

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD051420\
 Data File : VD065739.D
 Acq On : 14 May 2020 13:28
 Operator : VA/SY
 Sample : L2501-01
 Misc : 8.39G/5.00ml/MSVOA D/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 EO-01-051320

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\82D051320S.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	8.867	1169	1181	1195	rBV	616454	1242449	1.30%	0.148%
2	10.344	1423	1432	1437	rBV	1389578	2691995	2.83%	0.321%
3	10.408	1438	1443	1448	rVB	507258	1000218	1.05%	0.119%
4	10.526	1456	1463	1472	rVB	1543311	2718462	2.85%	0.324%
5	10.955	1527	1536	1542	rBV	617250	1072267	1.13%	0.128%
6	11.197	1569	1577	1581	rVV	952888	1829896	1.92%	0.218%
7	11.250	1581	1586	1592	rVB	1169595	2140412	2.25%	0.255%
8	11.355	1599	1604	1608	rBV2	709087	1255881	1.32%	0.150%
9	11.432	1608	1617	1628	rVB4	2986374	7525640	7.90%	0.897%
10	11.532	1628	1634	1640	rBV	2750474	4478970	4.70%	0.534%
11	11.650	1649	1654	1658	rBV	678857	1150239	1.21%	0.137%
12	11.750	1664	1671	1675	rBV	1606897	2918963	3.06%	0.348%
13	11.861	1680	1690	1699	rVV3	21033505	40244108	42.24%	4.798%
14	11.938	1699	1703	1709	rVB2	1935718	3449951	3.62%	0.411%
15	12.085	1720	1728	1733	rVB3	1012992	1949647	2.05%	0.232%
16	12.185	1733	1745	1753	rBV2	5401591	18245533	19.15%	2.175%
17	12.279	1753	1761	1765	rVB	4503408	7487131	7.86%	0.893%
18	12.361	1765	1775	1776	rBV3	2659613	5156143	5.41%	0.615%
19	12.397	1777	1781	1791	rVV5	4698764	11796168	12.38%	1.406%
20	12.485	1791	1796	1797	rVV3	1573338	2433391	2.55%	0.290%
21	12.567	1800	1810	1812	rVV4	6401759	16392371	17.20%	1.954%
22	12.591	1812	1814	1819	rVB	6485270	8230464	8.64%	0.981%
23	12.644	1819	1823	1825	rBV3	1398205	2130230	2.24%	0.254%
24	12.679	1825	1829	1833	rVV	5365671	7618226	8.00%	0.908%
25	12.726	1833	1837	1841	rVB4	892835	1327281	1.39%	0.158%
26	12.820	1846	1853	1858	rVB3	3178295	8186184	8.59%	0.976%
27	12.891	1858	1865	1868	rBV	9856979	16848453	17.68%	2.009%
28	12.961	1868	1877	1883	rVB2	54682305	95277538	100.00%	11.359%
29	13.014	1883	1886	1892	rVB2	3879086	6153884	6.46%	0.734%
30	13.126	1892	1905	1912	rBV4	5783399	21510787	22.58%	2.565%
31	13.191	1912	1916	1923	rVB2	12423450	19659468	20.63%	2.344%
32	13.279	1923	1931	1936	rBV	18915688	31901020	33.48%	3.803%
33	13.320	1936	1938	1943	rVB	2306589	2775782	2.91%	0.331%
34	13.373	1943	1947	1950	rBV2	2857935	4463404	4.68%	0.532%

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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

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Max Peaks: 100

Peak Location: TOP

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D051320S.M

Title : SW846 8260

35	13.473	1957	1964	1968	rBV3	10778326	20885744	21.92%	2.490%
36	13.514	1968	1971	1975	rVV2	8938369	13723514	14.40%	1.636%
37	13.549	1975	1977	1980	rVV	5319922	6811563	7.15%	0.812%
38	13.585	1980	1983	1985	rVV	8616350	11075884	11.62%	1.321%
39	13.614	1985	1988	1992	rVV	8165861	10881705	11.42%	1.297%
40	13.655	1992	1995	2001	rVB	7875962	10250161	10.76%	1.222%
41	13.732	2002	2008	2011	rBV5	2107152	4918295	5.16%	0.586%
42	13.779	2012	2016	2020	rBV2	7730626	10350834	10.86%	1.234%
43	13.820	2020	2023	2032	rVB3	6056720	13160644	13.81%	1.569%
44	13.908	2032	2038	2043	rBV	58126602	92152634	96.72%	10.987%
45	13.949	2043	2045	2048	rBV2	1473730	1892901	1.99%	0.226%
46	13.979	2048	2050	2053	rVB	1772525	1980637	2.08%	0.236%
47	14.067	2058	2065	2071	rVB2	7126874	14282899	14.99%	1.703%
48	14.126	2071	2075	2079	rVB2	6795335	9323471	9.79%	1.112%
49	14.167	2079	2082	2088	rVB3	4783543	7791547	8.18%	0.929%
50	14.238	2088	2094	2099	rBV3	8299241	14224273	14.93%	1.696%
51	14.296	2099	2104	2110	rBV6	3105544	6333489	6.65%	0.755%
52	14.391	2110	2120	2121	rBV3	7357194	15806689	16.59%	1.885%
53	14.426	2122	2126	2129	rVB3	9466305	13905014	14.59%	1.658%
54	14.467	2129	2133	2136	rBV	4329288	5443167	5.71%	0.649%
55	14.532	2141	2144	2148	rVB	7230975	8904737	9.35%	1.062%
56	14.579	2148	2152	2158	rVB4	4402996	6982554	7.33%	0.832%
57	14.638	2159	2162	2165	rBV2	1730722	2465611	2.59%	0.294%
58	14.673	2166	2168	2173	rVB2	1757754	2149871	2.26%	0.256%
59	14.761	2178	2183	2189	rVV2	32904489	50845395	53.37%	6.062%
60	14.832	2189	2195	2200	rVV4	12197797	28317344	29.72%	3.376%
61	14.879	2200	2203	2212	rVB2	9341072	17938974	18.83%	2.139%
62	14.996	2219	2223	2230	rVB	5333014	9723907	10.21%	1.159%
63	15.102	2237	2241	2246	rBV3	2959402	4481714	4.70%	0.534%
64	15.220	2254	2261	2269	rBV5	3036187	8477234	8.90%	1.011%
65	15.291	2270	2273	2282	rVB8	3723494	8793640	9.23%	1.048%
66	15.373	2282	2287	2296	rBV3	4196714	7886571	8.28%	0.940%
67	15.467	2299	2303	2308	rVB	1691084	2393822	2.51%	0.285%
68	15.520	2308	2312	2315	rBV2	1425874	2297450	2.41%	0.274%
69	15.608	2322	2327	2336	rVB	6037150	11604542	12.18%	1.384%
70	15.708	2337	2344	2352	rVV6	1642919	4600089	4.83%	0.548%
71	15.791	2353	2358	2364	rVB5	1245076	2329537	2.45%	0.278%

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Integration Parameters: RTEINT.P

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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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72	15.920	2375	2380	2390	rVB	2937016	5782201	6.07%	0.689%
73	16.238	2426	2434	2441	rBV	1216295	2867026	3.01%	0.342%
74	16.438	2463	2468	2475	rVB2	708511	1456288	1.53%	0.174%

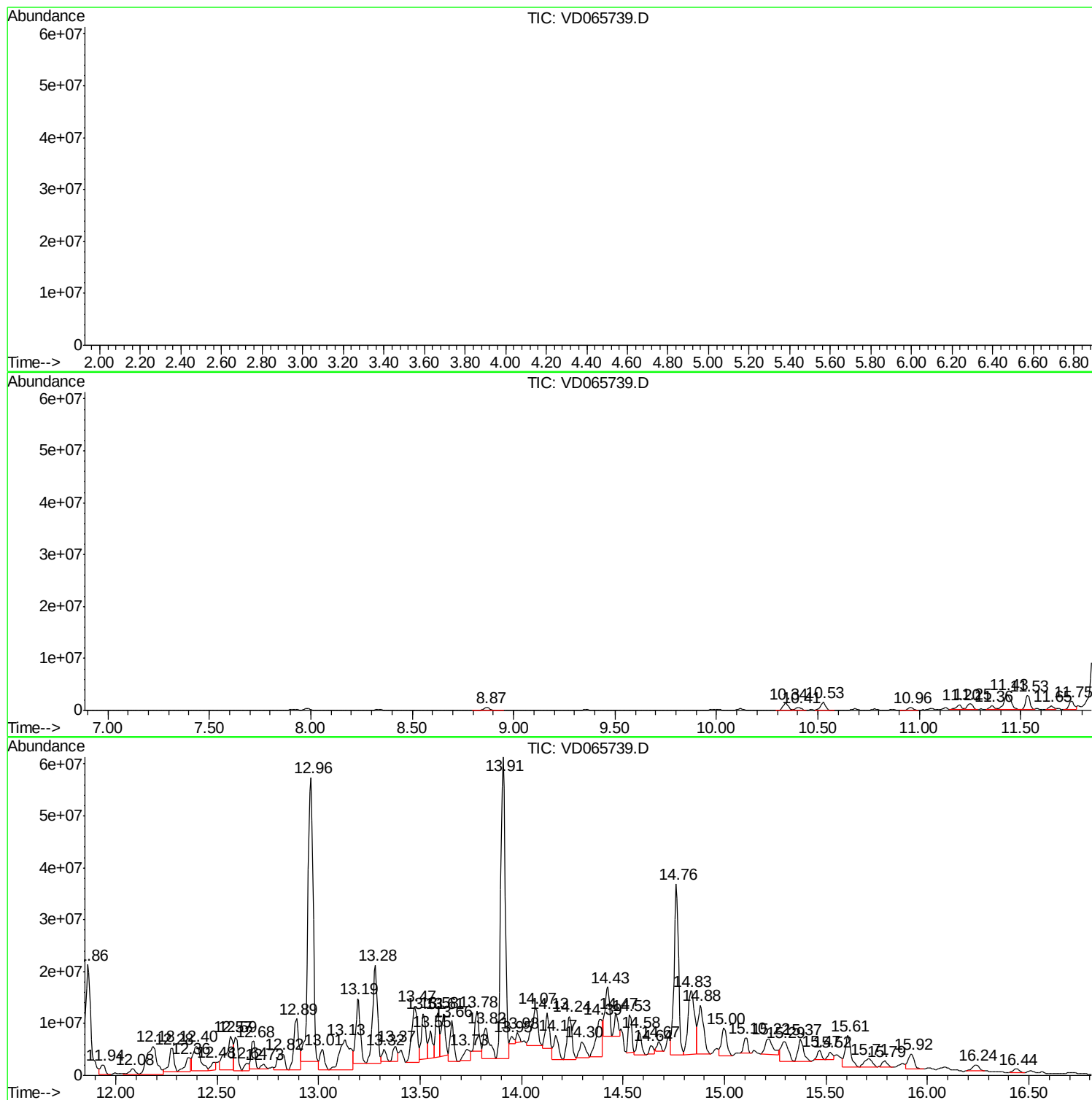
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D051320S.M
Quant Title : SW846 8260

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



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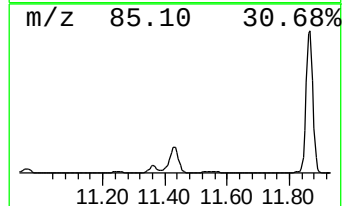
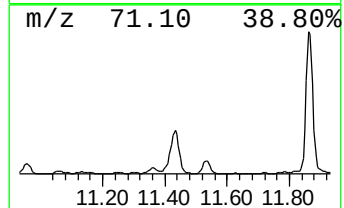
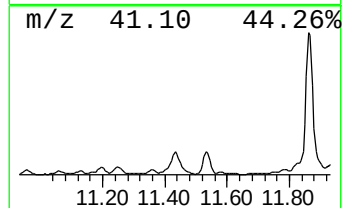
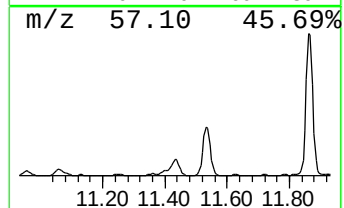
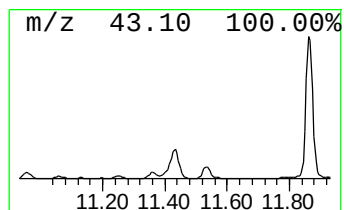
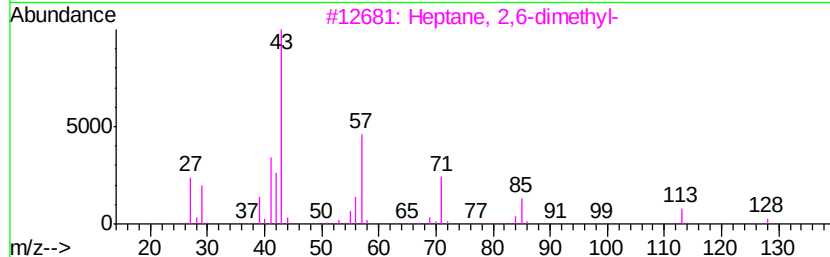
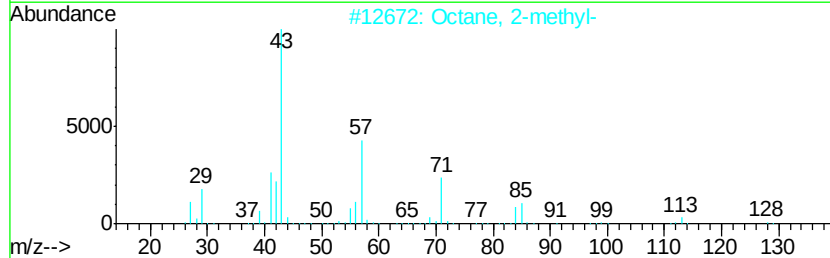
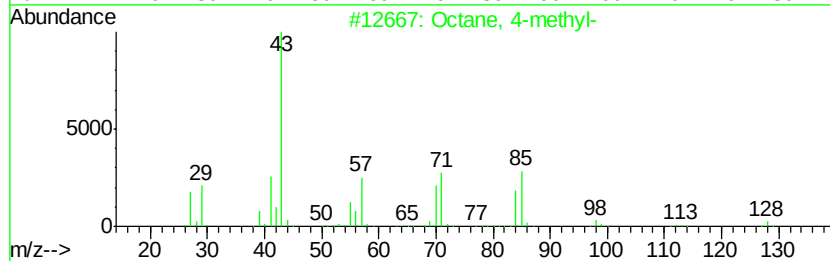
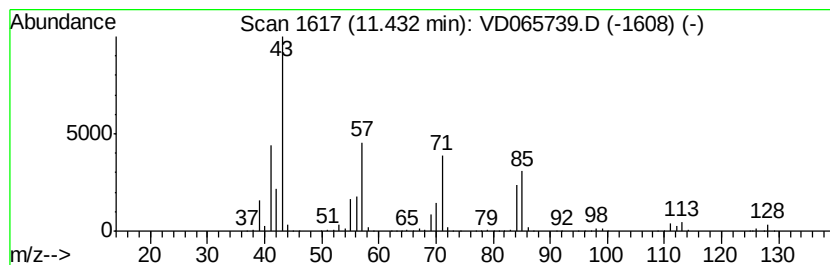
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D051320S.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	327.13 ug/l	7525640	Chlorobenzene-d5	11.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 4-methyl-	128	C9H20	002216-34-4	64
2		Octane, 2-methyl-	128	C9H20	003221-61-2	64
3		Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	53
4		Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	50
5		Ether, hexyl pentyl	172	C11H24O	032357-83-8	50



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Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

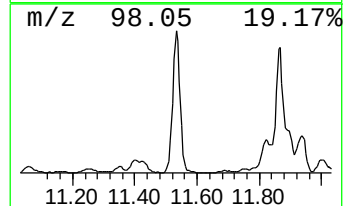
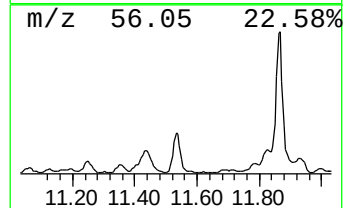
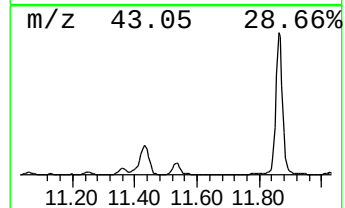
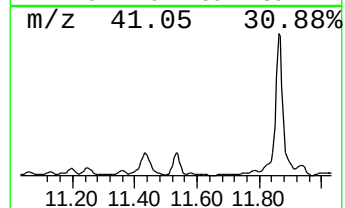
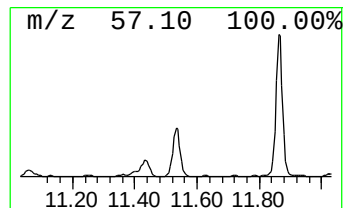
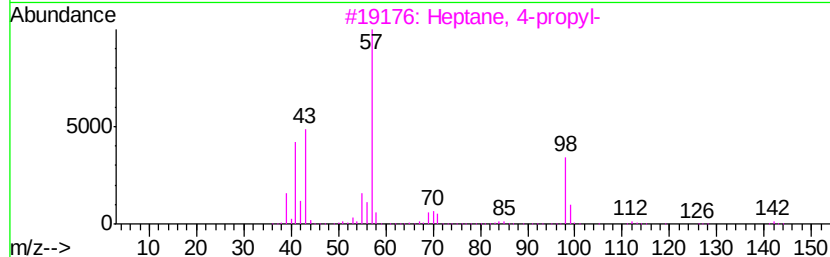
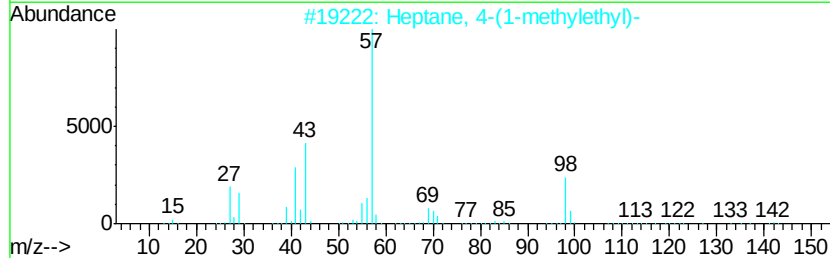
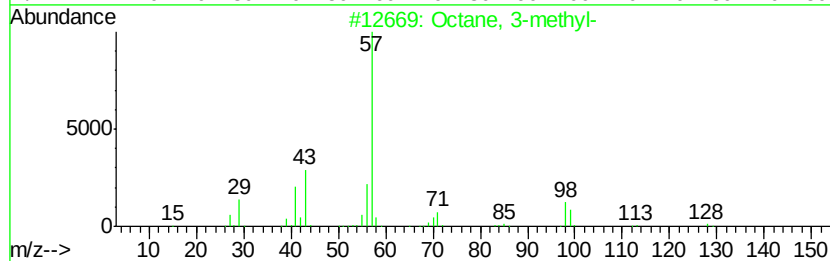
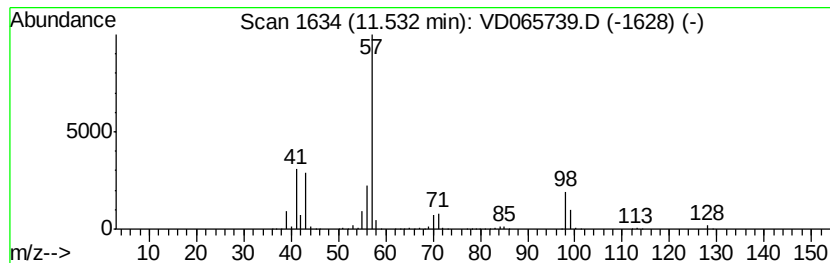
Instrument :
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 2 Octane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.53	194.70 ug/l	4478970	Chlorobenzene-d5	11.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 3-methyl-	128	C9H20	002216-33-3	91
2		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	64
3		Heptane, 4-propyl-	142	C10H22	003178-29-8	59
4		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	58
5		Heptane, 3-ethyl-	128	C9H20	015869-80-4	50



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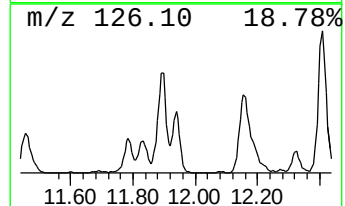
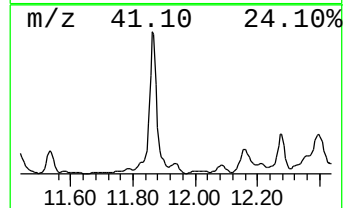
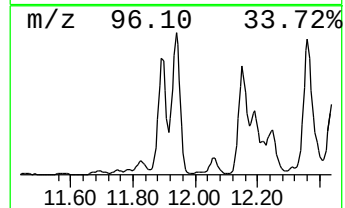
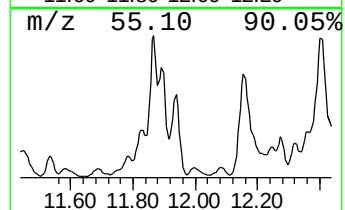
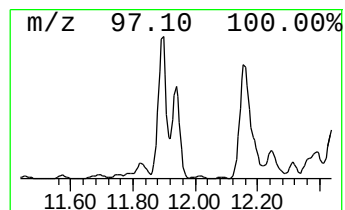
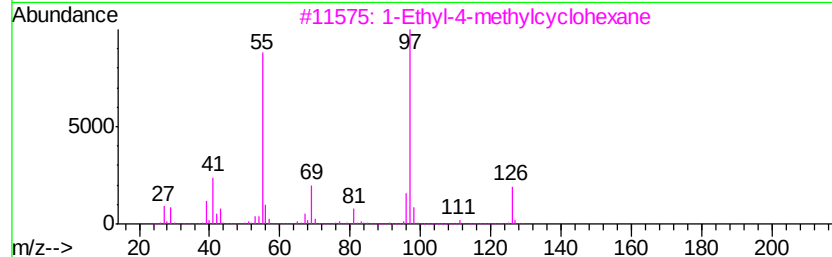
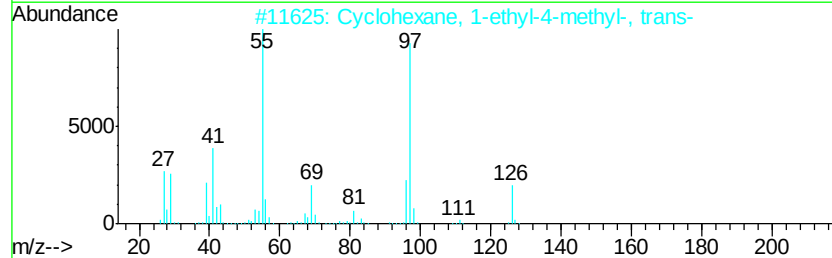
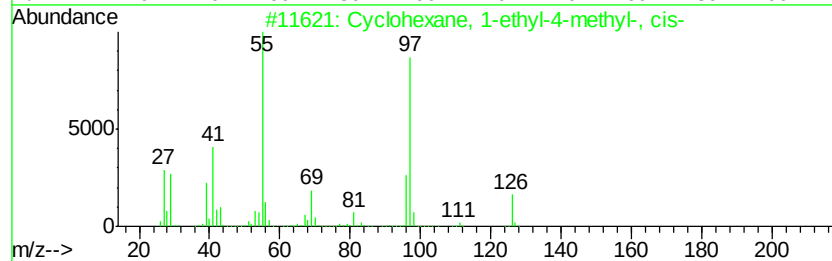
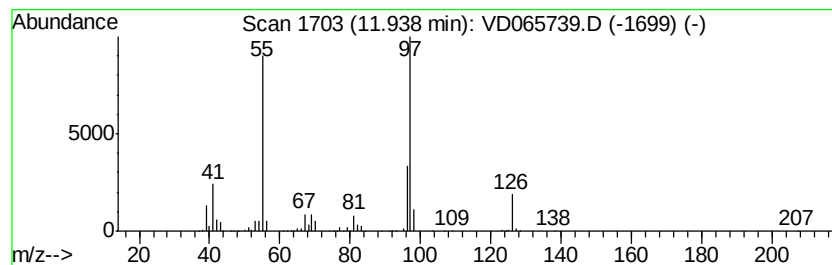
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1-ethyl-4-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.94	149.97 ug/l	3449950	Chlorobenzene-d5	11.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	86
2		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	83
3		1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	80
4		cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	80
5		1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	80



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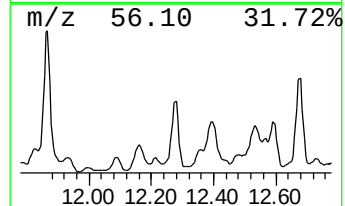
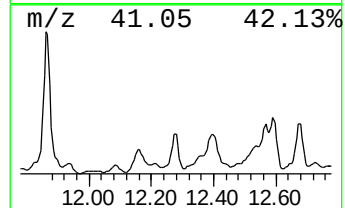
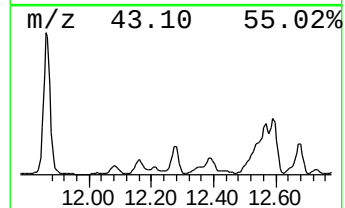
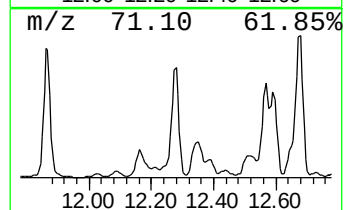
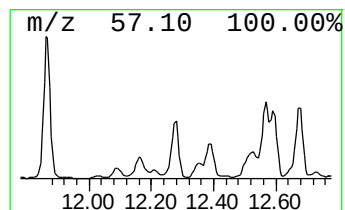
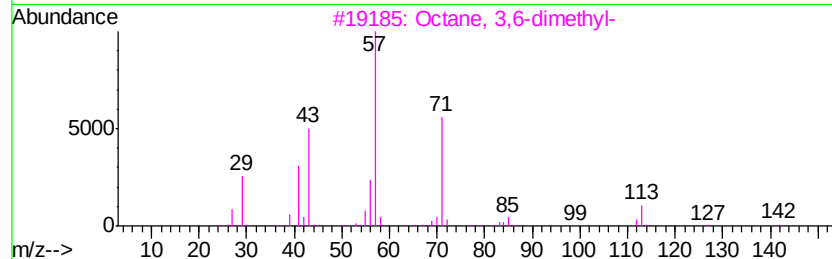
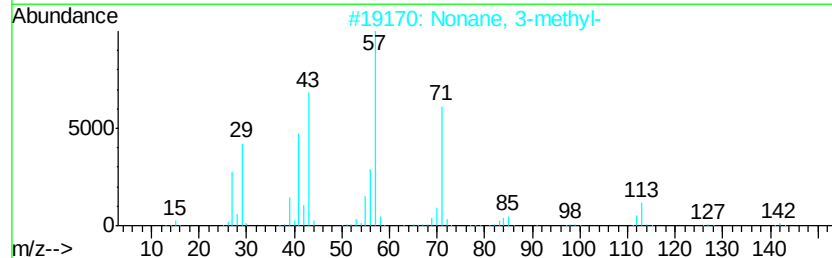
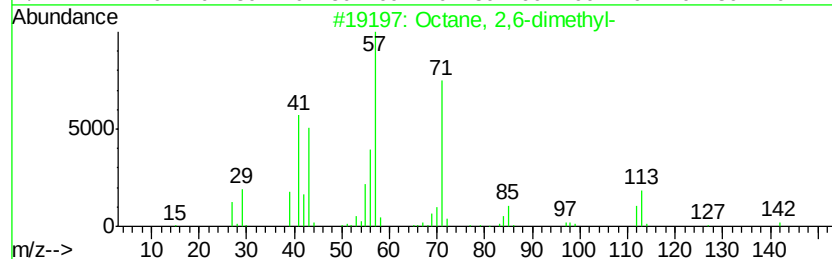
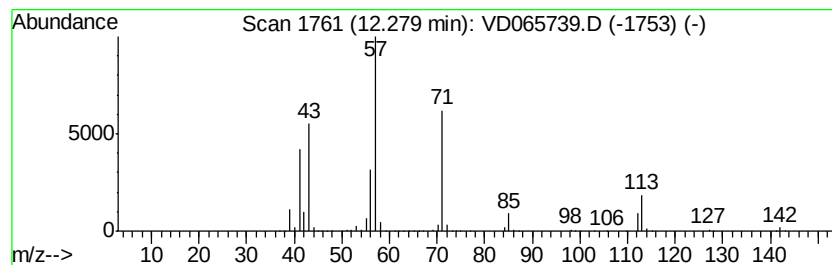
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Quant Title : SW846 8260

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

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	325.46 ug/l	7487130	Chlorobenzene-d5	11.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	90
3		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	86
4		Decane, 2,6,7-trimethyl-	184	C13H28	062108-25-2	64
5		Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1000309-19-2	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-01-051320

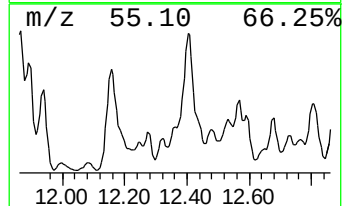
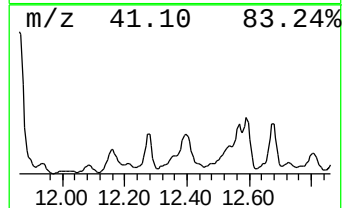
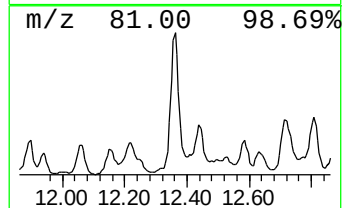
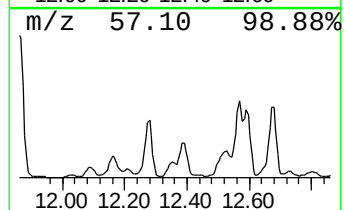
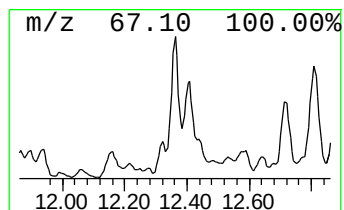
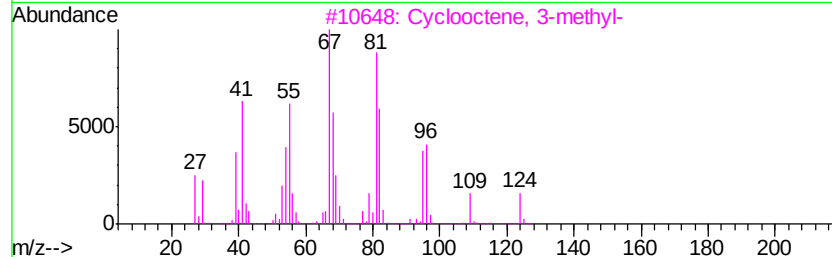
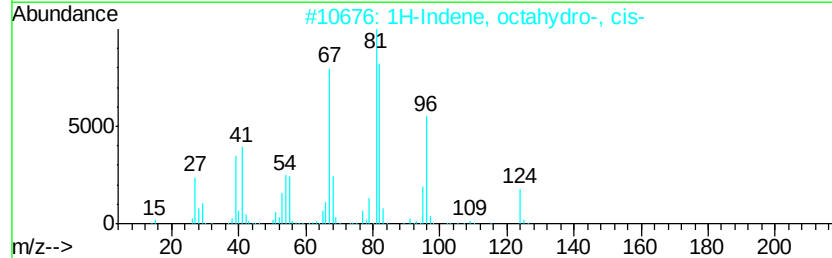
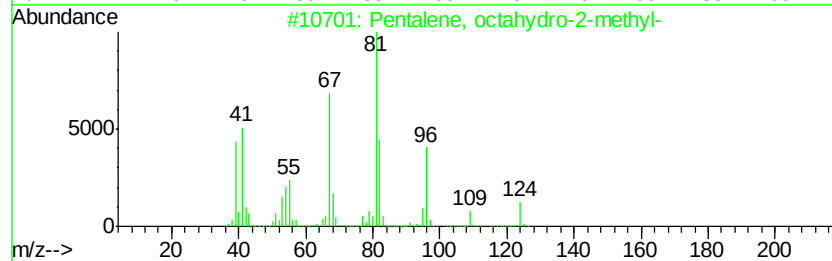
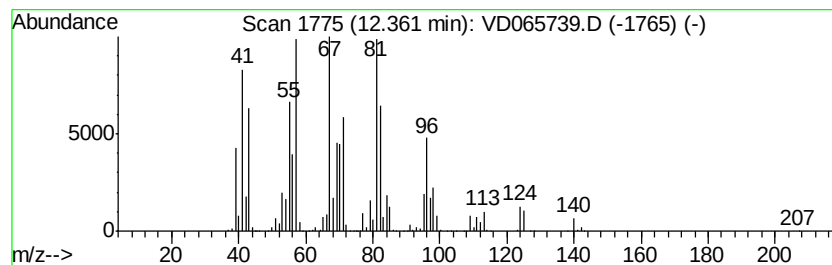
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Quant Title : SW846 8260

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

Peak Number 5 Pentalene, octahydro-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	224.13 ug/l	5156140	Chlorobenzene-d5	11.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	78
2		1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	55
3		Cyclooctene, 3-methyl-	124	C9H16	013152-05-1	49
4		Cyclopropane, (2-methylenebutyl)-	110	C8H14	074685-56-6	46
5		Cyclohexanone, 4-ethenyl-	124	C8H12O	001740-64-3	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

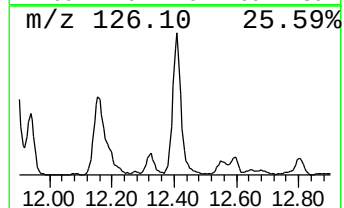
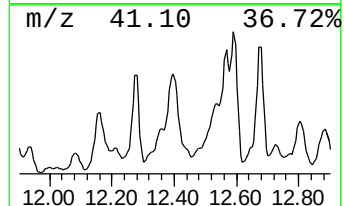
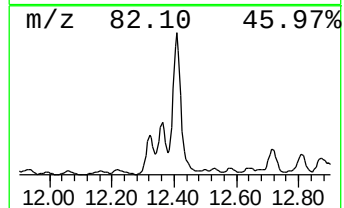
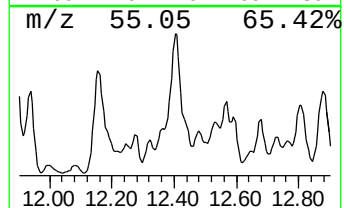
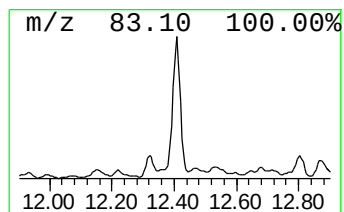
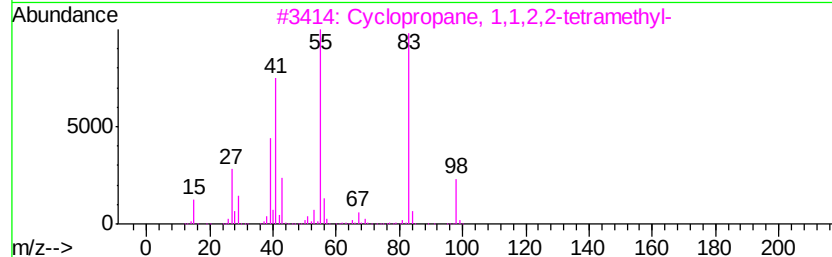
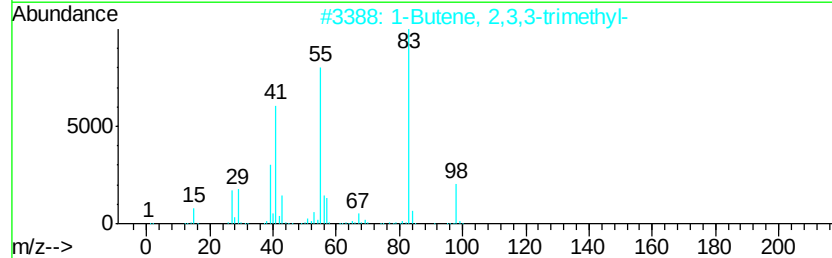
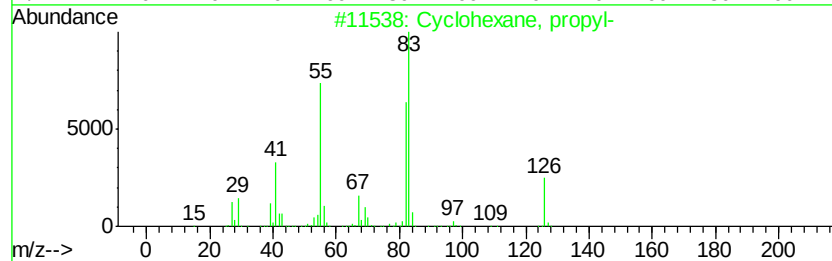
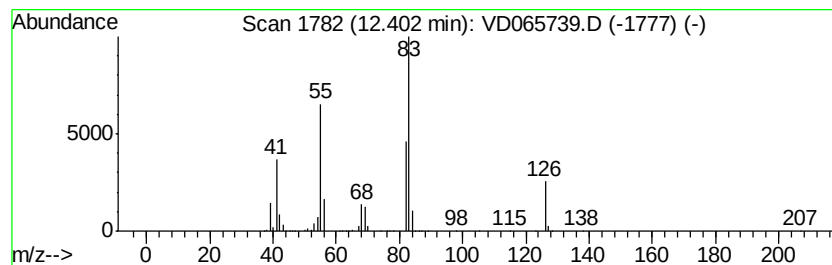
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ClientSampleId :
EO-01-051320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D051320S.M
Quant Title : SW846 8260

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

Peak Number 6 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
12.40	512.77 ug/l	11796200	Chlorobenzene-d5	11.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, propyl-	126	C9H18	001678-92-8	52
2		1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	49
3		Cyclopropane, 1,1,2,2-tetramethyl-	98	C7H14	004127-47-3	49
4		1-Tetradecanol, 14-chloro-	248	C14H29ClO	073937-05-0	47
5		Cyclohexane, methyl-	98	C7H14	000108-87-2	45



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-01-051320

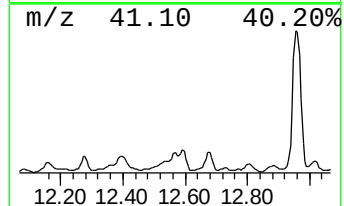
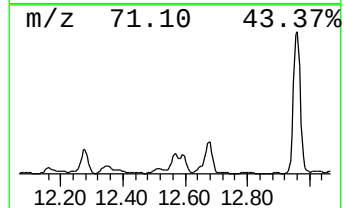
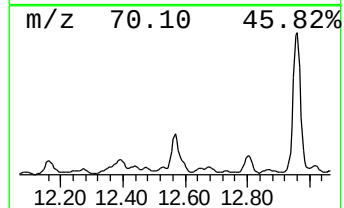
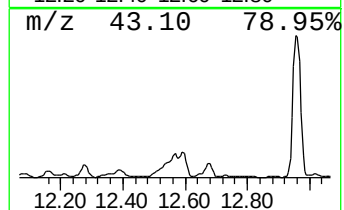
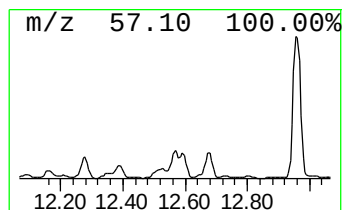
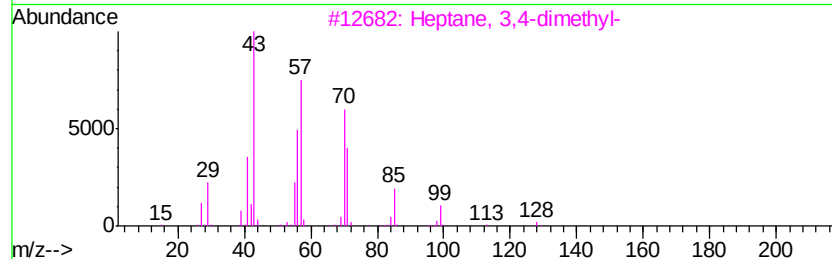
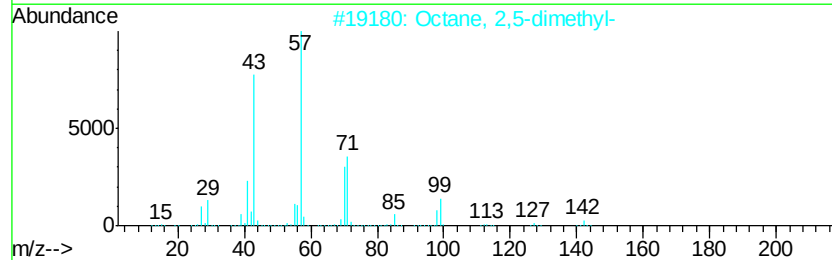
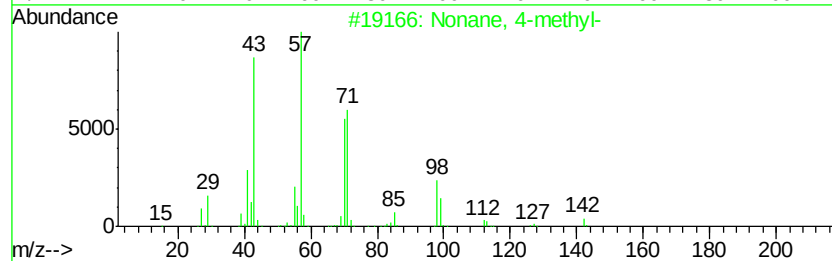
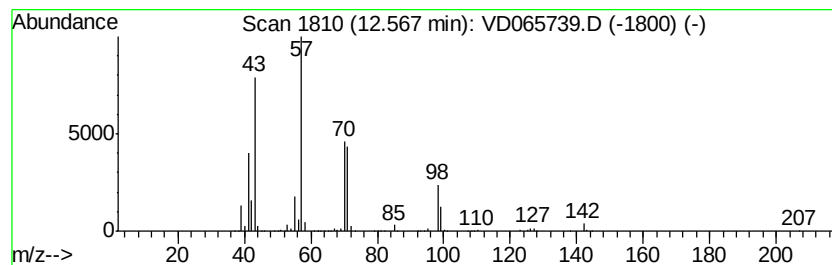
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Quant Title : SW846 8260

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

Peak Number 7 Nonane, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.57	712.56 ug/l	16392400	Chlorobenzene-d5	11.65

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 4-methyl-	142	C10H22	017301-94-9	91
2		Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	64
3		Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	59
4		Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5		Hexane, 3-methyl-	100	C7H16	000589-34-4	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-01-051320

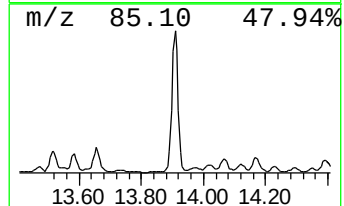
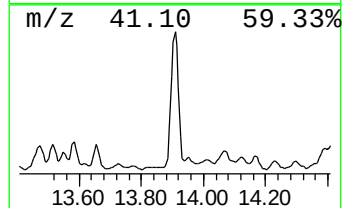
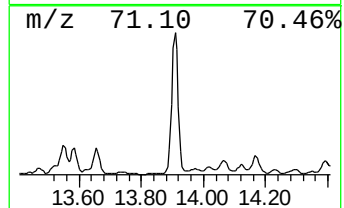
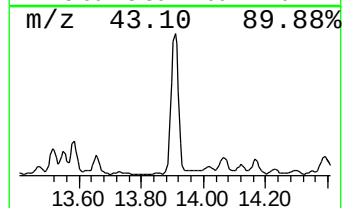
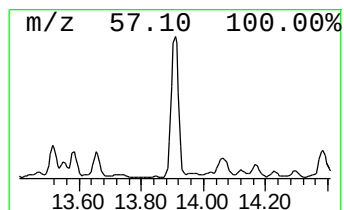
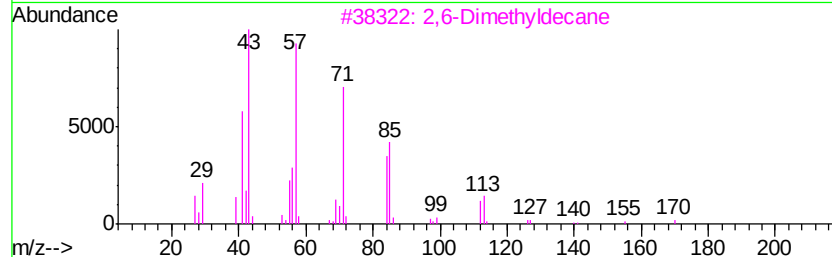
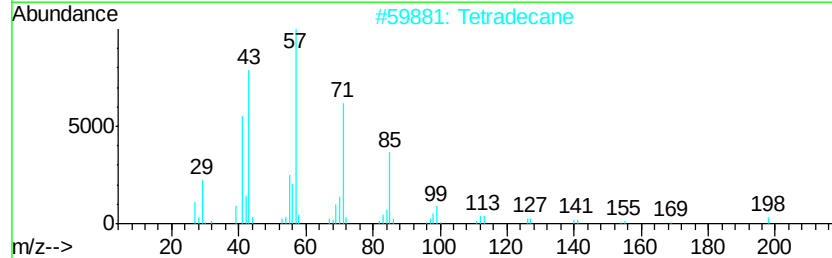
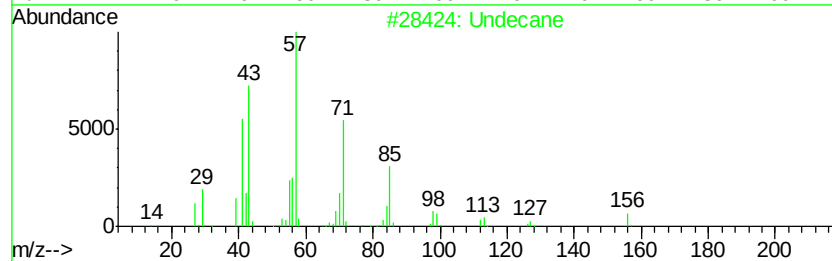
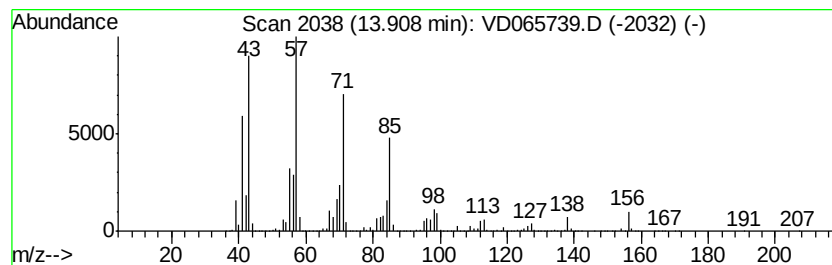
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Quant Title : SW846 8260

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

Peak Number 8 Undecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	416.01 ug/l	92152600	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	95
2	Tetradecane		198	C14H30	000629-59-4	80
3	2,6-Dimethyldecane		170	C12H26	013150-81-7	70
4	Undecane, 2-methyl-		170	C12H26	007045-71-8	68
5	Nonadecane		268	C19H40	000629-92-5	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

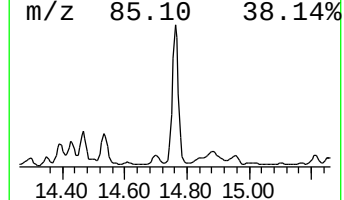
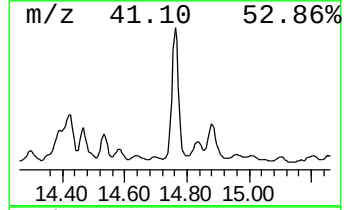
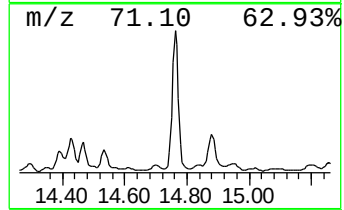
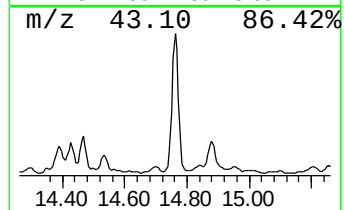
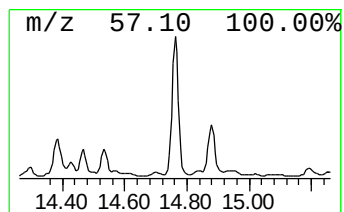
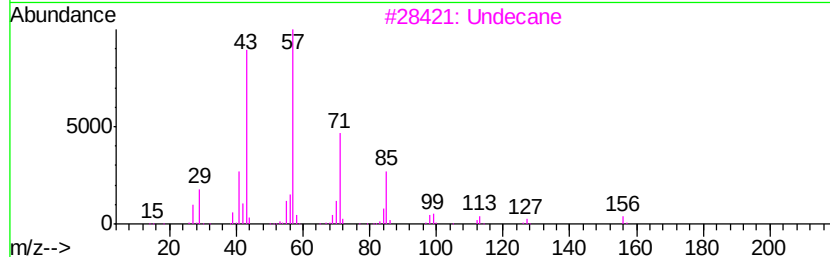
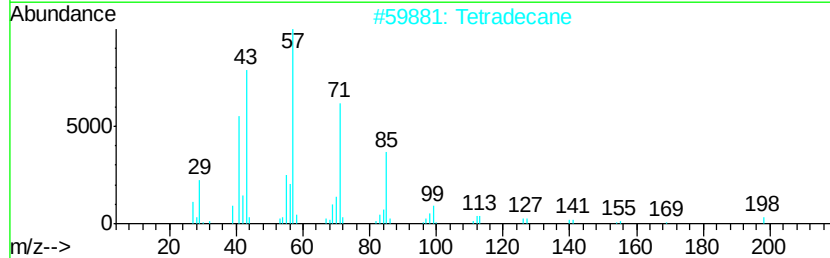
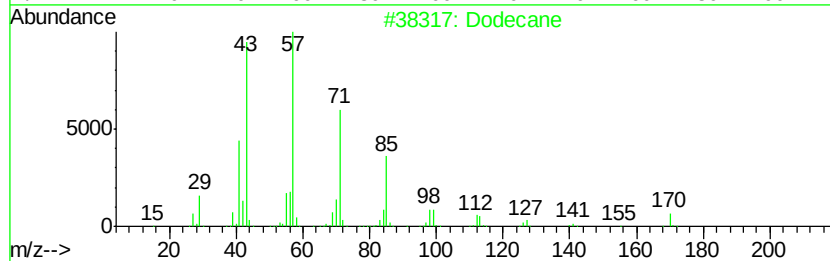
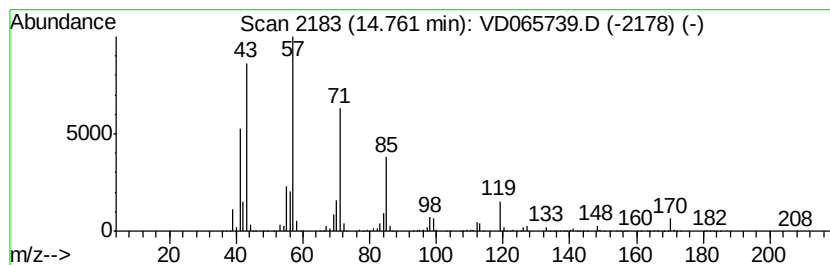
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D051320S.M
Quant Title : SW846 8260

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

Peak Number 9 Dodecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	229.53 ug/l	50845400	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Dodecane	170 C12H26	000112-40-3	96
2	Tetradecane	198 C14H30	000629-59-4	87
3	Undecane	156 C11H24	001120-21-4	80
4	Hexadecane	226 C16H34	000544-76-3	72
5	Dodecane, 2-methyl-	184 C13H28	001560-97-0	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-01-051320

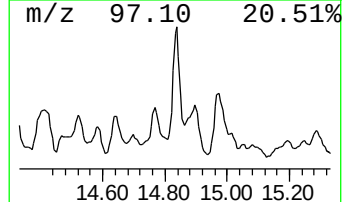
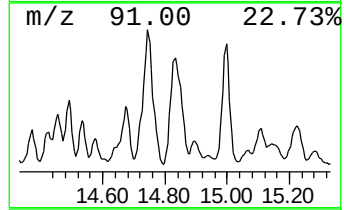
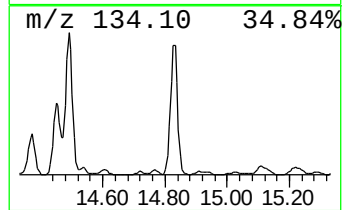
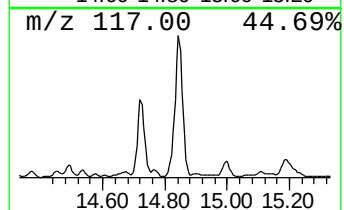
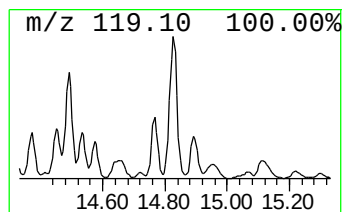
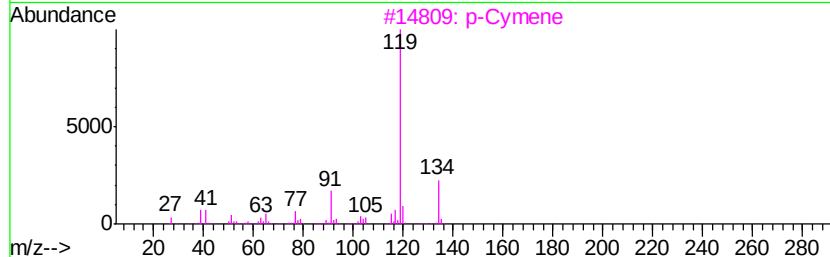
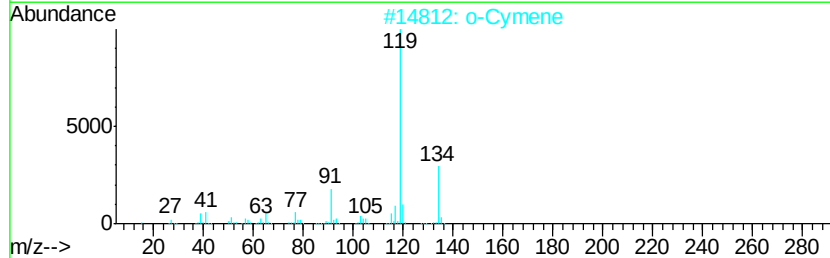
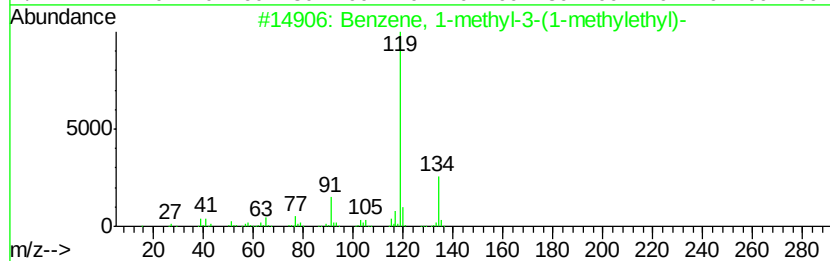
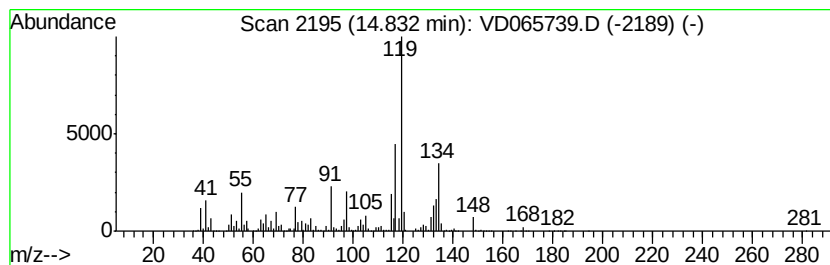
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-methyl-3-(1-meth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	127.83 ug/l	28317300	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	60
2		o-Cymene	134	C10H14	000527-84-4	60
3		p-Cymene	134	C10H14	000099-87-6	55
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	55
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	55



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD051420\
Data File : VD065739.D
Acq On : 14 May 2020 13:28
Operator : VA/SY
Sample : L2501-01
Misc : 8.39G/5.00ml/MSVOA_D/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-01-051320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D051320S.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
Octane, 4-methyl-	11.43	327.1	ug/l	7525640	3	11.65	1150240	50.0
Octane, 3-methyl-	11.53	194.7	ug/l	4478970	3	11.65	1150240	50.0
Cyclohexane, 1-et...	11.94	150.0	ug/l	3449950	3	11.65	1150240	50.0
Octane, 2,6-dimet...	12.28	325.5	ug/l	7487130	3	11.65	1150240	50.0
Pentalene, octahy...	12.36	224.1	ug/l	5156140	3	11.65	1150240	50.0
Cyclohexane, propyl-	12.40	512.8	ug/l	11796200	3	11.65	1150240	50.0
Nonane, 4-methyl-	12.57	712.6	ug/l	16392400	3	11.65	1150240	50.0
Undecane	13.91	416.0	ug/l	92152600	4	13.58	11075900	50.0
Dodecane	14.76	229.5	ug/l	50845400	4	13.58	11075900	50.0
Benzene, 1-methyl...	14.83	127.8	ug/l	28317300	4	13.58	11075900	50.0