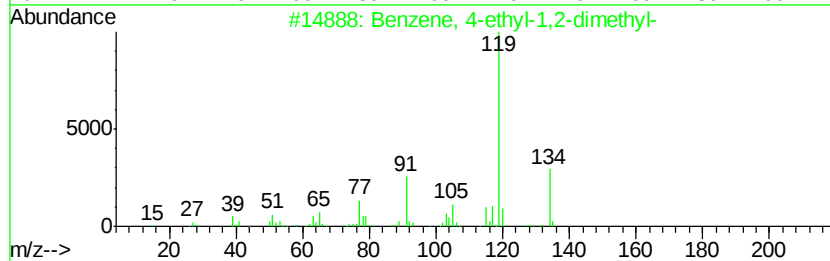
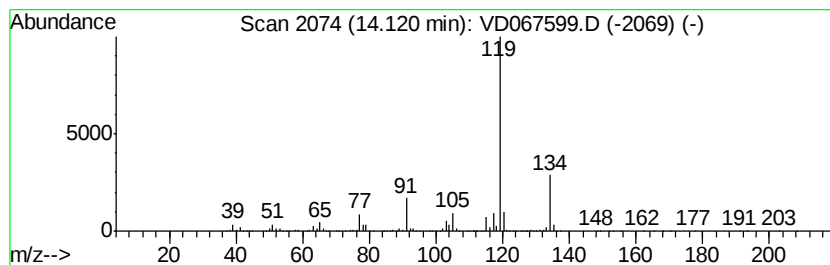
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.12	61.42 ug/l	2359670	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



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ClientSampleId :
M-99(0-10)

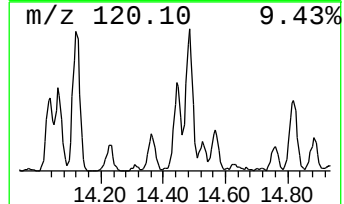
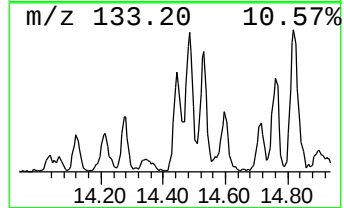
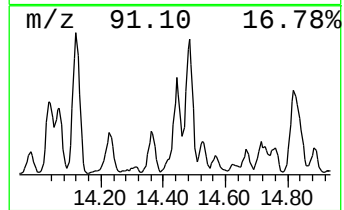
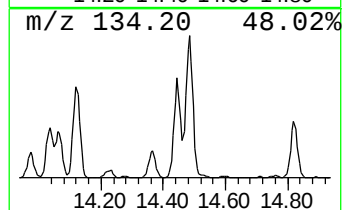
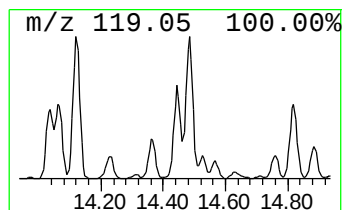
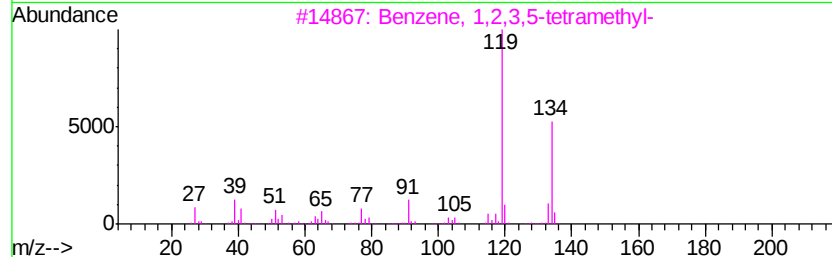
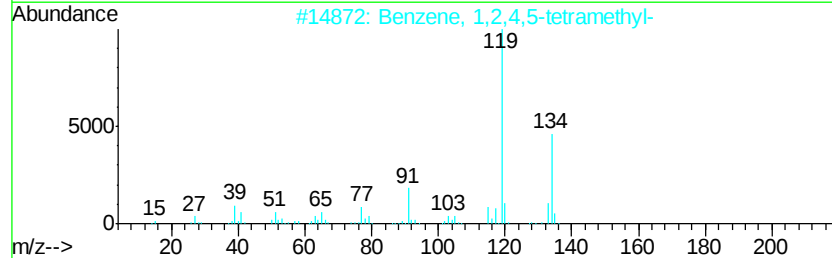
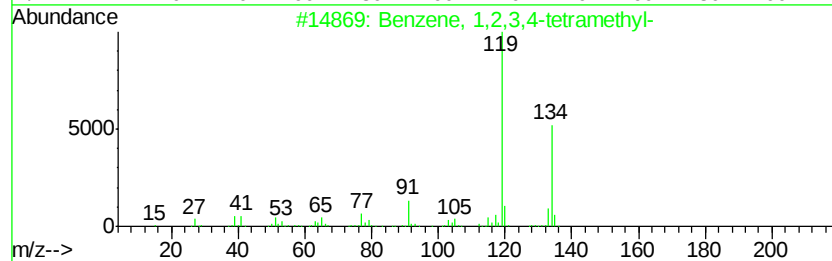
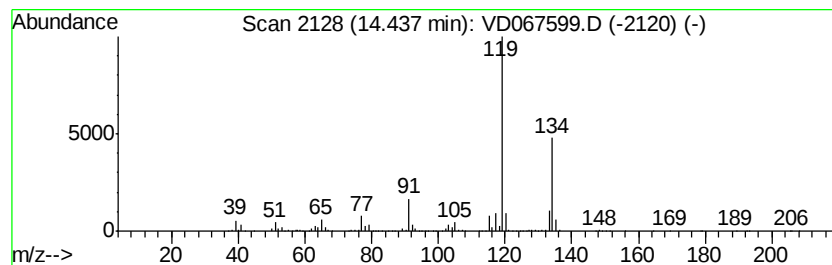
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.44	46.92 ug/l	1802450	1,4-Dichlorobenzene-d4	13.58

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	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5	o-Cymene	134	C10H14	000527-84-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111120\
Data File : VD067599.D
Acq On : 11 Nov 2020 13:46
Operator : VA/SY
Sample : L4702-04
Misc : 6.22G/5.00ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
M-99(0-10)

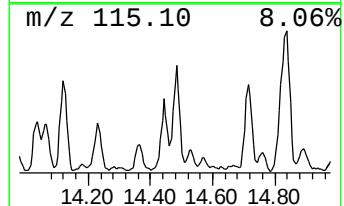
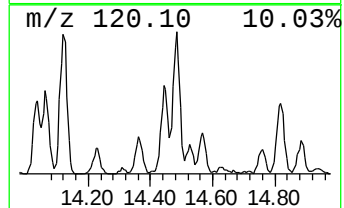
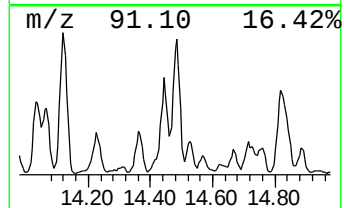
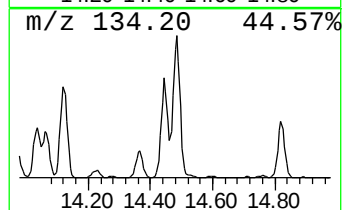
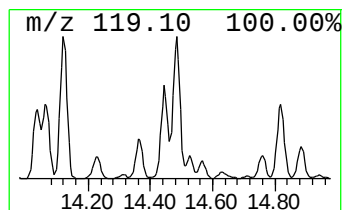
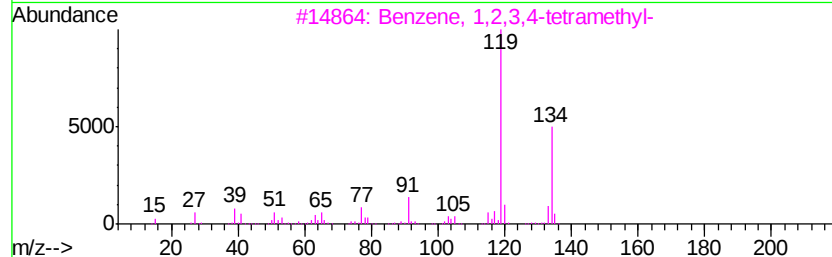
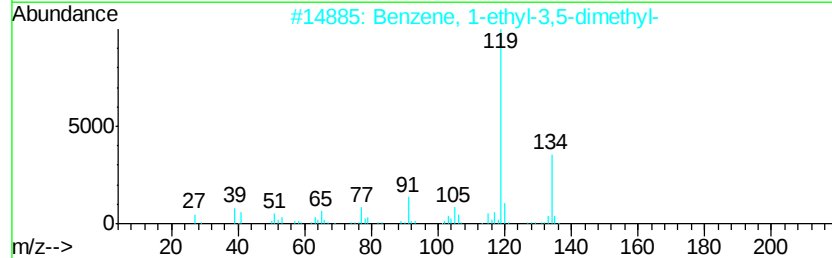
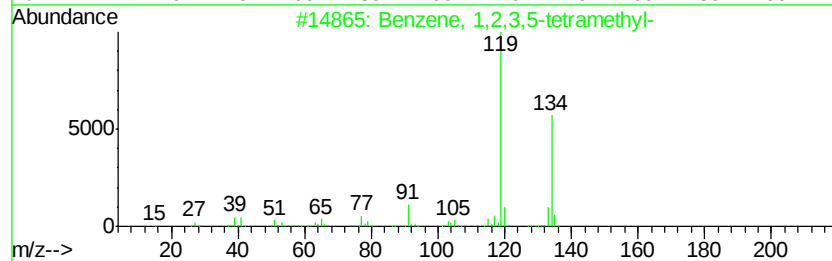
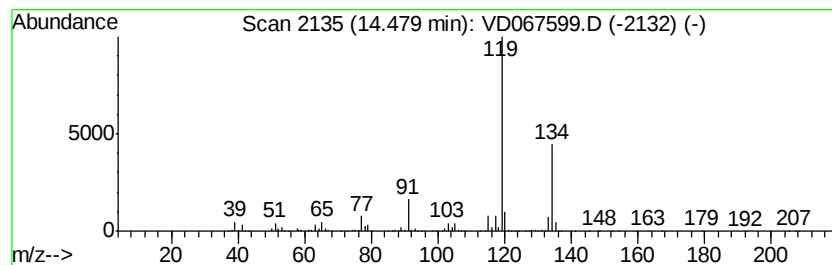
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	66.46 ug/l	2553330	1,4-Dichlorobenzene-d4	13.58

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
5	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111120\
Data File : VD067599.D
Acq On : 11 Nov 2020 13:46
Operator : VA/SY
Sample : L4702-04
Misc : 6.22G/5.00ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
M-99(0-10)

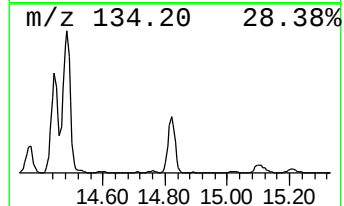
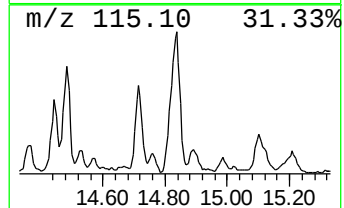
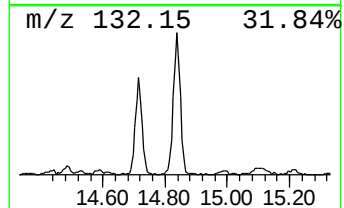
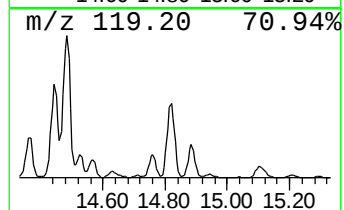
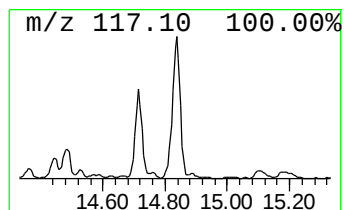
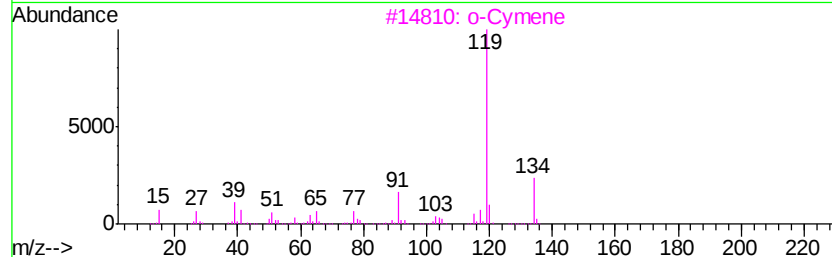
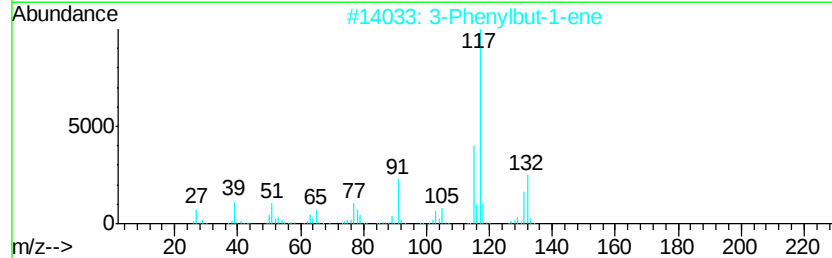
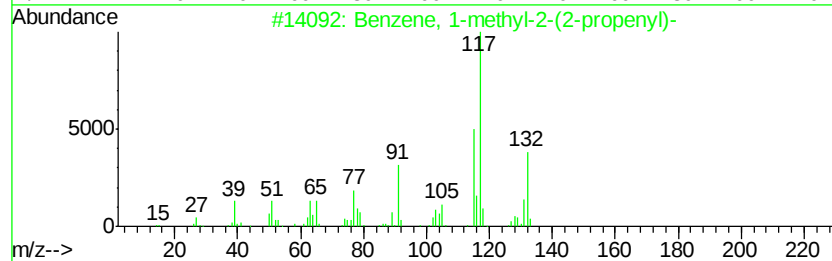
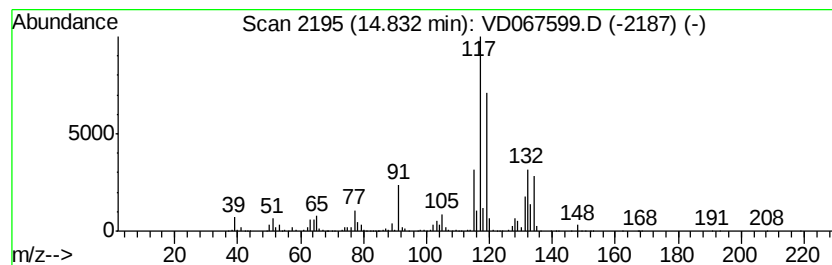
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	68.57 ug/l	2634190	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	80
3		o-Cymene	134	C10H14	000527-84-4	60
4		p-Cymene	134	C10H14	000099-87-6	60
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111120\
Data File : VD067599.D
Acq On : 11 Nov 2020 13:46
Operator : VA/SY
Sample : L4702-04
Misc : 6.22G/5.00ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
M-99(0-10)

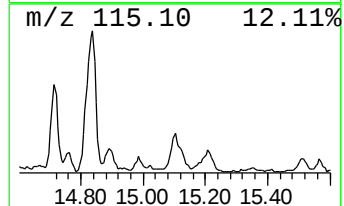
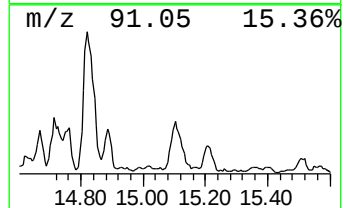
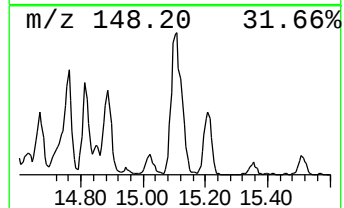
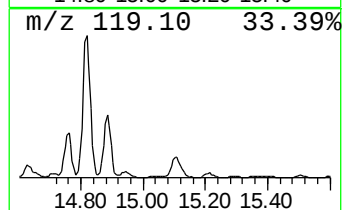
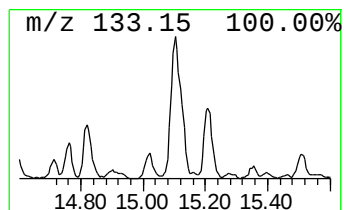
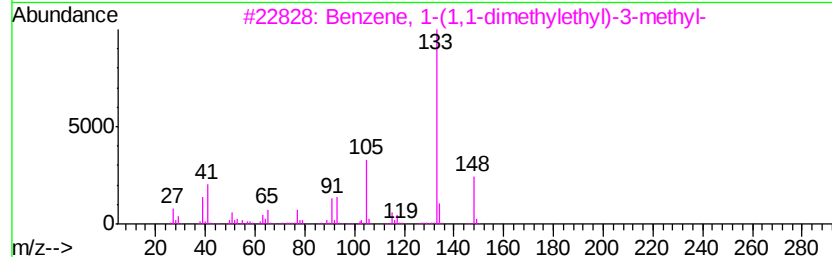
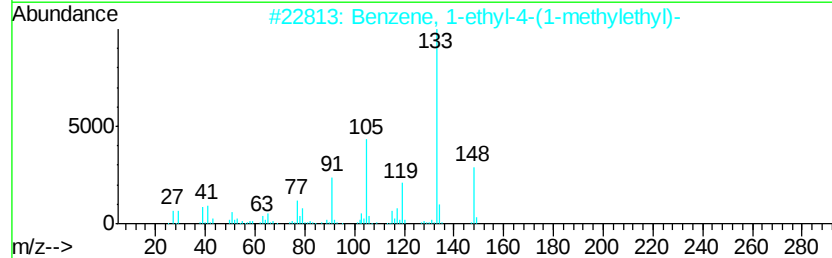
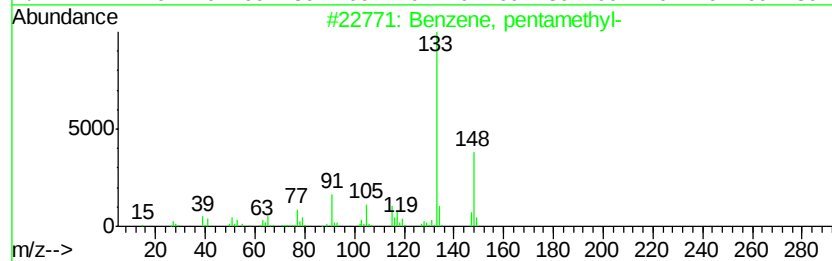
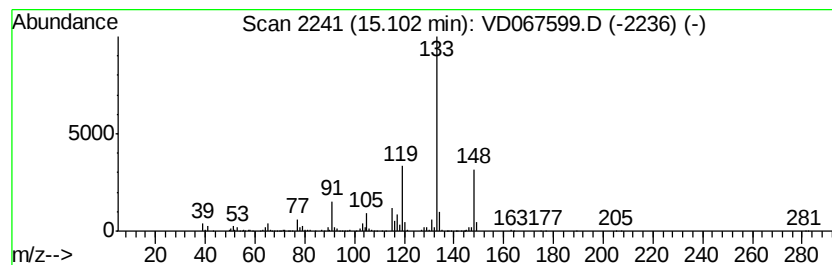
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, pentamethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	30.27 ug/l	1162730	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, pentamethyl-	148	C11H16	000700-12-9	91
2		Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	90
3		Benzene, 1-(1,1-dimethylethyl)-3-methyl-	148	C11H16	001075-38-3	81
4		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	80
5		Benzene, 2,4-dimethyl-1-(1-methylethyl)-	148	C11H16	004706-89-2	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111120\
Data File : VD067599.D
Acq On : 11 Nov 2020 13:46
Operator : VA/SY
Sample : L4702-04
Misc : 6.22G/5.00ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

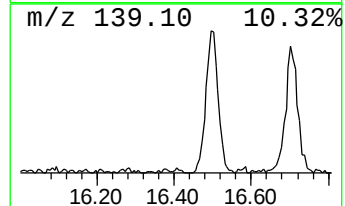
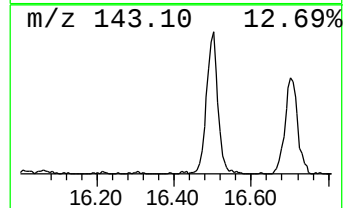
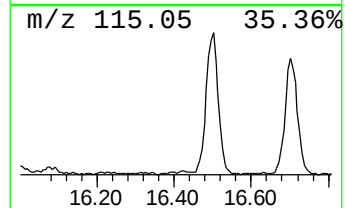
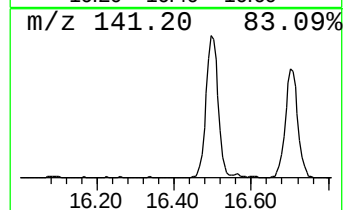
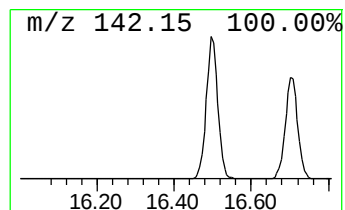
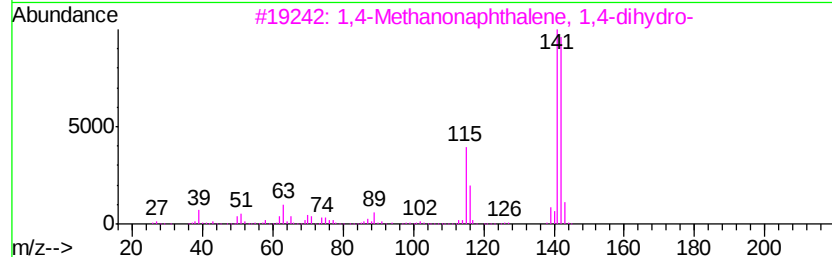
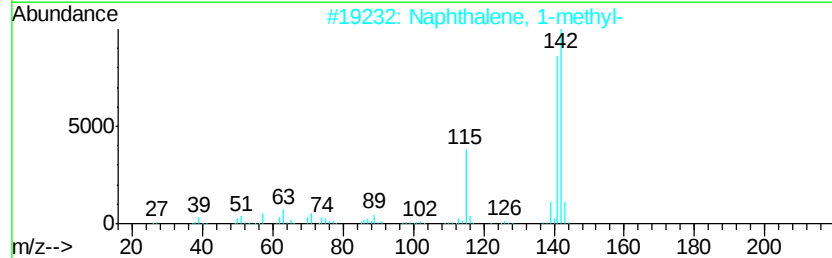
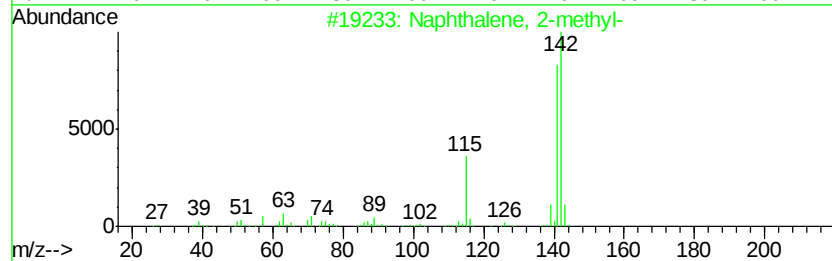
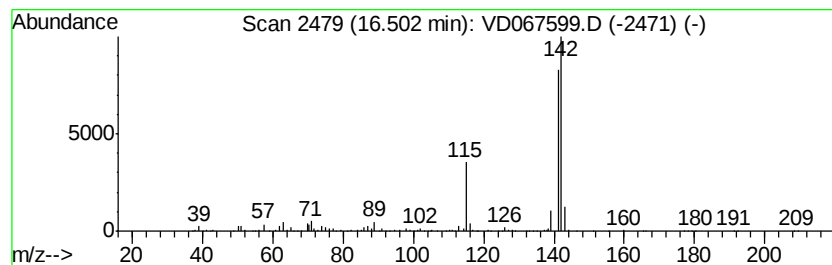
Instrument :
MSVOA_D
ClientSampleId :
M-99(0-10)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	24.39 ug/l	936914	1,4-Dichlorobenzene-d4	13.58
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		1,4-Methanonaphthalene, 1,4-dihydro...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 91
5		Naphthalene, 1-(2-hydroxypropyl)	186 C13H14O	027653-13-0 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD111120\
Data File : VD067599.D
Acq On : 11 Nov 2020 13:46
Operator : VA/SY
Sample : L4702-04
Misc : 6.22G/5.00ml/MSVOA D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
M-99(0-10)

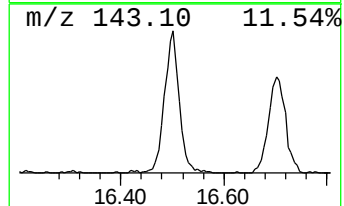
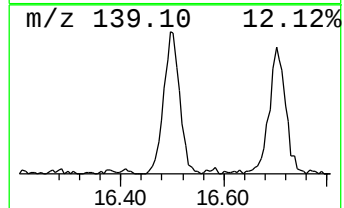
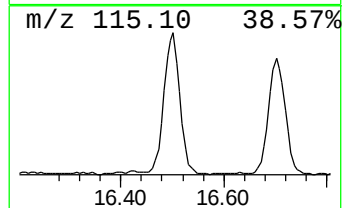
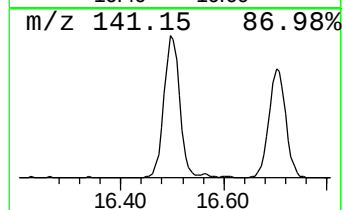
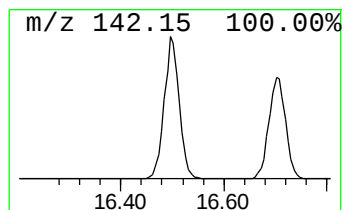
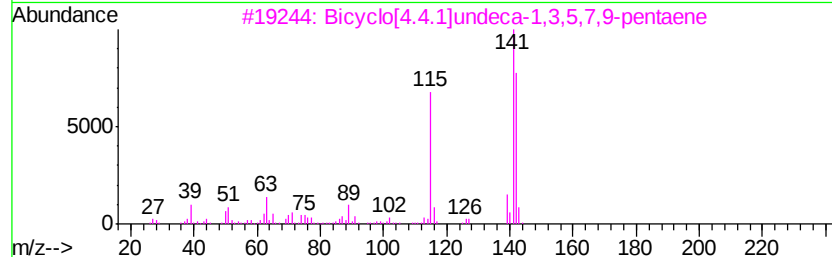
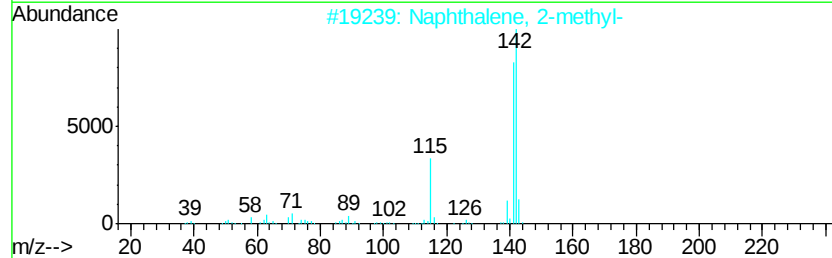
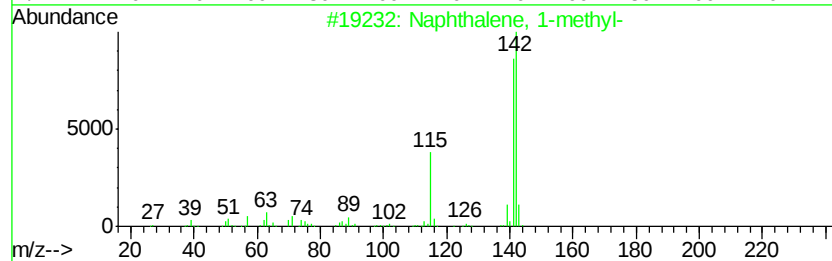
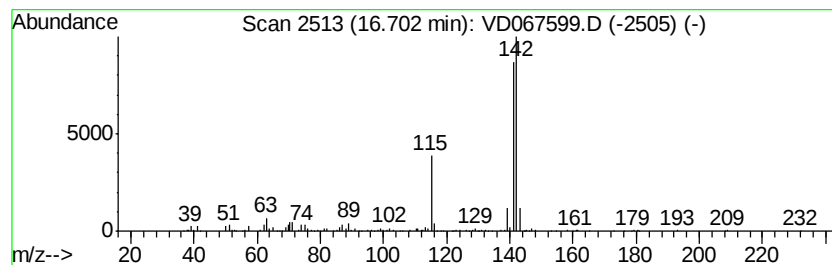
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.M
Quant Title : SW846 8260

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

Peak Number 10 Bicyclo[4.4.1]undeca-1,3,5,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	25.72 ug/l	988014	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	97
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
3		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	83



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD111120\
Data File : VD067599.D
Acq On : 11 Nov 2020 13:46
Operator : VA/SY
Sample : L4702-04
Misc : 6.22G/5.00ml/MSVOA_D/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
M-99(0-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110320S.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-methyl...	13.77	59.3	ug/l	2278620	4	13.58	1920900	50.0
Benzene, 2-ethyl...	14.04	35.1	ug/l	1350080	4	13.58	1920900	50.0
o-Cymene	14.06	29.9	ug/l	1149020	4	13.58	1920900	50.0
Benzene, 4-ethyl...	14.12	61.4	ug/l	2359670	4	13.58	1920900	50.0
Benzene, 1,2,3,4...	14.44	46.9	ug/l	1802450	4	13.58	1920900	50.0
Benzene, 1,2,3,5...	14.48	66.5	ug/l	2553330	4	13.58	1920900	50.0
Benzene, 1-methyl...	14.83	68.6	ug/l	2634190	4	13.58	1920900	50.0
Benzene, pentamet...	15.10	30.3	ug/l	1162730	4	13.58	1920900	50.0
Naphthalene, 1-me...	16.50	24.4	ug/l	936914	4	13.58	1920900	50.0
Bicyclo[4.4.1]und...	16.70	25.7	ug/l	988014	4	13.58	1920900	50.0