

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD092520\
 Data File : VD066894.D
 Acq On : 26 Sep 2020 02:11
 Operator : VA/SY
 Sample : L4154-04
 Misc : 5.01G/5.00ml/MSVOA D/SOIL
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :

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\82D092220S.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	4.967	505	518	531	rBV3	27761	99077	3.24%	0.210%
2	5.997	685	693	700	rBV4	10563	31380	1.03%	0.066%
3	7.191	887	896	905	rBV2	51390	146647	4.80%	0.310%
4	7.914	1009	1019	1024	rBV	352199	851630	27.88%	1.802%
5	7.979	1024	1030	1041	rVB	679254	1497811	49.04%	3.169%
6	8.326	1072	1089	1100	rBV2	302877	807164	26.43%	1.708%
7	8.861	1169	1180	1196	rBV	968314	1925763	63.05%	4.075%
8	9.350	1249	1263	1270	rBV5	35548	90510	2.96%	0.192%
9	9.973	1364	1369	1373	rBV4	19448	38875	1.27%	0.082%
10	10.114	1381	1393	1396	rBV6	21446	47172	1.54%	0.100%
11	10.338	1422	1431	1440	rBV	1551625	2837640	92.91%	6.004%
12	10.514	1455	1461	1469	rVB3	75544	150626	4.93%	0.319%
13	10.679	1483	1489	1492	rBV3	33100	63064	2.06%	0.133%
14	10.720	1492	1496	1500	rVV2	55984	104510	3.42%	0.221%
15	10.773	1501	1505	1511	rVV5	31826	63535	2.08%	0.134%
16	10.867	1515	1521	1525	rVB6	20498	40761	1.33%	0.086%
17	10.949	1530	1535	1540	rVB3	40154	64372	2.11%	0.136%
18	11.049	1547	1552	1554	rBV3	20664	34370	1.13%	0.073%
19	11.085	1554	1558	1561	rVB2	30058	47725	1.56%	0.101%
20	11.191	1572	1576	1579	rVB	53753	72129	2.36%	0.153%
21	11.244	1579	1585	1590	rVB	118081	202870	6.64%	0.429%
22	11.349	1598	1603	1607	rBV6	54729	91849	3.01%	0.194%
23	11.426	1608	1616	1628	rVB6	126317	416865	13.65%	0.882%
24	11.526	1628	1633	1639	rBV	108901	171558	5.62%	0.363%
25	11.644	1645	1653	1664	rBV	1446831	2569156	84.12%	5.436%
26	11.744	1664	1670	1673	rBV2	53586	93553	3.06%	0.198%
27	11.773	1673	1675	1679	rVB4	34943	40651	1.33%	0.086%
28	11.855	1679	1689	1699	rBV4	468847	1071386	35.08%	2.267%
29	11.991	1707	1712	1718	rBV9	24830	56982	1.87%	0.121%
30	12.079	1723	1727	1731	rVB2	32678	53318	1.75%	0.113%
31	12.149	1731	1739	1741	rBV4	188143	339103	11.10%	0.717%
32	12.267	1754	1759	1764	rVB	178904	271652	8.89%	0.575%
33	12.355	1764	1774	1775	rBV4	171087	318527	10.43%	0.674%
34	12.379	1775	1778	1785	rVB4	166875	365512	11.97%	0.773%

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35	12.473	1790	1794	1796	rBV4	41018	65567	2.15%	0.139%
36	12.555	1805	1808	1810	rBV2	67471	75868	2.48%	0.161%
37	12.579	1810	1812	1816	rVB2	136547	164827	5.40%	0.349%
38	12.632	1816	1821	1826	rBV	1247229	2019013	66.11%	4.272%
39	12.714	1831	1835	1840	rBV6	46449	72641	2.38%	0.154%
40	12.796	1844	1849	1856	rVB3	274355	531284	17.39%	1.124%
41	12.879	1856	1863	1867	rBV5	245363	504676	16.52%	1.068%
42	12.949	1867	1875	1883	rBV2	1166600	2002667	65.57%	4.237%
43	13.120	1896	1904	1911	rVV4	184118	699855	22.91%	1.481%
44	13.185	1911	1915	1922	rVV2	484867	821821	26.91%	1.739%
45	13.273	1925	1930	1934	rVV	298762	535334	17.53%	1.133%
46	13.314	1934	1937	1940	rVV3	109053	159223	5.21%	0.337%
47	13.361	1942	1945	1949	rVV2	101893	166588	5.45%	0.352%
48	13.402	1949	1952	1955	rVV3	78119	94275	3.09%	0.199%
49	13.461	1956	1962	1966	rVV5	344029	671680	21.99%	1.421%
50	13.508	1966	1970	1973	rVV2	265048	409560	13.41%	0.867%
51	13.538	1973	1975	1976	rVV2	137289	131809	4.32%	0.279%
52	13.579	1976	1982	1990	rVB2	1637365	3054242	100.00%	6.462%
53	13.649	1990	1994	1999	rVB4	192656	258907	8.48%	0.548%
54	13.726	2001	2007	2010	rBV7	133916	291924	9.56%	0.618%
55	13.773	2011	2015	2018	rVV2	221585	311784	10.21%	0.660%
56	13.808	2018	2021	2030	rVB2	265251	485335	15.89%	1.027%
57	13.896	2030	2036	2041	rBV	1477085	2214973	72.52%	4.686%
58	13.967	2046	2048	2052	rVB	105666	131993	4.32%	0.279%
59	14.061	2057	2064	2069	rVB4	234684	543793	17.80%	1.151%
60	14.114	2069	2073	2078	rVB	292656	444583	14.56%	0.941%
61	14.161	2078	2081	2087	rVB4	130326	207524	6.79%	0.439%
62	14.226	2087	2092	2098	rBV3	445780	744413	24.37%	1.575%
63	14.285	2098	2102	2109	rVB7	142473	336348	11.01%	0.712%
64	14.367	2109	2116	2119	rBV2	242015	534552	17.50%	1.131%
65	14.414	2120	2124	2127	rVB3	225064	318732	10.44%	0.674%
66	14.526	2139	2143	2146	rBV2	321872	430151	14.08%	0.910%
67	14.567	2146	2150	2157	rVB4	313647	520228	17.03%	1.101%
68	14.626	2158	2160	2161	rBV2	50754	43963	1.44%	0.093%
69	14.714	2172	2175	2177	rBV2	108772	138629	4.54%	0.293%
70	14.755	2177	2182	2187	rVB2	940445	1439461	47.13%	3.046%
71	14.826	2188	2194	2199	rBV3	550275	1390277	45.52%	2.942%

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72	14.873	2199	2202	2210	rVB3	388534	726553	23.79%	1.537%
73	14.990	2218	2222	2228	rVB4	220453	362428	11.87%	0.767%
74	15.096	2235	2240	2245	rBV5	242088	449438	14.72%	0.951%
75	15.208	2253	2259	2268	rBV4	264117	640048	20.96%	1.354%
76	15.290	2269	2273	2280	rVB6	226651	471957	15.45%	0.999%
77	15.367	2281	2286	2295	rBV3	417515	859698	28.15%	1.819%
78	15.461	2297	2302	2307	rBV	165397	323692	10.60%	0.685%
79	15.514	2307	2311	2313	rBV5	97279	157270	5.15%	0.333%
80	15.602	2321	2326	2335	rVB	669067	1227309	40.18%	2.597%
81	15.702	2336	2343	2349	rBV5	172947	441000	14.44%	0.933%
82	15.779	2352	2356	2363	rVB5	149384	279301	9.14%	0.591%
83	15.914	2375	2379	2386	rVB	228947	413714	13.55%	0.875%
84	16.079	2400	2407	2414	rBV6	117881	324204	10.61%	0.686%
85	16.232	2424	2433	2439	rBV4	213118	606062	19.84%	1.282%
86	16.355	2449	2454	2460	rVB	133797	219854	7.20%	0.465%
87	16.432	2462	2467	2473	rVV4	159325	353914	11.59%	0.749%
88	16.502	2474	2479	2484	rVV2	154148	310188	10.16%	0.656%
89	16.567	2484	2490	2496	rVB2	263356	495280	16.22%	1.048%
90	16.702	2506	2513	2527	rVB8	159285	555371	18.18%	1.175%

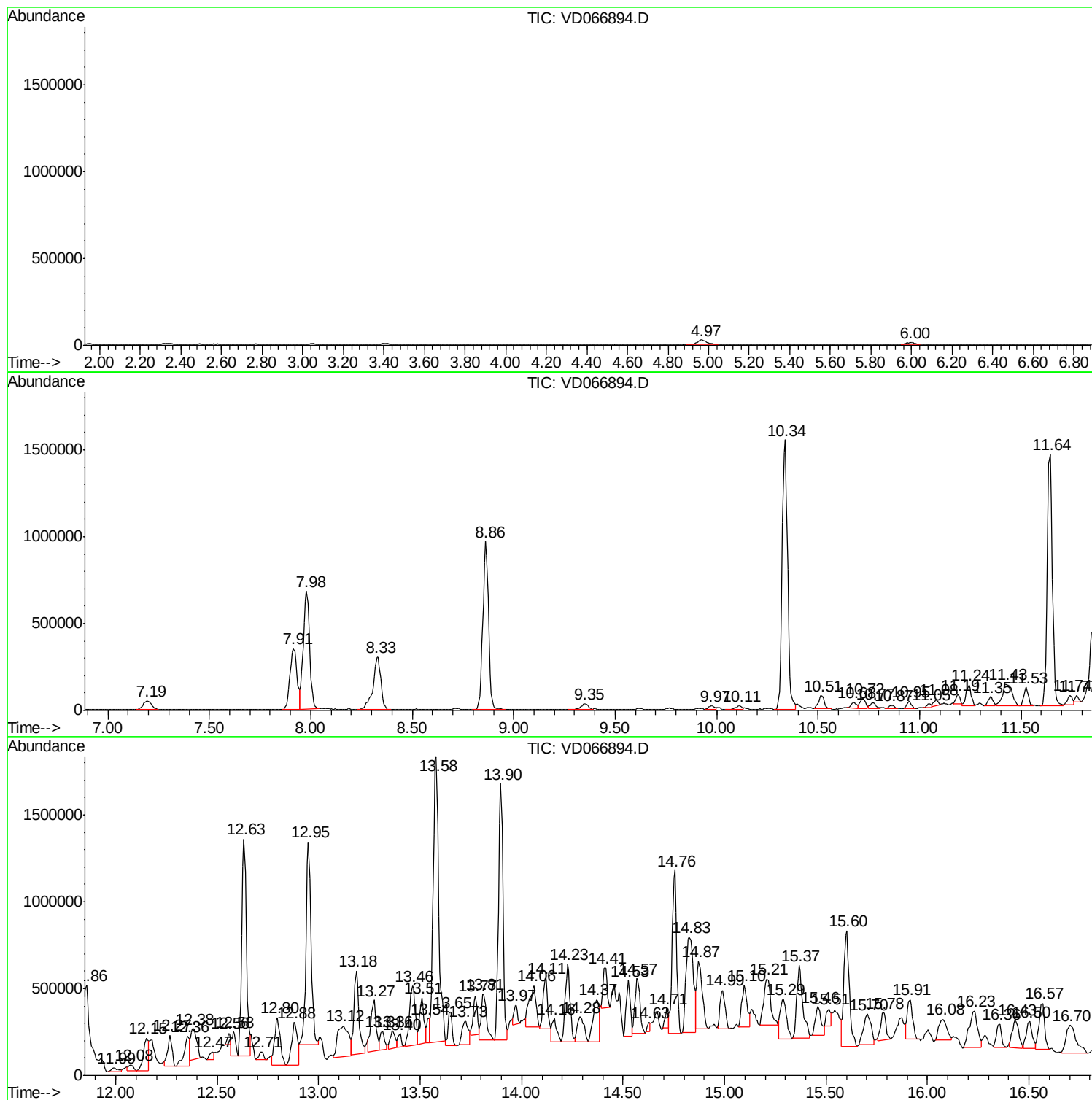
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.M
Quant Title : SW846 8260

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



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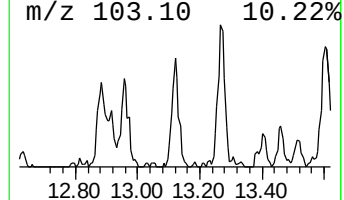
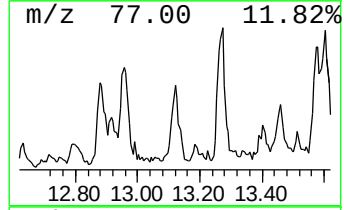
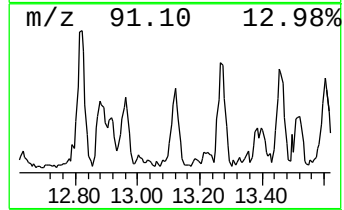
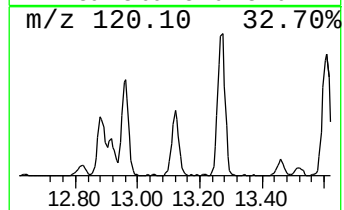
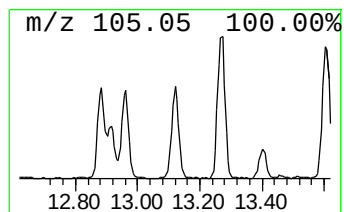
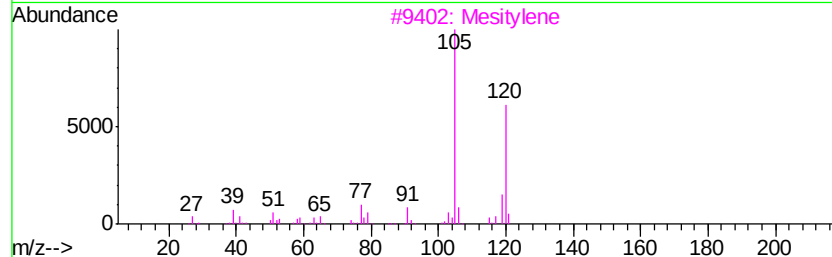
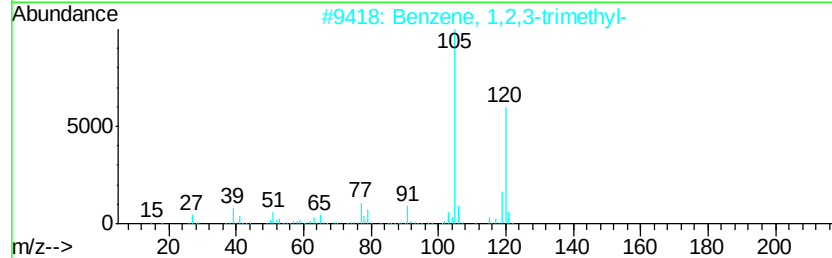
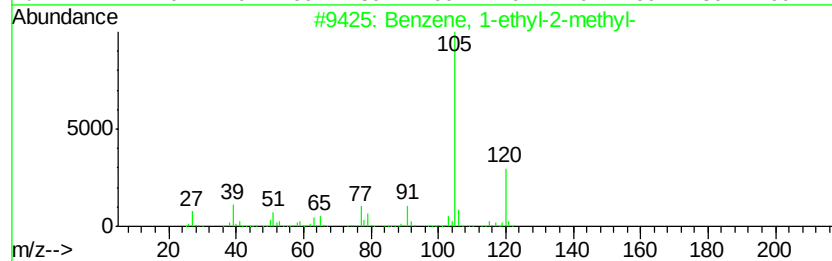
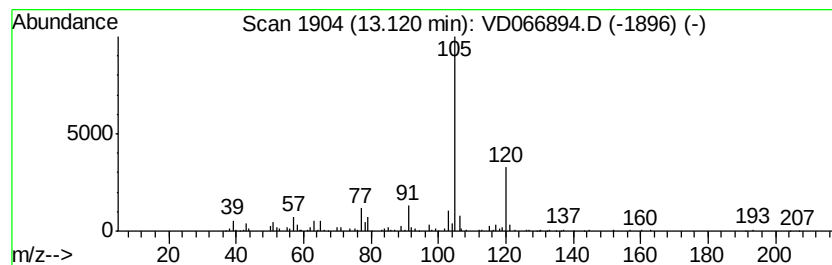
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	11.46 ug/l	699855	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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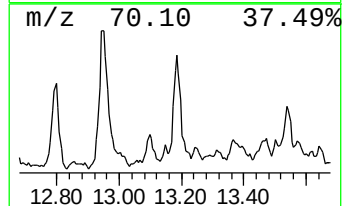
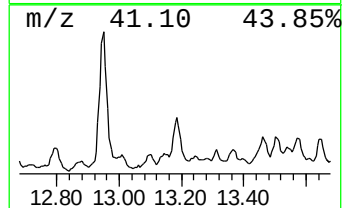
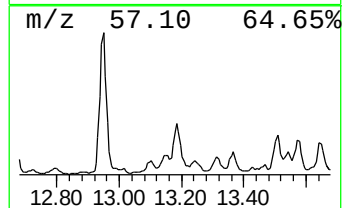
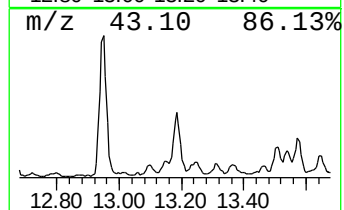
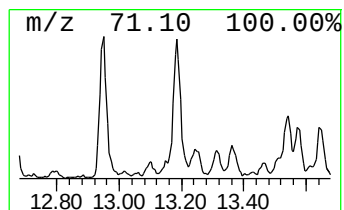
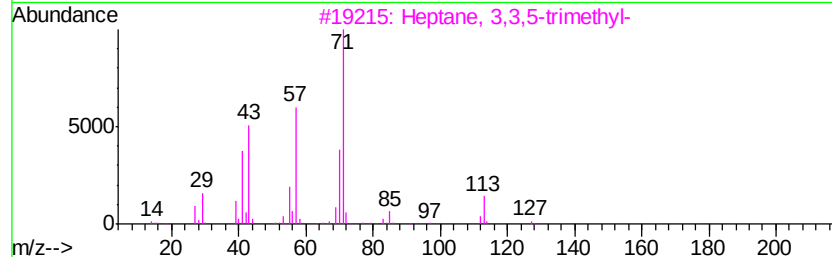
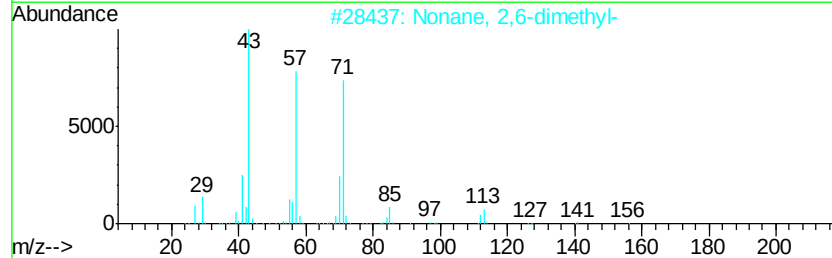
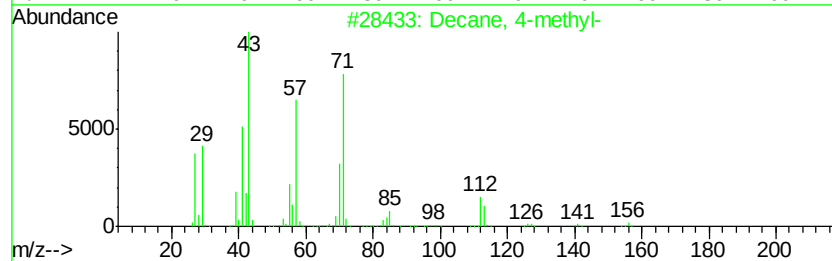
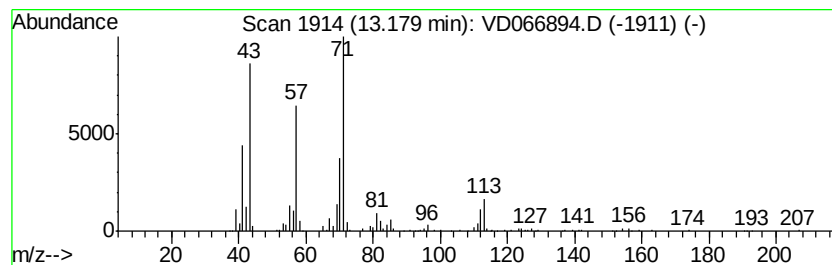
Instrument :
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Quant Title : SW846 8260

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

Peak Number 2 Decane, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.18	13.45 ug/l	821821	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Decane, 4-methyl-	156 C11H24	002847-72-5	83
2	Nonane, 2,6-dimethyl-	156 C11H24	017302-28-2	76
3	Heptane, 3,3,5-trimethyl-	142 C10H22	007154-80-5	64
4	Heptane, 3-ethyl-5-methyl-	142 C10H22	052896-90-9	64
5	Octane, 3,3-dimethyl-	142 C10H22	004110-44-5	59



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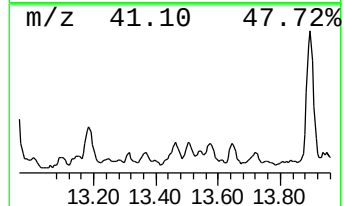
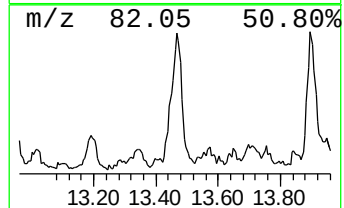
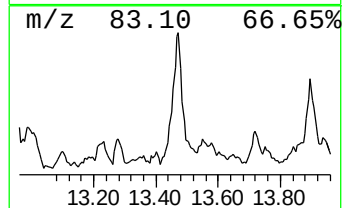
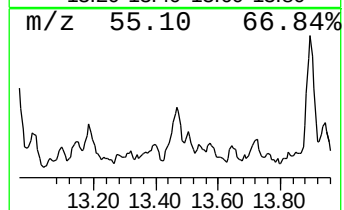
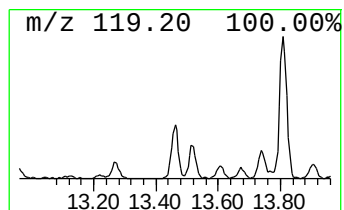
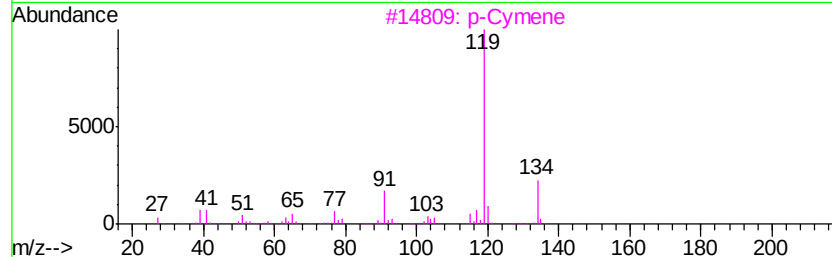
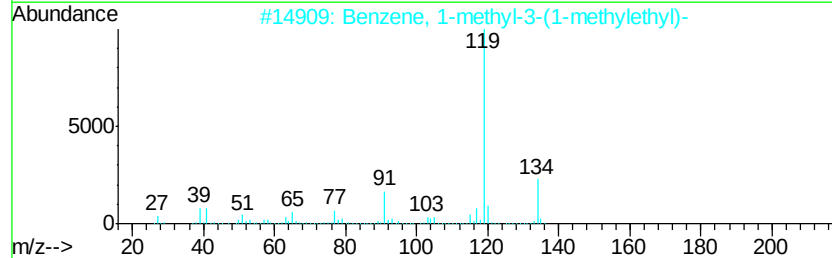
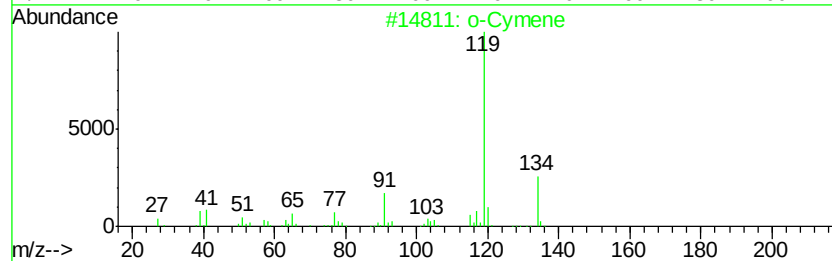
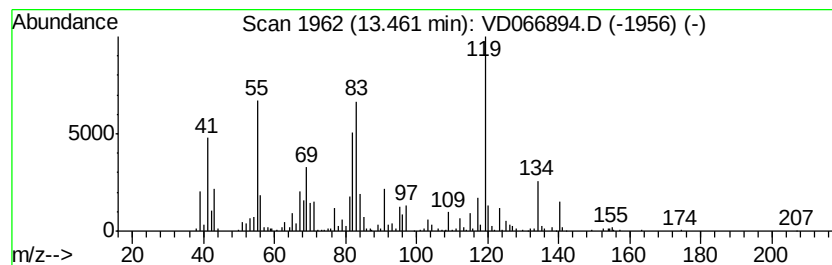
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.M
Quant Title : SW846 8260

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

Peak Number 3 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.46	11.00 ug/l	671680	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	86
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	86
3		p-Cymene	134	C10H14	000099-87-6	86
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	46
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	46



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ClientSampleId :

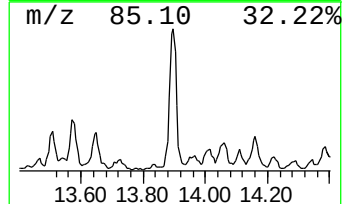
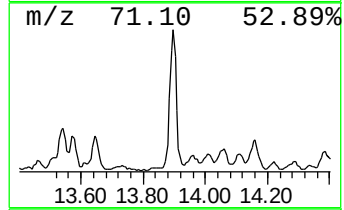
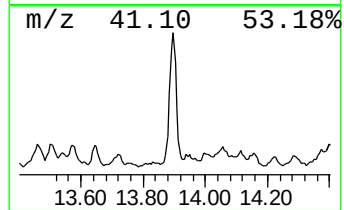
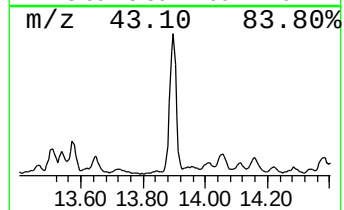
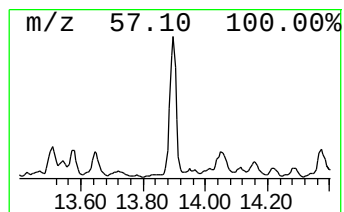
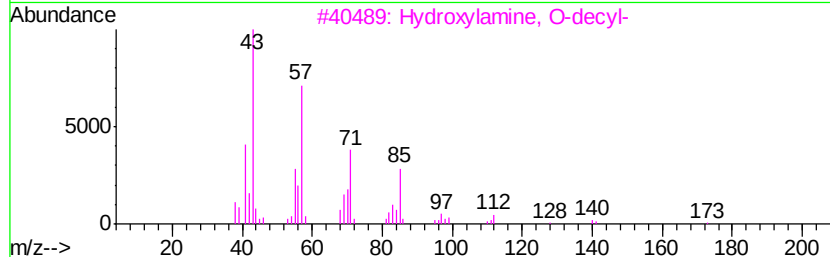
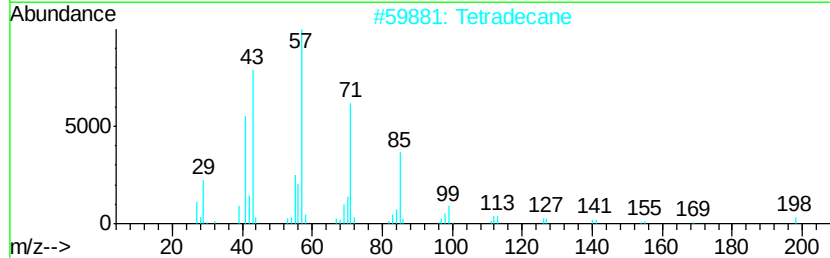
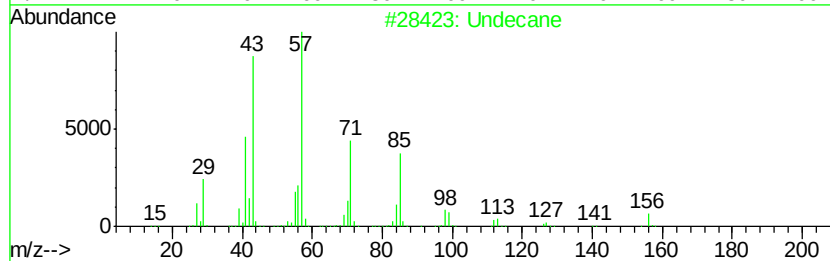
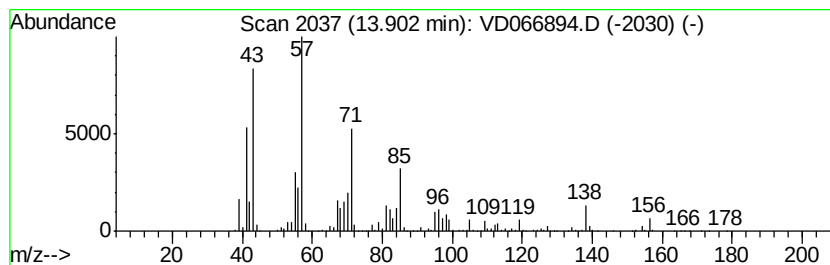
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Quant Title : SW846 8260

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

Peak Number 4 Undecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	36.26 ug/l	2214970	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	91
2	Tetradecane		198	C14H30	000629-59-4	59
3	Hvdroxylamine, O-decyl-		173	C10H23NO	029812-79-1	58
4	Oxalic acid, allyl tridecyl ester		312	C18H32O4	1000309-24-1	53
5	Ether, hexyl pentyl		172	C11H24O	032357-83-8	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092520\
Data File : VD066894.D
Acq On : 26 Sep 2020 02:11
Operator : VA/SY
Sample : L4154-04
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :

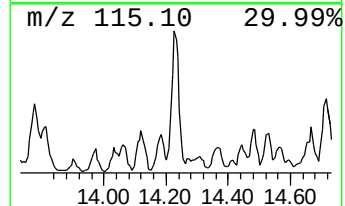
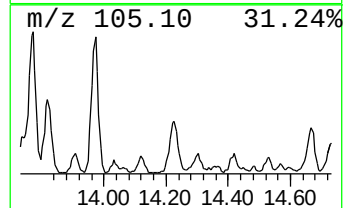
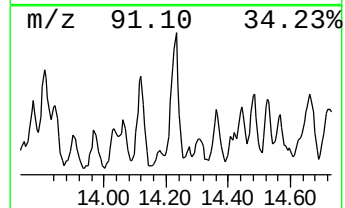
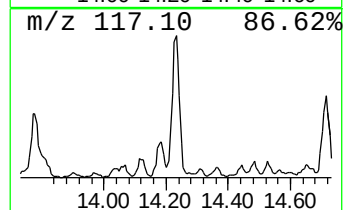
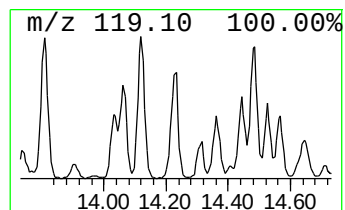
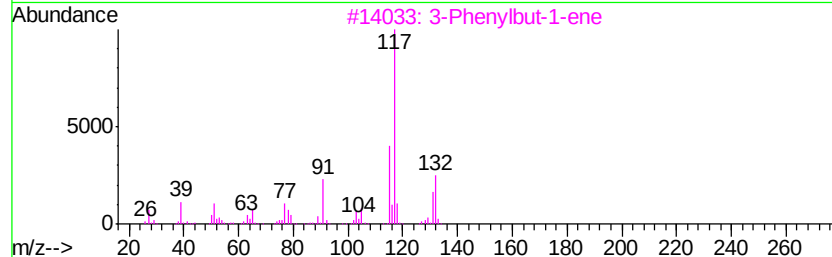
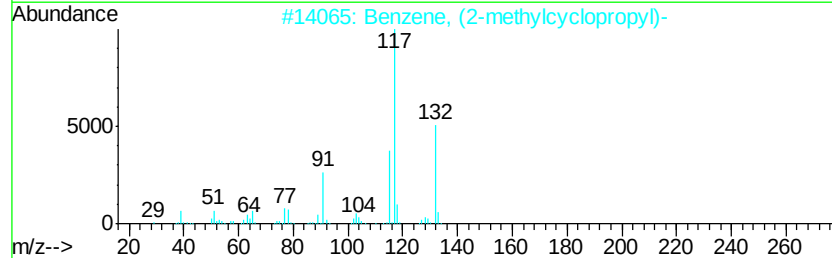
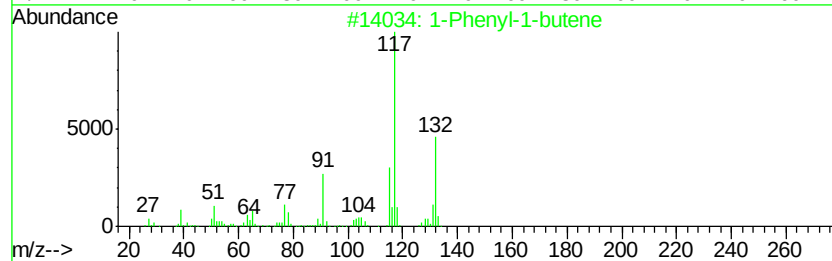
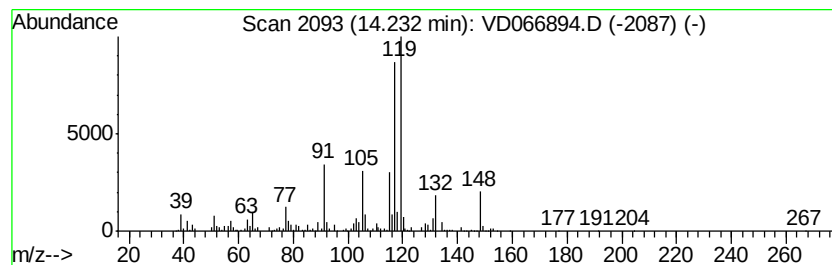
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Quant Title : SW846 8260

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

Peak Number 5 1-Phenyl-1-butene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	12.19 ug/l	744413	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	52
2		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	46
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	46
4		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	46
5		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092520\
 Data File : VD066894.D
 Acq On : 26 Sep 2020 02:11
 Operator : VA/SY
 Sample : L4154-04
 Misc : 5.01G/5.00ml/MSVOA D/SOIL
 ALS Vial : 36 Sample Multiplier: 1

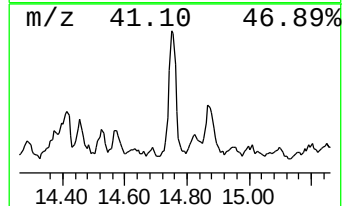
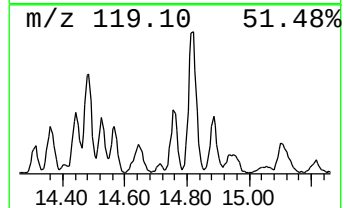
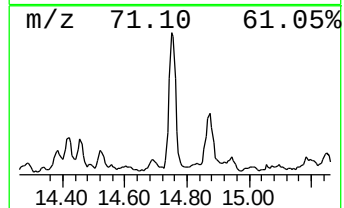
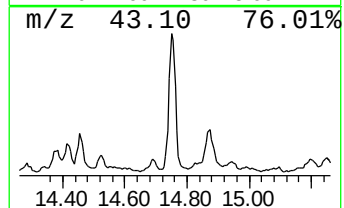
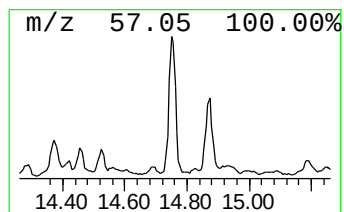
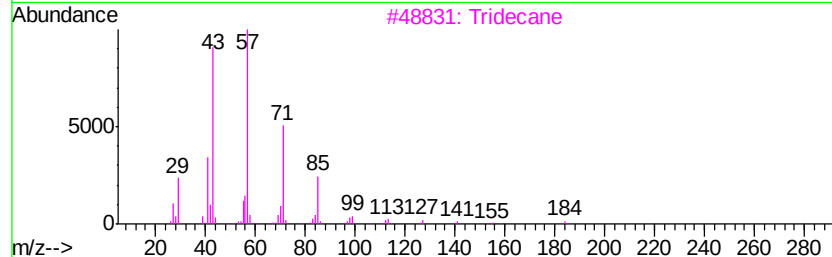
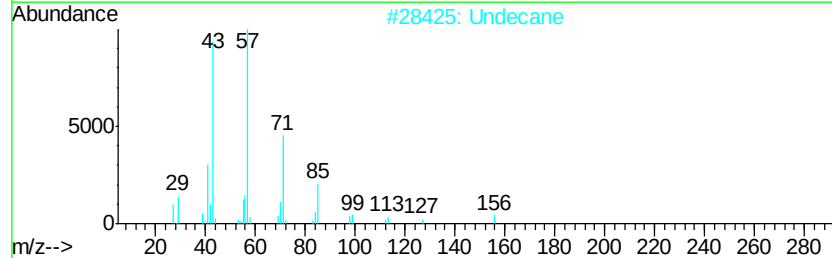
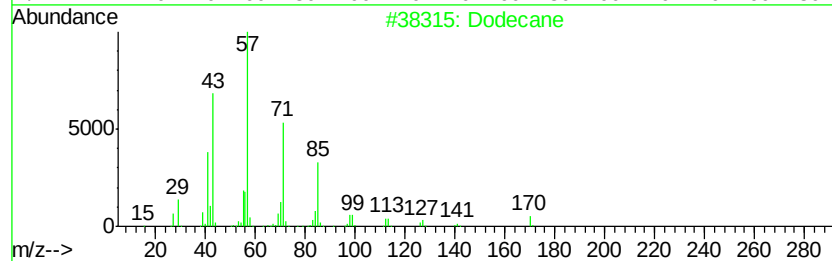
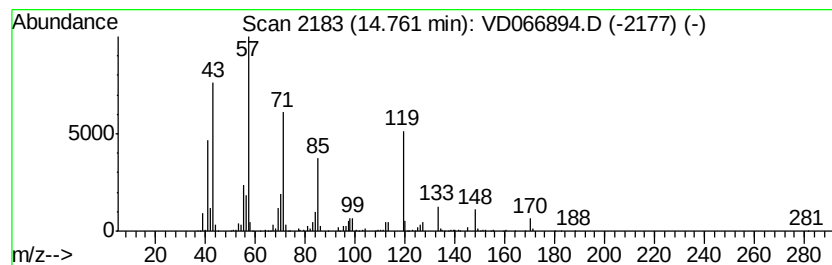
Instrument :
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 ClientSampleId :

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.M
 Quant Title : SW846 8260

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

 Peak Number 6 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	23.56 ug/l	1439460	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	96
2	Undecane	156 C11H24	001120-21-4	53
3	Tridecane	184 C13H28	000629-50-5	53
4	Hexadecane	226 C16H34	000544-76-3	52
5	Tetradecane	198 C14H30	000629-59-4	52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092520\
Data File : VD066894.D
Acq On : 26 Sep 2020 02:11
Operator : VA/SY
Sample : L4154-04
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :

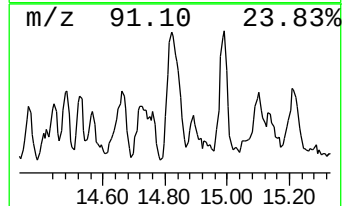
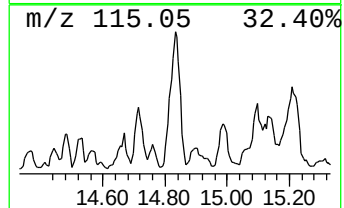
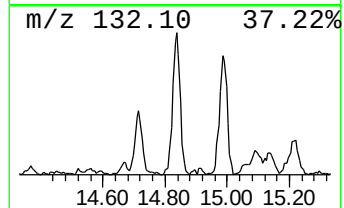
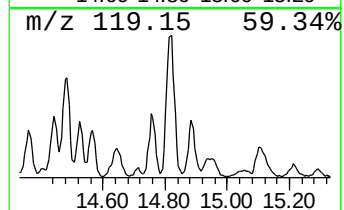
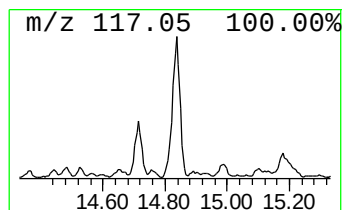
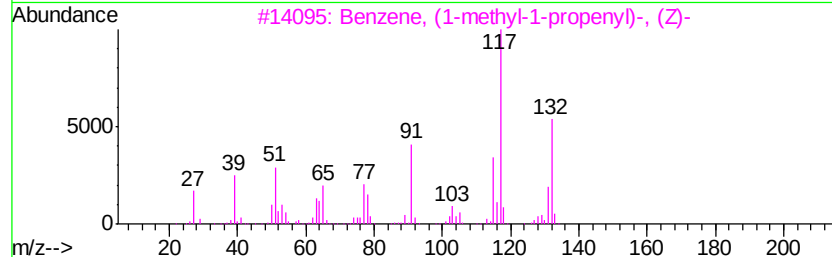
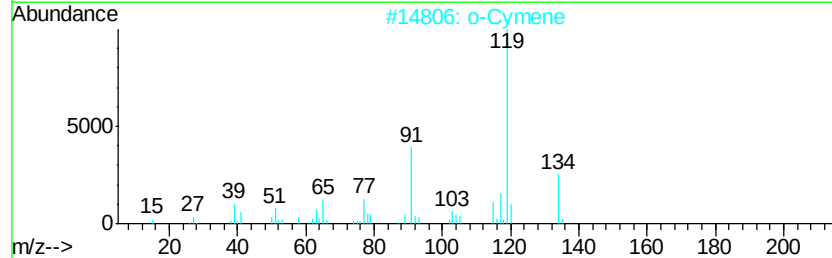
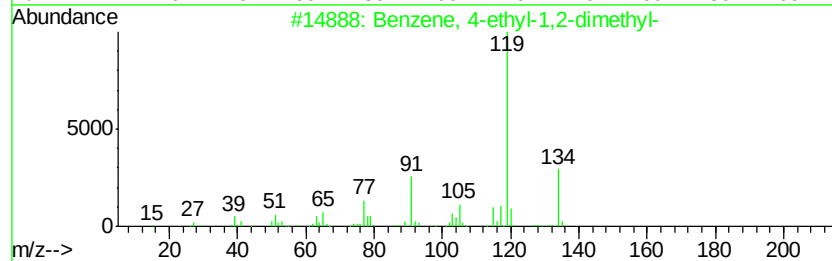
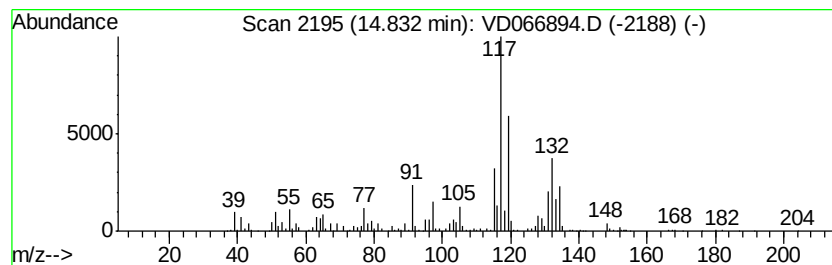
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	22.76 ug/l	1390280	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	50
2		o-Cymene	134	C10H14	000527-84-4	50
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	49
4		p-Cymene	134	C10H14	000099-87-6	46
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092520\
Data File : VD066894.D
Acq On : 26 Sep 2020 02:11
Operator : VA/SY
Sample : L4154-04
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :

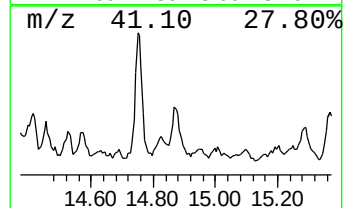
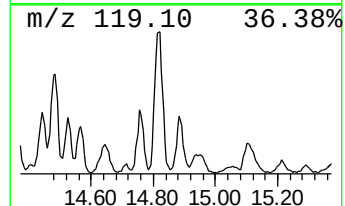
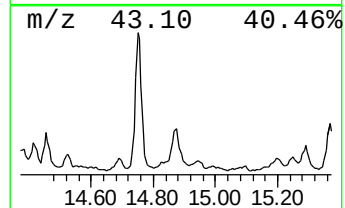
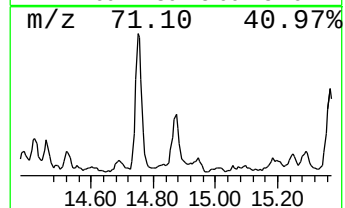
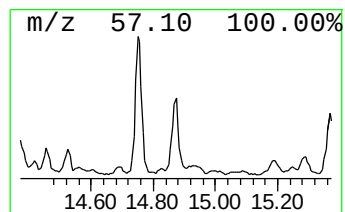
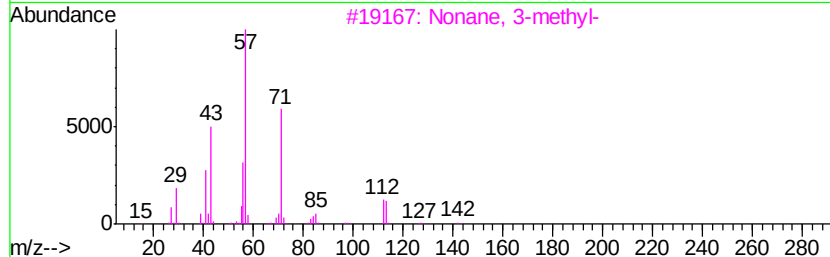
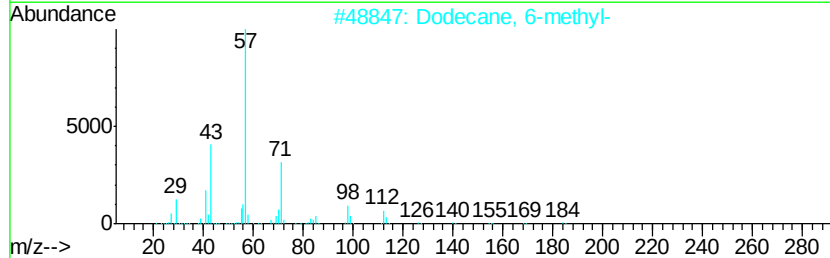
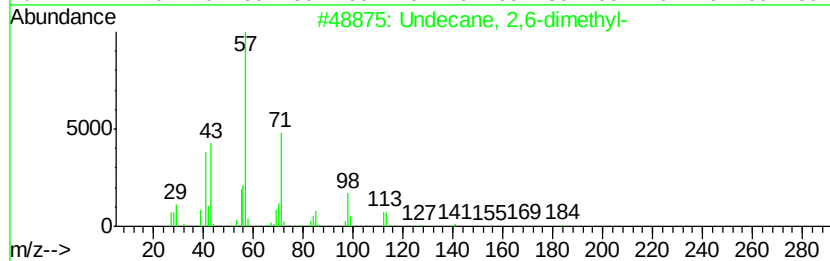
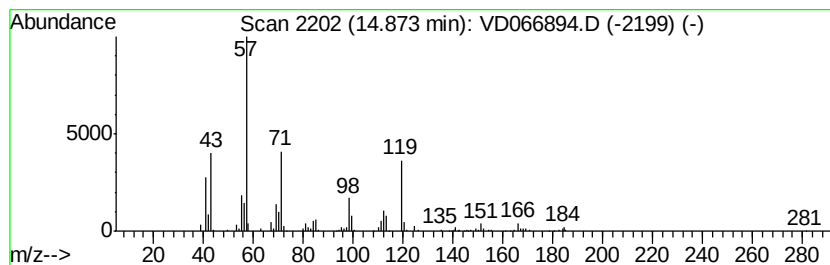
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Quant Title : SW846 8260

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

Peak Number 8 Undecane, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.87	11.89 ug/l	726553	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	93
2		Dodecane, 6-methyl-	184	C13H28	006044-71-9	53
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	53
4		Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	50
5		1-Decene, 4-methyl-	154	C11H22	013151-29-6	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092520\
Data File : VD066894.D
Acq On : 26 Sep 2020 02:11
Operator : VA/SY
Sample : L4154-04
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 36 Sample Multiplier: 1

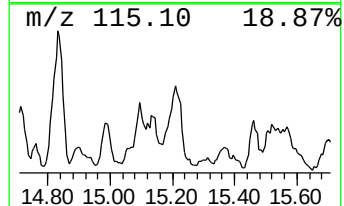
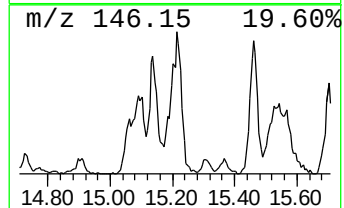
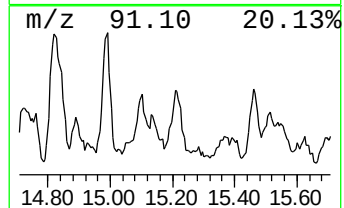
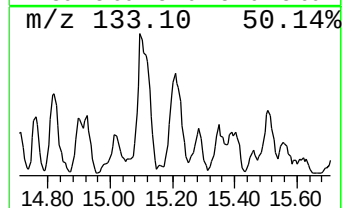
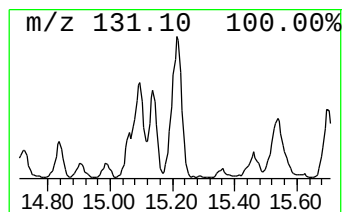
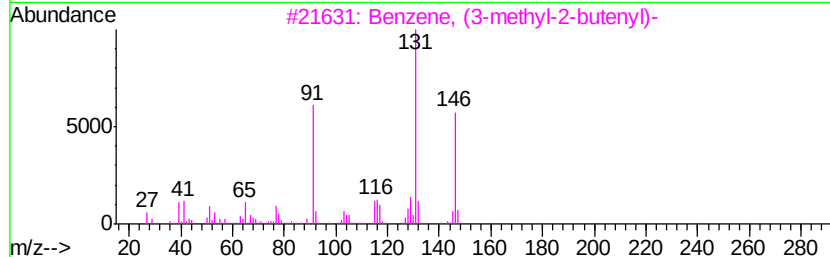
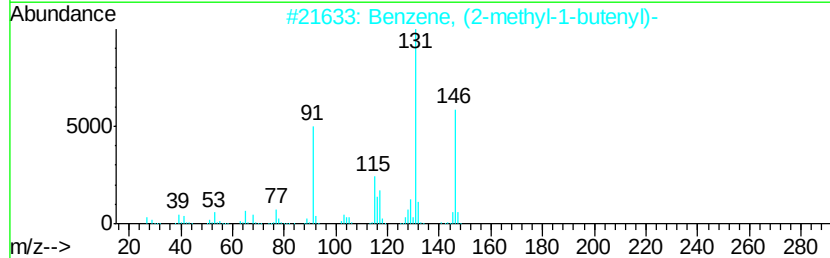
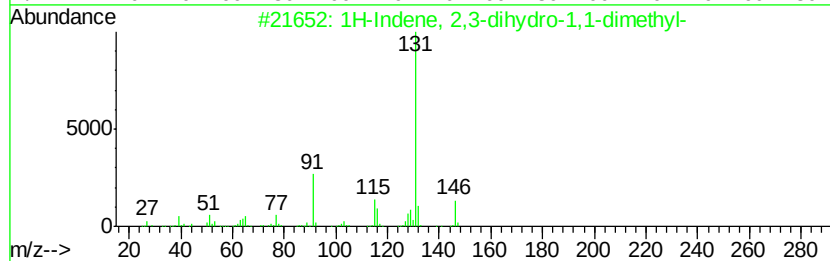
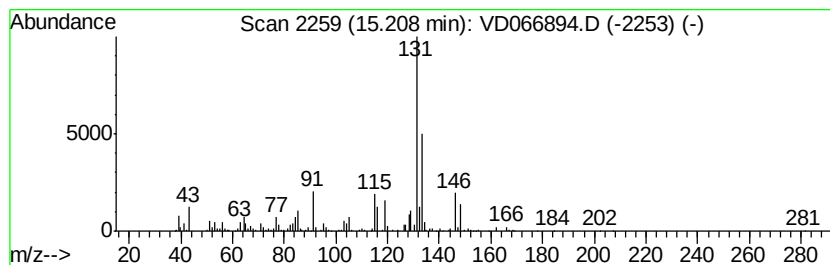
Instrument :
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ClientSampleId :

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.M
Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,1-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	10.48 ug/l	640048	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9	90
2	Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6	90
3	Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3	90
4	Benzene, 1-methyl-2-(1-methyl-2-...	146 C11H14	097664-19-2	64
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD092520\
Data File : VD066894.D
Acq On : 26 Sep 2020 02:11
Operator : VA/SY
Sample : L4154-04
Misc : 5.01G/5.00ml/MSVOA D/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :

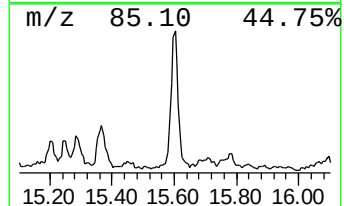
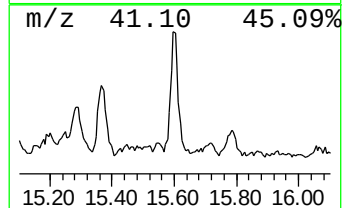
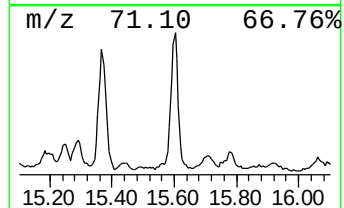
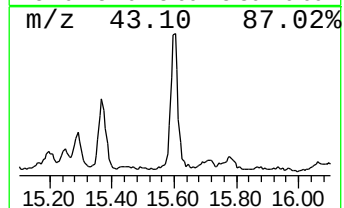
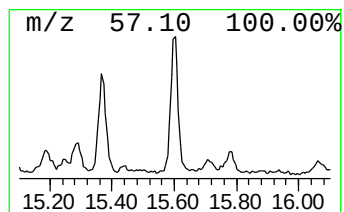
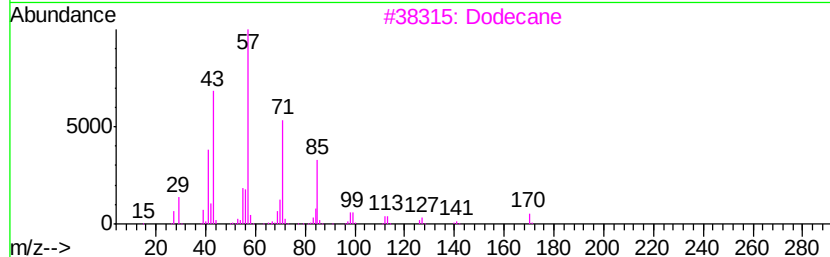
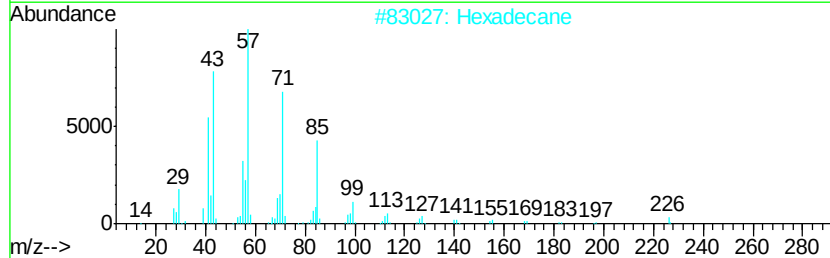
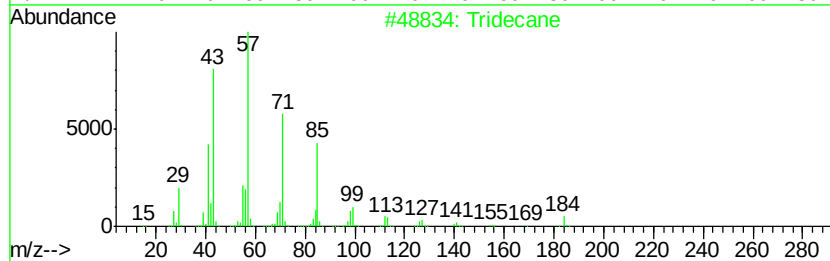
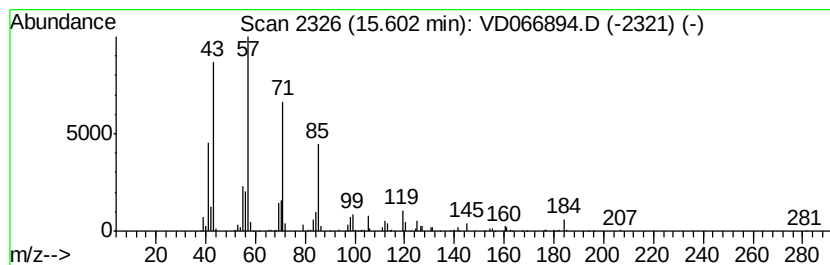
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Quant Title : SW846 8260

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

Peak Number 10 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.60	20.09 ug/l	1227310	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	96
2		Hexadecane	226	C16H34	000544-76-3	81
3		Dodecane	170	C12H26	000112-40-3	72
4		Pentadecane	212	C15H32	000629-62-9	72
5		Tetradecane	198	C14H30	000629-59-4	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD092520\
Data File : VD066894.D
Acq On : 26 Sep 2020 02:11
Operator : VA/SY
Sample : L4154-04
Misc : 5.01G/5.00ml/MSVOA_D/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D092220S.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-ethyl-...	13.12	11.5	ug/l	699855	4	13.58	3054240	50.0
Decane, 4-methyl-	13.18	13.4	ug/l	821821	4	13.58	3054240	50.0
o-Cymene	13.46	11.0	ug/l	671680	4	13.58	3054240	50.0
Undecane	13.90	36.3	ug/l	2214970	4	13.58	3054240	50.0
1-Phenyl-1-butene	14.23	12.2	ug/l	744413	4	13.58	3054240	50.0
Dodecane	14.76	23.6	ug/l	1439460	4	13.58	3054240	50.0
Benzene, 4-ethyl-...	14.83	22.8	ug/l	1390280	4	13.58	3054240	50.0
Undecane, 2,6-dim...	14.87	11.9	ug/l	726553	4	13.58	3054240	50.0
1H-Indene, 2,3-di...	15.21	10.5	ug/l	640048	4	13.58	3054240	50.0
Tridecane	15.60	20.1	ug/l	1227310	4	13.58	3054240	50.0