

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD022020\
 Data File : VD065253.D
 Acq On : 20 Feb 2020 13:35
 Operator : VA/SY
 Sample : L1601-05
 Misc : 5.68G/5.00ml/MSVOA D/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 SW-4

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\82D021820S.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	3.038	182	190	200	rBV	137117	295709	8.42%	0.472%
2	3.409	245	253	265	rVB2	100212	271730	7.74%	0.434%
3	3.879	323	333	343	rVB2	75635	200317	5.71%	0.320%
4	4.162	368	381	392	rVB3	25782	94623	2.70%	0.151%
5	5.015	501	526	543	rBV7	105899	689970	19.65%	1.102%
6	5.485	592	606	622	rBV3	107638	379464	10.81%	0.606%
7	5.809	647	661	670	rBV4	20940	71057	2.02%	0.114%
8	5.991	679	692	706	rBV	105339	328305	9.35%	0.524%
9	6.344	741	752	763	rVB3	58982	179305	5.11%	0.286%
10	6.479	765	775	785	rBV4	32002	100796	2.87%	0.161%
11	6.779	815	826	836	rBV3	42186	128312	3.65%	0.205%
12	6.926	839	851	858	rBV3	34643	103342	2.94%	0.165%
13	7.032	858	869	877	rBV2	212981	604624	17.22%	0.966%
14	7.220	892	901	914	rVB	31676	93255	2.66%	0.149%
15	7.738	981	989	996	rVB3	79961	191727	5.46%	0.306%
16	7.920	1009	1020	1025	rBV	434818	1136885	32.38%	1.816%
17	7.985	1025	1031	1039	rVV	995663	2392190	68.14%	3.821%
18	8.067	1039	1045	1056	rVB2	163498	410038	11.68%	0.655%
19	8.191	1057	1066	1073	rBV2	174205	409106	11.65%	0.654%
20	8.350	1082	1093	1103	rVB3	657864	1752678	49.92%	2.800%
21	8.509	1103	1120	1131	rVV	480671	1510099	43.01%	2.412%
22	8.609	1131	1137	1146	rVV5	61103	160067	4.56%	0.256%
23	8.726	1146	1157	1165	rVB2	185362	403714	11.50%	0.645%
24	8.867	1170	1181	1194	rBV	1298597	2839146	80.87%	4.535%
25	9.144	1223	1228	1239	rVB3	53142	133385	3.80%	0.213%
26	9.356	1250	1264	1269	rBV	354811	874324	24.90%	1.397%
27	9.409	1269	1273	1281	rVB	153766	303607	8.65%	0.485%
28	9.491	1281	1287	1291	rBV2	34071	65254	1.86%	0.104%
29	9.538	1291	1295	1301	rVB2	64586	116476	3.32%	0.186%
30	9.632	1301	1311	1317	rVB5	31090	83795	2.39%	0.134%
31	9.785	1330	1337	1346	rVB	436314	877841	25.01%	1.402%
32	9.879	1346	1353	1354	rBV	109704	185243	5.28%	0.296%
33	9.914	1354	1359	1366	rVB	530234	1023967	29.17%	1.636%
34	9.979	1366	1370	1374	rBV	210383	344572	9.82%	0.550%

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Integrator: RTE
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 Title : SW846 8260

35	10.120	1384	1394	1405	rBV	288648	688203	19.60%	1.099%
36	10.226	1405	1412	1414	rVV3	56791	100329	2.86%	0.160%
37	10.273	1414	1420	1425	rVV	347138	618321	17.61%	0.988%
38	10.344	1425	1432	1439	rVV	1928922	3510638	100.00%	5.608%
39	10.408	1439	1443	1449	rVV	232540	444572	12.66%	0.710%
40	10.456	1449	1451	1456	rVV3	46273	86788	2.47%	0.139%
41	10.526	1456	1463	1470	rVV2	331278	632677	18.02%	1.011%
42	10.726	1492	1497	1500	rVV2	81164	157574	4.49%	0.252%
43	10.785	1500	1507	1515	rVB7	116207	348411	9.92%	0.557%
44	10.867	1516	1521	1527	rVB3	80626	146405	4.17%	0.234%
45	10.961	1527	1537	1542	rBV2	73088	154824	4.41%	0.247%
46	11.061	1542	1554	1561	rBV	130381	364657	10.39%	0.583%
47	11.161	1568	1571	1574	rVV5	36111	57187	1.63%	0.091%
48	11.361	1600	1605	1608	rBV2	78061	115822	3.30%	0.185%
49	11.432	1608	1617	1628	rVB2	401667	950721	27.08%	1.519%
50	11.532	1628	1634	1645	rBV	307429	545467	15.54%	0.871%
51	11.650	1645	1654	1661	rBV	1626978	2936323	83.64%	4.691%
52	11.720	1662	1666	1667	rVV	110679	154847	4.41%	0.247%
53	11.750	1667	1671	1680	rVB	325878	551773	15.72%	0.881%
54	11.855	1680	1689	1701	rBV	766989	1646882	46.91%	2.631%
55	12.091	1724	1729	1735	rVV3	80733	163544	4.66%	0.261%
56	12.185	1736	1745	1754	rVV2	241555	589861	16.80%	0.942%
57	12.273	1754	1760	1765	rVB2	80835	139302	3.97%	0.223%
58	12.355	1767	1774	1778	rBV6	53643	120951	3.45%	0.193%
59	12.485	1791	1796	1801	rBV	241666	420291	11.97%	0.671%
60	12.638	1812	1822	1832	rVB2	1275492	2771220	78.94%	4.427%
61	12.779	1841	1846	1849	rBV	63888	105208	3.00%	0.168%
62	12.826	1849	1854	1860	rVB	919668	1425640	40.61%	2.277%
63	12.891	1860	1865	1868	rBV	130141	200921	5.72%	0.321%
64	12.920	1868	1870	1873	rVV	50190	57217	1.63%	0.091%
65	12.967	1873	1878	1895	rVB4	193324	488456	13.91%	0.780%
66	13.126	1897	1905	1914	rBV	147957	323419	9.21%	0.517%
67	13.220	1914	1921	1925	rVB3	36151	74730	2.13%	0.119%
68	13.279	1925	1931	1936	rVB	322711	488769	13.92%	0.781%
69	13.402	1944	1952	1956	rBV3	133048	350040	9.97%	0.559%
70	13.585	1975	1983	2005	rVB2	1504249	3053719	86.98%	4.878%
71	13.749	2005	2011	2013	rBV	438202	656717	18.71%	1.049%

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72	13.785	2013	2017	2022	rVB	1298299	1964342	55.95%	3.138%
73	13.832	2022	2025	2026	rBV2	86511	89254	2.54%	0.143%
74	13.914	2034	2039	2044	rBV	84930	139294	3.97%	0.223%
75	13.979	2045	2050	2057	rVB	391934	618935	17.63%	0.989%
76	14.044	2057	2061	2063	rBV2	40012	61117	1.74%	0.098%
77	14.132	2069	2076	2082	rBV	1452154	2250533	64.11%	3.595%
78	14.191	2082	2086	2089	rVV	111876	171669	4.89%	0.274%
79	14.238	2089	2094	2101	rVB2	444137	766678	21.84%	1.225%
80	14.308	2101	2106	2112	rBV2	34038	64682	1.84%	0.103%
81	14.455	2121	2131	2136	rBV	1006233	1639653	46.71%	2.619%
82	14.538	2140	2145	2149	rBV	138372	190236	5.42%	0.304%
83	14.608	2154	2157	2159	rVV	38892	55530	1.58%	0.089%
84	14.632	2159	2161	2165	rVV2	46125	71456	2.04%	0.114%
85	14.726	2171	2177	2182	rVV	604494	1015153	28.92%	1.622%
86	14.767	2182	2184	2189	rVB	92111	111039	3.16%	0.177%
87	14.844	2189	2197	2203	rBV3	1072893	2351211	66.97%	3.756%
88	14.896	2203	2206	2213	rVB2	197632	335464	9.56%	0.536%
89	14.996	2218	2223	2228	rBV2	64280	105505	3.01%	0.169%
90	15.108	2236	2242	2247	rBV	279637	442688	12.61%	0.707%
91	15.220	2254	2261	2268	rVB4	204866	452818	12.90%	0.723%
92	15.414	2281	2294	2301	rBV	1749855	3139431	89.43%	5.015%
93	15.520	2307	2312	2317	rBV3	67984	122181	3.48%	0.195%
94	15.579	2317	2322	2331	rVB2	51683	105018	2.99%	0.168%
95	15.708	2338	2344	2352	rBV	91775	186075	5.30%	0.297%
96	15.890	2369	2375	2383	rVV	48430	148208	4.22%	0.237%
97	16.008	2391	2395	2401	rVV	64649	117359	3.34%	0.187%
98	16.090	2402	2409	2416	rVB4	36209	88355	2.52%	0.141%
99	16.514	2472	2481	2496	rVB	444945	977254	27.84%	1.561%
100	16.720	2507	2516	2525	rBV	173194	419827	11.96%	0.671%

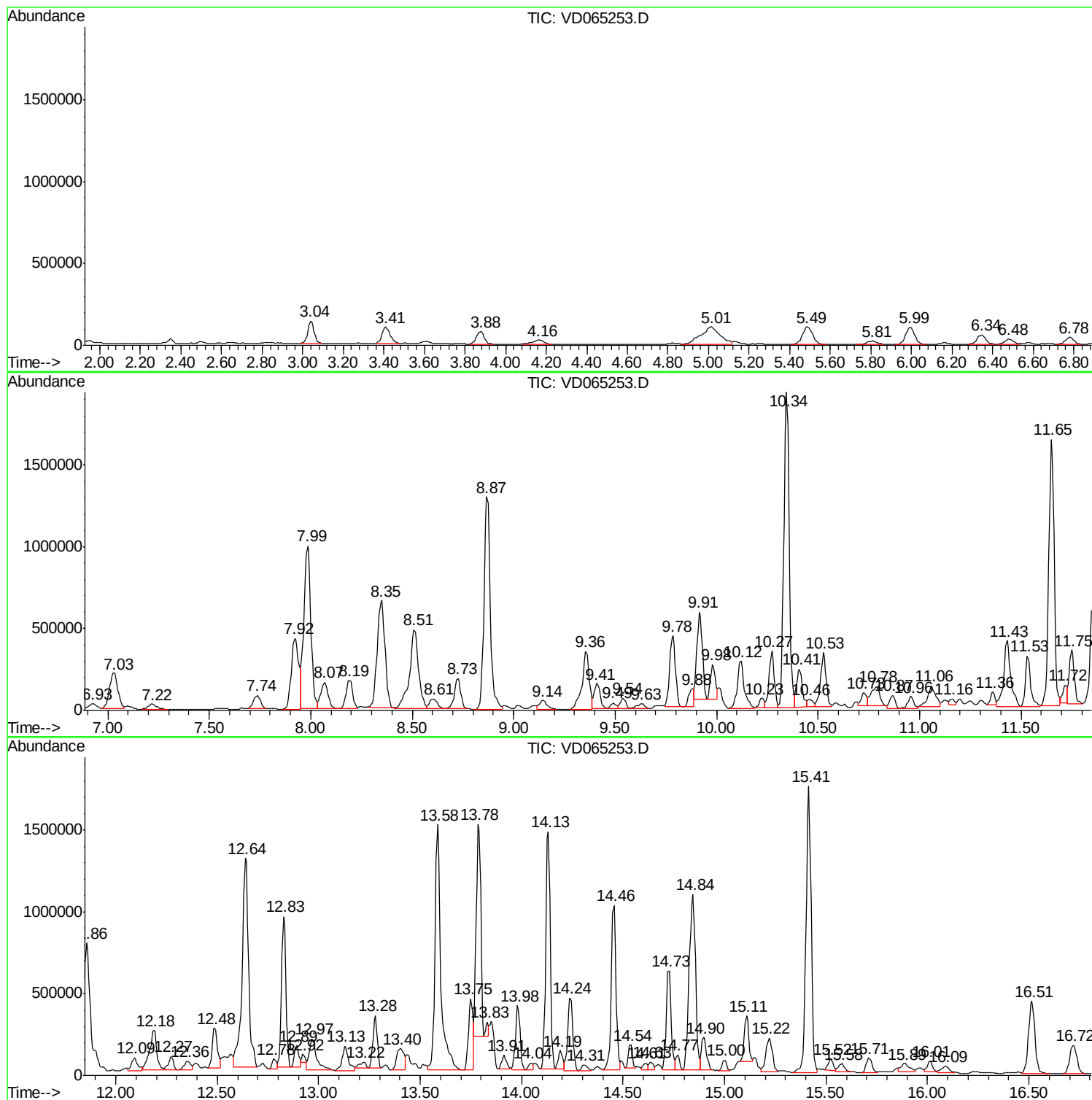
Sum of corrected areas: 62598314

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SW-4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_D
ClientSampled :
SW-4

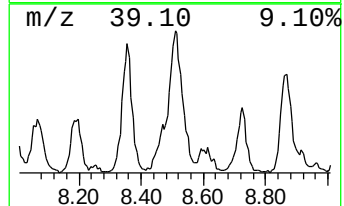
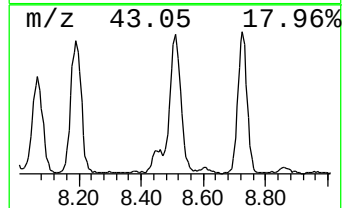
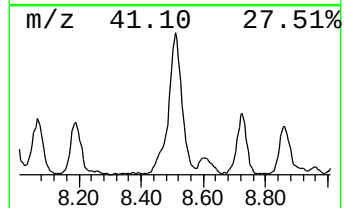
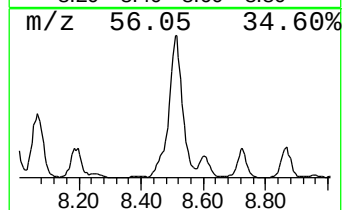
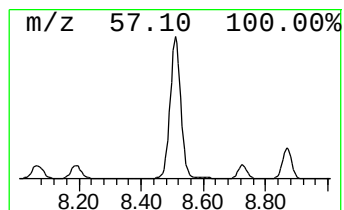
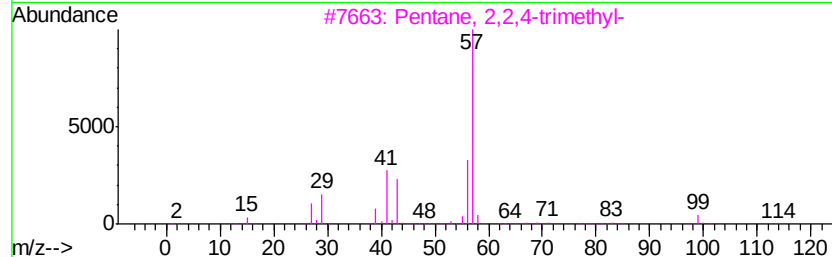
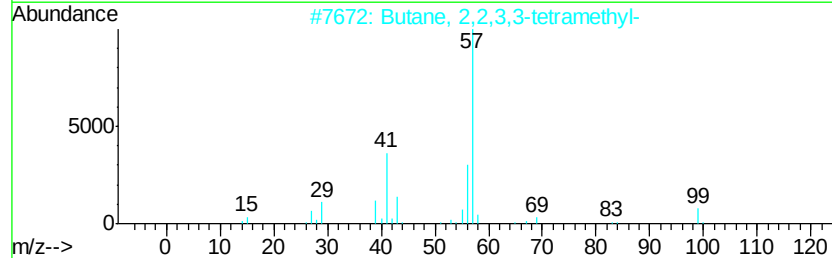
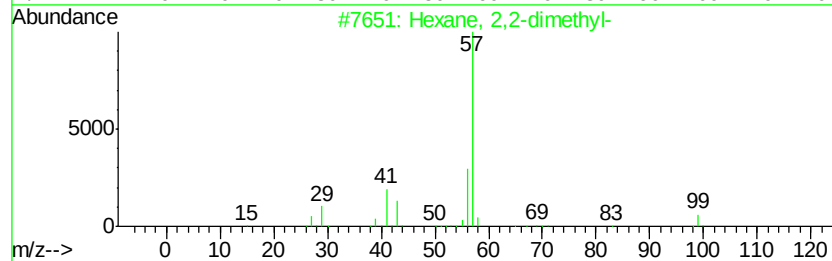
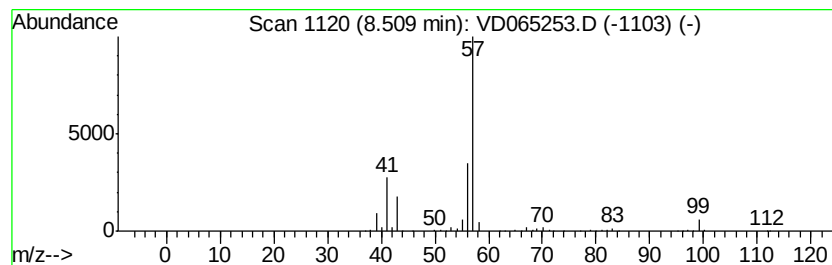
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 2,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.51	26.59 ug/l	1510100	1,4-Difluorobenzene	8.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59



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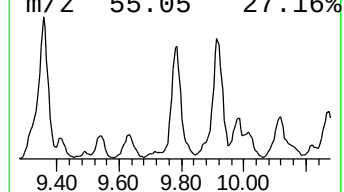
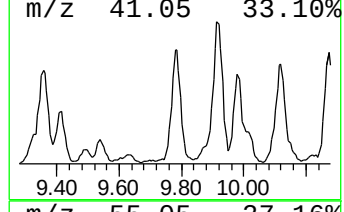
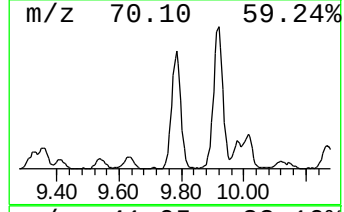
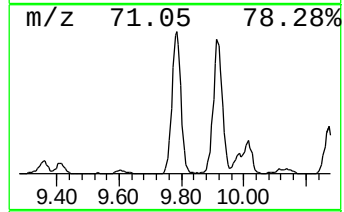
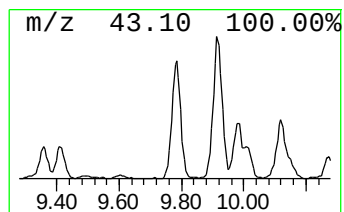
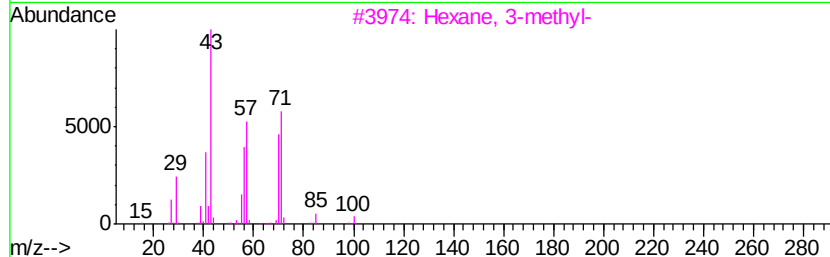
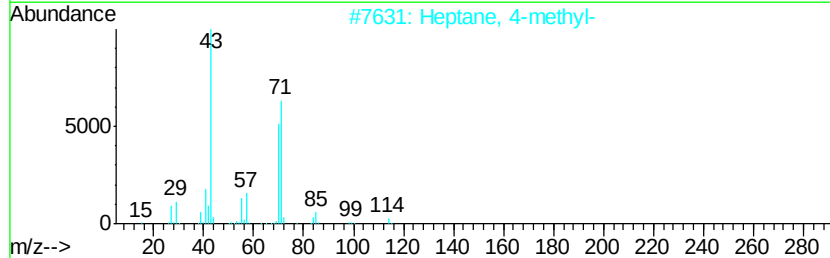
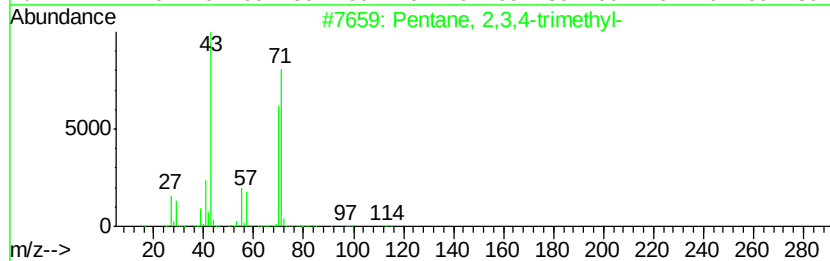
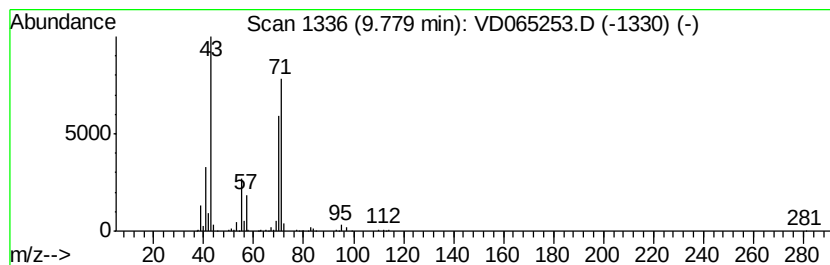
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	15.46 ug/l	877841	1,4-Difluorobenzene	8.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3		Hexane, 3-methyl-	100	C7H16	000589-34-4	78
4		Hexanoic acid, 1,1-dimethylpropyl...	186	C11H22O2	116423-69-9	78
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78



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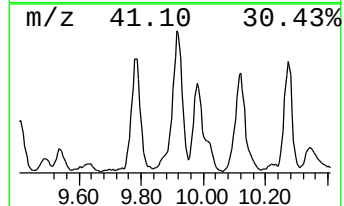
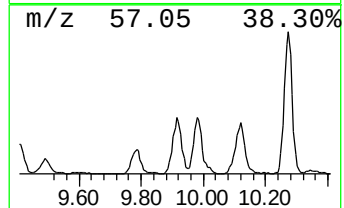
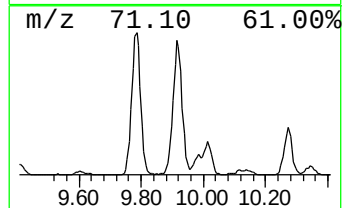
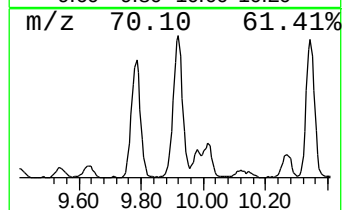
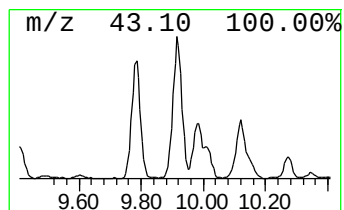
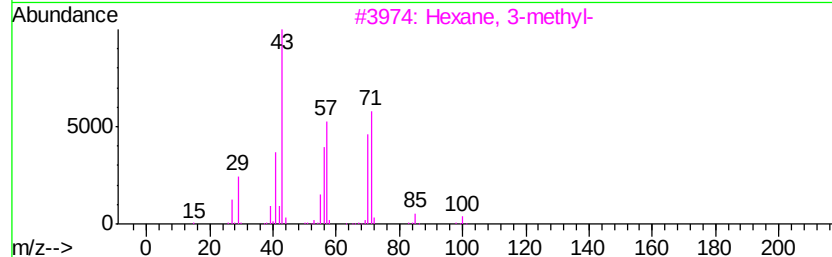
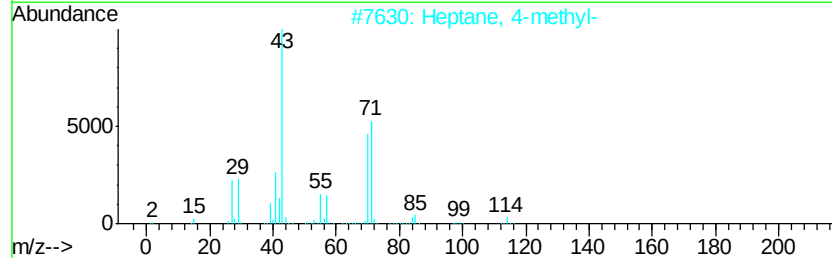
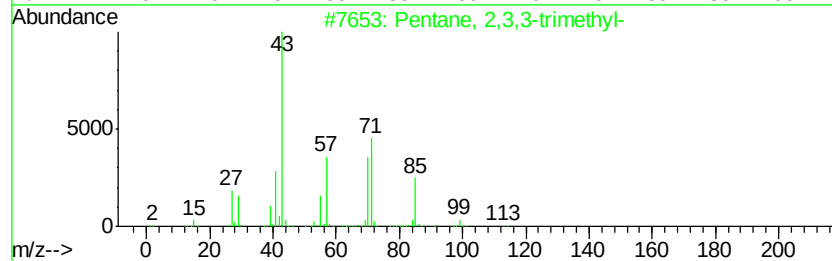
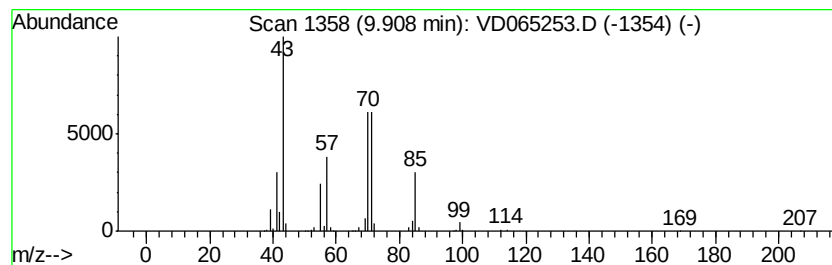
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3,3-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.91	18.03 ug/l	1023970	1,4-Difluorobenzene	8.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	80
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3		Hexane, 3-methyl-	100	C7H16	000589-34-4	64
4		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5		Pentane, 3-ethyl-	100	C7H16	000617-78-7	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SW-4

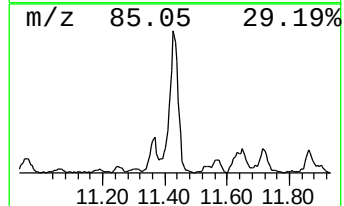
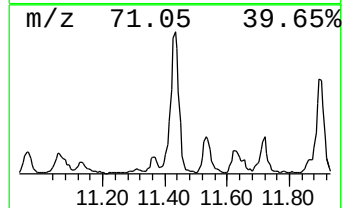
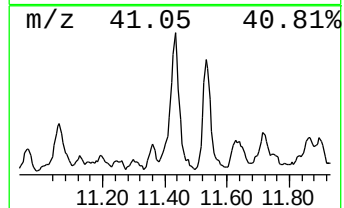
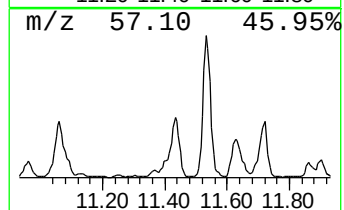
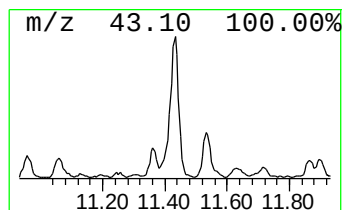
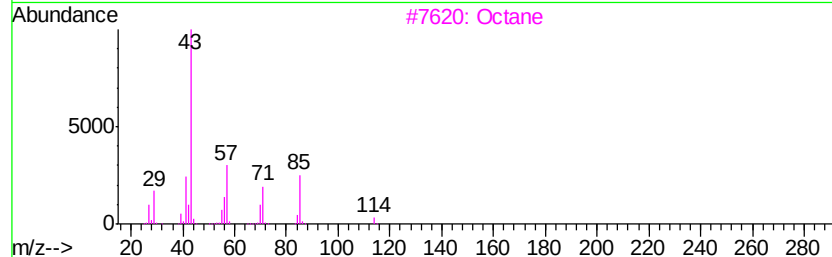
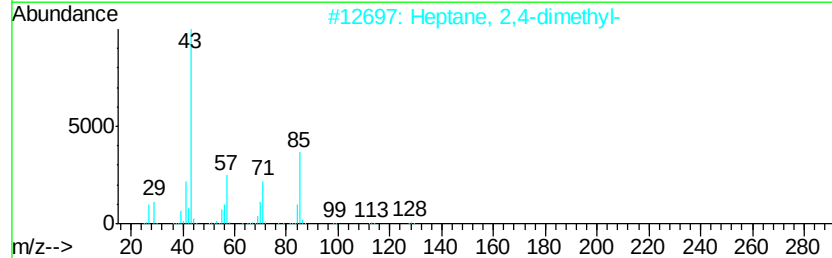
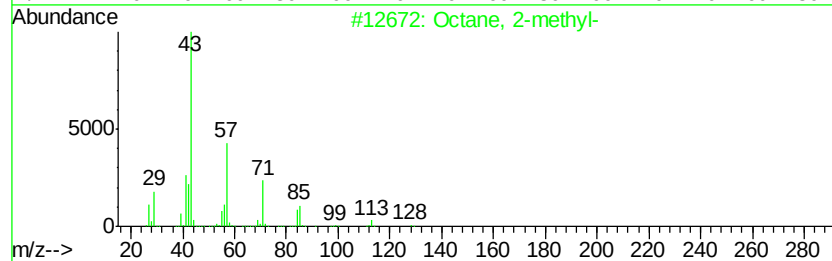
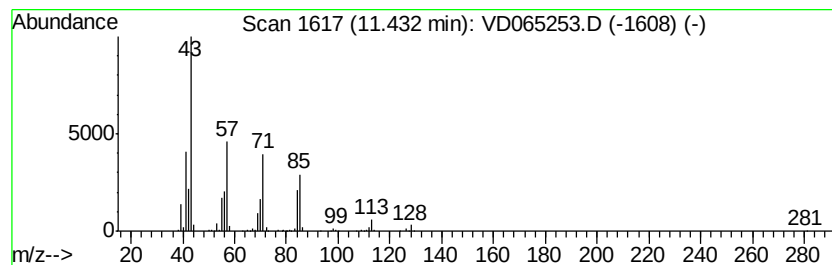
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Quant Title : SW846 8260

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

Peak Number 4 Octane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.43	16.19 ug/l	950721	Chlorobenzene-d5	11.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 2-methyl-	128	C9H20	003221-61-2	83
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3		Octane	114	C8H18	000111-65-9	59
4		Octane, 4-methyl-	128	C9H20	002216-34-4	53
5		Ether, hexyl pentyl	172	C11H24O	032357-83-8	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SW-4

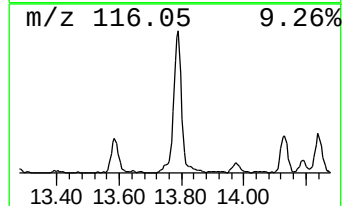
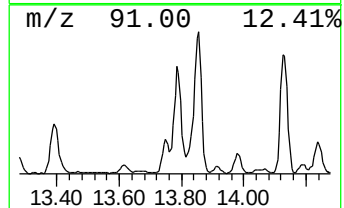
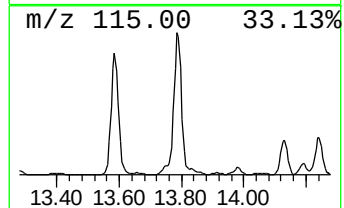
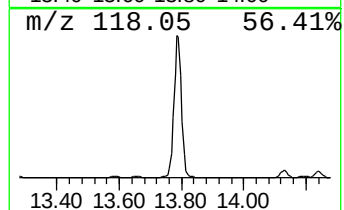
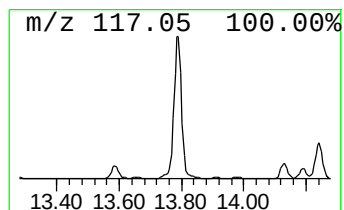
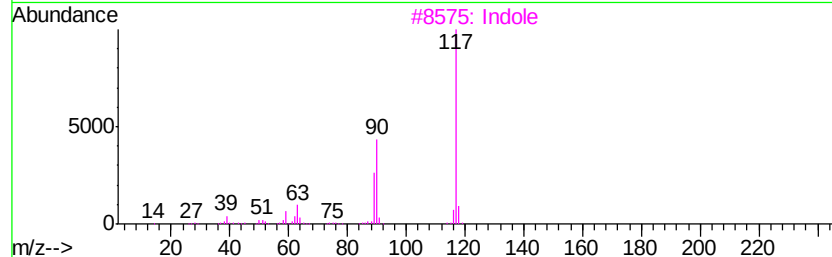
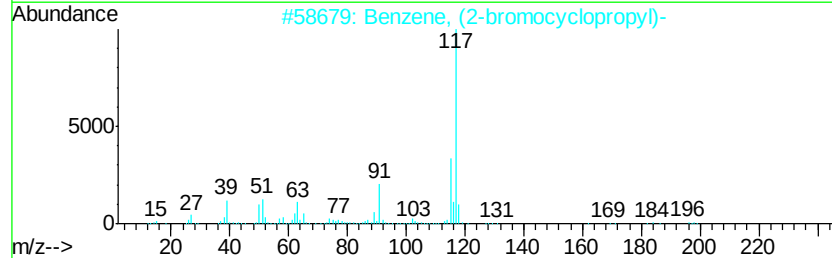
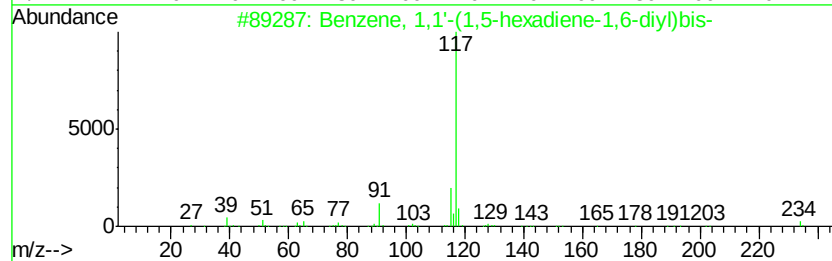
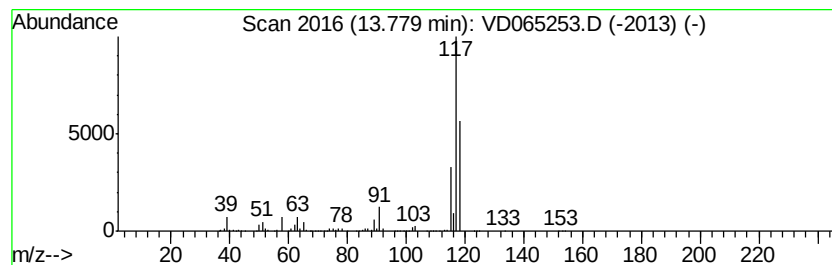
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	32.16 ug/l	1964340	1,4-Dichlorobenzene-d4	13.58

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
3		Indole	117	C8H7N	000120-72-9	43
4		m-Aminophenylacetylene	117	C8H7N	054060-30-9	43
5		trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SW-4

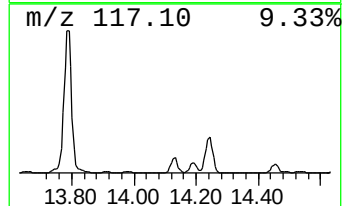
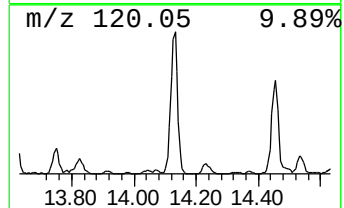
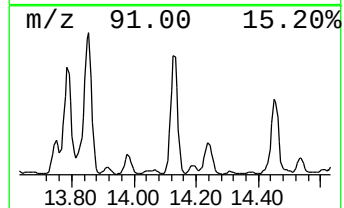
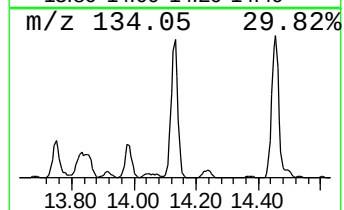
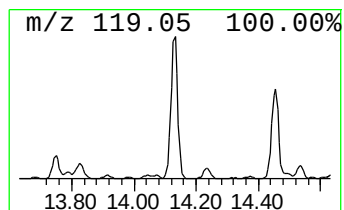
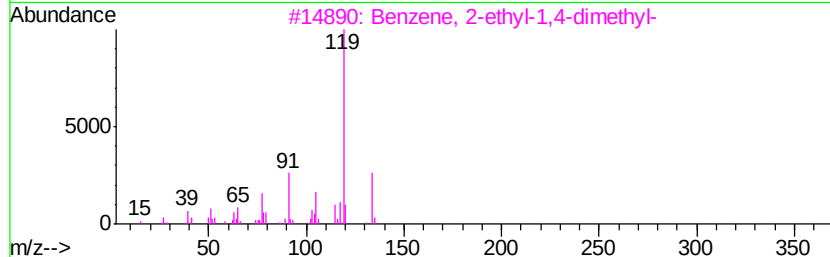
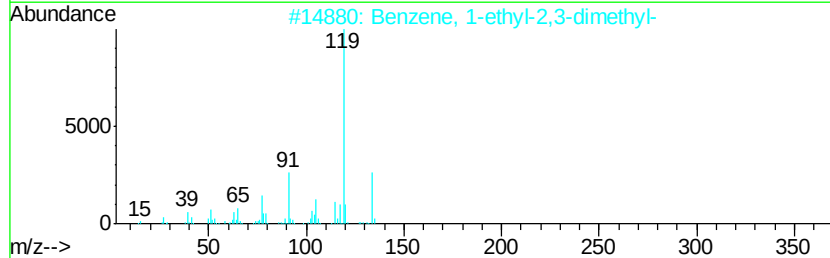
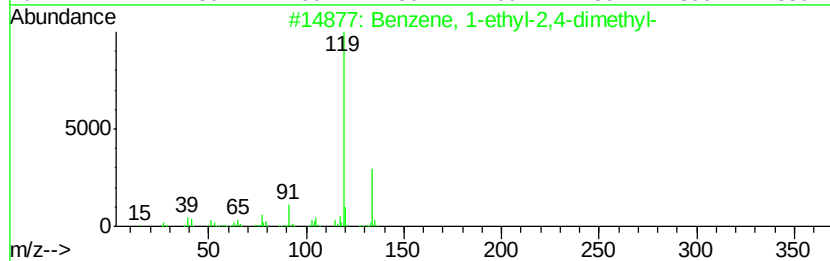
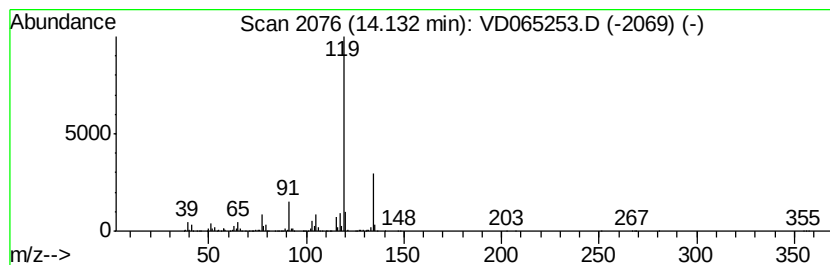
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	36.85 ug/l	2250530	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	97
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

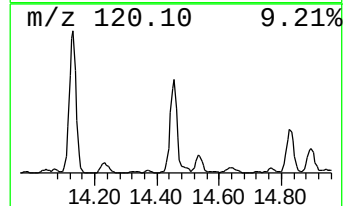
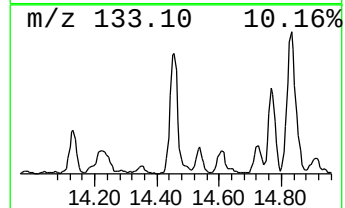
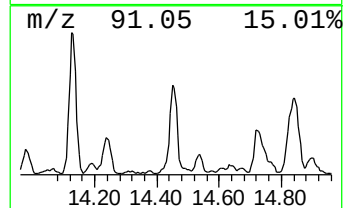
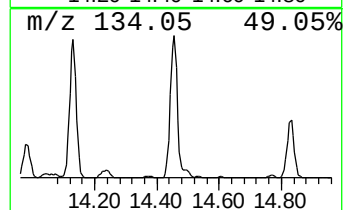
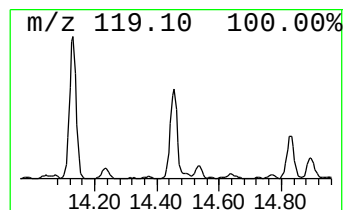
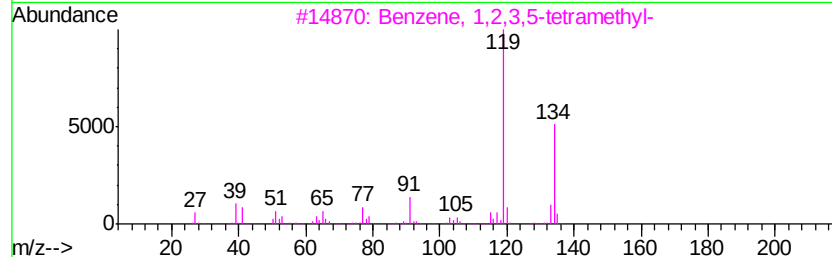
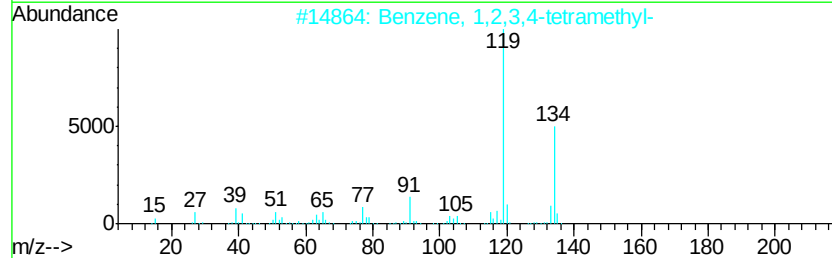
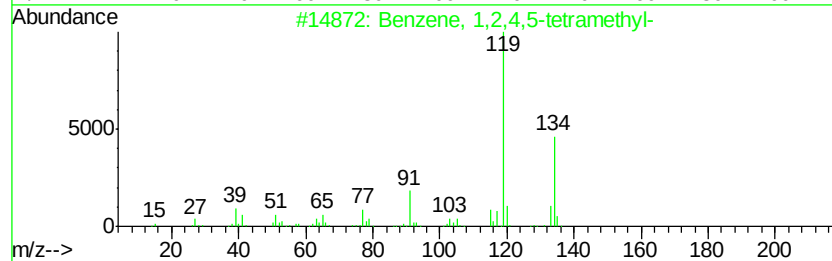
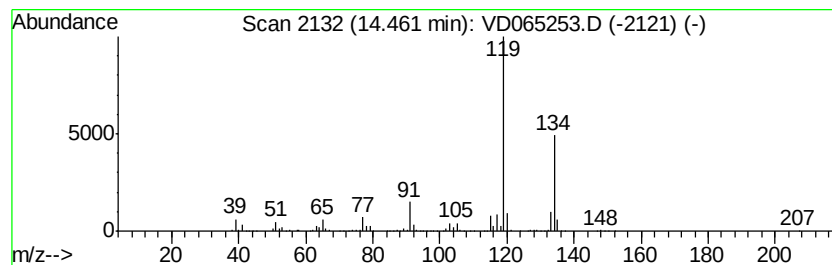
Instrument :
MSVOA_D
ClientSampleId :
SW-4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	26.85 ug/l	1639650	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7	95
4	o-Cymene	134 C10H14	000527-84-4	94
5	Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

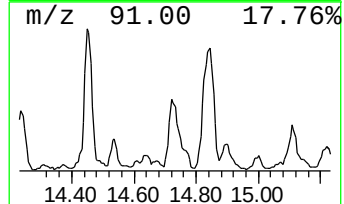
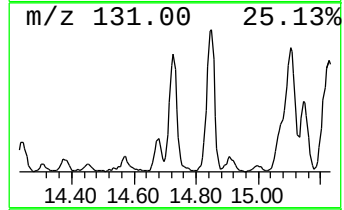
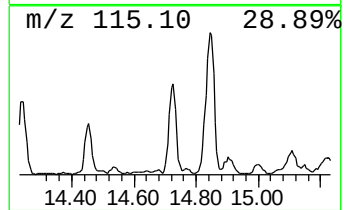
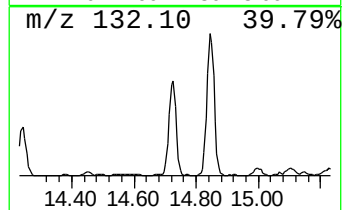
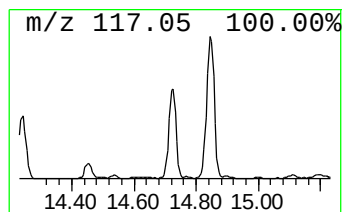
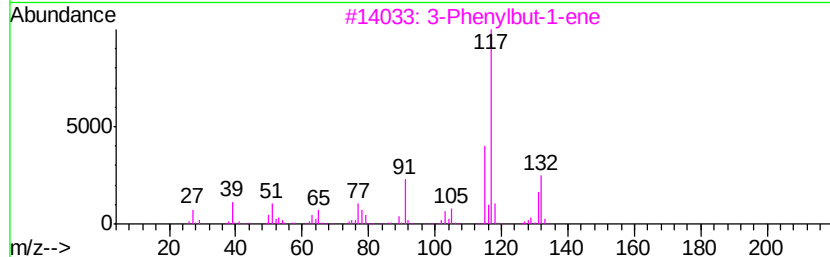
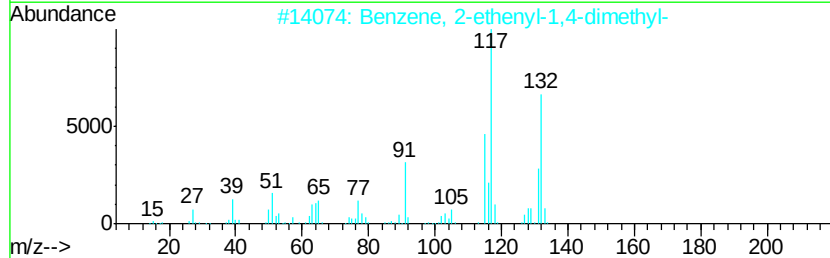
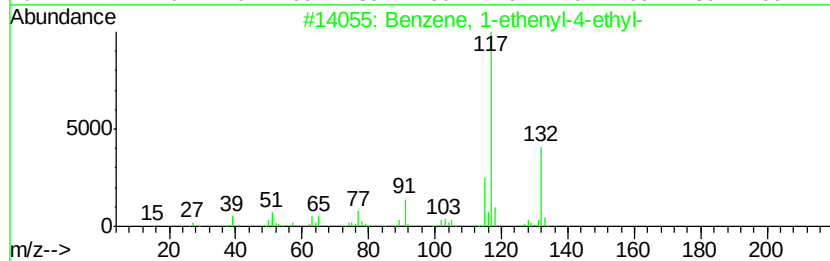
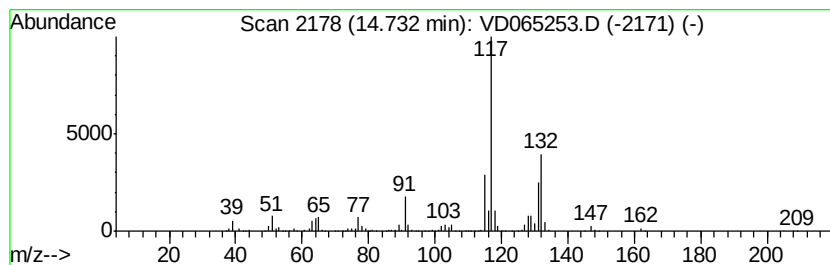
Instrument :
MSVOA_D
ClientSampled :
SW-4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	16.62 ug/l	1015150	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7 94
2		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 93
3		3-Phenylbut-1-ene	132 C10H12	000934-10-1 91
4		1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1 90
5		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA D/SOIL
ALS Vial : 8 Sample Multiplier: 1

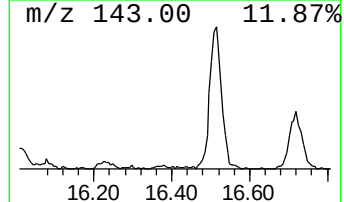
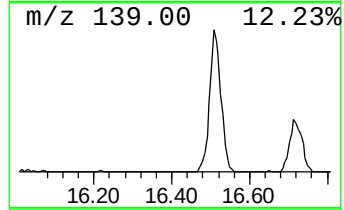
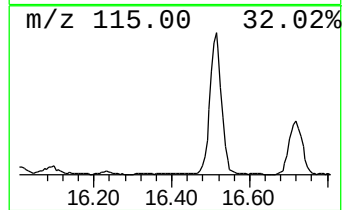
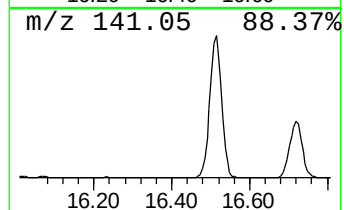
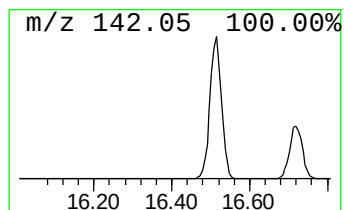
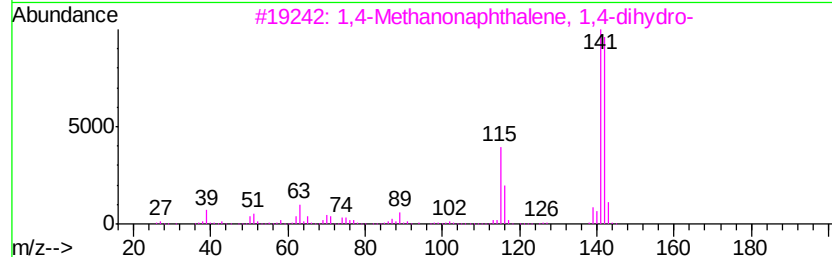
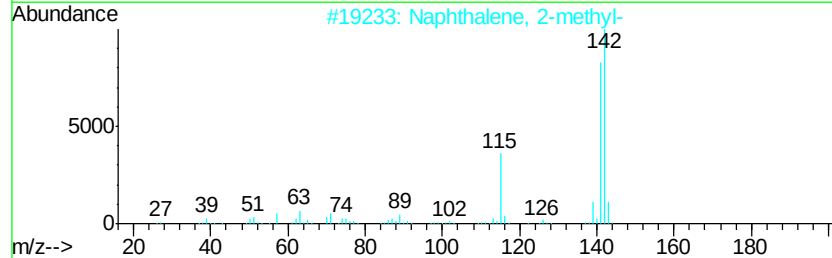
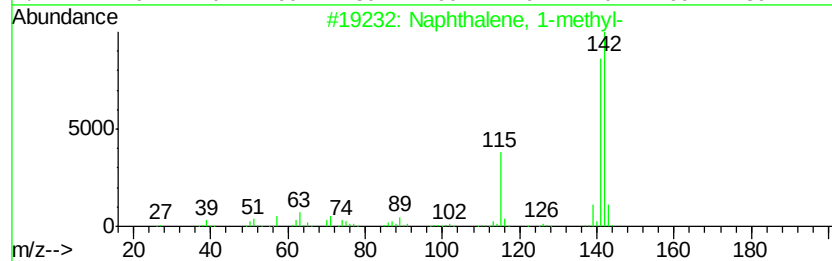
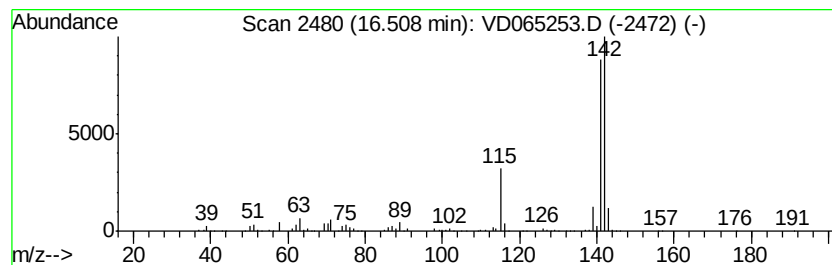
Instrument :
MSVOA_D
ClientSampled :
SW-4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.51	16.00 ug/l	977254	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
2		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
3		1,4-Methanonaphthalene, 1,4-dihv...	142 C11H10	004453-90-1 91
4		Benzocycloheptatriene	142 C11H10	000264-09-5 87
5		Naphthalene, 1-[(2-propenyloxy)m...	198 C14H14O	074685-39-5 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_D\DATA\VD022020\
Data File : VD065253.D
Acq On : 20 Feb 2020 13:35
Operator : VA/SY
Sample : L1601-05
Misc : 5.68G/5.00ml/MSVOA_D/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SW-4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D021820S.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
Hexane, 2,2-dimet...	8.51	26.6	ug/l	1510100	2	8.87	2839150	50.0
Pentane, 2,3,4-tr...	9.78	15.5	ug/l	877841	2	8.87	2839150	50.0
Pentane, 2,3,3-tr...	9.91	18.0	ug/l	1023970	2	8.87	2839150	50.0
Octane, 2-methyl-	11.43	16.2	ug/l	950721	3	11.65	2936320	50.0
Benzene, 1,1'-(1,...	13.78	32.2	ug/l	1964340	4	13.58	3053720	50.0
Benzene, 1-ethyl-...	14.13	36.9	ug/l	2250530	4	13.58	3053720	50.0
Benzene, 1,2,4,5-...	14.46	26.9	ug/l	1639650	4	13.58	3053720	50.0
Benzene, 1-etheny...	14.73	16.6	ug/l	1015150	4	13.58	3053720	50.0
Naphthalene, 1-me...	16.51	16.0	ug/l	977254	4	13.58	3053720	50.0