

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
 Data File : VY001275.D
 Acq On : 17 Jan 2020 19:02
 Operator : SY/MD
 Sample : L1006-01
 Misc : 5.00G/5ML/MSVOA Y/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 HD-01-011520

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_Y\METHODS\82Y011020S.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	4.767	485	497	509	rBV	17163	55022	4.21%	0.332%
2	7.761	978	988	994	rBV	169688	412100	31.56%	2.484%
3	7.828	994	999	1011	rVB	295215	661347	50.65%	3.986%
4	8.181	1047	1057	1068	rBV	177663	392213	30.04%	2.364%
5	8.724	1136	1146	1155	rBV	496376	957499	73.33%	5.771%
6	10.205	1380	1389	1397	rBV	738016	1305693	100.00%	7.869%
7	11.108	1531	1537	1543	rBV2	19883	35354	2.71%	0.213%
8	11.217	1549	1555	1559	rBV3	7888	15188	1.16%	0.092%
9	11.309	1566	1570	1579	rVB2	16036	28398	2.17%	0.171%
10	11.504	1595	1602	1613	rBV	687731	1122605	85.98%	6.766%
11	11.644	1620	1625	1628	rBV2	11440	15184	1.16%	0.092%
12	11.693	1628	1633	1637	rBV6	14789	30097	2.31%	0.181%
13	11.748	1638	1642	1646	rVV	28582	51084	3.91%	0.308%
14	11.796	1646	1650	1655	rVB2	24287	42282	3.24%	0.255%
15	11.857	1655	1660	1664	rBV3	15680	32397	2.48%	0.195%
16	11.943	1670	1674	1679	rVB2	18882	32192	2.47%	0.194%
17	12.016	1679	1686	1692	rBV5	83350	193805	14.84%	1.168%
18	12.071	1692	1695	1697	rVV2	22108	33796	2.59%	0.204%
19	12.101	1697	1700	1708	rVB6	26620	57777	4.43%	0.348%
20	12.211	1708	1718	1720	rBV4	85982	173865	13.32%	1.048%
21	12.248	1720	1724	1730	rVV2	154438	325628	24.94%	1.962%
22	12.333	1734	1738	1741	rVV4	46570	94076	7.21%	0.567%
23	12.382	1741	1746	1753	rVB4	79237	187353	14.35%	1.129%
24	12.491	1758	1764	1770	rBV	551723	902487	69.12%	5.439%
25	12.583	1770	1779	1782	rBV7	48413	124818	9.56%	0.752%
26	12.656	1786	1791	1798	rVB	158674	290616	22.26%	1.751%
27	12.735	1798	1804	1809	rBV3	86389	207548	15.90%	1.251%
28	12.821	1809	1818	1822	rBV4	188789	453249	34.71%	2.732%
29	12.869	1823	1826	1831	rVB2	123364	190621	14.60%	1.149%
30	12.924	1831	1835	1837	rBV3	29539	44951	3.44%	0.271%
31	12.961	1837	1841	1844	rBV3	92687	136385	10.45%	0.822%
32	13.004	1844	1848	1852	rVV5	39661	88314	6.76%	0.532%
33	13.046	1852	1855	1862	rVB2	132179	223760	17.14%	1.349%
34	13.107	1863	1865	1868	rBV2	19268	24557	1.88%	0.148%

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Integration Parameters: RTEINT.P

Integrator: RTE
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35	13.174	1873	1876	1881	rBV3	79590	111193	8.52%	0.670%
36	13.229	1881	1885	1888	rBV2	93197	134575	10.31%	0.811%
37	13.327	1893	1901	1905	rBV3	256134	534416	40.93%	3.221%
38	13.436	1913	1919	1932	rVB	652244	1157822	88.67%	6.978%
39	13.577	1938	1942	1945	rBV3	71708	117498	9.00%	0.708%
40	13.680	1955	1959	1965	rBV4	44919	96768	7.41%	0.583%
41	13.753	1966	1971	1976	rBV4	304874	563153	43.13%	3.394%
42	13.802	1976	1979	1983	rVV2	68428	92646	7.10%	0.558%
43	13.863	1986	1989	1992	rVB5	42602	60295	4.62%	0.363%
44	13.973	2004	2007	2012	rVB3	114650	191059	14.63%	1.151%
45	14.022	2012	2015	2020	rVB3	65373	95494	7.31%	0.576%
46	14.083	2020	2025	2030	rVB3	129543	218566	16.74%	1.317%
47	14.150	2031	2036	2042	rBV5	73461	153614	11.76%	0.926%
48	14.205	2042	2045	2049	rBV3	44411	66315	5.08%	0.400%
49	14.265	2050	2055	2060	rBV3	228515	438009	33.55%	2.640%
50	14.387	2071	2075	2078	rBV	73940	100947	7.73%	0.608%
51	14.424	2078	2081	2088	rVV5	139531	203516	15.59%	1.227%
52	14.485	2089	2091	2093	rVV3	21331	19282	1.48%	0.116%
53	14.619	2110	2113	2118	rVB3	55839	73375	5.62%	0.442%
54	14.686	2118	2124	2130	rVV3	158091	339047	25.97%	2.043%
55	14.741	2130	2133	2139	rVB4	100680	165630	12.69%	0.998%
56	14.845	2146	2150	2154	rVB2	56727	87525	6.70%	0.527%
57	14.948	2163	2167	2171	rBV2	93203	144089	11.04%	0.868%
58	14.985	2171	2173	2180	rVV4	50251	81215	6.22%	0.489%
59	15.058	2180	2185	2192	rVB2	107725	228330	17.49%	1.376%
60	15.217	2206	2211	2214	rBV6	50574	94033	7.20%	0.567%
61	15.247	2214	2216	2220	rVV4	39382	69144	5.30%	0.417%
62	15.296	2220	2224	2228	rVV	97130	162353	12.43%	0.978%
63	15.369	2228	2236	2244	rVV6	69578	267173	20.46%	1.610%
64	15.436	2244	2247	2255	rVB5	39615	79407	6.08%	0.479%
65	15.527	2255	2262	2269	rBV4	54914	140693	10.78%	0.848%
66	15.607	2271	2275	2280	rVB5	23309	40881	3.13%	0.246%
67	15.692	2280	2289	2291	rBV4	63549	147064	11.26%	0.886%
68	15.729	2292	2295	2303	rVB2	82019	140423	10.75%	0.846%
69	15.814	2305	2309	2315	rVB5	29104	57037	4.37%	0.344%
70	15.893	2315	2322	2327	rBV4	50658	124633	9.55%	0.751%
71	16.033	2338	2345	2351	rBV3	104557	222083	17.01%	1.338%

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72	16.229	2369	2377	2383	rBV3	121048	260285	19.93%	1.569%
73	16.296	2383	2388	2393	rBV	43309	81265	6.22%	0.490%
74	16.497	2413	2421	2426	rBV6	42696	131664	10.08%	0.794%
75	16.546	2427	2429	2435	rVB2	33251	49850	3.82%	0.300%
76	16.625	2437	2442	2450	rVB4	44395	108076	8.28%	0.651%

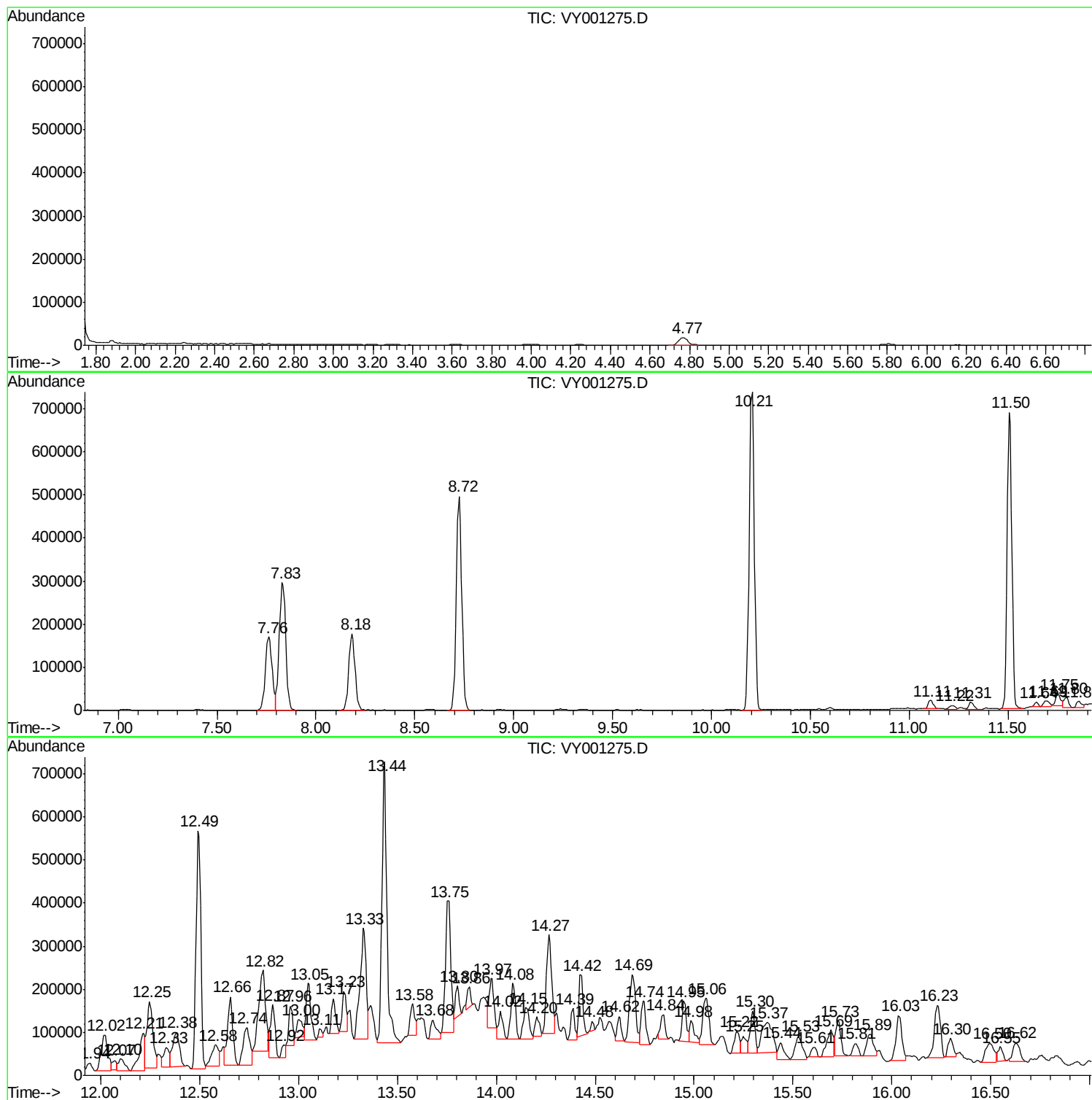
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y011020S.M
Quant Title : SW846 8260

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



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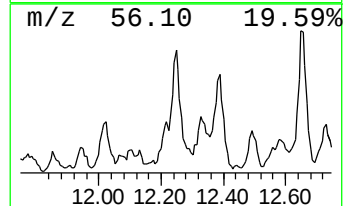
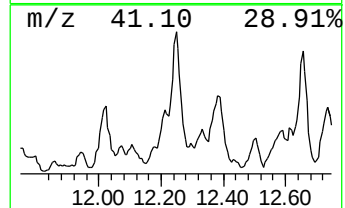
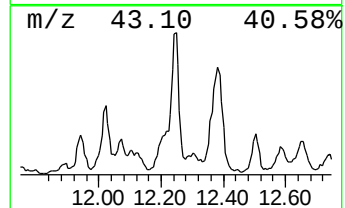
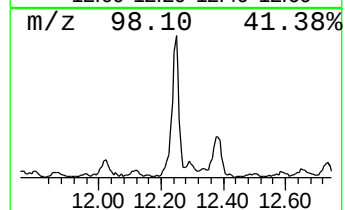
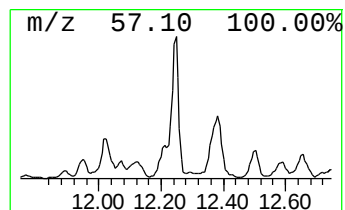
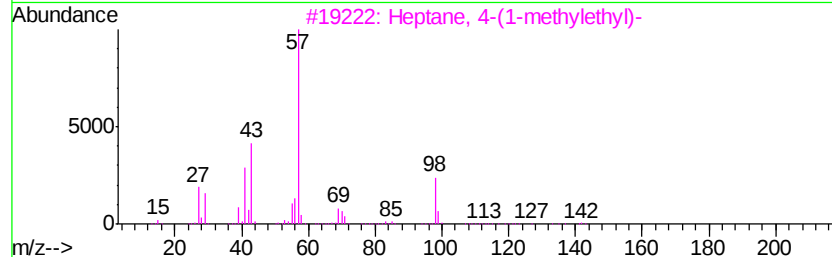
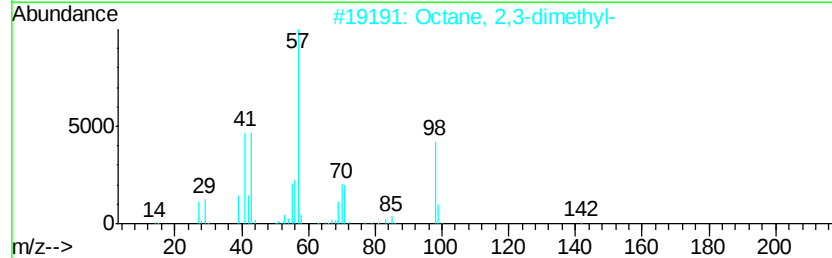
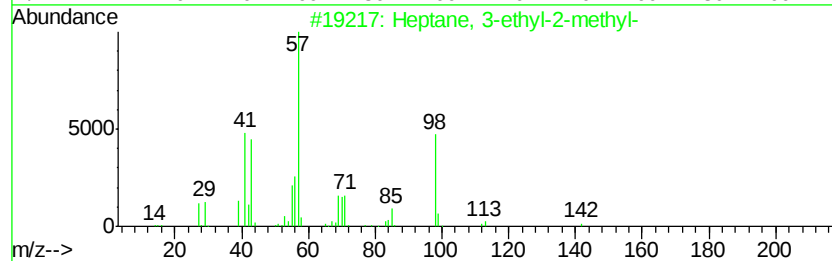
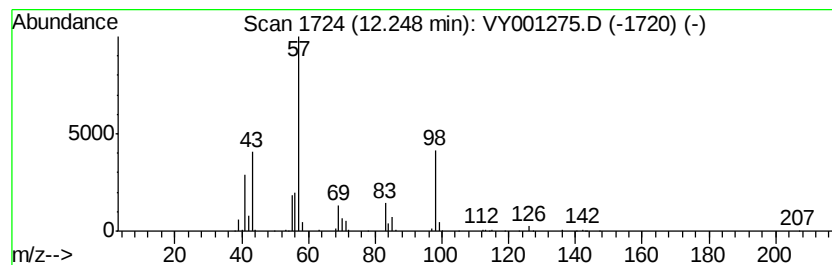
Instrument :
MSVOA_Y
ClientSampleId :
HD-01-011520

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y011020S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.25	14.50 ug/l	325628	Chlorobenzene-d5	11.50		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	83
2		Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	64
3		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	47
4		Cyclohexanone, 4-(1,1-dimethylet...	154	C10H18O	000098-53-3	42
5		Ether, tert-butyl isopropylidene...	154	C10H18O	024524-56-9	42



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ALS Vial : 12 Sample Multiplier: 1

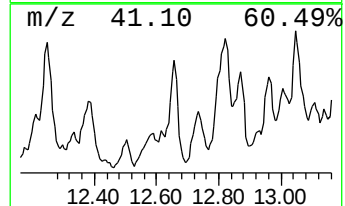
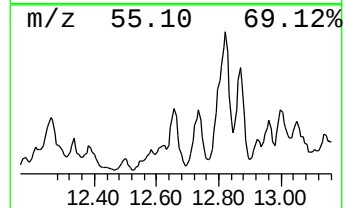
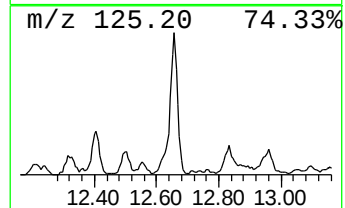
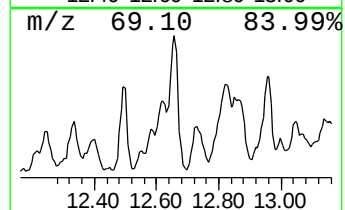
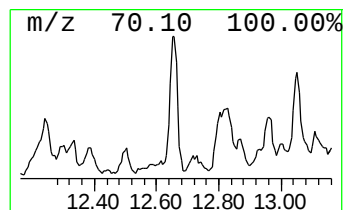
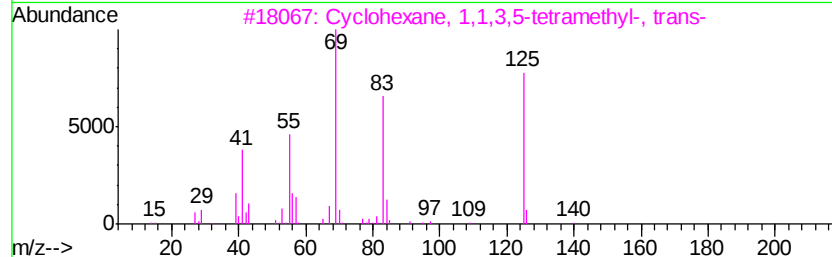
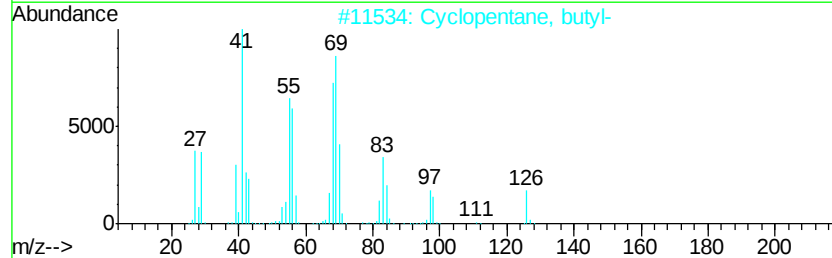
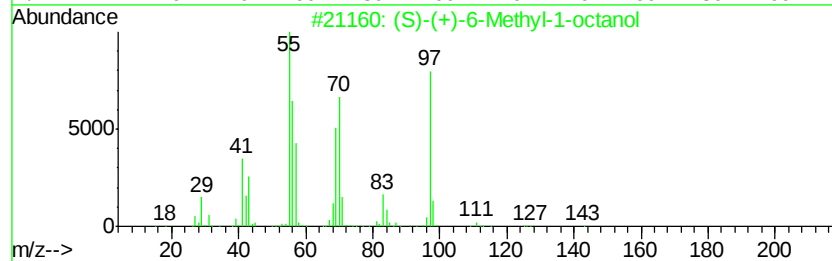
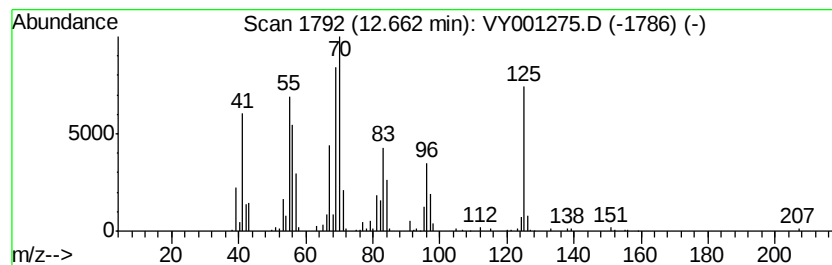
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (S)-(+)-6-Methyl-1-octanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.66	12.55 ug/l	290616	1,4-Dichlorobenzene-d4	13.44		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(S)-(+)-6-Methyl-1-octanol	144	C9H20O	110453-78-6	53	
2	Cyclopentane, butyl-	126	C9H18	002040-95-1	50	
3	Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	43	
4	Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	43	
5	Octane, 2-chloro-	148	C8H17Cl	000628-61-5	38	



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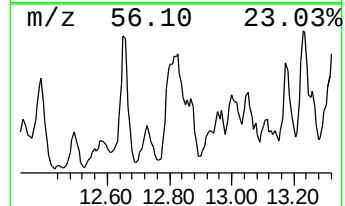
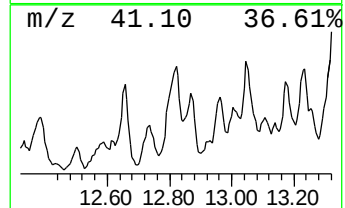
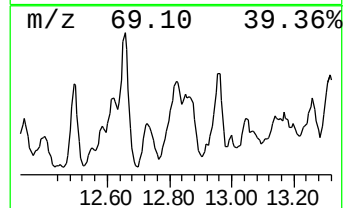
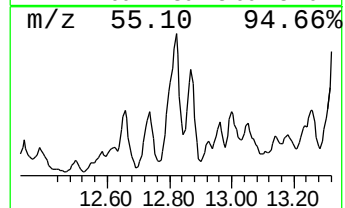
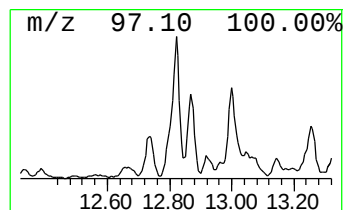
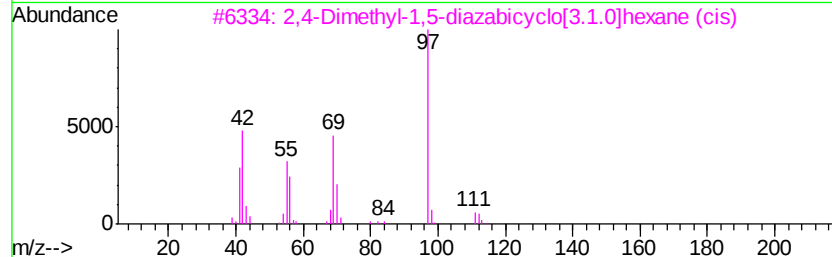
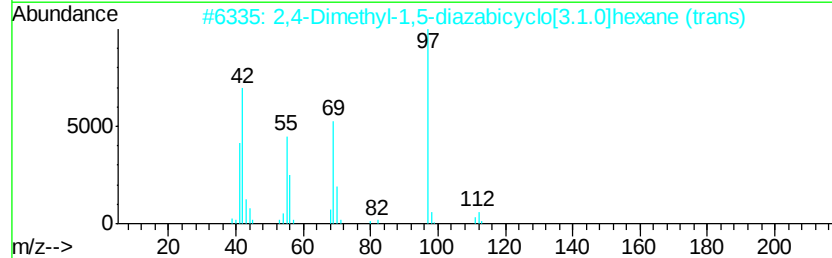
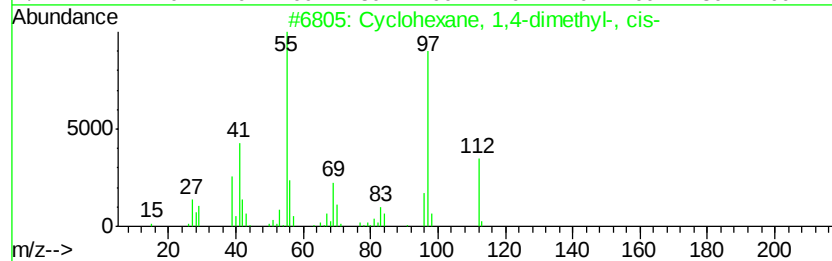
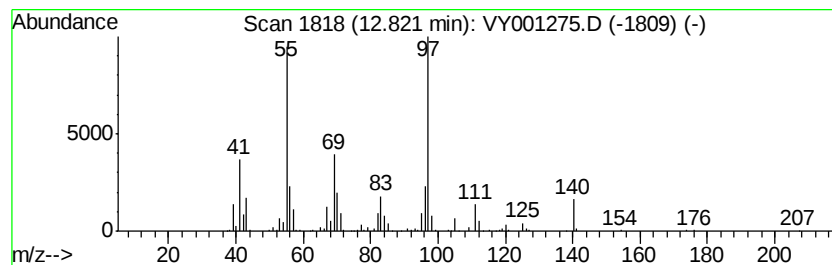
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1,4-dimethyl-,... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.82	19.57 ug/l	453249	1,4-Dichlorobenzene-d4	13.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	72
2		2,4-Dimethyl-1,5-diazabicyclo[3.1.0]hexane (trans)	112	C6H12N2	1000283-20-9	53
3		2,4-Dimethyl-1,5-diazabicyclo[3.1.0]hexane (cis)	112	C6H12N2	100463-01-2	52
4		Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	49
5		Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	49



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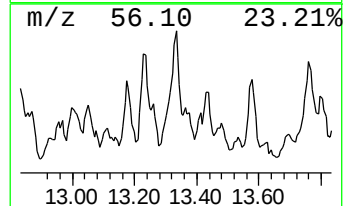
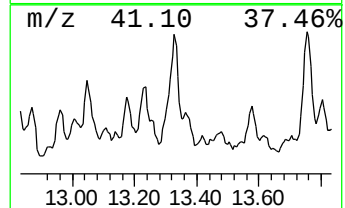
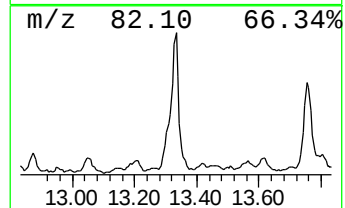
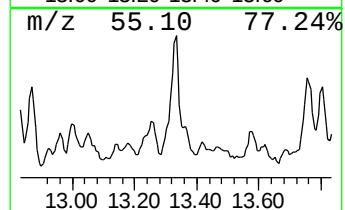
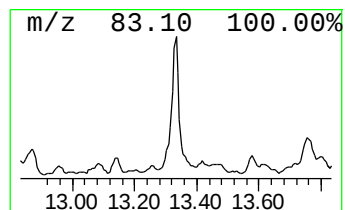
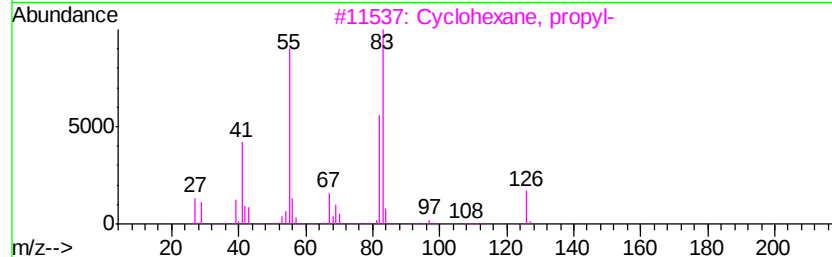
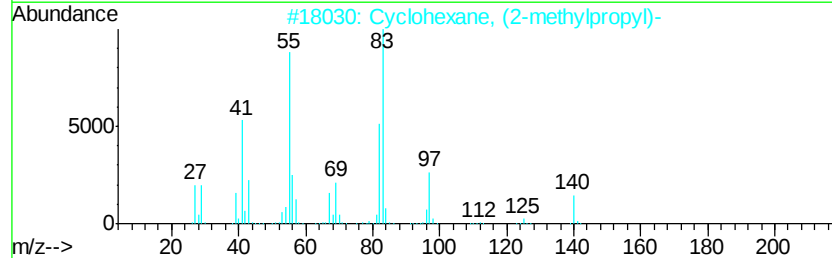
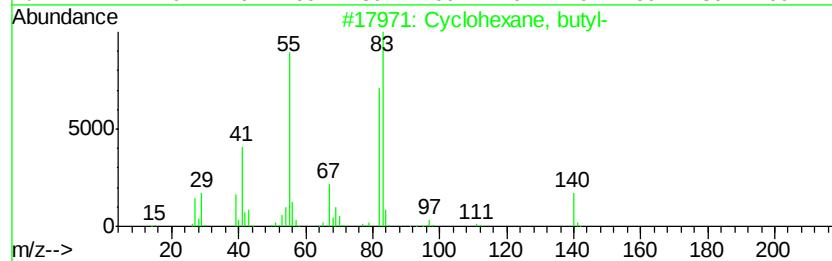
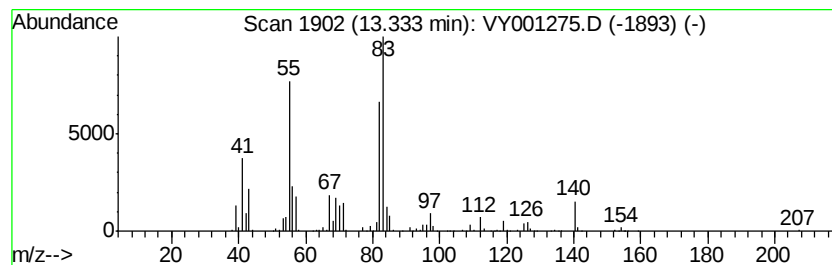
Instrument :
MSVOA_Y
ClientSampleId :
HD-01-011520

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y011020S.M
Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, butyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	23.08 ug/l	534416	1,4-Dichlorobenzene-d4	13.44
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, butyl-	140 C10H20	001678-93-9 74
2		Cyclohexane, (2-methylpropyl)-	140 C10H20	001678-98-4 74
3		Cyclohexane, propyl-	126 C9H18	001678-92-8 64
4		Cyclohexane, pentyl-	154 C11H22	004292-92-6 64
5		Cyclohexane, (1-methylethyl)-	126 C9H18	000696-29-7 53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HD-01-011520

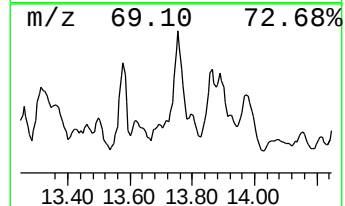
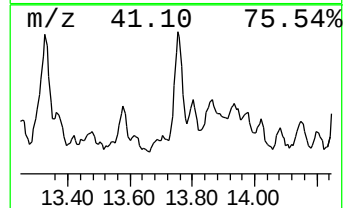
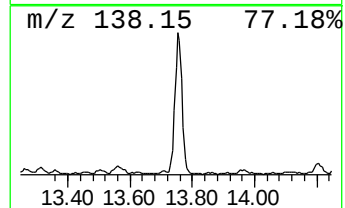
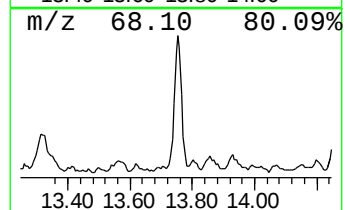
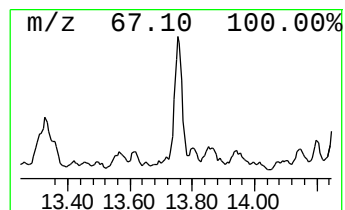
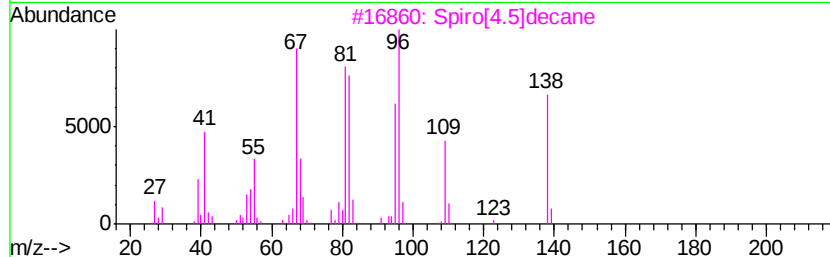
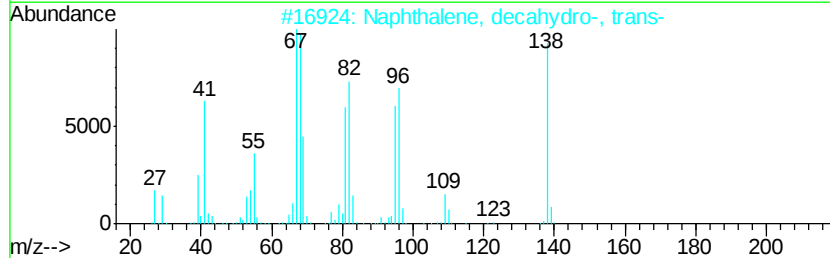
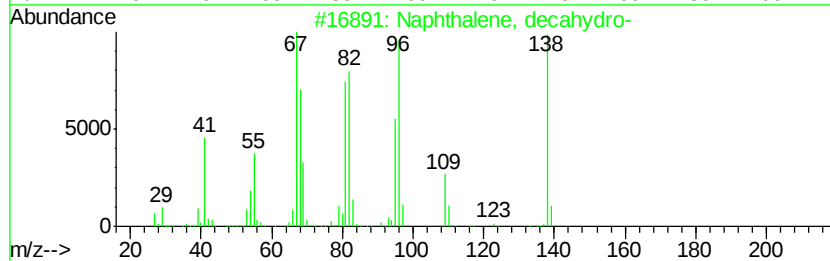
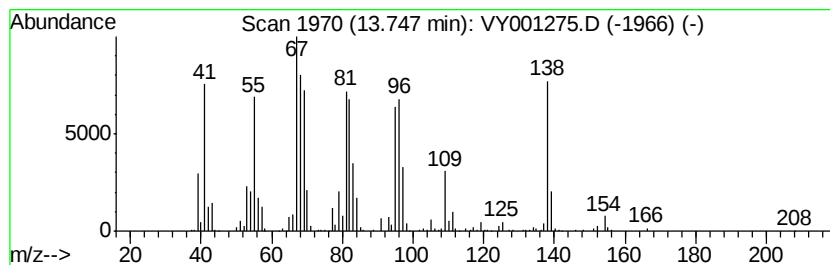
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, decahydro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	24.32 ug/l	563153	1,4-Dichlorobenzene-d4	13.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, decahydro-	138	C10H18	000091-17-8	91
2		Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	81
3		Spiro[4.5]decane	138	C10H18	000176-63-6	76
4		Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	70
5		Cyclodecene	138	C10H18	003618-12-0	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

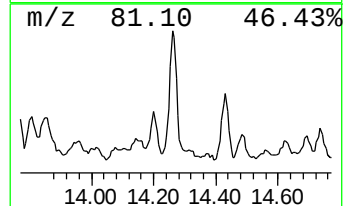
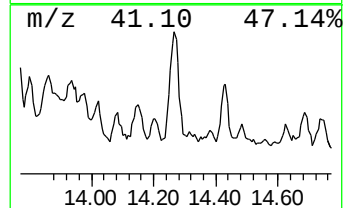
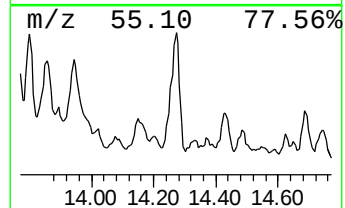
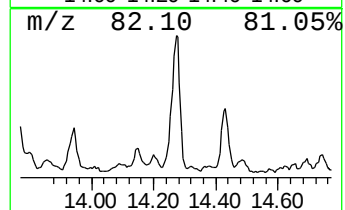
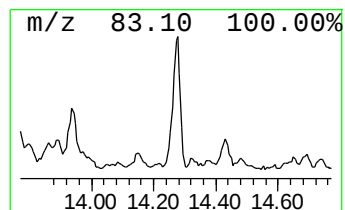
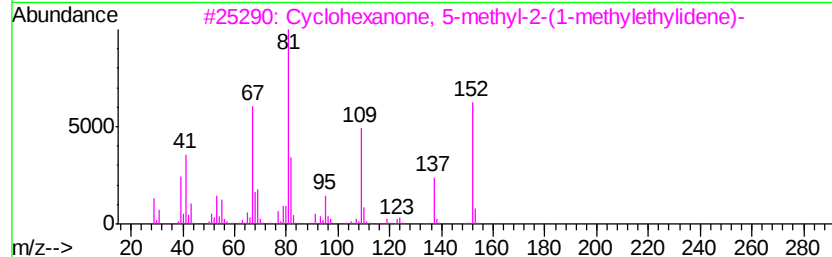
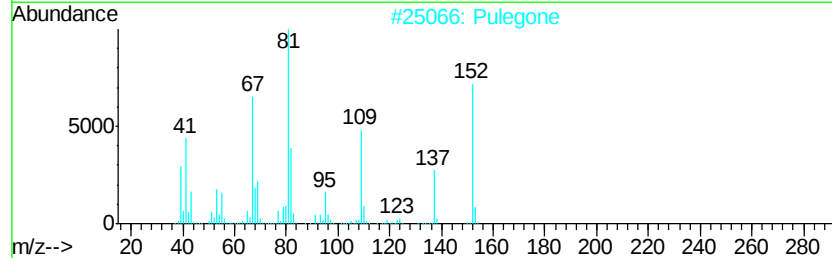
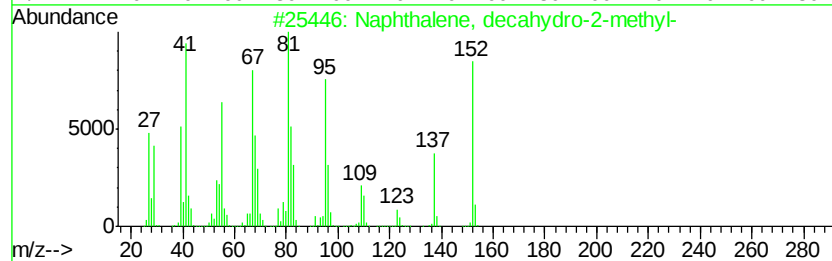
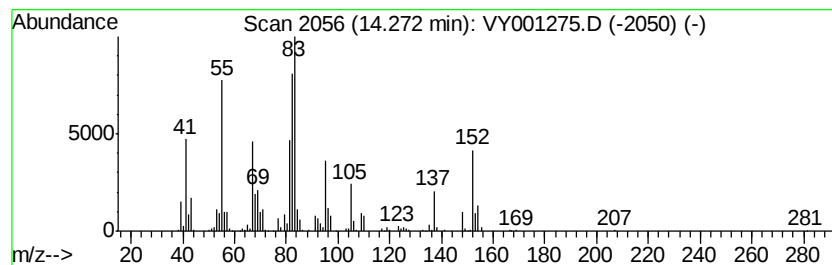
Instrument :
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ClientSampleId :
HD-01-011520

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y011020S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, decahydro-2-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	18.92 ug/l	438009	1,4-Dichlorobenzene-d4	13.44
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 81
2		Pulegone	152 C10H16O	000089-82-7 55
3		Cyclohexanone, 5-methyl-2-(1-met...	152 C10H16O	015932-80-6 49
4		Bicyclo[3.1.1]heptane, 2,6,6-tri...	138 C10H18	006876-13-7 46
5		Cyclohexanone, 5-methyl-2-(1-met...	152 C10H16O	000529-00-0 45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

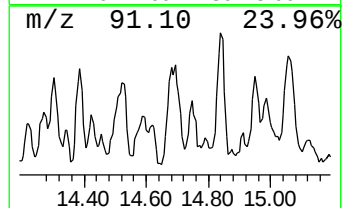
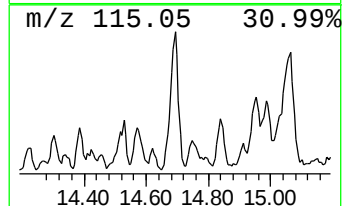
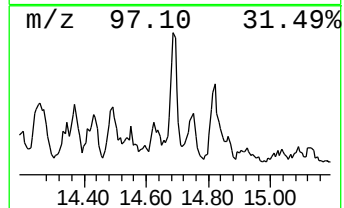
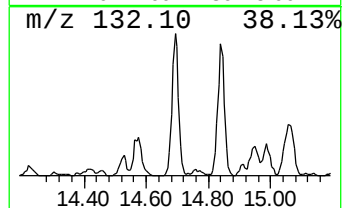
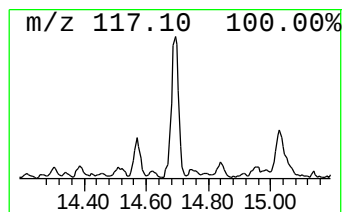
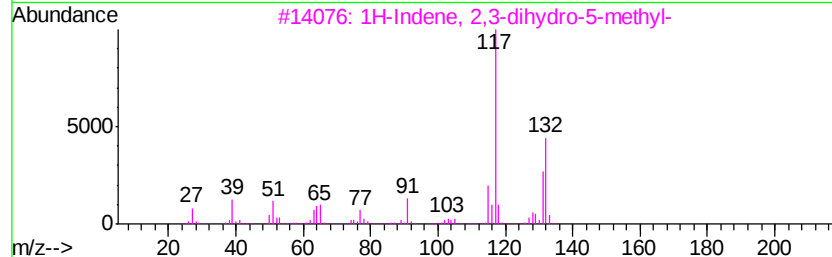
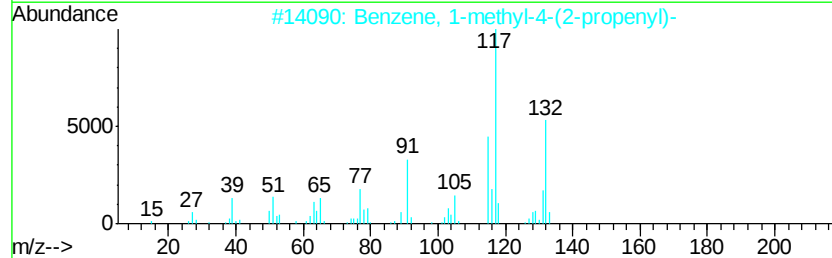
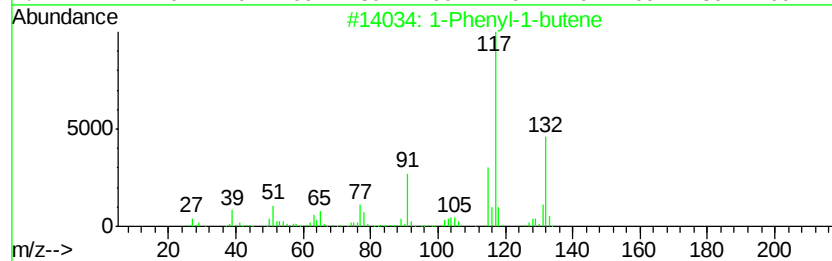
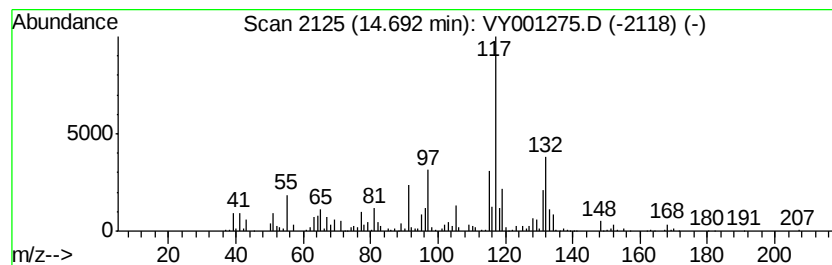
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ClientSampleId :
HD-01-011520

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y011020S.M
Quant Title : SW846 8260

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

Peak Number 7 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	14.64 ug/l	339047	1,4-Dichlorobenzene-d4	13.44
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Phenyl-1-butene	132 C10H12	000824-90-8	91
2	Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9	83
3	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	60
4	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000768-00-3	60
5	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HD-01-011520

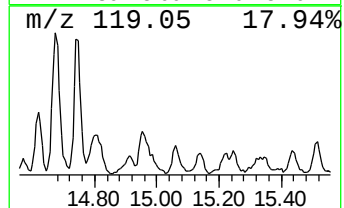
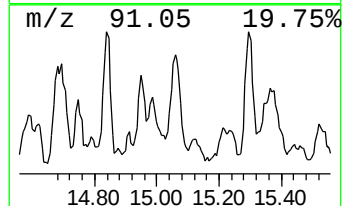
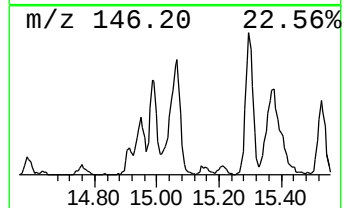
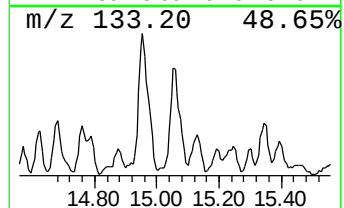
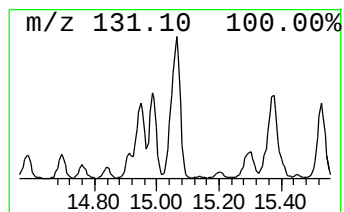
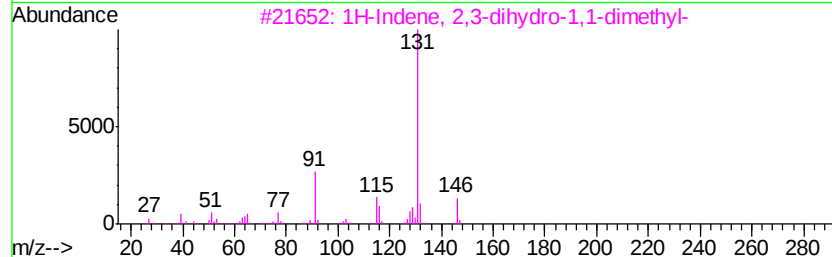
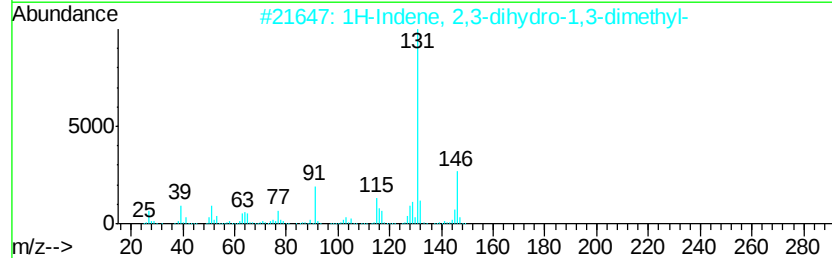
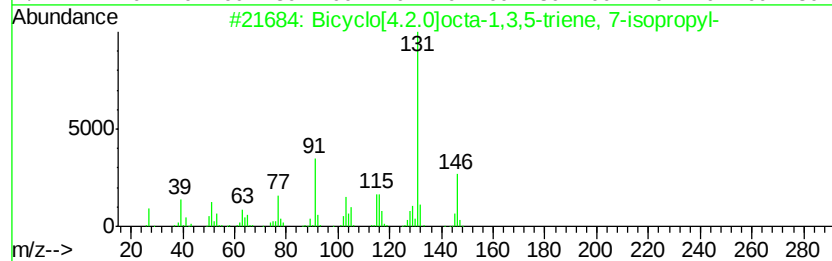
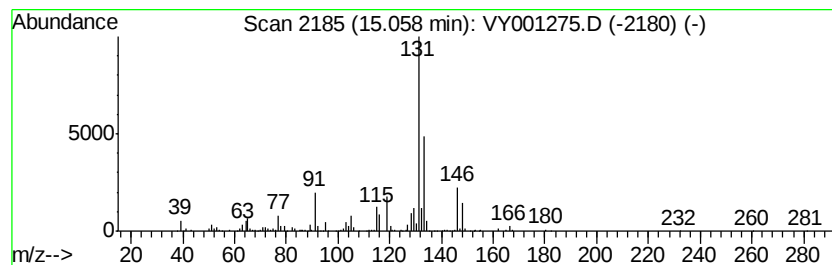
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Quant Title : SW846 8260

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

Peak Number 8 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	9.86 ug/l	228330	1,4-Dichlorobenzene-d4	13.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	64
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
3		1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HD-01-011520

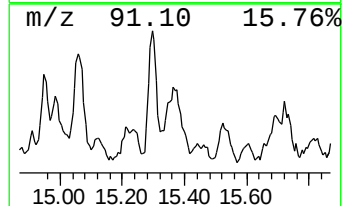
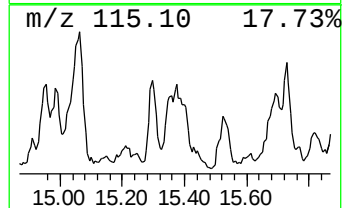
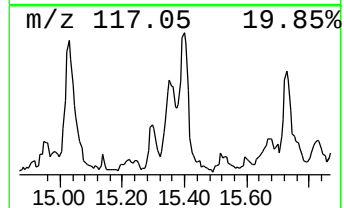
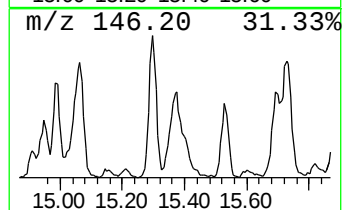
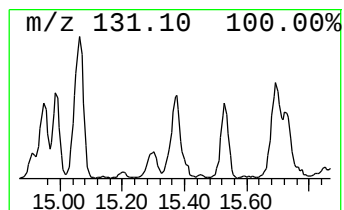
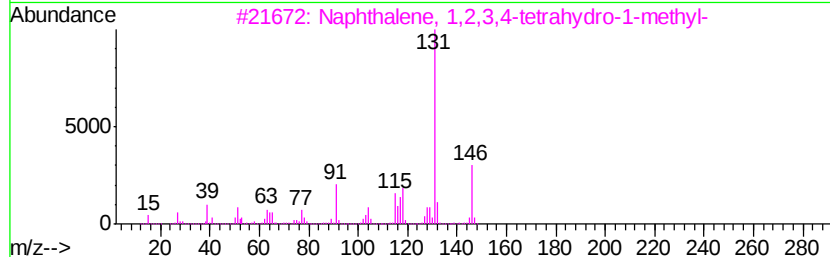
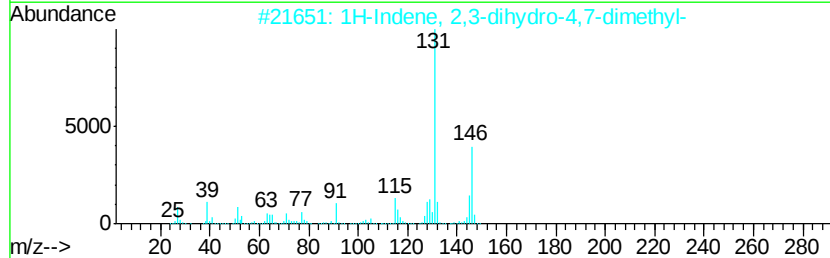
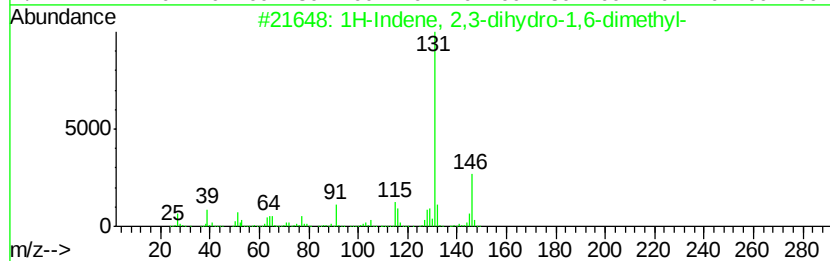
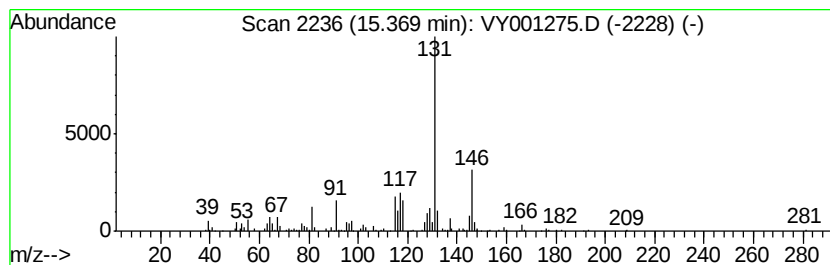
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	11.54 ug/l	267173	1,4-Dichlorobenzene-d4	13.44

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94	
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93	
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	90	
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90	
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	87	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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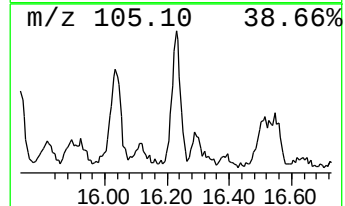
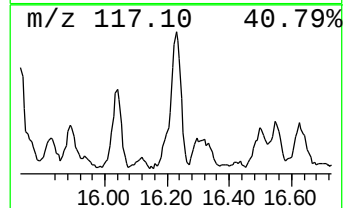
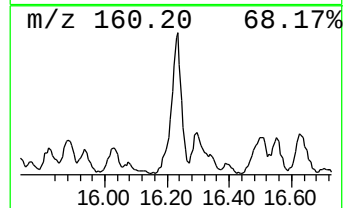
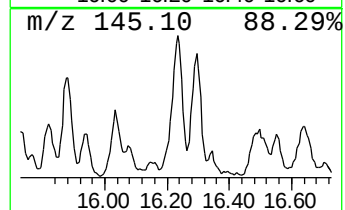
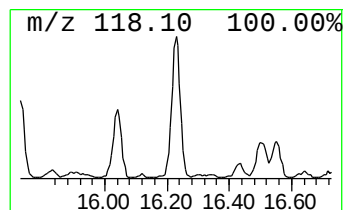
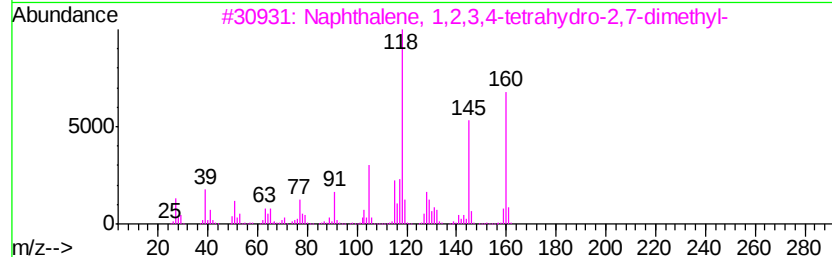
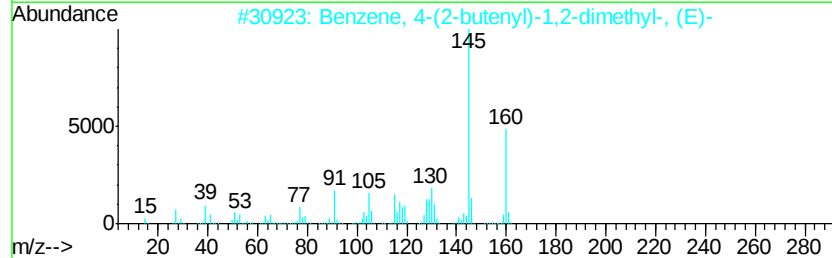
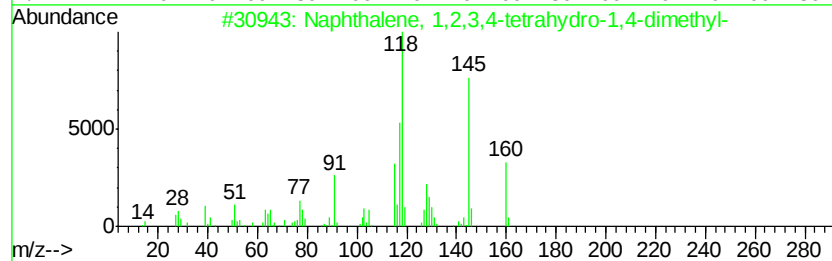
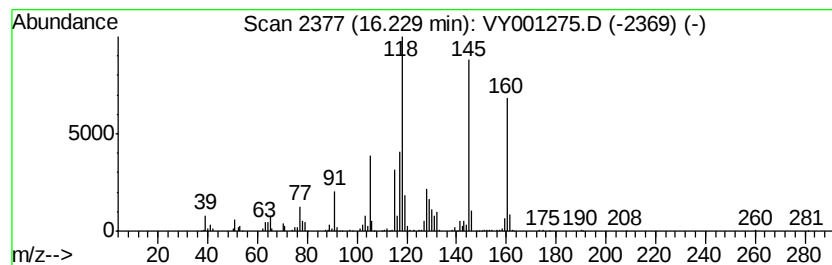
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	11.24 ug/l	260285	1,4-Dichlorobenzene-d4	13.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	94
2		Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	83
3		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	76
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	70
5		Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_Y\DATA\VY011720\
Data File : VY001275.D
Acq On : 17 Jan 2020 19:02
Operator : SY/MD
Sample : L1006-01
Misc : 5.00G/5ML/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HD-01-011520

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y011020S.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
Heptane, 3-ethyl-...	12.25	14.5	ug/l	325628	3	11.50	1122610	50.0
(S)-(+)-6-Methyl-...	12.66	12.6	ug/l	290616	4	13.44	1157820	50.0
Cyclohexane, 1,4-...	12.82	19.6	ug/l	453249	4	13.44	1157820	50.0
Cyclohexane, butyl-	13.33	23.1	ug/l	534416	4	13.44	1157820	50.0
Naphthalene, deca...	13.75	24.3	ug/l	563153	4	13.44	1157820	50.0
Naphthalene, deca...	14.27	18.9	ug/l	438009	4	13.44	1157820	50.0
1-Phenyl-1-butene	14.69	14.6	ug/l	339047	4	13.44	1157820	50.0
Bicyclo[4.2.0]oct...	15.06	9.9	ug/l	228330	4	13.44	1157820	50.0
1H-Indene, 2,3-di...	15.37	11.5	ug/l	267173	4	13.44	1157820	50.0
Naphthalene, 1,2,...	16.23	11.2	ug/l	260285	4	13.44	1157820	50.0