

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY060320\
 Data File : VY002826.D
 Acq On : 03 Jun 2020 12:29
 Operator : SY/MD
 Sample : L2840-03
 Misc : 5.34G/5ML/MSVOA Y/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 DRUM-OIL

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\82Y052720S.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	7.724	973	982	988	rBV	179452	430963	1.65%	0.145%
2	7.797	988	994	1005	rVB	327281	715888	2.74%	0.241%
3	8.151	1042	1052	1064	rVB2	160910	356608	1.37%	0.120%
4	8.693	1133	1141	1154	rBV	463145	922546	3.53%	0.311%
5	10.181	1378	1385	1390	rBV	794242	1370306	5.25%	0.461%
6	10.242	1390	1395	1402	rVB	496685	868899	3.33%	0.292%
7	11.278	1557	1565	1576	rVB5	216987	565337	2.16%	0.190%
8	11.382	1576	1582	1586	rBV	193100	312099	1.19%	0.105%
9	11.492	1594	1600	1606	rBV	633511	1025110	3.92%	0.345%
10	11.589	1609	1616	1625	rVV	1452411	2473138	9.47%	0.832%
11	11.699	1625	1634	1644	rVV	6555991	12195201	46.68%	4.105%
12	12.028	1676	1688	1696	rBV	4763740	7875811	30.15%	2.651%
13	12.199	1712	1716	1719	rBV	185422	297101	1.14%	0.100%
14	12.241	1719	1723	1728	rVV4	236420	446184	1.71%	0.150%
15	12.327	1732	1737	1742	rBV	634641	988880	3.79%	0.333%
16	12.479	1758	1762	1767	rVV	409177	601313	2.30%	0.202%
17	12.668	1785	1793	1798	rBV	1871177	3003241	11.50%	1.011%
18	12.735	1798	1804	1807	rVV	9128140	13440578	51.45%	4.524%
19	12.766	1807	1809	1813	rVV	3447434	5003405	19.15%	1.684%
20	12.808	1813	1816	1822	rVV	5034316	7686074	29.42%	2.587%
21	12.973	1830	1843	1850	rBV	4004202	6373152	24.40%	2.145%
22	13.040	1850	1854	1861	rVB	397874	623800	2.39%	0.210%
23	13.119	1861	1867	1873	rBV	17307020	26124562	100.00%	8.793%
24	13.253	1881	1889	1893	rBV2	544105	1088657	4.17%	0.366%
25	13.314	1893	1899	1904	rBV2	1117125	1732001	6.63%	0.583%
26	13.369	1904	1908	1911	rBV2	469211	662824	2.54%	0.223%
27	13.455	1911	1922	1928	rVV	5491943	8895688	34.05%	2.994%
28	13.595	1940	1945	1946	rBV	1215941	1590603	6.09%	0.535%
29	13.625	1946	1950	1954	rVV	6565363	10201933	39.05%	3.434%
30	13.668	1954	1957	1966	rVB3	3812893	6930543	26.53%	2.333%
31	13.753	1966	1971	1976	rBV3	2056902	3031399	11.60%	1.020%
32	13.820	1976	1982	1987	rBV	1431178	2198835	8.42%	0.740%
33	13.887	1988	1993	1995	rBV	2356991	3465216	13.26%	1.166%
34	13.918	1995	1998	2002	rVB	2449489	3335239	12.77%	1.123%

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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
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Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
 Title : SW846 8260

35	13.973	2002	2007	2012	rVB	3803931	5335688	20.42%	1.796%
36	14.034	2012	2017	2020	rBV	912683	1402012	5.37%	0.472%
37	14.083	2020	2025	2030	rVB	3990344	6150668	23.54%	2.070%
38	14.156	2030	2037	2041	rBV5	508727	1266531	4.85%	0.426%
39	14.217	2041	2047	2050	rBV2	1360668	2327222	8.91%	0.783%
40	14.259	2051	2054	2056	rVV4	425327	651258	2.49%	0.219%
41	14.296	2056	2060	2063	rVV	2002936	3155367	12.08%	1.062%
42	14.339	2063	2067	2071	rVB	3662072	5195547	19.89%	1.749%
43	14.381	2071	2074	2078	rBV	1545420	1987404	7.61%	0.669%
44	14.424	2078	2081	2085	rVB2	1071011	1291664	4.94%	0.435%
45	14.521	2092	2097	2100	rBV2	1312406	1776269	6.80%	0.598%
46	14.570	2100	2105	2108	rBV	3210611	4809449	18.41%	1.619%
47	14.613	2108	2112	2117	rVB	4597461	6310304	24.15%	2.124%
48	14.686	2117	2124	2129	rBV4	7035275	15847306	60.66%	5.334%
49	14.735	2129	2132	2139	rVB3	1633484	2759262	10.56%	0.929%
50	14.832	2139	2148	2154	rBV	4903514	7795289	29.84%	2.624%
51	14.912	2157	2161	2162	rBV2	865803	1135773	4.35%	0.382%
52	14.948	2162	2167	2170	rVV3	2449324	4328217	16.57%	1.457%
53	14.979	2170	2172	2177	rVV	1418187	1905793	7.30%	0.641%
54	15.058	2178	2185	2190	rVV4	2777269	5700647	21.82%	1.919%
55	15.137	2195	2198	2204	rVV4	1314899	2237961	8.57%	0.753%
56	15.235	2204	2214	2219	rVV2	2554033	7496143	28.69%	2.523%
57	15.296	2219	2224	2228	rVV	2377346	3774761	14.45%	1.271%
58	15.351	2228	2233	2239	rVV4	1841096	5539325	21.20%	1.865%
59	15.442	2243	2248	2255	rVB	5435377	8777795	33.60%	2.955%
60	15.527	2255	2262	2269	rBV3	2009476	4851180	18.57%	1.633%
61	15.607	2270	2275	2280	rVB3	755673	1437569	5.50%	0.484%
62	15.692	2280	2289	2291	rBV3	1417231	3235675	12.39%	1.089%
63	15.729	2291	2295	2303	rVB	4808813	8916708	34.13%	3.001%
64	15.820	2306	2310	2315	rVB3	584767	1043741	4.00%	0.351%
65	15.887	2315	2321	2333	rBV3	1313853	3441759	13.17%	1.158%
66	16.039	2338	2346	2352	rBV	3241800	7018529	26.87%	2.362%
67	16.161	2362	2366	2369	rBV2	328716	503485	1.93%	0.169%
68	16.235	2370	2378	2383	rVV3	2265377	4856763	18.59%	1.635%
69	16.296	2383	2388	2395	rVV	3668620	7189610	27.52%	2.420%
70	16.363	2396	2399	2409	rVB	1011768	1977277	7.57%	0.666%
71	16.497	2413	2421	2427	rBV2	2134979	5243580	20.07%	1.765%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Title : SW846 8260

72	16.631	2438	2443	2453	rVB3	601220	1585868	6.07%	0.534%
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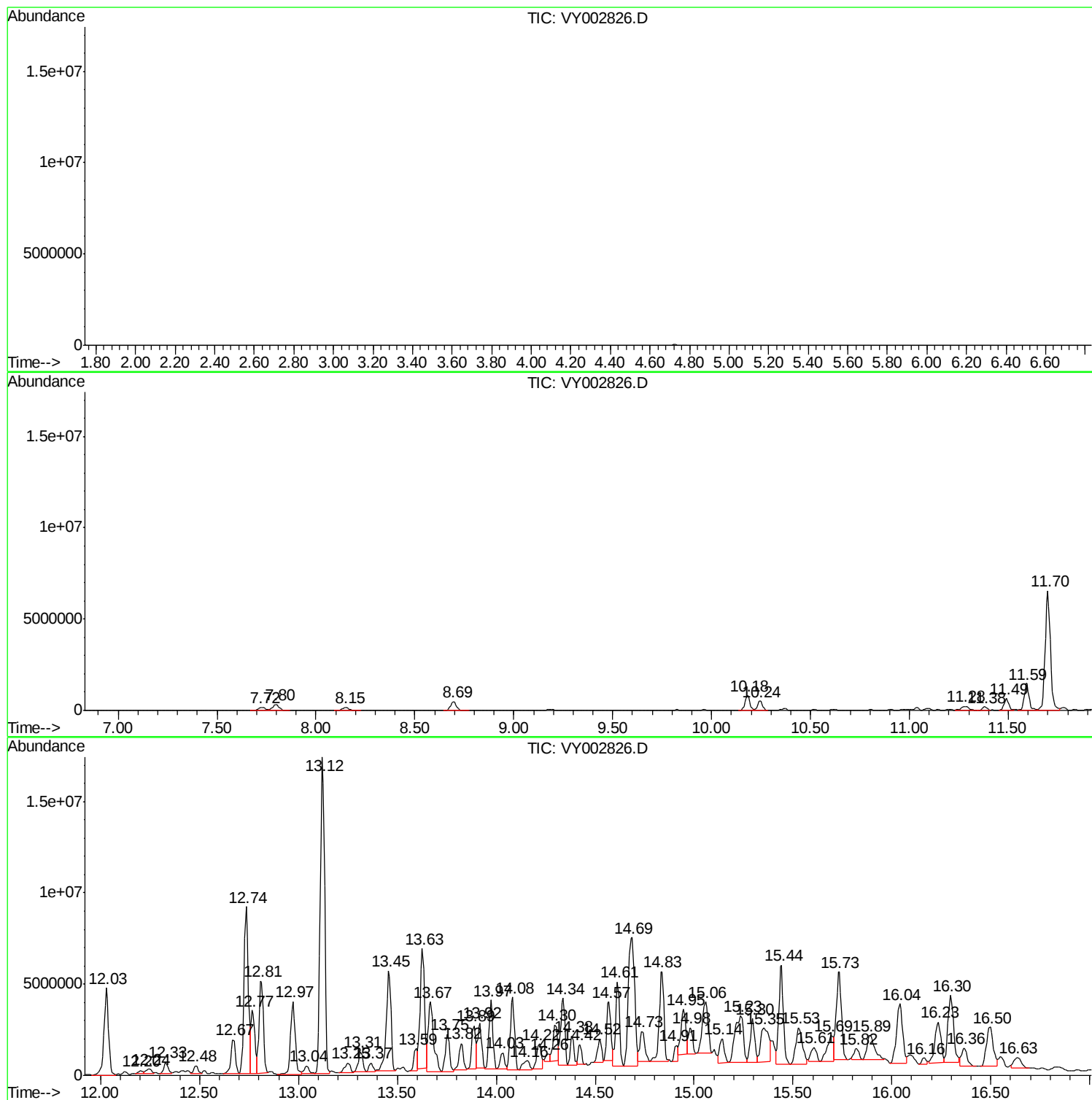
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Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY060320\
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Instrument :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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



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Instrument :
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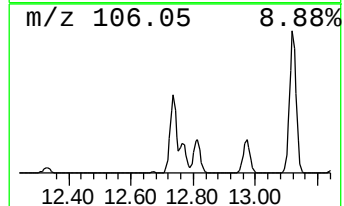
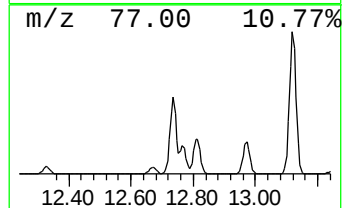
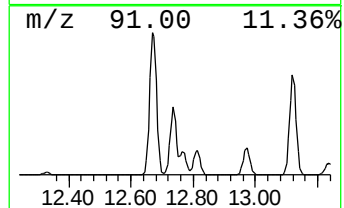
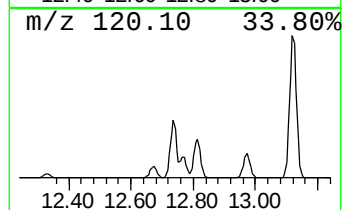
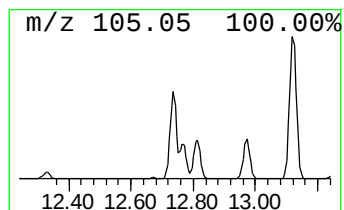
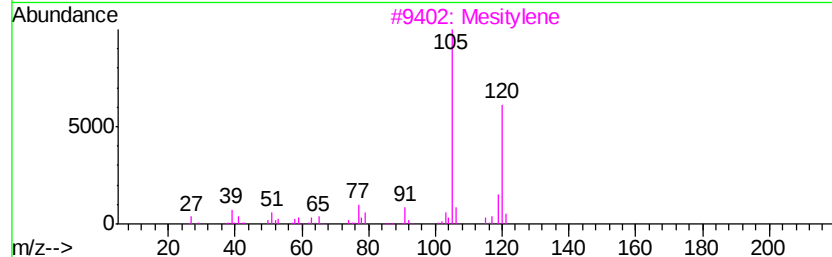
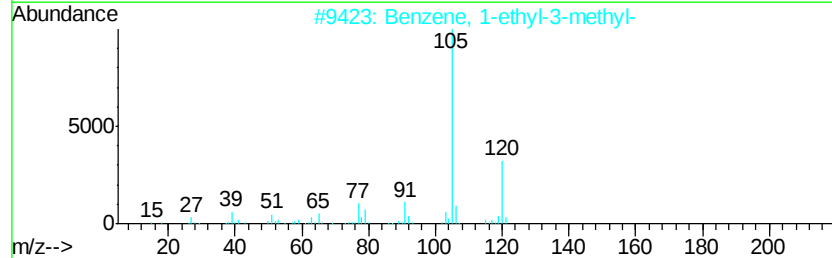
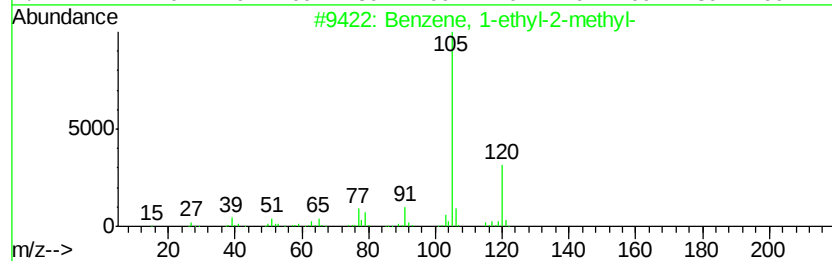
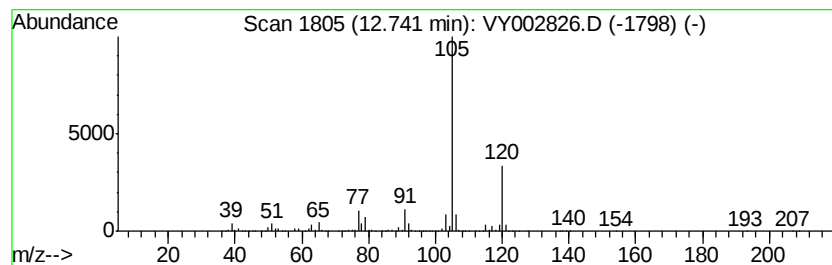
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	75.55 ug/l	13440600	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Misc : 5.34G/5ML/MSVOA Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

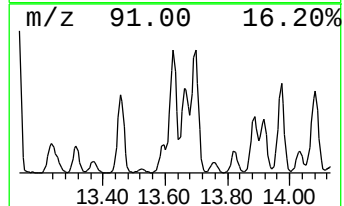
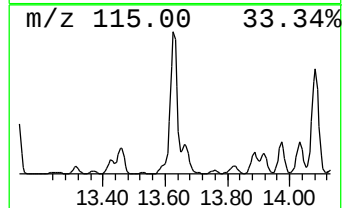
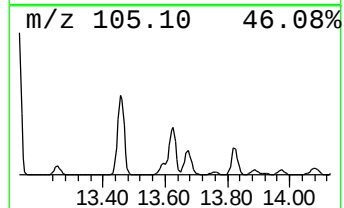
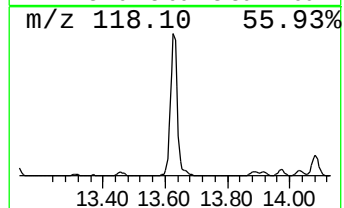
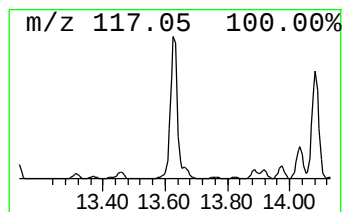
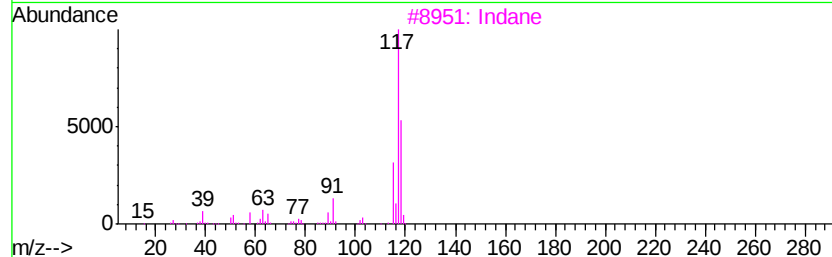
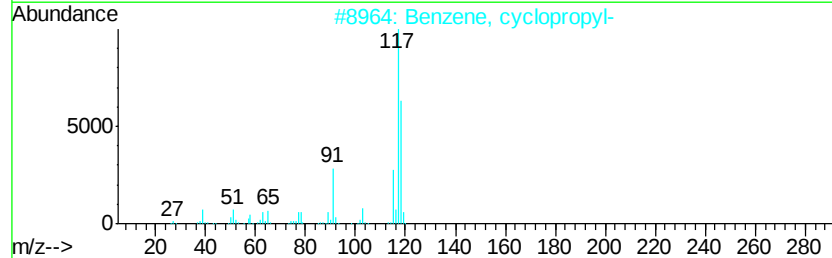
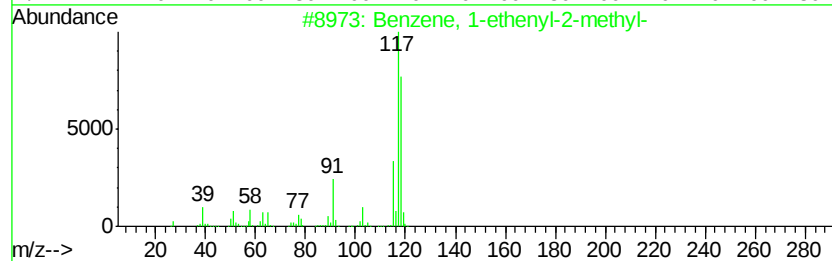
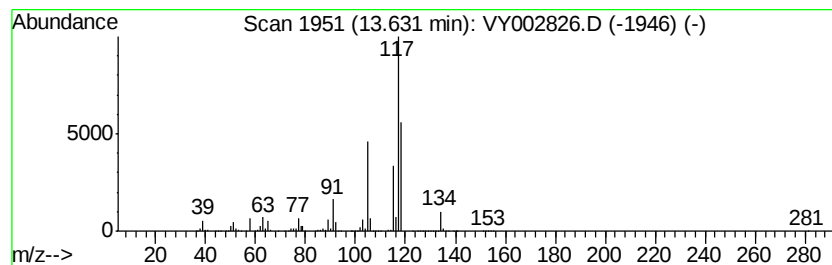
Instrument :
MSVOA_Y
ClientSampleId :
DRUM-OIL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.63	57.34 ug/l	10201900	1,4-Dichlorobenzene-d4	13.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 70
2		Benzene, cyclopropyl-	118 C9H10	000873-49-4 70
3		Indane	118 C9H10	000496-11-7 70
4		Benzene, 2-propenyl-	118 C9H10	000300-57-2 70
5		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 58



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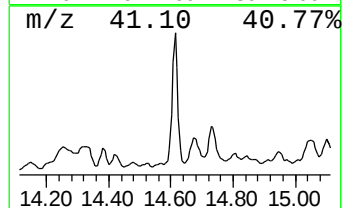
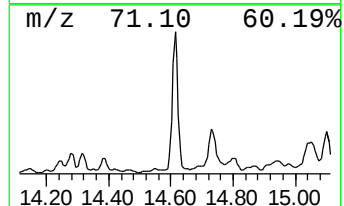
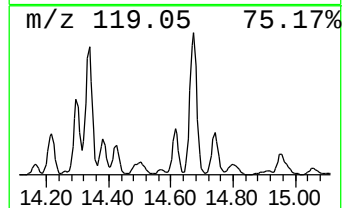
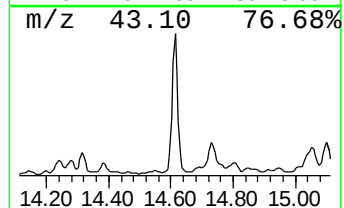
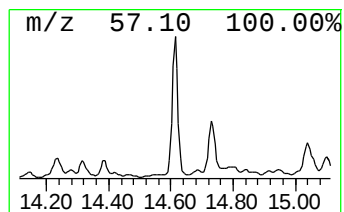
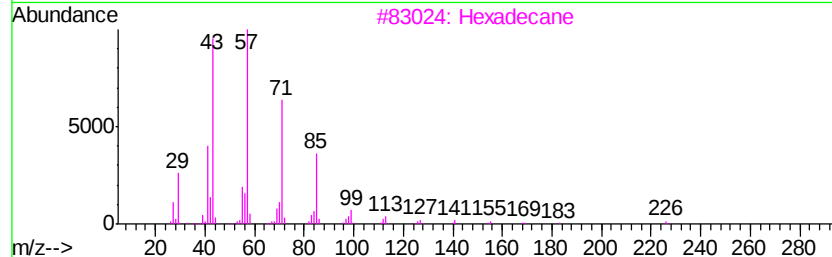
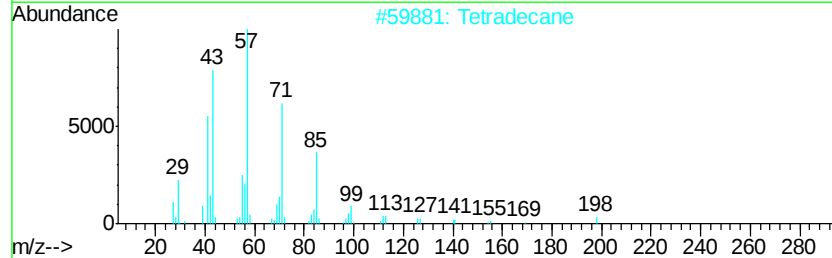
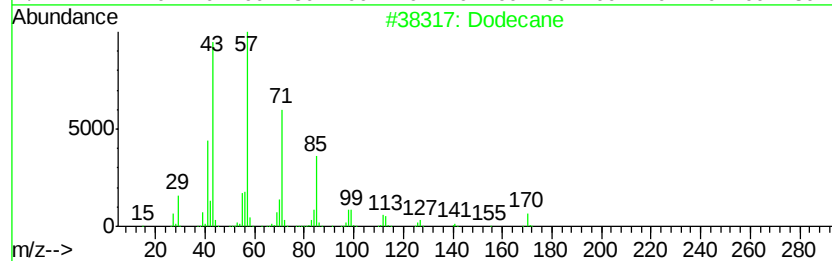
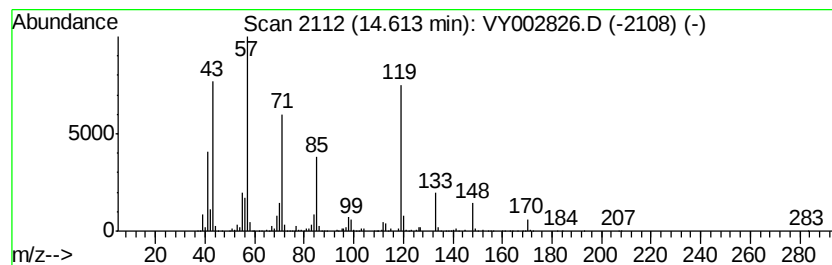
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Quant Title : SW846 8260

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

Peak Number 4 Dodecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	35.47 ug/l	6310300	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	94
2		Tetradecane	198	C14H30	000629-59-4	30
3		Hexadecane	226	C16H34	000544-76-3	30
4		Tridecane	184	C13H28	000629-50-5	27
5		Pentadecane	212	C15H32	000629-62-9	27



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Misc : 5.34G/5ML/MSVOA Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

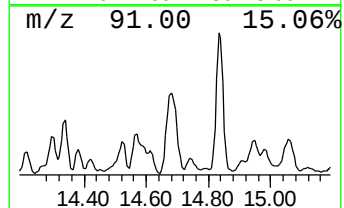
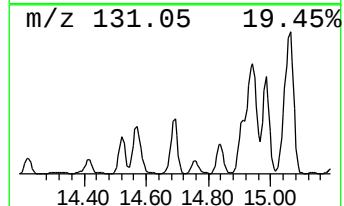
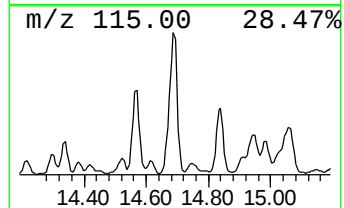
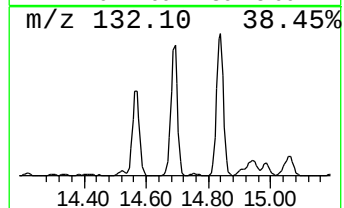
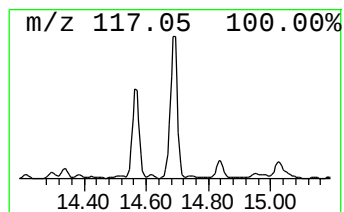
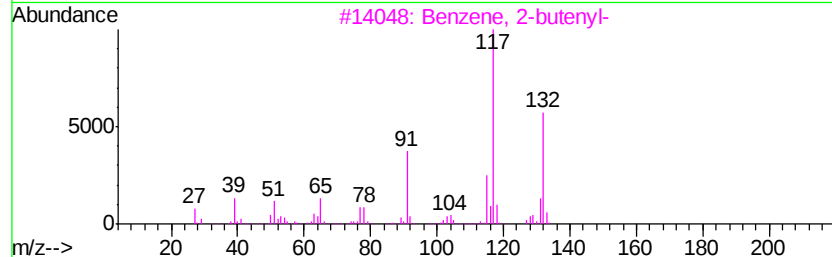
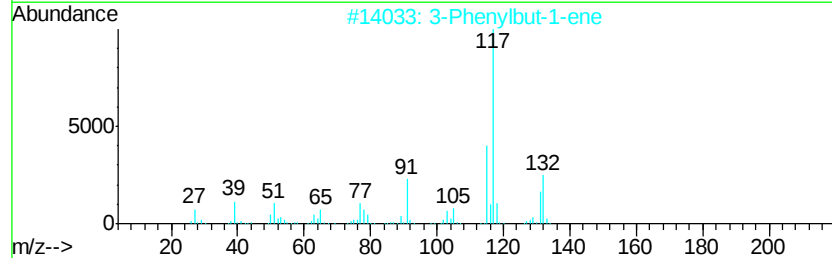
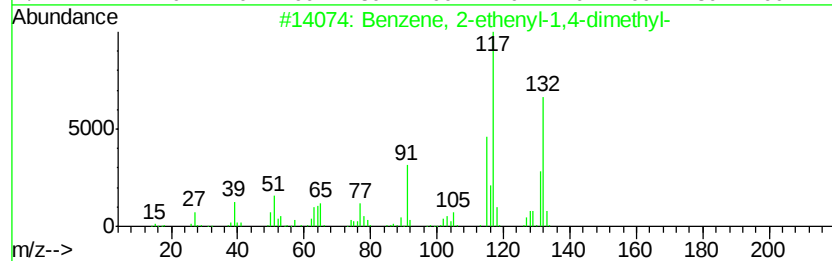
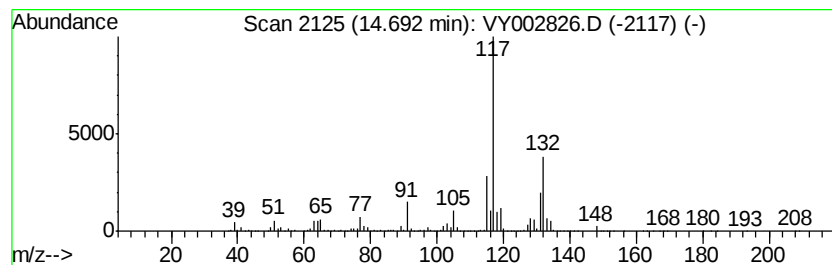
Instrument :
MSVOA_Y
ClientSampleId :
DRUM-OIL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	89.07 ug/l	15847300	1,4-Dichlorobenzene-d4	13.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 92
2		3-Phenylbut-1-ene	132 C10H12	000934-10-1 81
3		Benzene, 2-butenyl-	132 C10H12	001560-06-1 81
4		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 76
5		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY060320\
Data File : VY002826.D
Acq On : 03 Jun 2020 12:29
Operator : SY/MD
Sample : L2840-03
Misc : 5.34G/5ML/MSVOA Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
DRUM-OIL

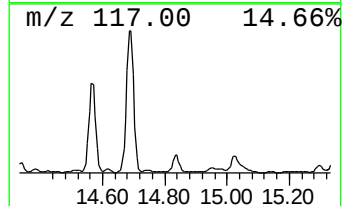
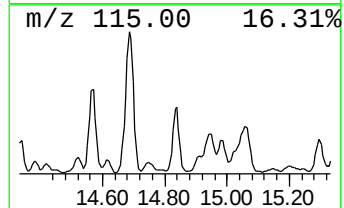
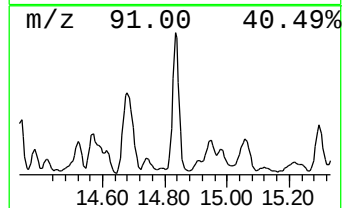
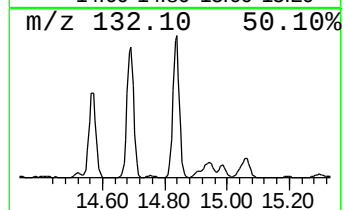
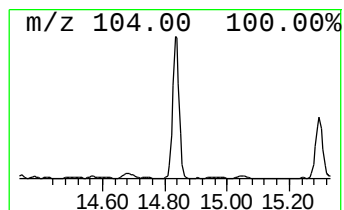
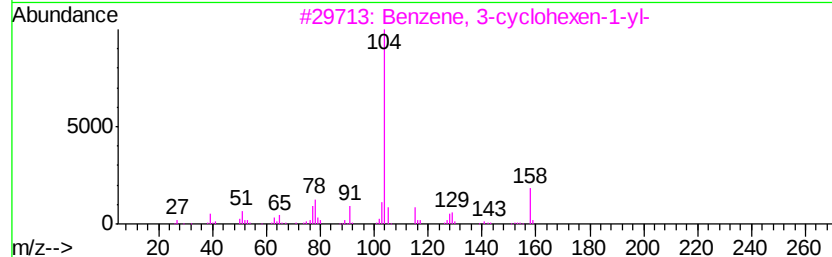
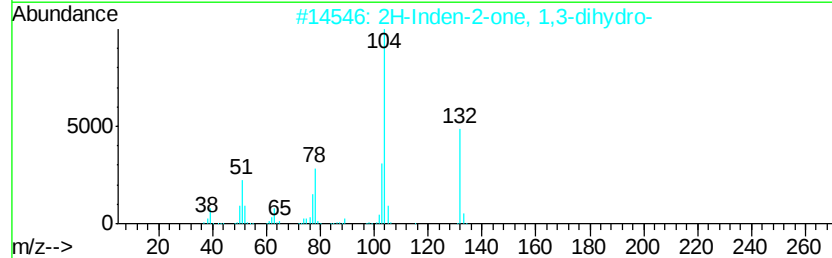
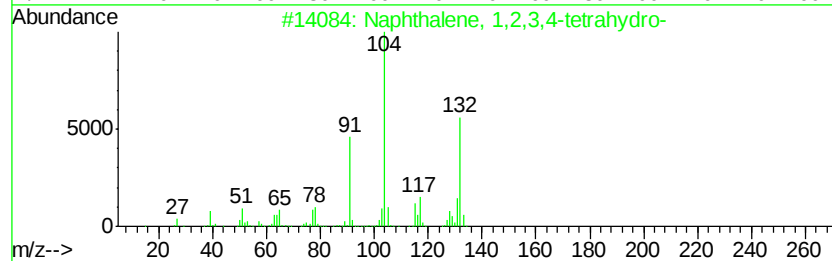
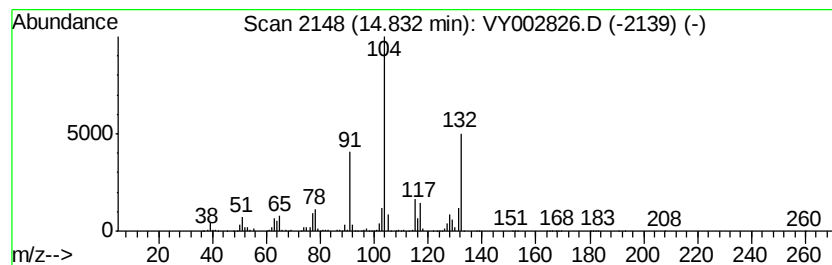
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	43.82 ug/l	7795290	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
3		Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	47
4		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46
5		Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY060320\
Data File : VY002826.D
Acq On : 03 Jun 2020 12:29
Operator : SY/MD
Sample : L2840-03
Misc : 5.34G/5ML/MSVOA Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

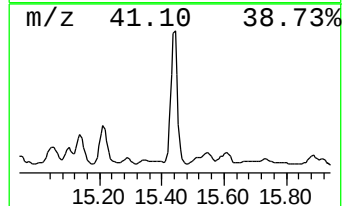
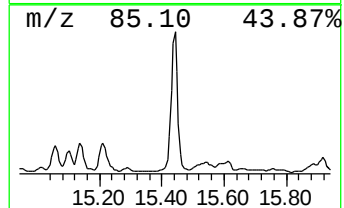
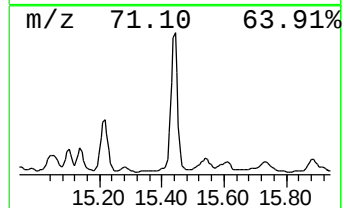
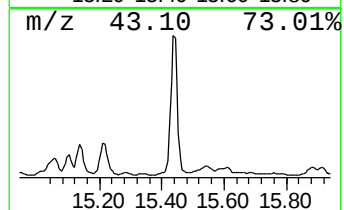
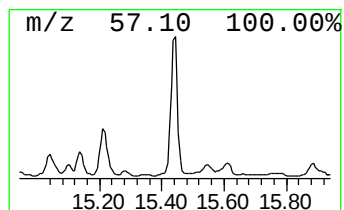
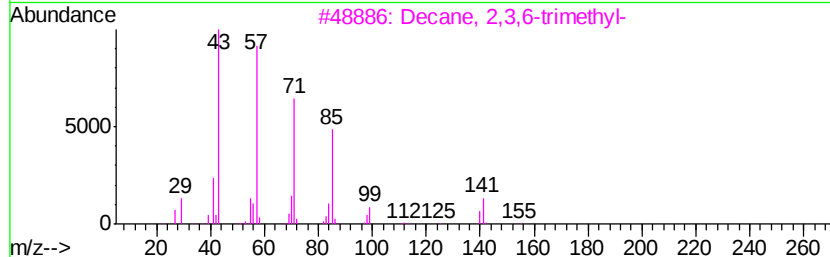
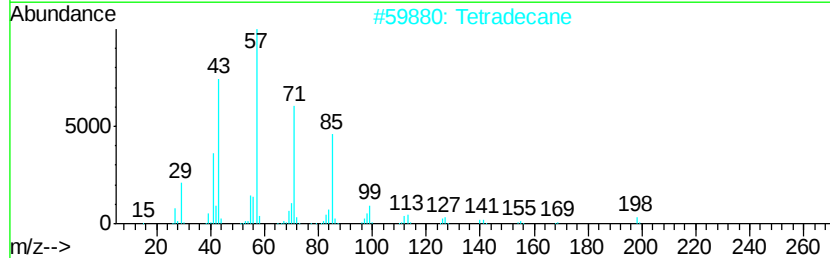
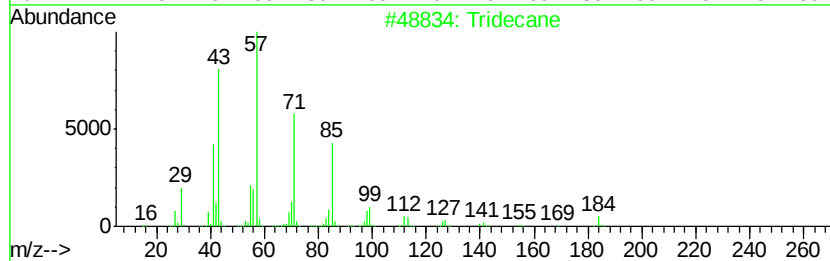
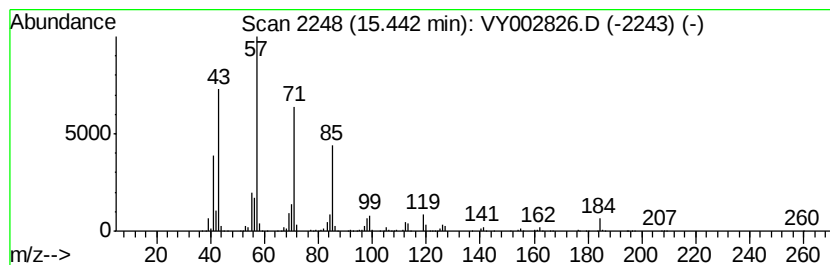
Instrument :
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ClientSampleId :
DRUM-OIL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

Peak Number 7 Tridecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.44	49.34 ug/l	8777800	1,4-Dichlorobenzene-d4	13.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	98
2	Tetradecane	198	C14H30	000629-59-4	87
3	Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	80
4	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	72
5	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY060320\
Data File : VY002826.D
Acq On : 03 Jun 2020 12:29
Operator : SY/MD
Sample : L2840-03
Misc : 5.34G/5ML/MSVOA Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

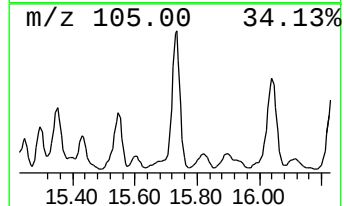
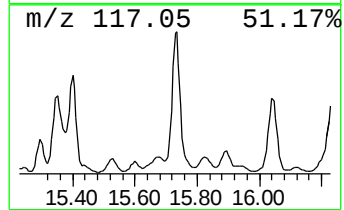
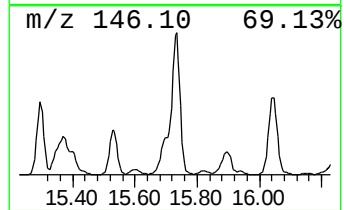
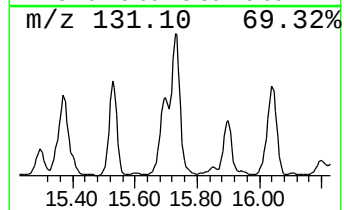
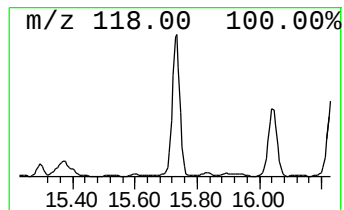
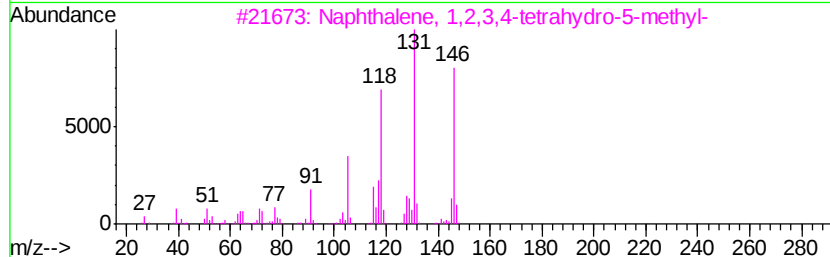
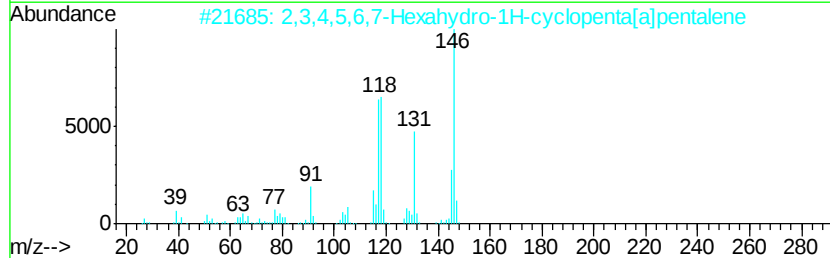
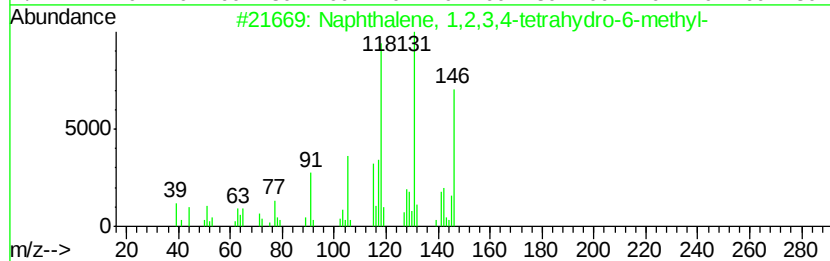
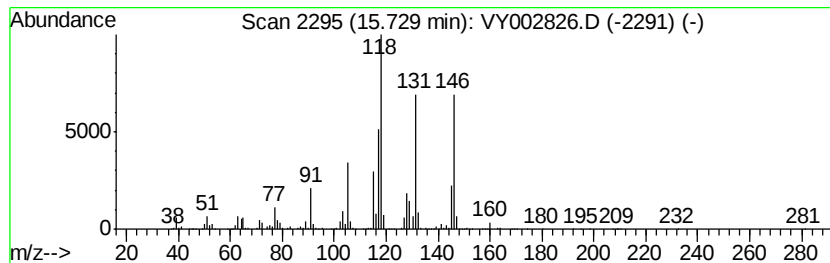
Instrument :
MSVOA_Y
ClientSampleId :
DRUM-OIL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	50.12 ug/l	8916710	1,4-Dichlorobenzene-d4	13.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 86
2		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 76
3		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 64
4		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 55
5		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	013065-07-1 52



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA Y\DATA\VY060320\
Data File : VY002826.D
Acq On : 03 Jun 2020 12:29
Operator : SY/MD
Sample : L2840-03
Misc : 5.34G/5ML/MSVOA Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
DRUM-OIL

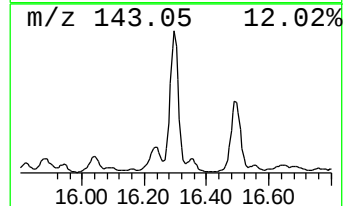
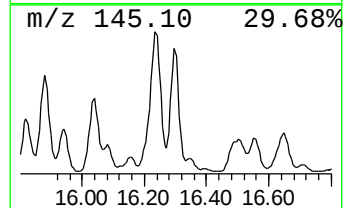
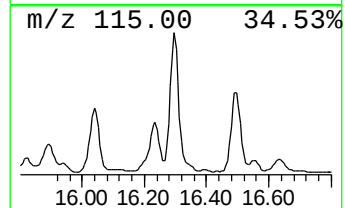
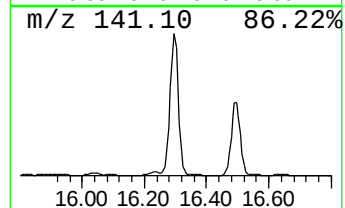
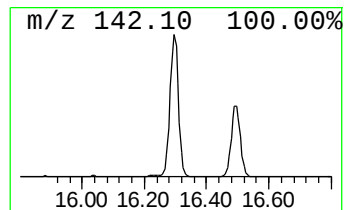
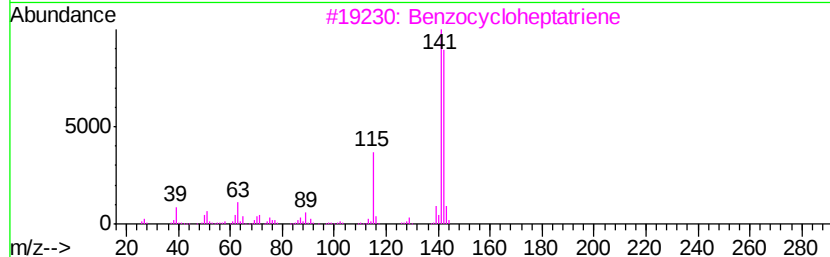
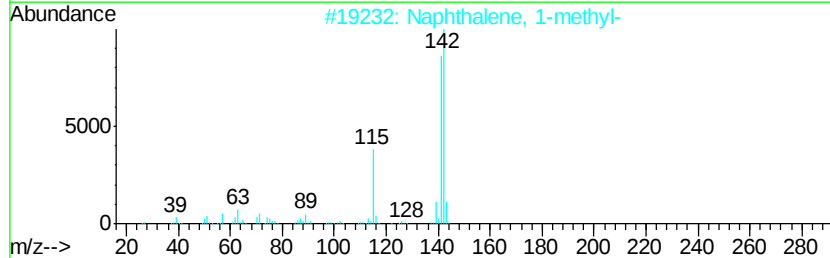
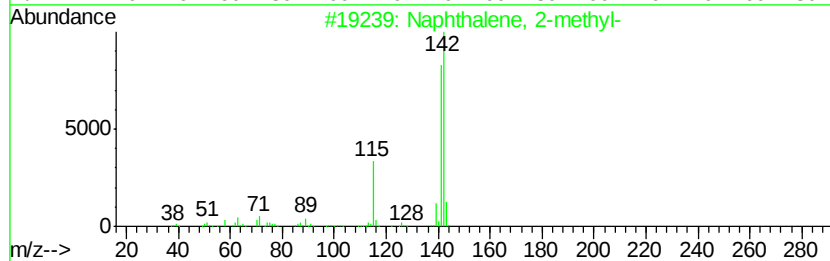
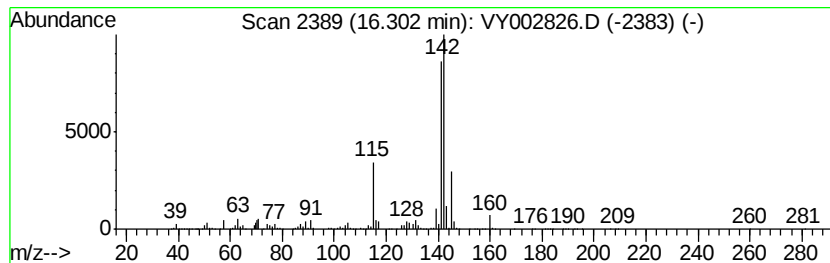
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.30	40.41 ug/l	7189610	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	87
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	81
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_Y\DATA\VY060320\
Data File : VY002826.D
Acq On : 03 Jun 2020 12:29
Operator : SY/MD
Sample : L2840-03
Misc : 5.34G/5ML/MSVOA_Y/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
DRUM-OIL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y052720S.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-...	12.74	75.5	ug/l	13440600	4	13.42	8895690	50.0
Benzene, 1-etheny...	13.63	57.3	ug/l	10201900	4	13.42	8895690	50.0
Dodecane	14.61	35.5	ug/l	6310300	4	13.42	8895690	50.0
Benzene, 2-etheny...	14.69	89.1	ug/l	15847300	4	13.42	8895690	50.0
Naphthalene, 1,2,...	14.83	43.8	ug/l	7795290	4	13.42	8895690	50.0
Tridecane	15.44	49.3	ug/l	8777800	4	13.42	8895690	50.0
Naphthalene, 1,2,...	15.73	50.1	ug/l	8916710	4	13.42	8895690	50.0
Naphthalene, 1-me...	16.30	40.4	ug/l	7189610	4	13.42	8895690	50.0