

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD073020\
 Data File : VD066119.D
 Acq On : 30 Jul 2020 18:48
 Operator : VA/SY
 Sample : L3507-01
 Misc : 8.17G/5.00ml/MSVOA D/SOIL
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SU-01-073020

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\82D072120S.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	1.944	3	4	12	rVB2	42000	41136	3.21%	0.338%
2	2.315	62	67	78	rBV3	15897	36987	2.88%	0.303%
3	3.403	246	252	267	rVB4	10731	31852	2.48%	0.261%
4	4.150	370	379	389	rBV	31478	88047	6.87%	0.722%
5	4.238	389	394	400	rVB3	7140	16542	1.29%	0.136%
6	4.961	505	517	529	rBV2	34928	112397	8.77%	0.922%
7	7.208	891	899	910	rBV	13830	39299	3.06%	0.322%
8	7.914	1008	1019	1024	rBV	163797	396115	30.89%	3.250%
9	7.979	1024	1030	1040	rVB	328353	723390	56.42%	5.936%
10	8.332	1081	1090	1102	rBV	133397	294282	22.95%	2.415%
11	8.720	1148	1156	1164	rBV5	6946	16087	1.25%	0.132%
12	8.861	1172	1180	1192	rBV	445977	900811	70.25%	7.391%
13	9.308	1250	1256	1259	rBV4	6008	12964	1.01%	0.106%
14	9.779	1327	1336	1343	rVB2	18312	35384	2.76%	0.290%
15	9.914	1353	1359	1364	rVB3	19699	38601	3.01%	0.317%
16	9.979	1364	1370	1373	rBV4	12686	19196	1.50%	0.158%
17	10.114	1386	1393	1397	rBV3	18471	37500	2.92%	0.308%
18	10.232	1406	1413	1417	rBV	85103	158958	12.40%	1.304%
19	10.267	1417	1419	1423	rVB3	24612	31929	2.49%	0.262%
20	10.338	1423	1431	1440	rBV	698660	1282215	100.00%	10.521%
21	10.526	1455	1463	1470	rVB10	14925	30799	2.40%	0.253%
22	10.585	1470	1473	1479	rBV6	6964	13312	1.04%	0.109%
23	10.791	1501	1508	1514	rVV7	11322	30733	2.40%	0.252%
24	10.867	1517	1521	1525	rVB4	18971	30638	2.39%	0.251%
25	10.955	1530	1536	1543	rBV3	23313	44107	3.44%	0.362%
26	11.055	1547	1553	1561	rVB2	37427	81171	6.33%	0.666%
27	11.249	1579	1586	1592	rBV5	17044	34435	2.69%	0.283%
28	11.355	1600	1604	1608	rBV6	23219	42416	3.31%	0.348%
29	11.426	1608	1616	1629	rVV4	128090	344183	26.84%	2.824%
30	11.532	1629	1634	1646	rVB	103942	203751	15.89%	1.672%
31	11.649	1646	1654	1661	rBV	708125	1273226	99.30%	10.447%
32	11.843	1683	1687	1689	rBV5	11548	19927	1.55%	0.164%
33	11.890	1692	1695	1700	rVB5	27141	40000	3.12%	0.328%
34	12.038	1715	1720	1722	rBV6	6998	13367	1.04%	0.110%

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 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SU-01-073020

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

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Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
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35	12.079	1722	1727	1733	rBV3	27906	51664	4.03%	0.424%
36	12.155	1735	1740	1746	rBV4	61366	121735	9.49%	0.999%
37	12.273	1753	1760	1766	rVB3	67457	134660	10.50%	1.105%
38	12.349	1767	1773	1777	rBV3	35564	72174	5.63%	0.592%
39	12.496	1792	1798	1801	rBV3	71865	142429	11.11%	1.169%
40	12.567	1807	1810	1811	rBV2	32310	32377	2.53%	0.266%
41	12.590	1812	1814	1816	rVV2	44591	49444	3.86%	0.406%
42	12.638	1816	1822	1827	rVB	545949	874220	68.18%	7.173%
43	12.726	1833	1837	1841	rVB7	20322	34922	2.72%	0.287%
44	12.779	1841	1846	1849	rBV5	21788	41770	3.26%	0.343%
45	12.967	1869	1878	1881	rBV6	52868	141126	11.01%	1.158%
46	13.049	1889	1892	1897	rVB5	71810	115412	9.00%	0.947%
47	13.102	1898	1901	1905	rVB5	26783	35066	2.73%	0.288%
48	13.149	1906	1909	1913	rBV3	27285	37714	2.94%	0.309%
49	13.190	1913	1916	1918	rBV3	20652	24975	1.95%	0.205%
50	13.320	1934	1938	1944	rVB3	45634	76259	5.95%	0.626%
51	13.437	1944	1958	1963	rBV5	107661	378612	29.53%	3.107%
52	13.579	1974	1982	1988	rBV	717816	1202683	93.80%	9.868%
53	13.649	1988	1994	2004	rVB3	75537	216071	16.85%	1.773%
54	13.743	2006	2010	2014	rBV2	28409	47960	3.74%	0.394%
55	13.790	2014	2018	2021	rBV5	16851	23924	1.87%	0.196%
56	13.902	2033	2037	2043	rBV8	23136	46502	3.63%	0.382%
57	14.055	2059	2063	2069	rVB7	40049	76343	5.95%	0.626%
58	14.126	2069	2075	2080	rBV	111672	178372	13.91%	1.464%
59	14.232	2088	2093	2100	rVB5	46787	81674	6.37%	0.670%
60	14.314	2100	2107	2113	rBV8	20065	47912	3.74%	0.393%
61	14.384	2114	2119	2122	rBV4	20493	38499	3.00%	0.316%
62	14.455	2127	2131	2136	rVB	69323	112495	8.77%	0.923%
63	14.532	2140	2144	2149	rBV3	44670	66058	5.15%	0.542%
64	14.720	2172	2176	2181	rBV3	42127	72530	5.66%	0.595%
65	14.767	2181	2184	2188	rVB4	16633	26445	2.06%	0.217%
66	14.837	2188	2196	2201	rBV5	73277	185509	14.47%	1.522%
67	14.890	2201	2205	2212	rVB2	33138	63190	4.93%	0.518%
68	15.055	2229	2233	2236	rBV6	13410	23413	1.83%	0.192%
69	15.102	2236	2241	2246	rVV2	56305	105652	8.24%	0.867%
70	15.149	2246	2249	2255	rVV7	15863	29220	2.28%	0.240%
71	15.214	2255	2260	2270	rVB6	43555	112118	8.74%	0.920%

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Instrument :
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 ClientSampleId :
 SU-01-073020

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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72	15.296	2270	2274	2278	rVB6	9798	16158	1.26%	0.133%
73	15.373	2282	2287	2291	rBV7	12952	22661	1.77%	0.186%
74	15.520	2306	2312	2316	rVV8	19569	41629	3.25%	0.342%
75	15.561	2316	2319	2329	rVB9	18241	36655	2.86%	0.301%
76	15.708	2338	2344	2351	rBV3	29664	71709	5.59%	0.588%
77	15.873	2362	2372	2376	rBV3	20947	56865	4.43%	0.467%
78	16.008	2388	2395	2400	rBV8	15521	32801	2.56%	0.269%
79	16.078	2400	2407	2415	rVB4	18537	45810	3.57%	0.376%
80	16.231	2426	2433	2436	rBV7	9551	21579	1.68%	0.177%
81	16.508	2473	2480	2494	rVB2	43866	117331	9.15%	0.963%
82	16.714	2506	2515	2521	rBV	37199	87422	6.82%	0.717%

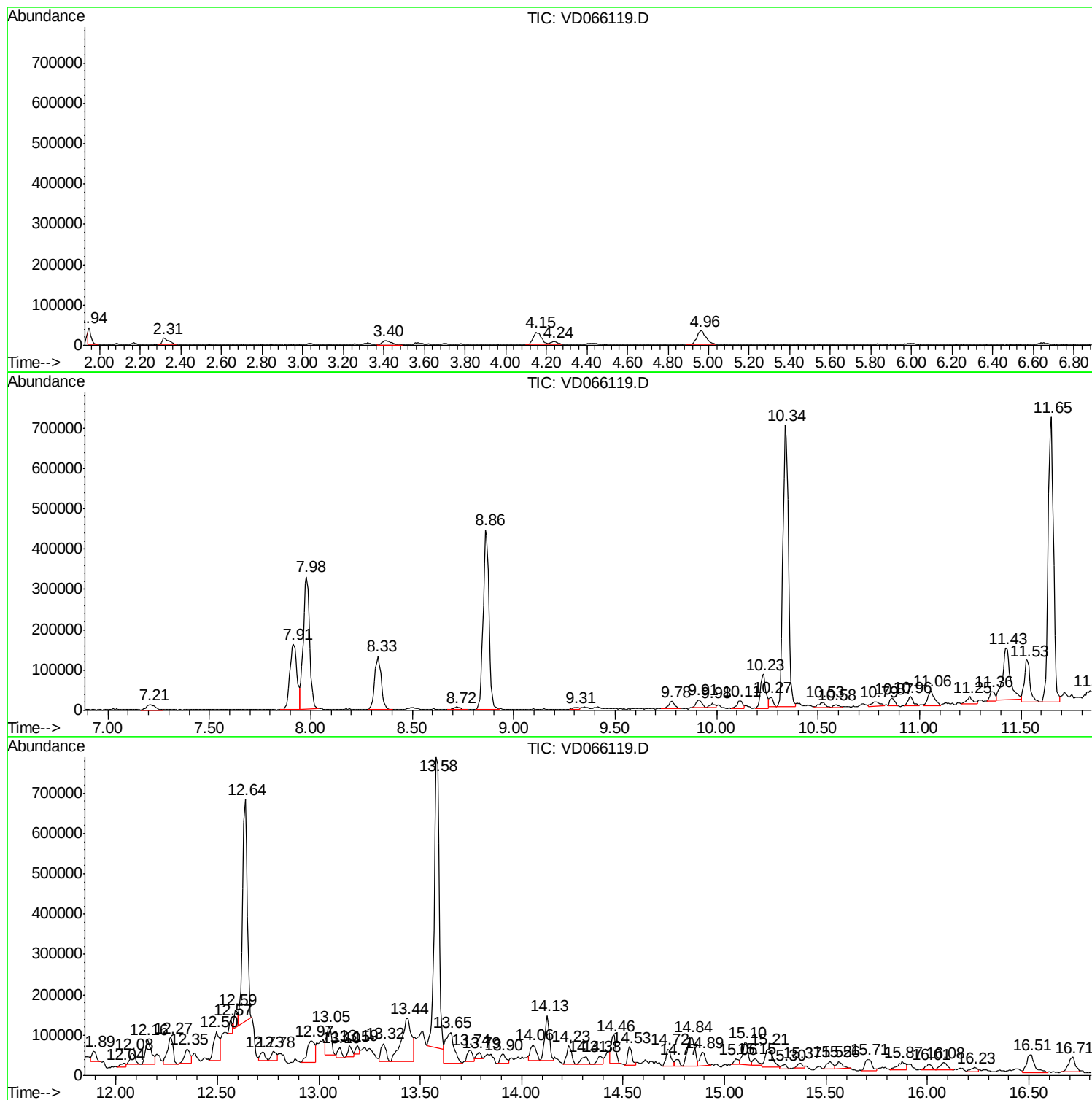
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Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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



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Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

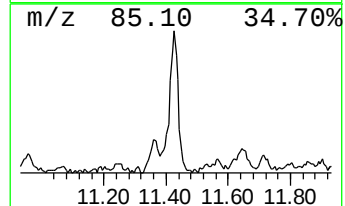
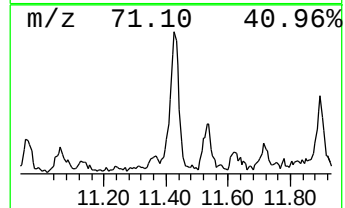
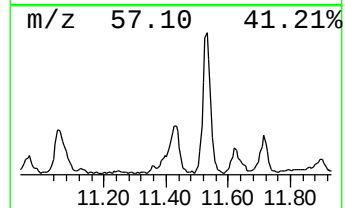
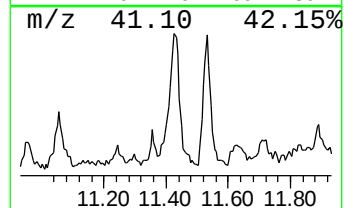
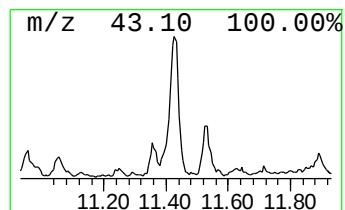
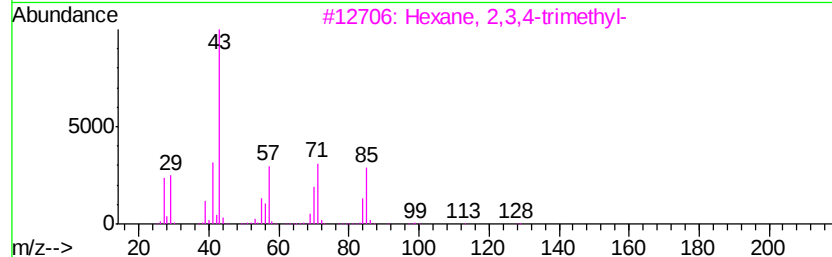
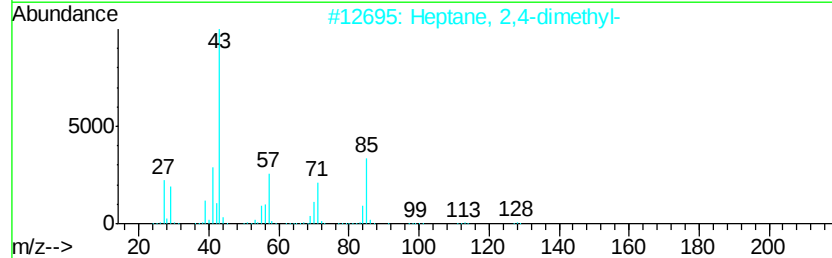
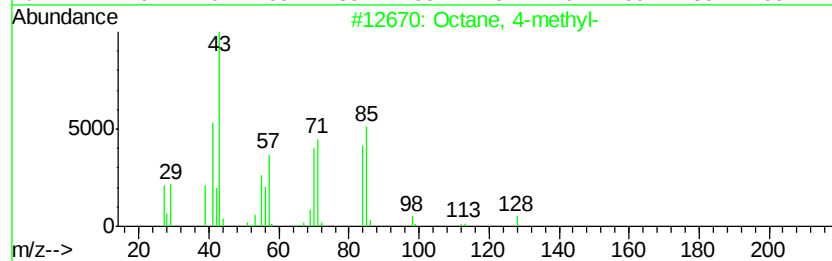
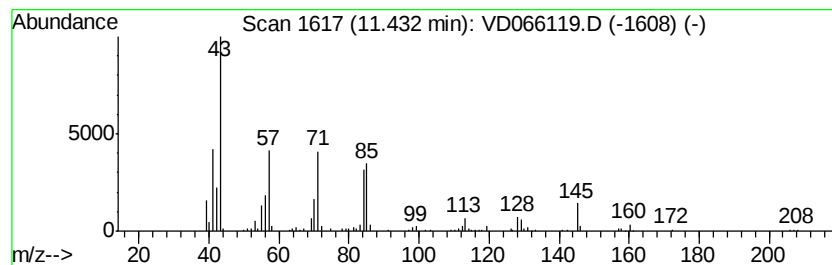
Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.43	13.52 ug/l	344183	Chlorobenzene-d5	11.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-methyl-		128	C9H20	002216-34-4	68
2	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	64
3	Hexane, 2,3,4-trimethyl-		128	C9H20	000921-47-1	64
4	Hexane, 3-ethyl-		114	C8H18	000619-99-8	59
5	Ether, hexyl pentyl		172	C11H24O	032357-83-8	53



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Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

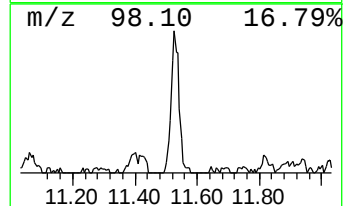
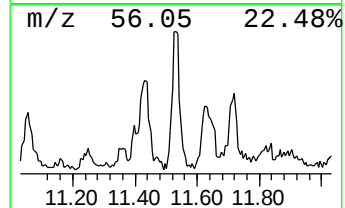
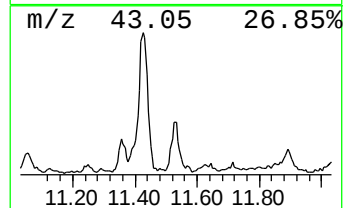
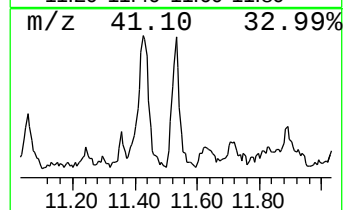
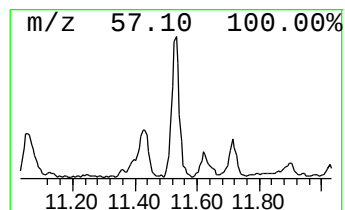
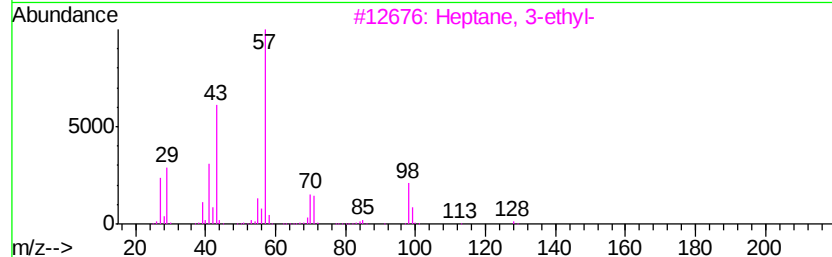
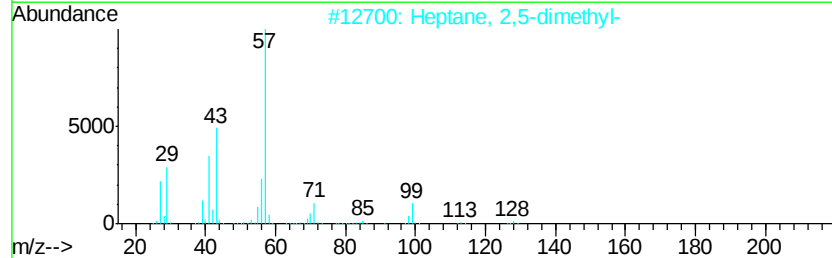
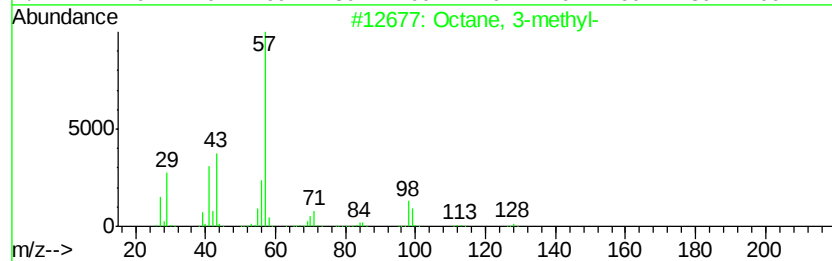
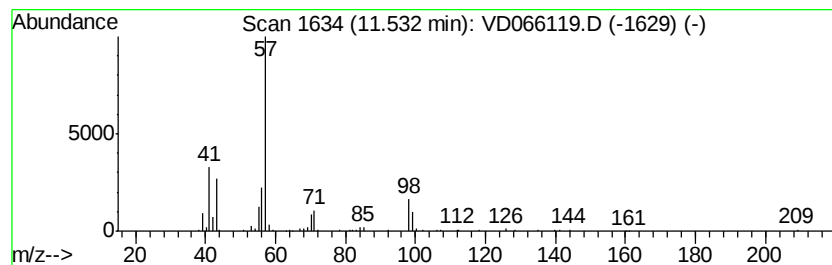
Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 2 Octane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.53	8.00 ug/l	203751	Chlorobenzene-d5	11.65		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-methyl-		128	C9H20	002216-33-3	87
2	Heptane, 2,5-dimethyl-		128	C9H20	002216-30-0	76
3	Heptane, 3-ethyl-		128	C9H20	015869-80-4	58
4	Undecane, 6-methyl-		170	C12H26	017302-33-9	53
5	Heptane, 3-ethyl-2-methyl-		142	C10H22	014676-29-0	50



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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SU-01-073020

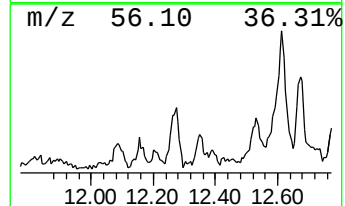
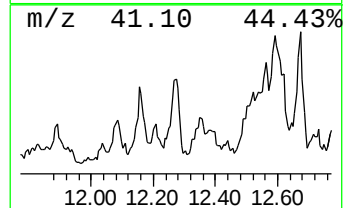
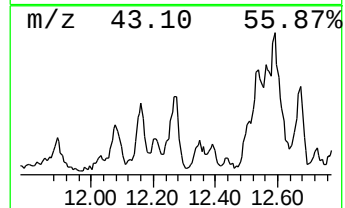
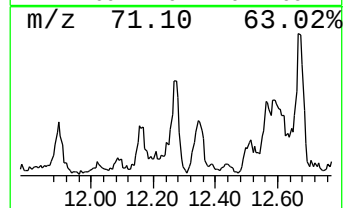
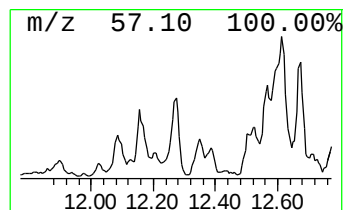
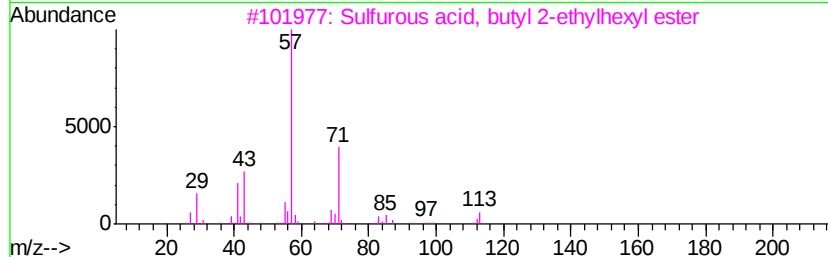
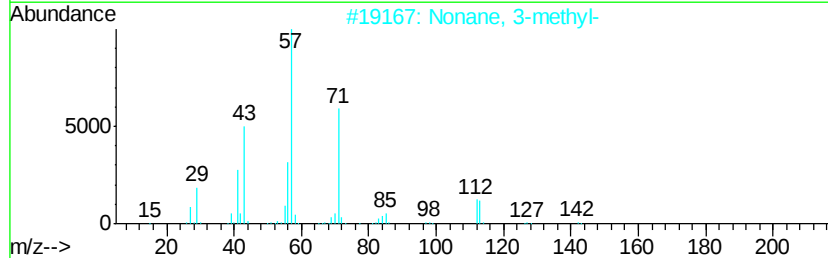
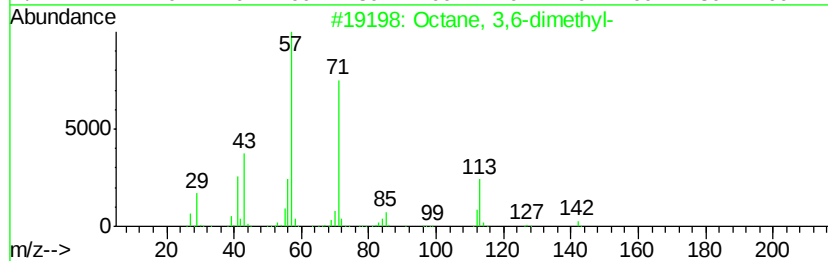
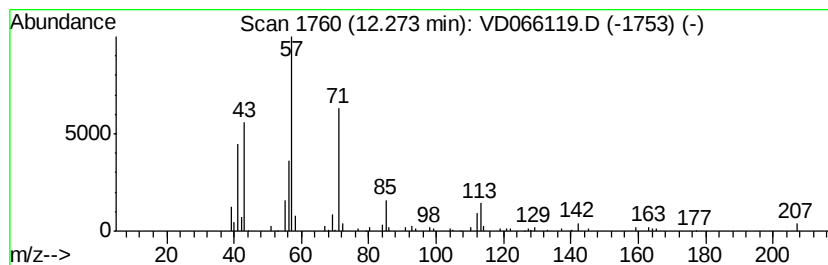
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 3 Octane, 3,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.27	5.29 ug/l	134660	Chlorobenzene-d5	11.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	78
2		Nonane, 3-methyl-	142	C10H22	005911-04-6	72
3		Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	59
4		Ethanol, 2-(octyloxy)-	174	C10H22O2	010020-43-6	59
5		Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	59



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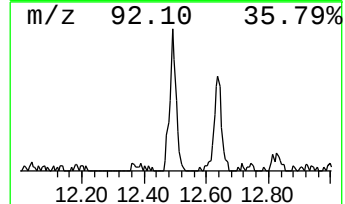
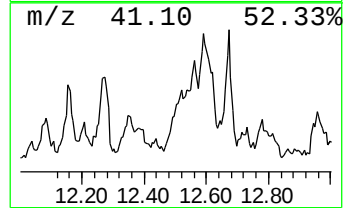
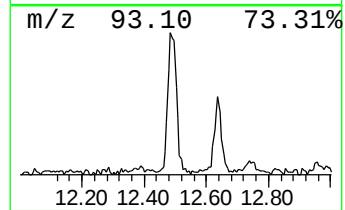
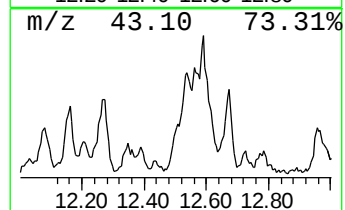
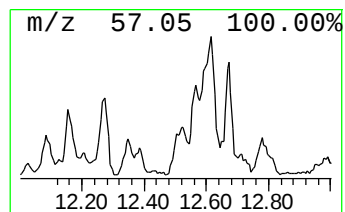
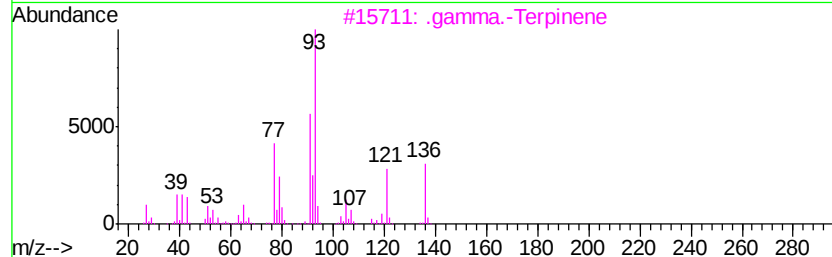
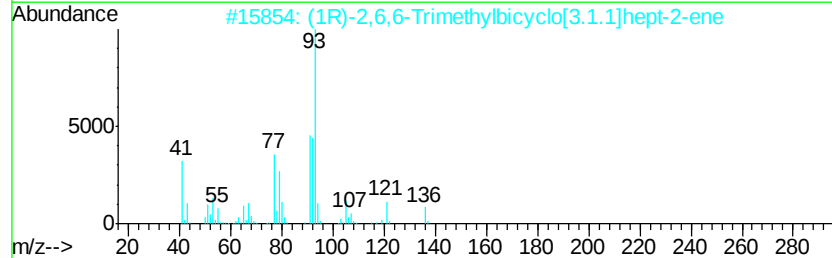
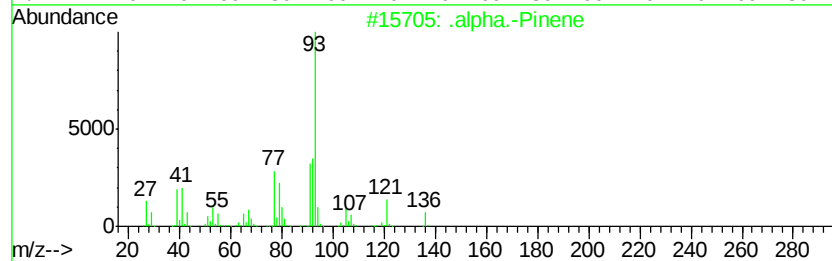
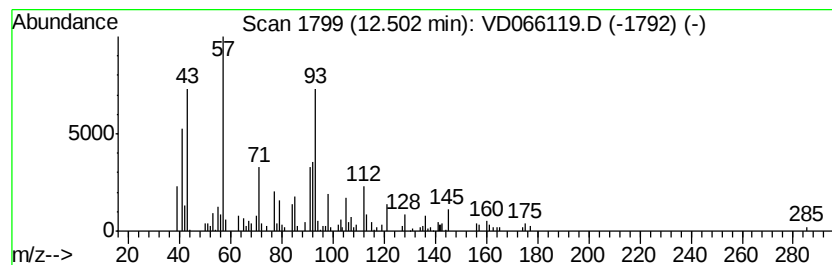
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MSVOA_D
ClientSampleId :
SU-01-073020

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 4 .alpha.-Pinene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.50	5.59 ug/l	142429	Chlorobenzene-d5	11.65	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	81
2	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	72
3	.gamma.-Terpinene	136	C10H16	000099-85-4	70
4	Bicyclo[3.1.0]hex-2-ene, 2-methy...	136	C10H16	002867-05-2	64
5	.beta.-Ocimene	136	C10H16	013877-91-3	62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD073020\
Data File : VD066119.D
Acq On : 30 Jul 2020 18:48
Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

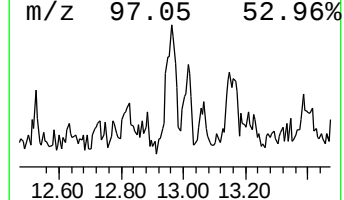
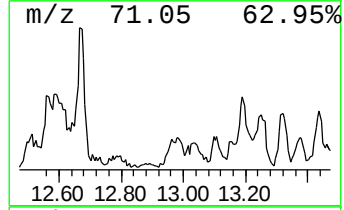
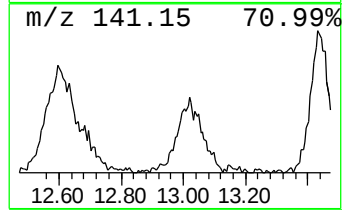
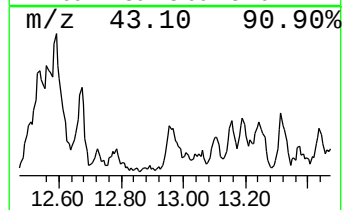
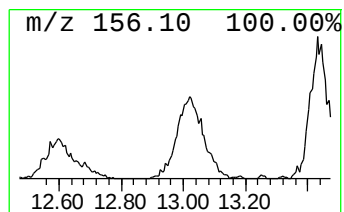
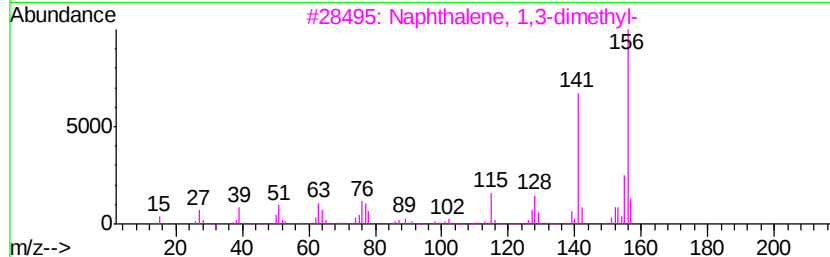
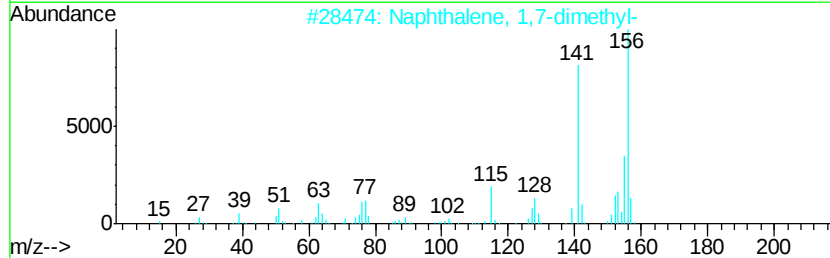
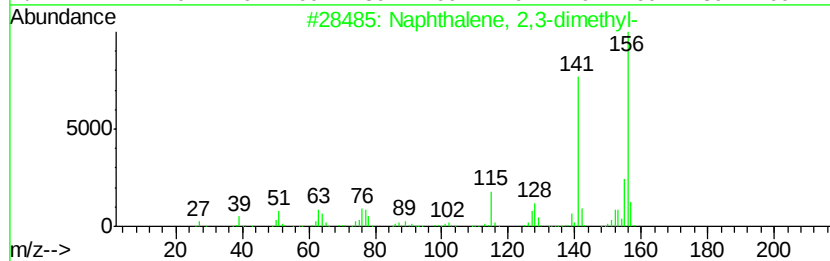
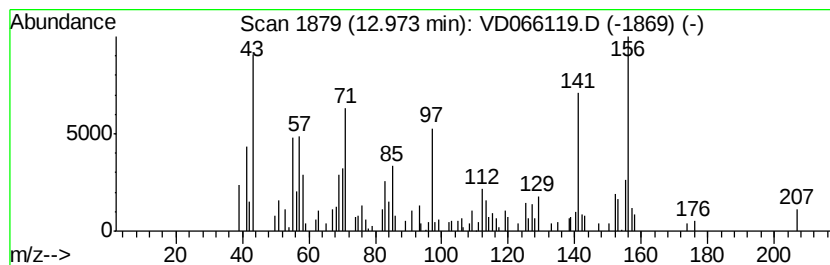
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Quant Title : SW846 8260

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

Peak Number 5 unknown12.97 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	5.87 ug/l	141126	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	38
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	38
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	30
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	30
5		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD073020\
Data File : VD066119.D
Acq On : 30 Jul 2020 18:48
Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

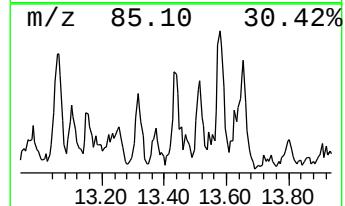
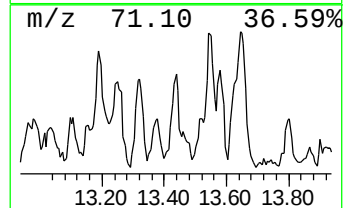
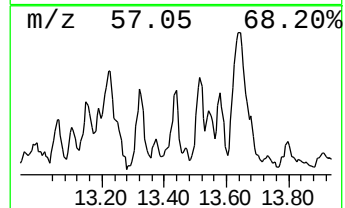
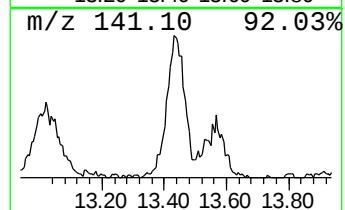
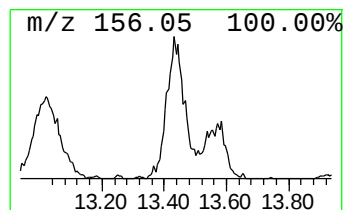
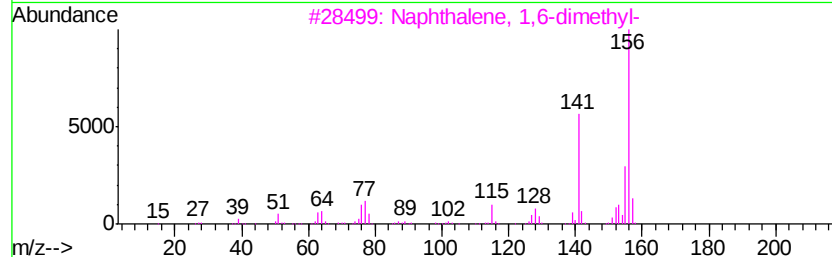
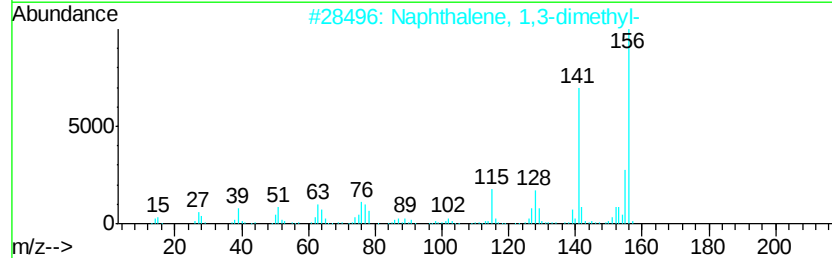
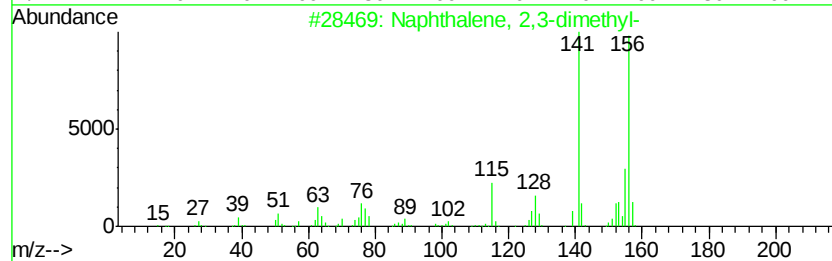
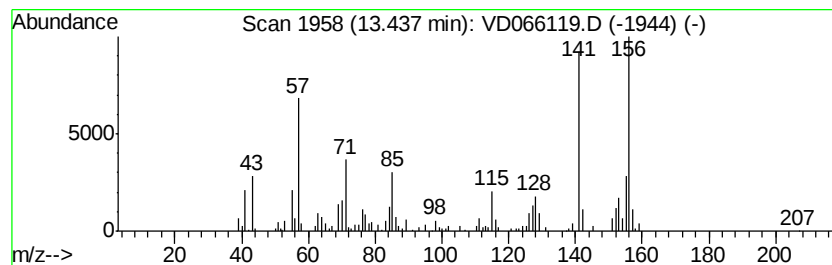
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	15.74 ug/l	378612	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	93
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	91
4		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	91
5		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD073020\
Data File : VD066119.D
Acq On : 30 Jul 2020 18:48
Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

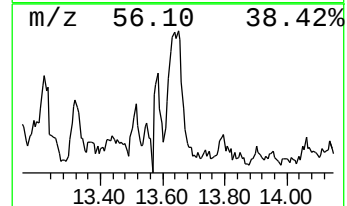
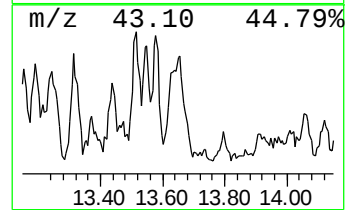
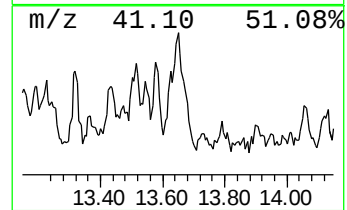
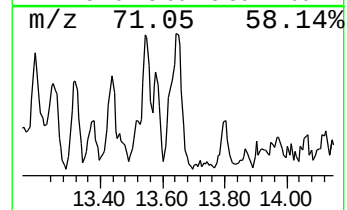
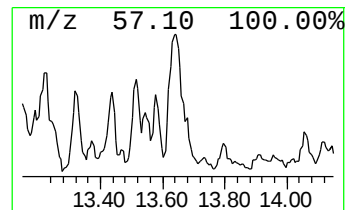
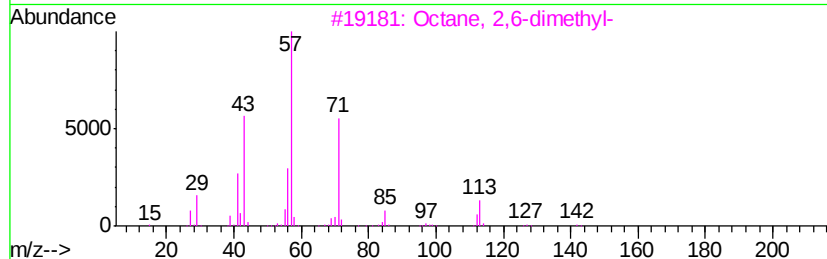
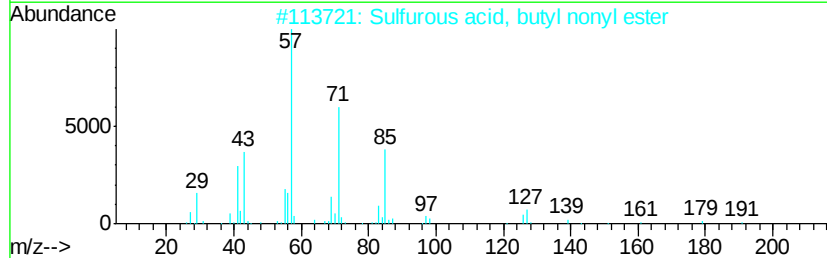
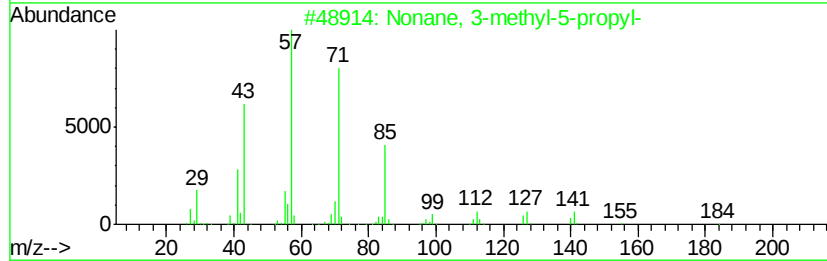
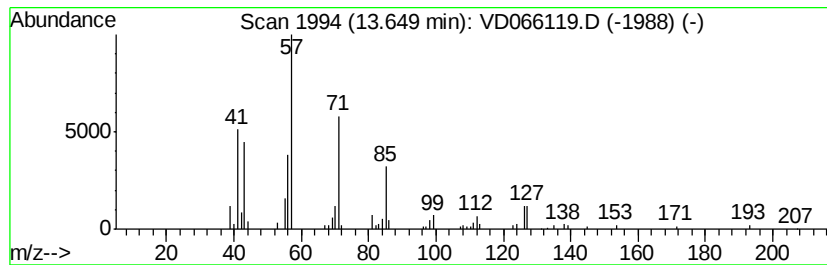
Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 7 Nonane, 3-methyl-5-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	8.98 ug/l	216071	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Nonane, 3-methyl-5-propyl-	184 C13H28	031081-18-2	53
2	Sulfurous acid, butyl nonyl ester	264 C13H28O3S	1000309-17-6	53
3	Octane, 2,6-dimethyl-	142 C10H22	002051-30-1	50
4	Tridecane	184 C13H28	000629-50-5	50
5	Tritetracontane	605 C43H88	007098-21-7	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD073020\
Data File : VD066119.D
Acq On : 30 Jul 2020 18:48
Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

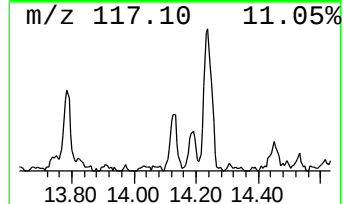
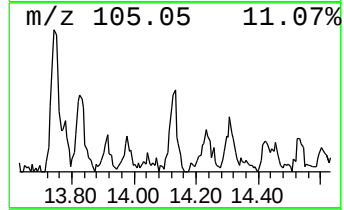
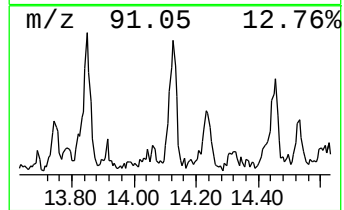
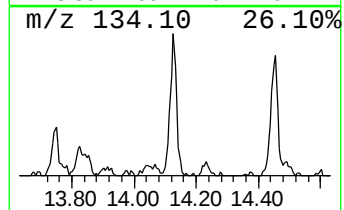
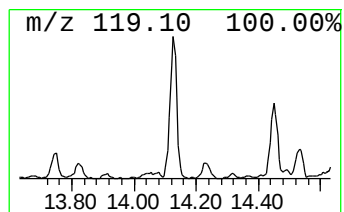
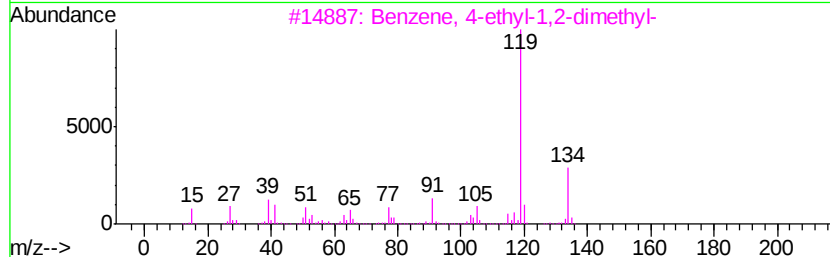
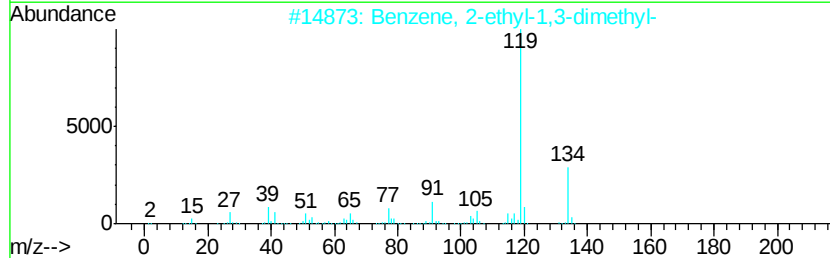
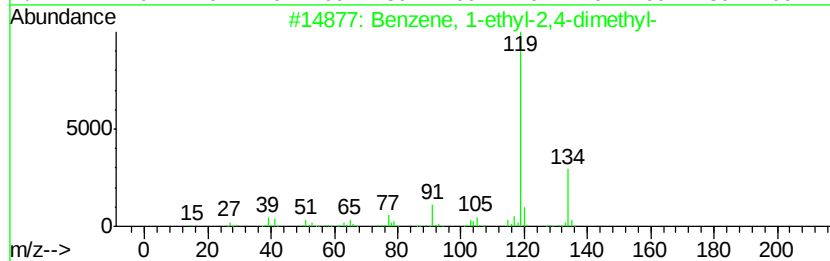
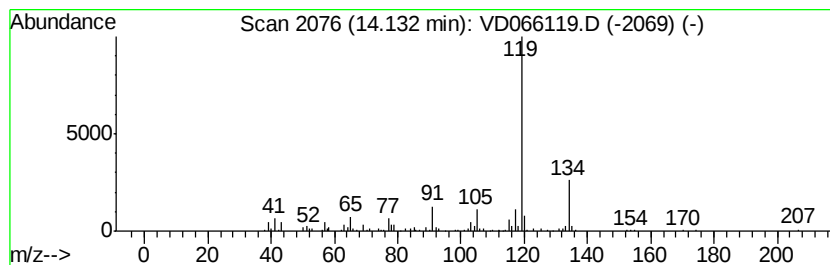
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	7.42 ug/l	178372	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD073020\
Data File : VD066119.D
Acq On : 30 Jul 2020 18:48
Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

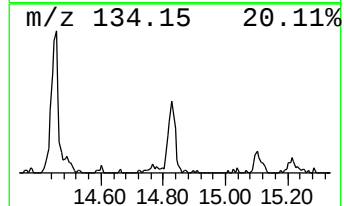
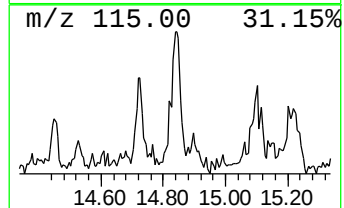
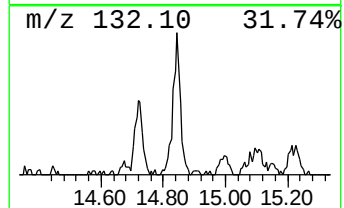
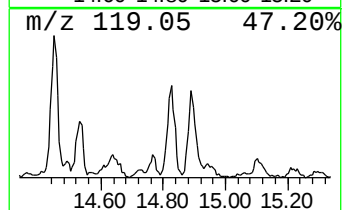
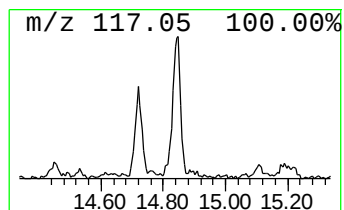
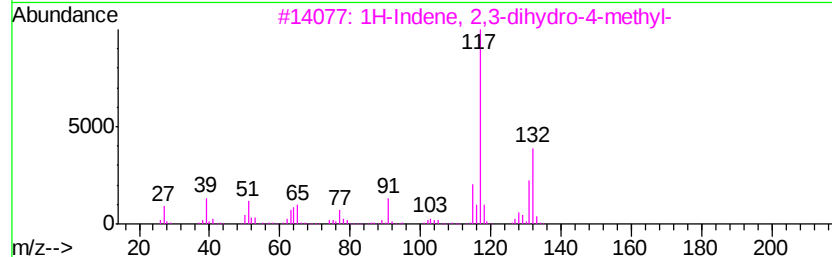
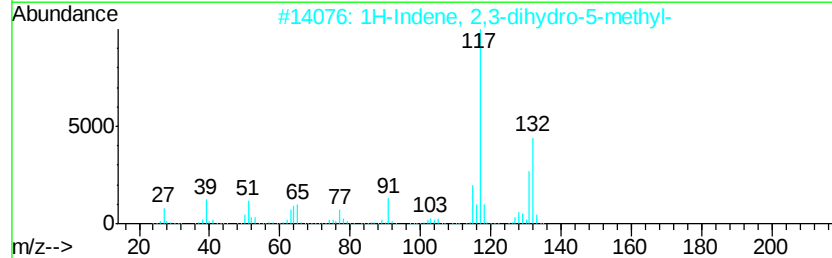
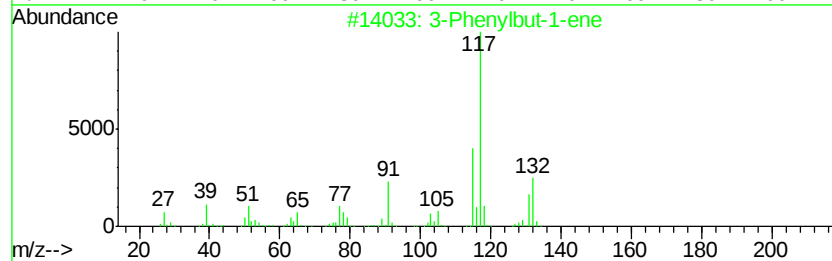
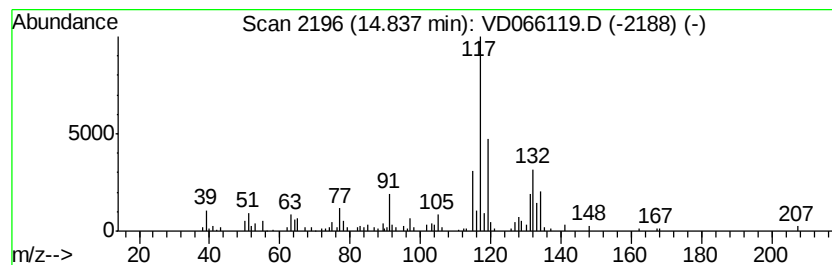
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.M
Quant Title : SW846 8260

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

Peak Number 9 3-Phenylbut-1-ene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	7.71 ug/l	185509	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	64
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD073020\
Data File : VD066119.D
Acq On : 30 Jul 2020 18:48
Operator : VA/SY
Sample : L3507-01
Misc : 8.17G/5.00ml/MSVOA_D/SOIL
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-01-073020

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D072120S.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	13.5	ug/l	344183	3	11.65	1273230	50.0
Octane, 3-methyl-	11.53	8.0	ug/l	203751	3	11.65	1273230	50.0
Octane, 3,6-dimet...	12.27	5.3	ug/l	134660	3	11.65	1273230	50.0
.alpha.-Pinene	12.50	5.6	ug/l	142429	3	11.65	1273230	50.0
unknown12.97	12.97	5.9	ug/l	141126	4	13.58	1202680	50.0
Naphthalene, 2,3-...	13.44	15.7	ug/l	378612	4	13.58	1202680	50.0
Nonane, 3-methyl-...	13.65	9.0	ug/l	216071	4	13.58	1202680	50.0
Benzene, 1-ethyl-...	14.13	7.4	ug/l	178372	4	13.58	1202680	50.0
3-Phenylbut-1-ene	14.84	7.7	ug/l	185509	4	13.58	1202680	50.0