

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD090420\
 Data File : VD066558.D
 Acq On : 04 Sep 2020 13:00
 Operator : VA/SY
 Sample : L3877-01
 Misc : 6.78G/5.00ml/MSVOA D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-1

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\82D090320S.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	7.726	970	987	999	rBV	1191248	2972175	1.08%	0.146%
2	7.979	1022	1030	1038	rVB5	1086757	2955932	1.08%	0.145%
3	8.355	1081	1094	1108	rBV3	70584089	207420023	75.67%	10.184%
4	9.350	1248	1263	1270	rBV	4144193	9277278	3.38%	0.455%
5	9.973	1364	1369	1383	rVB2	2326099	5500047	2.01%	0.270%
6	10.114	1383	1393	1404	rBV	1850289	4507063	1.64%	0.221%
7	10.214	1404	1410	1416	rBV	1738014	3233710	1.18%	0.159%
8	10.332	1423	1430	1435	rBV2	4196260	8593064	3.13%	0.422%
9	10.402	1435	1442	1448	rVB	4354436	9872472	3.60%	0.485%
10	10.520	1454	1462	1470	rVB2	4194876	8299684	3.03%	0.408%
11	10.679	1483	1489	1497	rVB	1974130	3806699	1.39%	0.187%
12	10.949	1525	1535	1542	rBV	1914170	3454192	1.26%	0.170%
13	11.049	1542	1552	1559	rBV2	1252007	3659198	1.33%	0.180%
14	11.191	1567	1576	1580	rVV	2905623	5335286	1.95%	0.262%
15	11.244	1580	1585	1590	rVB	2802007	4691047	1.71%	0.230%
16	11.432	1608	1617	1628	rVB4	3009674	8057956	2.94%	0.396%
17	11.532	1628	1634	1644	rBV	2330159	4424002	1.61%	0.217%
18	11.749	1663	1671	1680	rBV3	82692116	183024073	66.77%	8.986%
19	11.861	1680	1690	1699	rBV2	56614709	118887484	43.37%	5.837%
20	12.179	1732	1744	1752	rBV	13951282	29176169	10.64%	1.433%
21	12.267	1755	1759	1764	rVB	1806933	2820859	1.03%	0.138%
22	12.355	1764	1774	1777	rBV4	1826459	4630638	1.69%	0.227%
23	12.396	1777	1781	1786	rVB4	1727015	3336715	1.22%	0.164%
24	12.479	1789	1795	1800	rBV	6217727	9978297	3.64%	0.490%
25	12.820	1845	1853	1858	rBV	9701315	17401521	6.35%	0.854%
26	12.885	1858	1864	1867	rVV	28603761	48070344	17.54%	2.360%
27	12.920	1867	1870	1873	rVV	21915663	35062539	12.79%	1.722%
28	12.961	1873	1877	1884	rVV2	40181327	70198590	25.61%	3.447%
29	13.126	1896	1905	1912	rBV	19960453	39789796	14.52%	1.954%
30	13.191	1912	1916	1922	rVB	4717444	6845480	2.50%	0.336%
31	13.279	1922	1931	1937	rBV3	72345392	141872596	51.76%	6.966%
32	13.396	1942	1951	1956	rBV4	2182053	6870963	2.51%	0.337%
33	13.467	1956	1963	1967	rBV2	6911374	13443714	4.90%	0.660%
34	13.514	1967	1971	1979	rVB5	5223892	11339845	4.14%	0.557%

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35	13.614	1979	1988	1992	rBV2	40713914	76078085	27.75%	3.735%
36	13.649	1992	1994	2002	rVB2	9827105	13280419	4.84%	0.652%
37	13.743	2003	2010	2011	rBV2	5005509	7024071	2.56%	0.345%
38	13.779	2011	2016	2020	rVV2	30914271	57252321	20.89%	2.811%
39	13.814	2020	2022	2032	rVB3	20302132	34231951	12.49%	1.681%
40	13.902	2032	2037	2042	rBV	25446336	37697525	13.75%	1.851%
41	13.967	2042	2048	2053	rBV3	6023883	12494768	4.56%	0.613%
42	14.038	2055	2060	2062	rVV	10541058	16308297	5.95%	0.801%
43	14.067	2062	2065	2070	rVV2	13662157	20942713	7.64%	1.028%
44	14.126	2070	2075	2079	rVB	19563706	30054722	10.96%	1.476%
45	14.232	2088	2093	2098	rVB2	11143214	17750225	6.48%	0.872%
46	14.290	2098	2103	2110	rVB8	3216661	6932747	2.53%	0.340%
47	14.373	2110	2117	2121	rBV4	9010202	19626991	7.16%	0.964%
48	14.420	2121	2125	2127	rVV3	12202345	20501810	7.48%	1.007%
49	14.449	2127	2130	2133	rVV2	17826443	30171334	11.01%	1.481%
50	14.485	2133	2136	2141	rVV	22218492	38706400	14.12%	1.900%
51	14.532	2141	2144	2148	rVV3	10334169	15830509	5.78%	0.777%
52	14.573	2148	2151	2159	rVB6	5871471	11109761	4.05%	0.545%
53	14.673	2164	2168	2171	rBV	2779122	3661294	1.34%	0.180%
54	14.720	2171	2176	2179	rVV	8785636	14346143	5.23%	0.704%
55	14.761	2179	2183	2188	rVB	22060772	31566975	11.52%	1.550%
56	14.826	2188	2194	2200	rBV3	24057332	53606195	19.56%	2.632%
57	14.885	2200	2204	2212	rVB3	9459529	22278672	8.13%	1.094%
58	14.996	2217	2223	2229	rVB	14742970	24442429	8.92%	1.200%
59	15.108	2238	2242	2246	rBV3	3229583	4663224	1.70%	0.229%
60	15.220	2257	2261	2269	rVB4	3470135	8179916	2.98%	0.402%
61	15.296	2270	2274	2281	rVB4	2675256	6540214	2.39%	0.321%
62	15.426	2282	2296	2305	rVV2	93005582	274110524	100.00%	13.458%
63	15.514	2305	2311	2322	rVV	9654813	20006294	7.30%	0.982%
64	15.602	2322	2326	2335	rVB2	4357791	7833610	2.86%	0.385%
65	15.708	2336	2344	2351	rBV8	1277712	3635784	1.33%	0.179%
66	16.508	2472	2480	2495	rVB	17623649	39988681	14.59%	1.963%
67	16.714	2506	2515	2527	rVB	7290045	17063431	6.23%	0.838%

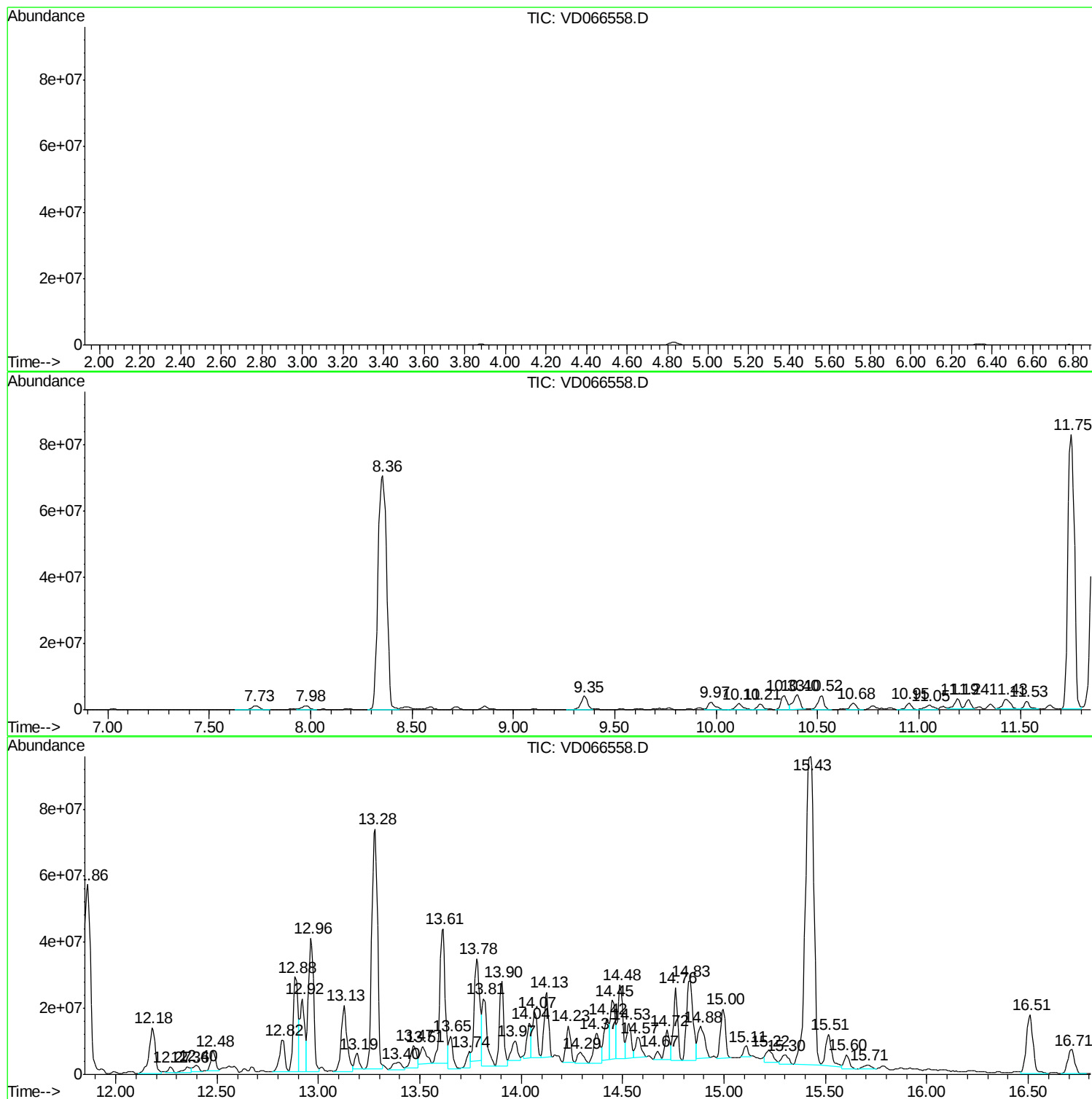
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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



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Instrument :
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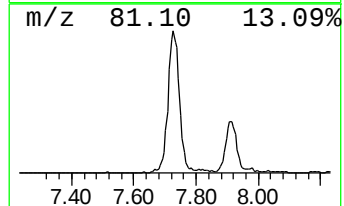
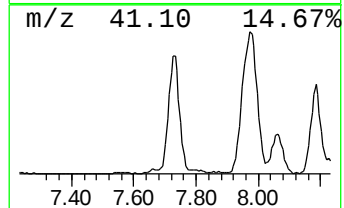
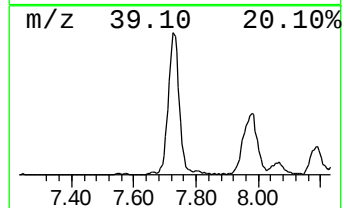
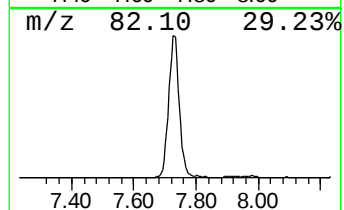
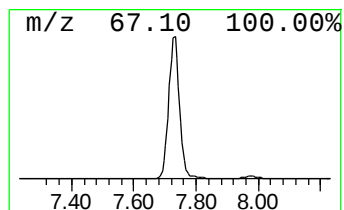
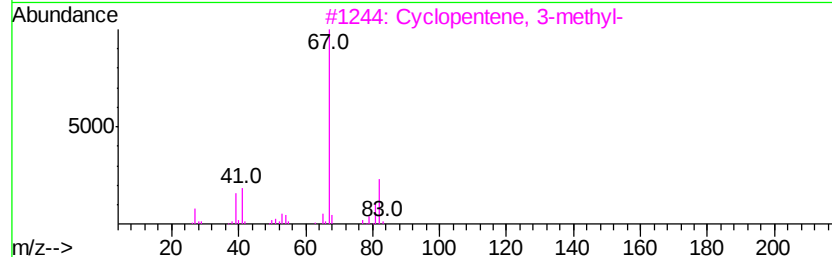
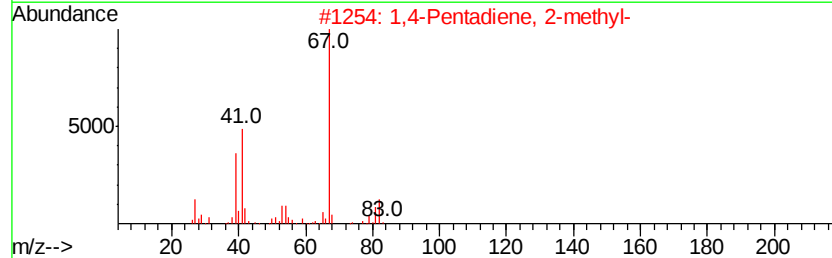
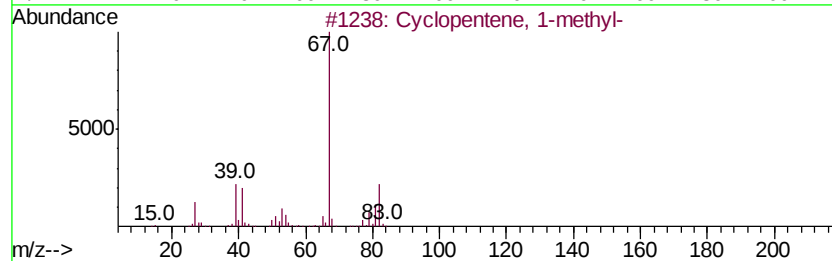
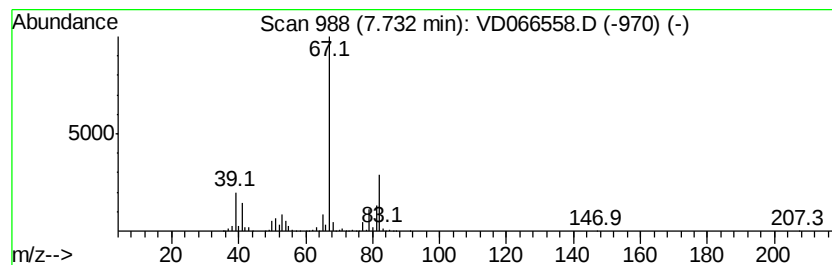
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.73	50.27 ug/l	2972180	Pentafluorobenzene	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	90
2		1,4-Pentadiene, 2-methyl-	82	C6H10	000763-30-4	90
3		Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	86
4		Cyclopentane, methylene-	82	C6H10	001528-30-9	81
5		1,4-Hexadiene	82	C6H10	000592-45-0	80



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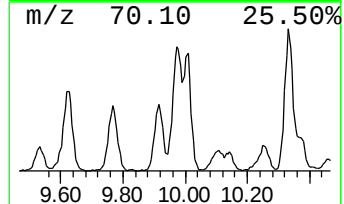
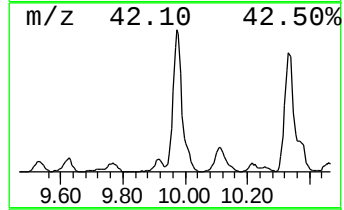
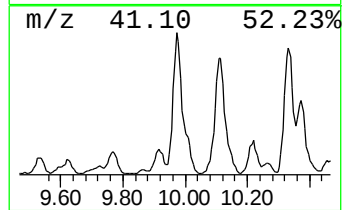
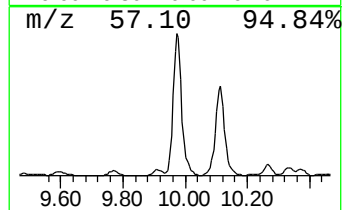
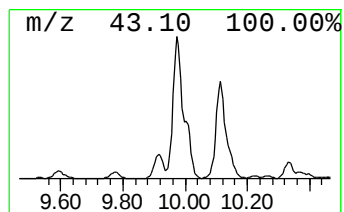
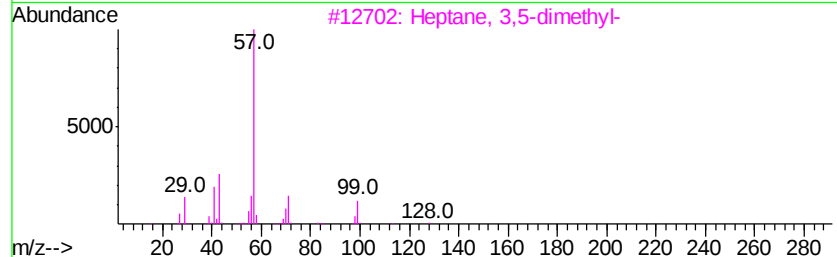
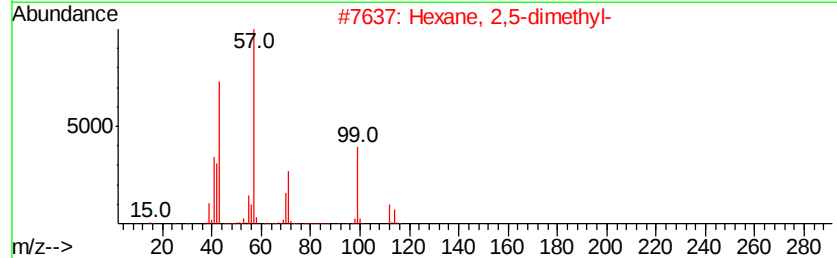
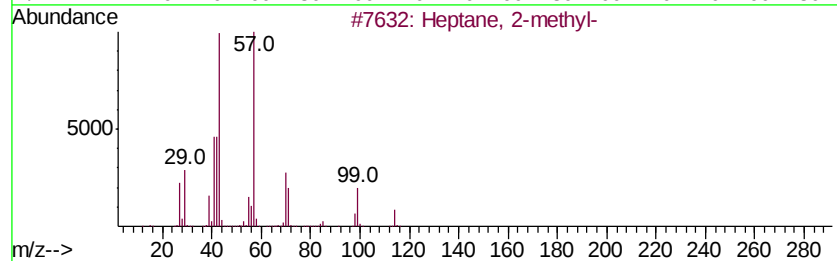
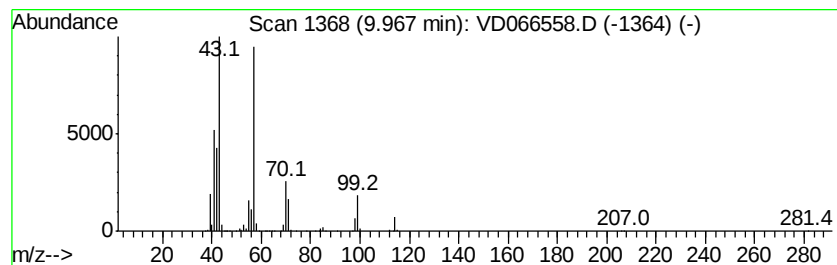
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Heptane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.97	29.64 ug/l	5500050	1,4-Difluorobenzene	8.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 2-methyl-	114	C8H18	000592-27-8	91
2		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	87
3		Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	38
4		1-Pentanol, 2,2,4-trimethyl-	130	C8H18O	000123-44-4	38
5		Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	38



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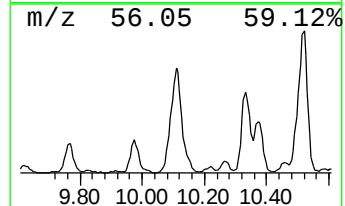
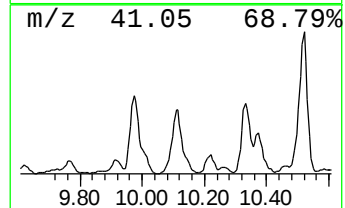
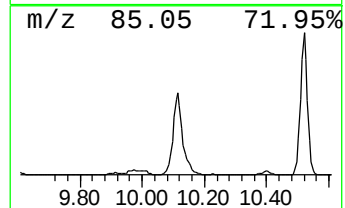
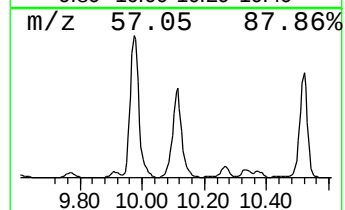
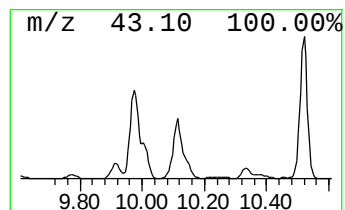
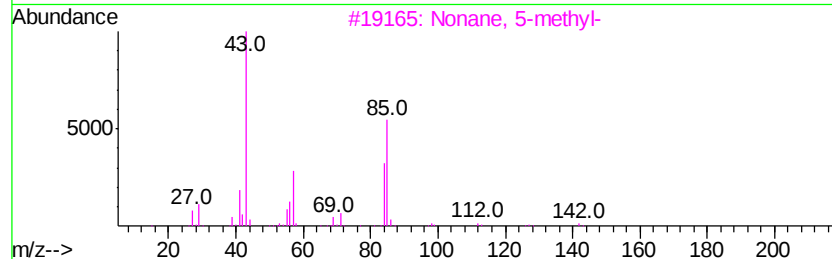
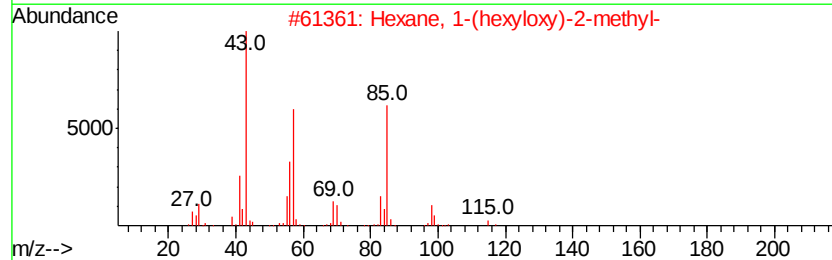
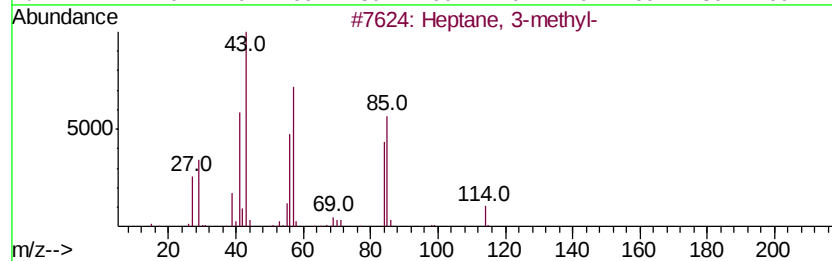
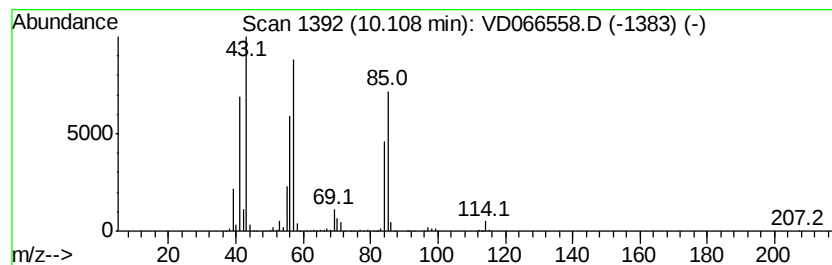
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Peak Number 3 Heptane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.11	24.29 ug/l	4507060	1,4-Difluorobenzene	8.86
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heptane, 3-methyl-	114 C8H18	000589-81-1	94
2	Hexane, 1-(hexyloxy)-2-methyl-	200 C13H28O	074421-17-3	64
3	Nonane, 5-methyl-	142 C10H22	015869-85-9	64
4	Heptane, 3,4,5-trimethyl-	142 C10H22	020278-89-1	59
5	Hexane, 2,4-dimethyl-	114 C8H18	000589-43-5	53



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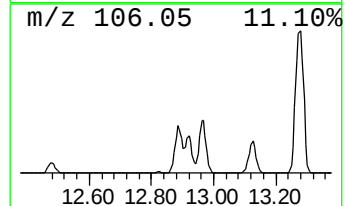
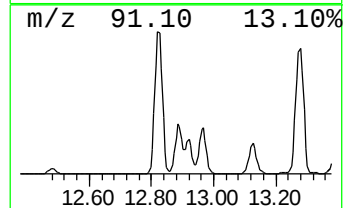
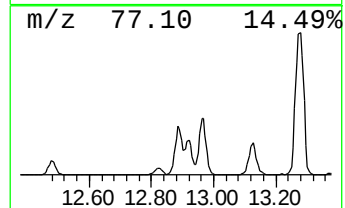
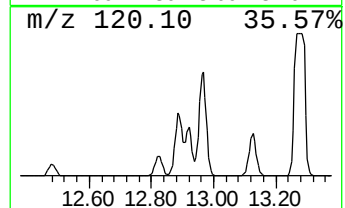
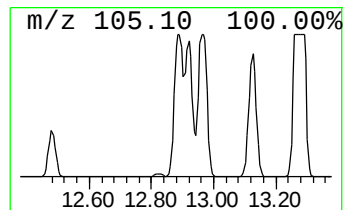
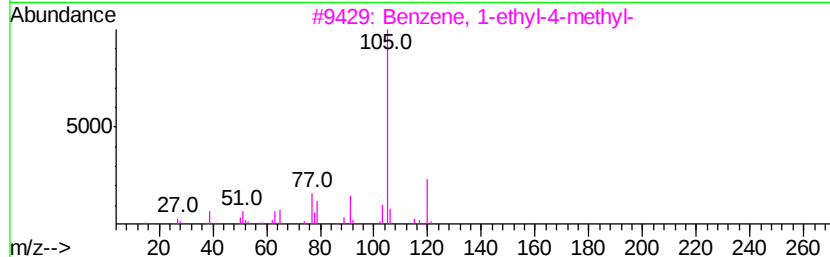
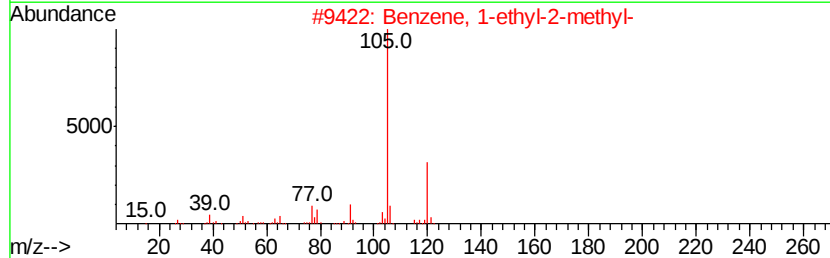
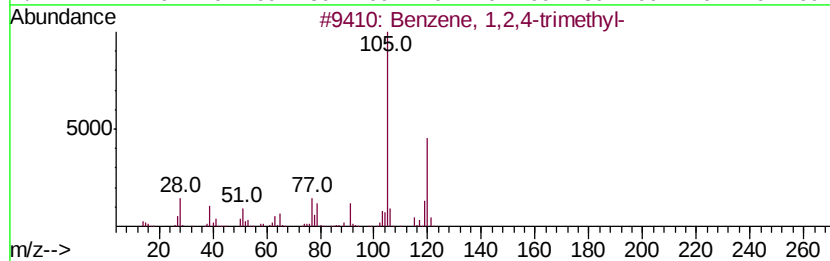
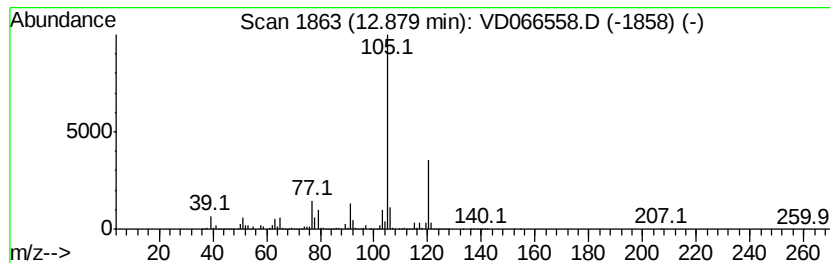
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Peak Number 4 Benzene, 1-ethyl-4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.88	31.59 ug/l	48070300	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
2		Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TP-1

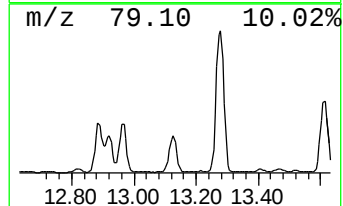
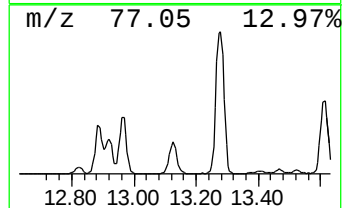
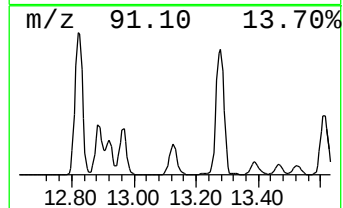
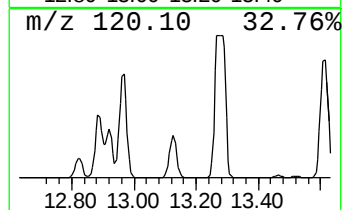
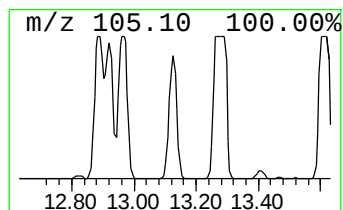
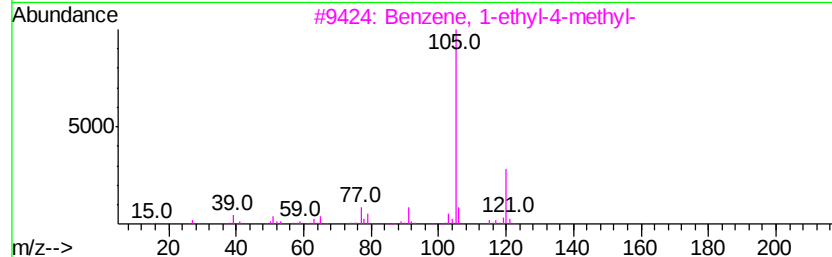
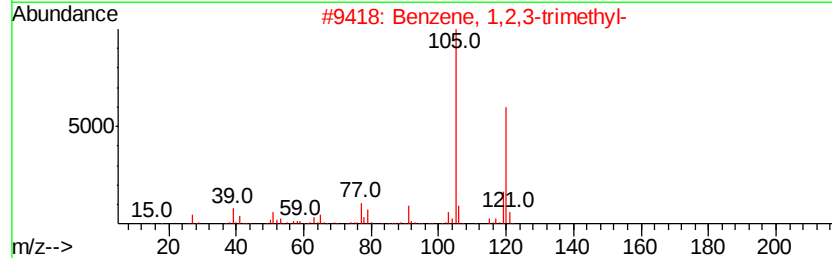
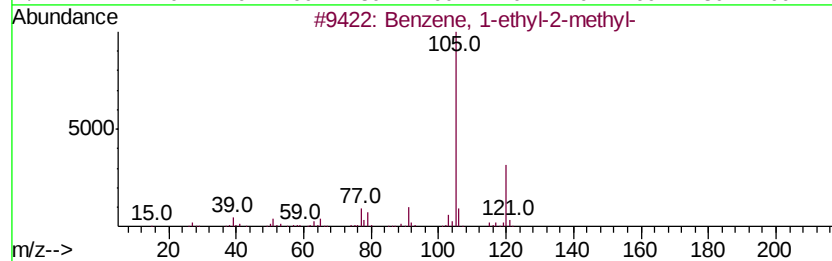
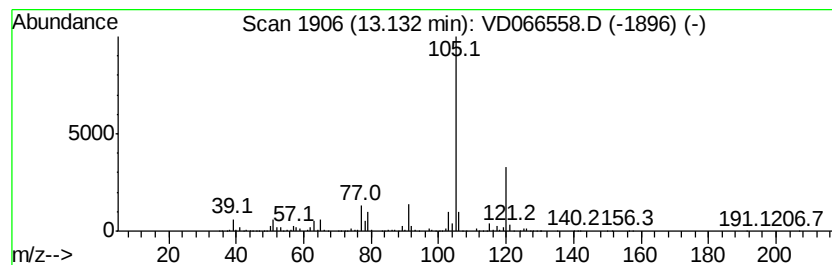
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	26.15 ug/l	39789800	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	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

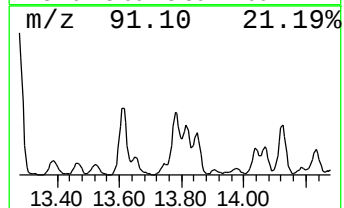
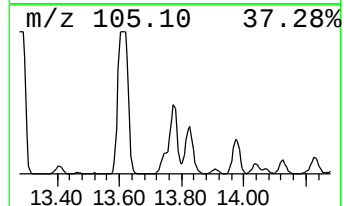
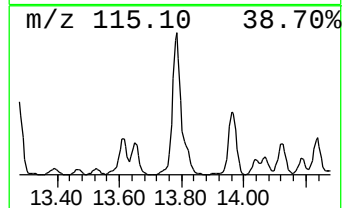
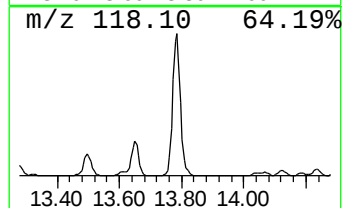
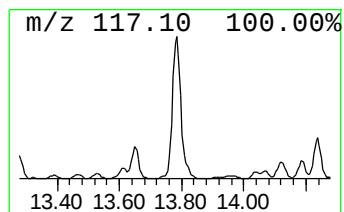
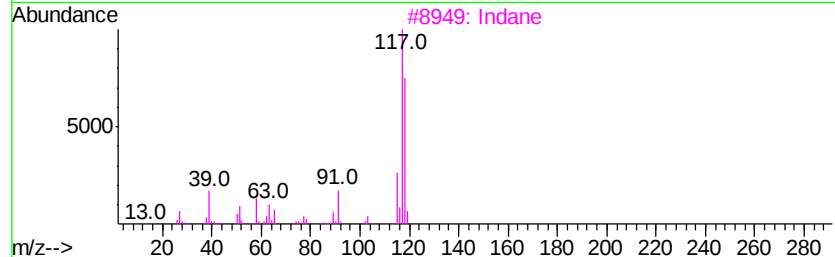
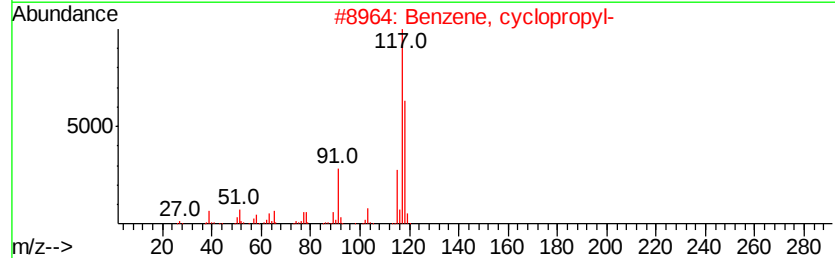
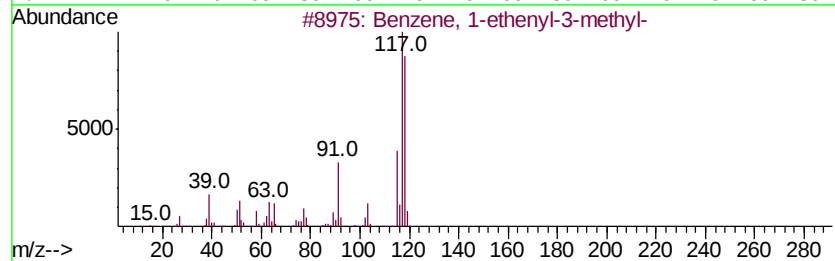
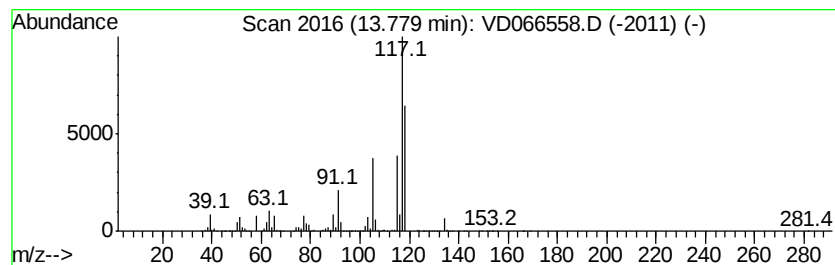
Instrument :
MSVOA_D
ClientSampleId :
TP-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	37.63 ug/l	57252300	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-3-methyl-	118 C9H10	000100-80-1 83
2		Benzene, cyclopropyl-	118 C9H10	000873-49-4 64
3		Indane	118 C9H10	000496-11-7 64
4		Benzene, 2-propenyl-	118 C9H10	000300-57-2 64
5		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

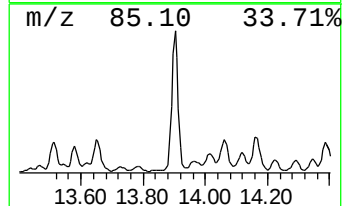
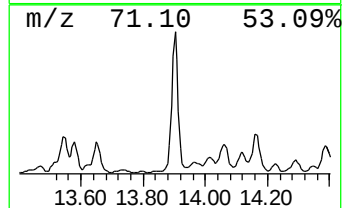
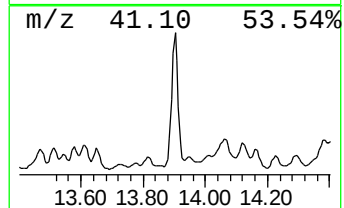
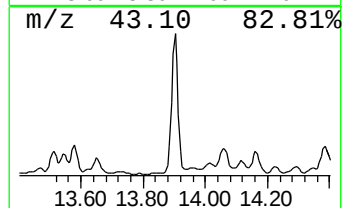
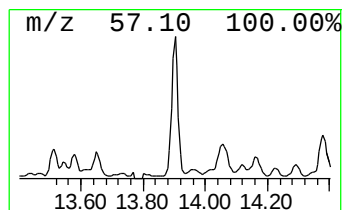
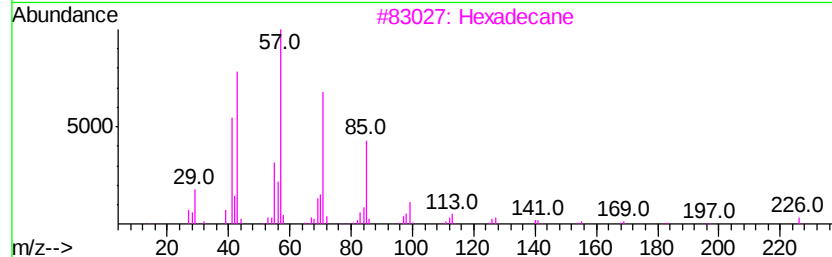
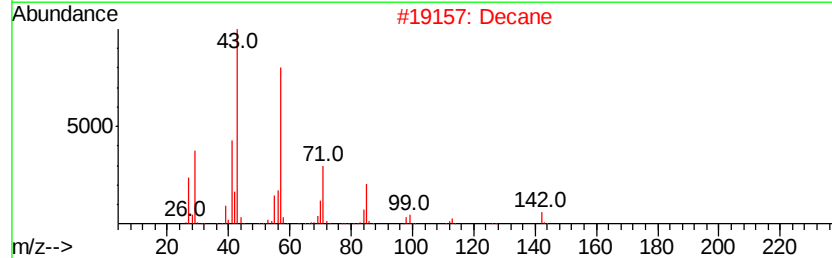
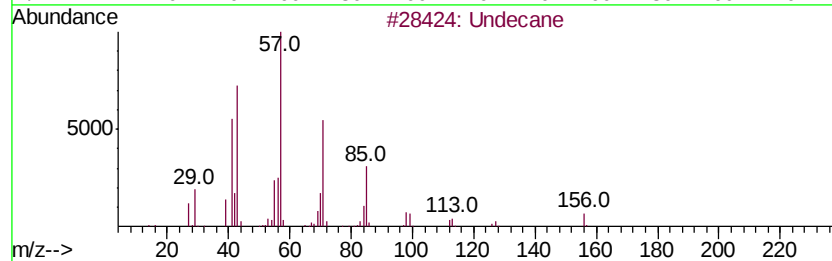
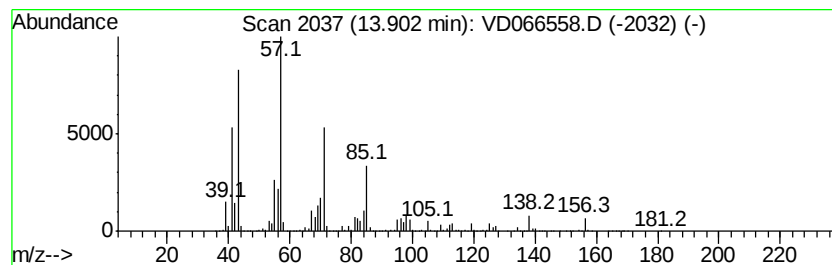
Instrument :
MSVOA_D
ClientSampleId :
TP-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 7 Undecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	24.78 ug/l	37697500	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Undecane	156 C11H24	001120-21-4	96
2	Decane	142 C10H22	000124-18-5	80
3	Hexadecane	226 C16H34	000544-76-3	80
4	Tetradecane	198 C14H30	000629-59-4	80
5	Heptadecane, 2,6,10,14-tetramethyl-	296 C21H44	018344-37-1	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampled :
TP-1

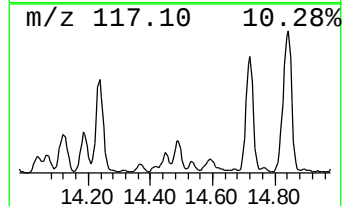
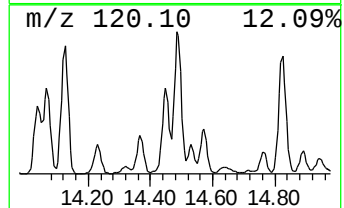
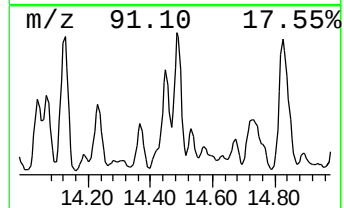
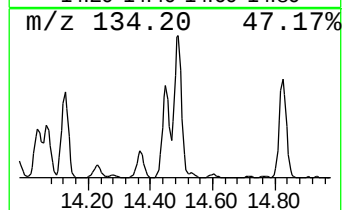
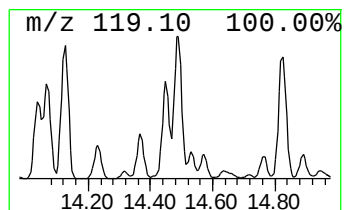
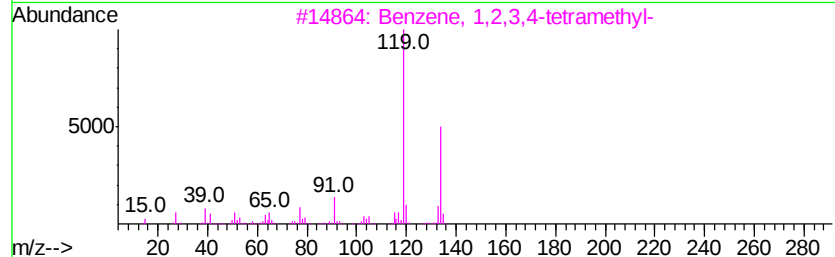
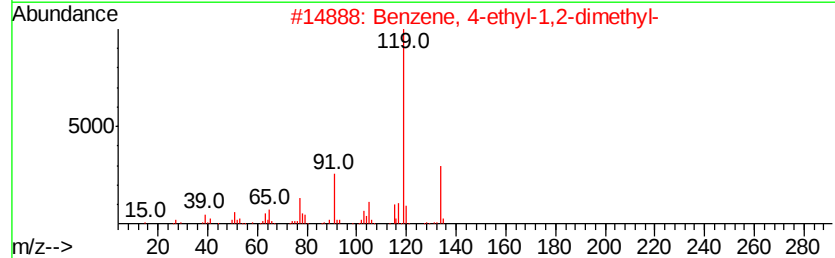
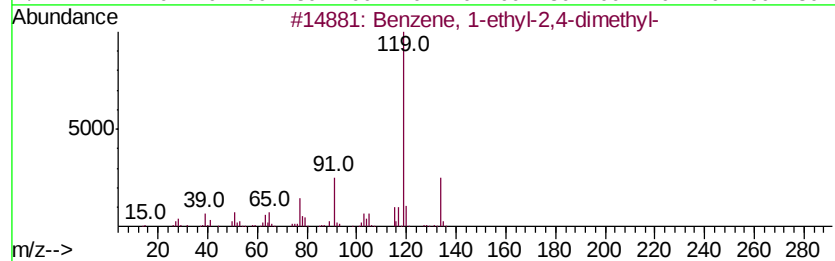
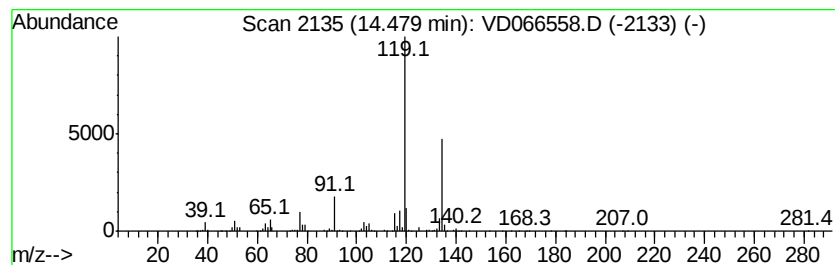
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	25.44 ug/l	38706400	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	93
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

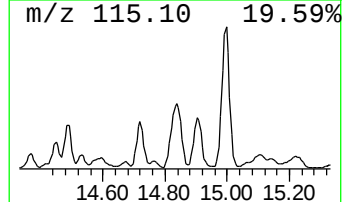
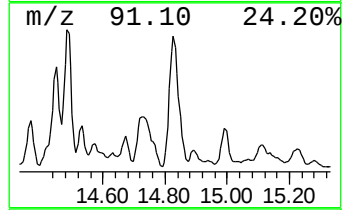
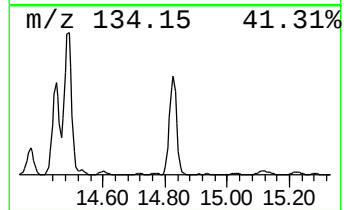
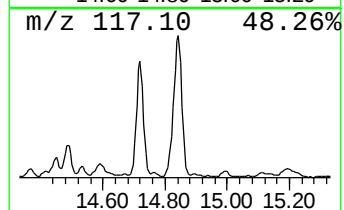
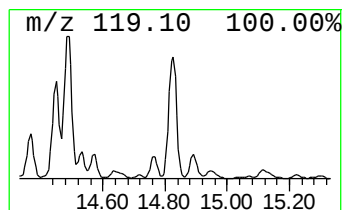
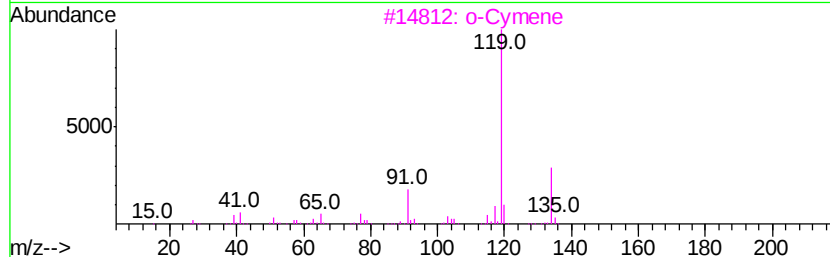
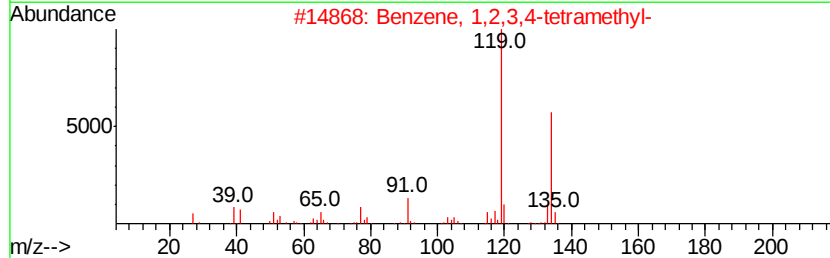
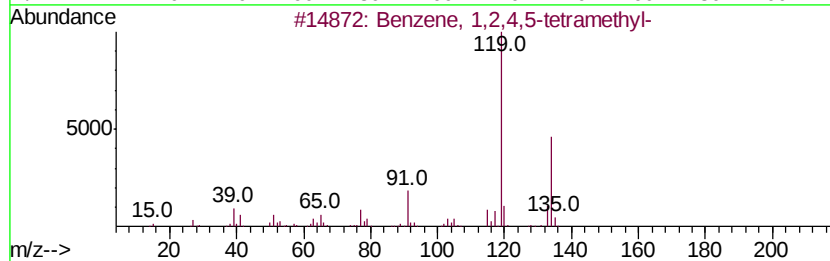
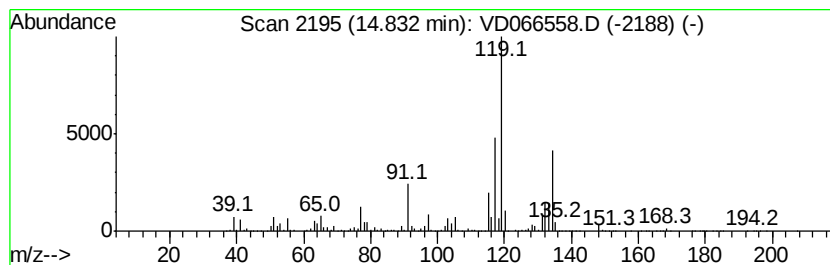
Instrument :
MSVOA_D
ClientSampleId :
TP-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	35.23 ug/l	53606200	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 95
2		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 81
3		o-Cymene	134 C10H14	000527-84-4 81
4		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 81
5		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TP-1

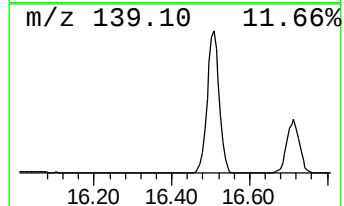
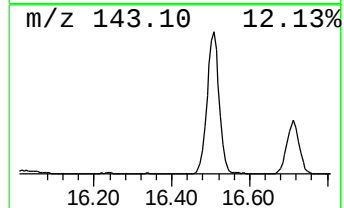
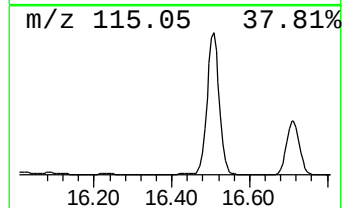
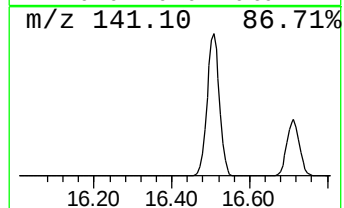
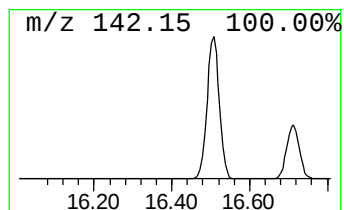
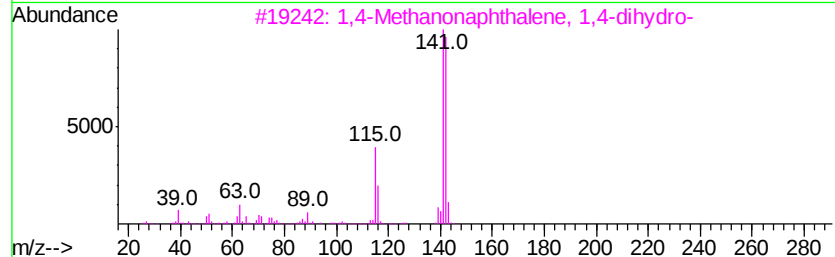
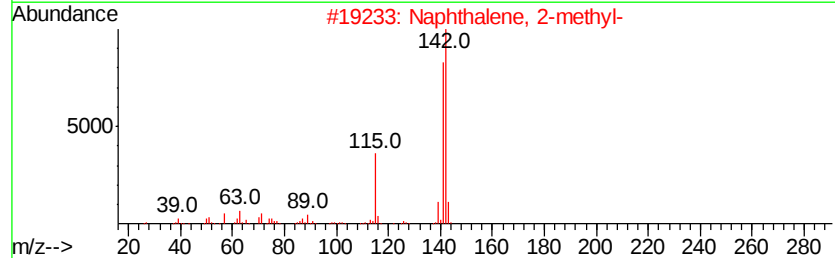
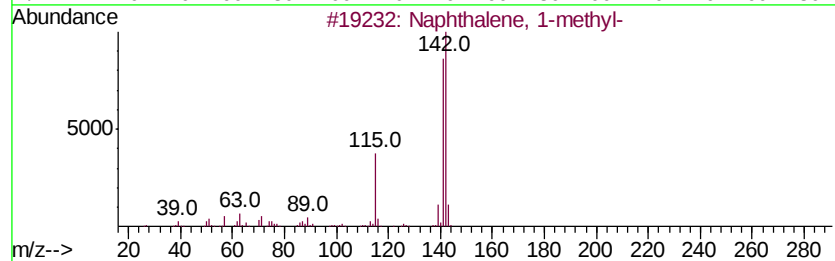
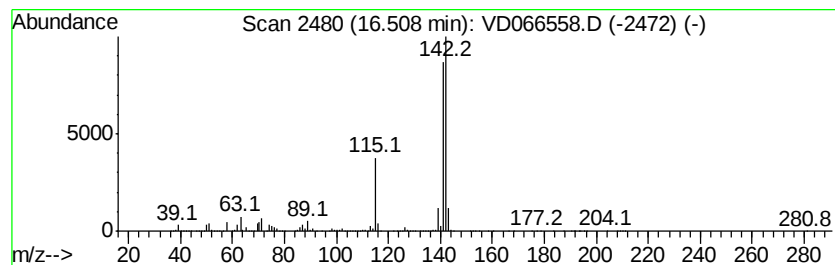
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	26.28 ug/l	39988700	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		1,4-Methanonaphthalene, 1,4-dihv...	142	C11H10	004453-90-1	91
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD090420\
Data File : VD066558.D
Acq On : 04 Sep 2020 13:00
Operator : VA/SY
Sample : L3877-01
Misc : 6.78G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
TP-1

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.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
Cyclopentene, 1-m...	7.73	50.3	ug/l	2972180	1	7.98	2955930	50.0
Heptane, 2-methyl-	9.97	29.6	ug/l	5500050	2	8.86	9277280	50.0
Heptane, 3-methyl-	10.11	24.3	ug/l	4507060	2	8.86	9277280	50.0
Benzene, 1-ethyl-...	12.88	31.6	ug/l	48070300	4	13.58	76078100	50.0
Benzene, 1-ethyl-...	13.13	26.1	ug/l	39789800	4	13.58	76078100	50.0
Benzene, 1-etheny...	13.78	37.6	ug/l	57252300	4	13.58	76078100	50.0
Undecane	13.90	24.8	ug/l	37697500	4	13.58	76078100	50.0
Benzene, 1-ethyl-...	14.48	25.4	ug/l	38706400	4	13.58	76078100	50.0
Benzene, 1,2,4,5-...	14.83	35.2	ug/l	53606200	4	13.58	76078100	50.0
Naphthalene, 1-me...	16.51	26.3	ug/l	39988700	4	13.58	76078100	50.0