

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW101718\
 Data File : VW006140.D
 Acq On : 17 Oct 2018 21:11
 Operator : VA/AP
 Sample : J5515-20
 Misc : 5.76G/5ML/MSVOA W/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 AOC1-02A

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_W\METHOD\82W101518S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.958	478	501	530	rBV5	45261	269678	13.66%	0.694%
2	5.434	565	579	598	rBV	42362	153029	7.75%	0.394%
3	5.940	650	662	673	rVB2	22828	69026	3.50%	0.178%
4	6.873	803	815	824	rBV	47436	150385	7.62%	0.387%
5	6.982	824	833	843	rVV	67803	194816	9.87%	0.501%
6	7.159	853	862	876	rVB4	19156	67755	3.43%	0.174%
7	7.891	970	982	985	rBV	172856	422984	21.43%	1.089%
8	7.952	985	992	999	rVV2	416442	1093846	55.42%	2.815%
9	8.037	999	1006	1018	rVV2	152384	390434	19.78%	1.005%
10	8.159	1018	1026	1033	rVV	176105	392758	19.90%	1.011%
11	8.256	1033	1042	1044	rVV3	18742	57427	2.91%	0.148%
12	8.311	1044	1051	1062	rVB	155864	371306	18.81%	0.956%
13	8.482	1062	1079	1090	rBV	479122	1740334	88.17%	4.479%
14	8.573	1090	1094	1105	rVB2	73796	176752	8.95%	0.455%
15	8.695	1105	1114	1127	rVB2	58420	145621	7.38%	0.375%
16	8.848	1127	1139	1150	rBV	556538	1246268	63.14%	3.208%
17	9.073	1170	1176	1180	rBV	36907	77505	3.93%	0.199%
18	9.122	1180	1184	1201	rVB2	32772	95072	4.82%	0.245%
19	9.335	1205	1219	1224	rBV	430495	1104881	55.97%	2.844%
20	9.390	1224	1228	1235	rVB	162650	311023	15.76%	0.801%
21	9.470	1235	1241	1245	rBV2	41432	91138	4.62%	0.235%
22	9.518	1245	1249	1255	rVB	75519	139128	7.05%	0.358%
23	9.610	1255	1264	1271	rVB3	86177	214107	10.85%	0.551%
24	9.762	1282	1289	1300	rVB	493369	1038538	52.61%	2.673%
25	9.896	1300	1311	1318	rBV	609417	1405843	71.22%	3.618%
26	9.957	1318	1321	1325	rVV	138973	266990	13.53%	0.687%
27	9.994	1325	1327	1335	rVB3	122249	211678	10.72%	0.545%
28	10.098	1335	1344	1357	rBV2	245499	661634	33.52%	1.703%
29	10.250	1364	1369	1375	rVB2	220504	404152	20.47%	1.040%
30	10.329	1375	1382	1394	rBV	779619	1591464	80.63%	4.096%
31	10.445	1394	1401	1404	rBV4	52734	99058	5.02%	0.255%
32	10.506	1404	1411	1418	rVB7	130047	371533	18.82%	0.956%
33	10.671	1433	1438	1445	rBV	110180	218994	11.09%	0.564%
34	10.756	1445	1452	1462	rVB4	183435	511495	25.91%	1.317%

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35	10.847	1463	1467	1473	rVB3	79697	145447	7.37%	0.374%
36	10.939	1473	1482	1488	rBV3	83214	188681	9.56%	0.486%
37	11.043	1488	1499	1506	rBV4	114854	366638	18.57%	0.944%
38	11.110	1506	1510	1513	rBV4	38363	58708	2.97%	0.151%
39	11.146	1513	1516	1519	rVV3	49981	78667	3.99%	0.202%
40	11.183	1519	1522	1526	rVV3	78847	124966	6.33%	0.322%
41	11.231	1526	1530	1535	rVV	100675	169421	8.58%	0.436%
42	11.292	1535	1540	1544	rVB2	49277	77635	3.93%	0.200%
43	11.341	1544	1548	1552	rBV2	65766	90625	4.59%	0.233%
44	11.414	1552	1560	1572	rVB4	176142	590894	29.94%	1.521%
45	11.518	1572	1577	1582	rBV	162035	282869	14.33%	0.728%
46	11.634	1589	1596	1603	rBV	609562	1047347	53.06%	2.696%
47	11.731	1609	1612	1616	rVV4	61884	99821	5.06%	0.257%
48	11.847	1621	1631	1633	rBV6	56611	159220	8.07%	0.410%
49	11.878	1633	1636	1641	rVB3	71472	111461	5.65%	0.287%
50	12.073	1664	1668	1673	rVB4	37535	67587	3.42%	0.174%
51	12.140	1673	1679	1686	rBV4	99537	244304	12.38%	0.629%
52	12.256	1694	1698	1702	rVB	74261	121429	6.15%	0.313%
53	12.341	1702	1712	1715	rBV5	81464	211151	10.70%	0.543%
54	12.469	1728	1733	1738	rBV	119047	219346	11.11%	0.565%
55	12.621	1751	1758	1767	rVB3	465231	934173	47.33%	2.404%
56	12.707	1767	1772	1777	rVB6	38781	69190	3.51%	0.178%
57	12.810	1777	1789	1794	rVB	268126	513126	26.00%	1.321%
58	12.871	1794	1799	1802	rBV5	52154	88862	4.50%	0.229%
59	12.951	1807	1812	1818	rVB6	69467	139534	7.07%	0.359%
60	13.109	1830	1838	1845	rBV2	110044	275301	13.95%	0.709%
61	13.170	1845	1848	1856	rVV5	65999	137285	6.95%	0.353%
62	13.256	1858	1862	1866	rVB	92295	135525	6.87%	0.349%
63	13.304	1866	1870	1874	rVB2	47547	66988	3.39%	0.172%
64	13.384	1875	1883	1891	rBV4	103490	328321	16.63%	0.845%
65	13.566	1908	1913	1922	rBV2	417129	725497	36.75%	1.867%
66	13.725	1933	1939	1942	rBV	509163	792742	40.16%	2.040%
67	13.768	1942	1946	1950	rVV	782822	1262323	63.95%	3.249%
68	13.810	1950	1953	1960	rVB2	182747	357207	18.10%	0.919%
69	13.890	1960	1966	1970	rBV2	151312	255812	12.96%	0.658%
70	14.036	1985	1990	1995	rBV2	98131	170653	8.65%	0.439%
71	14.103	1995	2001	2007	rBV	1018295	1519044	76.96%	3.910%

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72	14.164	2007	2011	2014	rVV	97421	142107	7.20%	0.366%
73	14.213	2014	2019	2025	rVB2	595456	1013717	51.36%	2.609%
74	14.280	2025	2030	2036	rBV4	61488	143574	7.27%	0.370%
75	14.353	2038	2042	2045	rBV3	42128	61923	3.14%	0.159%
76	14.396	2045	2049	2050	rVV2	96028	121882	6.17%	0.314%
77	14.426	2050	2054	2059	rVB	292561	461385	23.37%	1.188%
78	14.505	2063	2067	2071	rBV	179977	269118	13.63%	0.693%
79	14.554	2071	2075	2081	rVV3	45054	101297	5.13%	0.261%
80	14.694	2094	2098	2103	rBV2	313951	507867	25.73%	1.307%
81	14.743	2103	2106	2110	rVB	122604	168211	8.52%	0.433%
82	14.816	2110	2118	2123	rBV2	1053724	1973906	100.00%	5.081%
83	14.865	2123	2126	2133	rVB3	189503	355843	18.03%	0.916%
84	15.072	2149	2160	2164	rBV2	275108	644778	32.67%	1.660%
85	15.115	2164	2167	2172	rVV	163976	280139	14.19%	0.721%
86	15.188	2172	2179	2186	rVB2	364708	813929	41.23%	2.095%
87	15.341	2198	2204	2207	rBV4	138420	245180	12.42%	0.631%
88	15.432	2214	2219	2222	rBV2	60951	114924	5.82%	0.296%
89	15.481	2223	2227	2231	rVV2	113091	216912	10.99%	0.558%
90	15.536	2232	2236	2250	rVB	154318	411638	20.85%	1.060%
91	15.670	2251	2258	2264	rBV	223055	456033	23.10%	1.174%
92	15.743	2266	2270	2275	rBV4	51470	83485	4.23%	0.215%
93	15.834	2278	2285	2289	rBV2	132786	308394	15.62%	0.794%
94	15.962	2302	2306	2310	rVB3	58830	92189	4.67%	0.237%
95	16.035	2311	2318	2324	rBV4	113847	284044	14.39%	0.731%
96	16.188	2335	2343	2348	rBV2	115737	290678	14.73%	0.748%
97	16.304	2358	2362	2366	rVB	56385	82221	4.17%	0.212%
98	16.389	2372	2376	2381	rVB2	87172	169983	8.61%	0.438%
99	16.450	2382	2386	2392	rBV2	91860	172896	8.76%	0.445%
100	16.657	2412	2420	2433	rVB	345711	879235	44.54%	2.263%

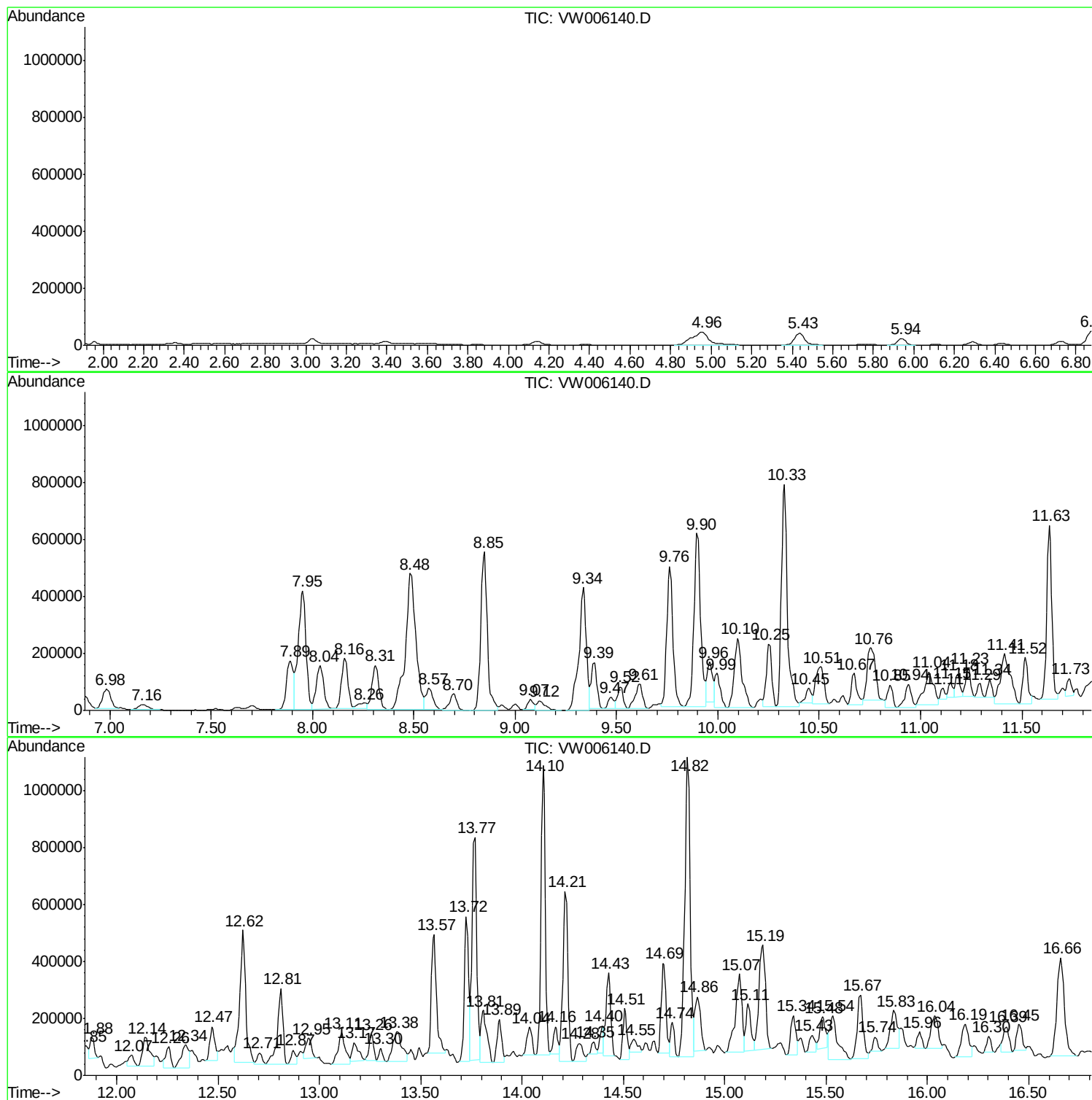
Sum of corrected areas: 38851940

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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.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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AOC1-02A

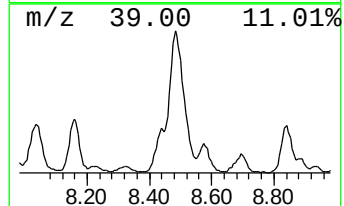
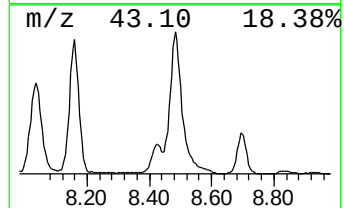
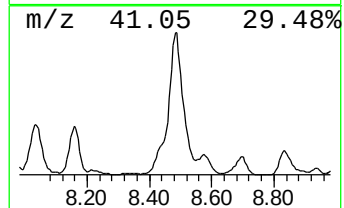
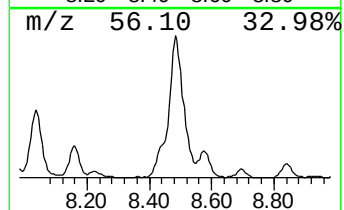
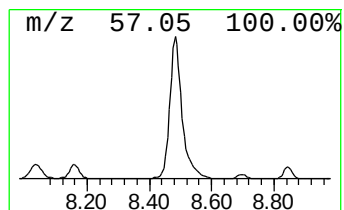
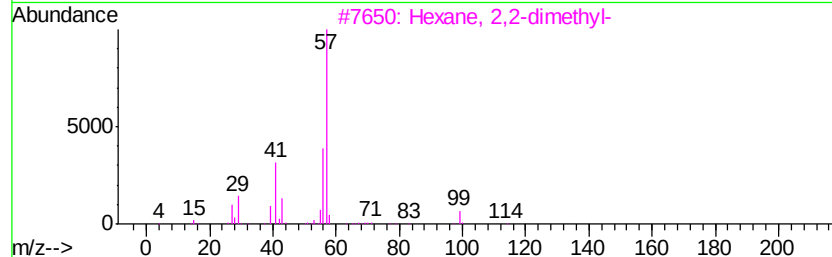
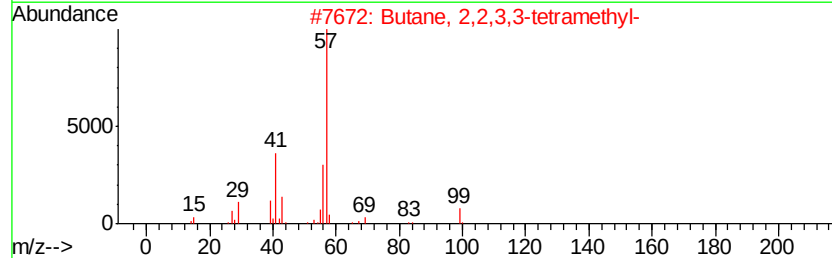
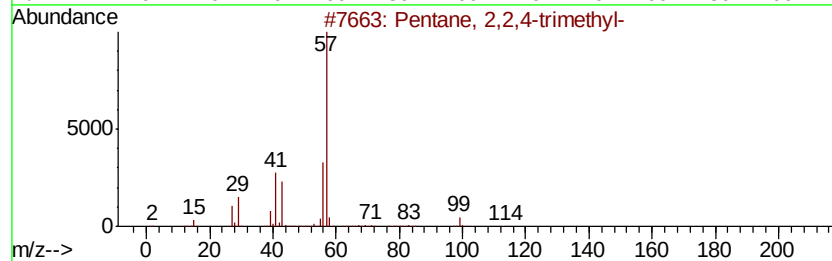
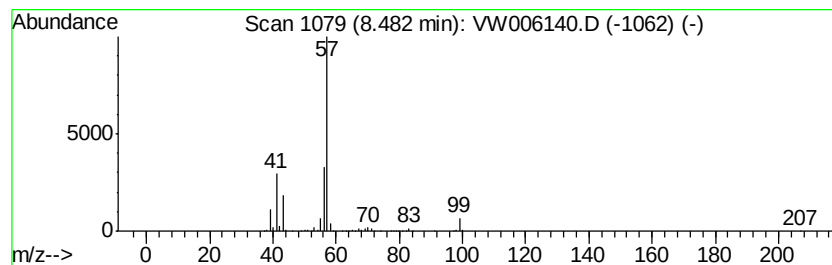
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 2,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.48	69.82 ug/l	1740330	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
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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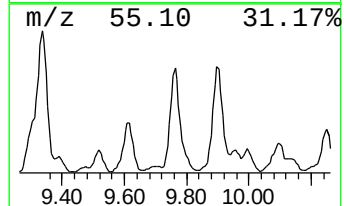
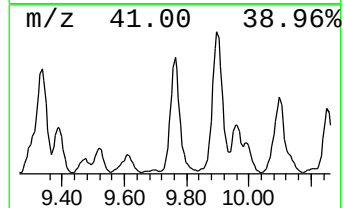
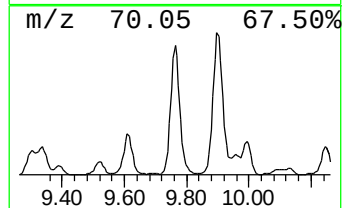
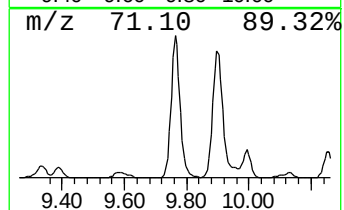
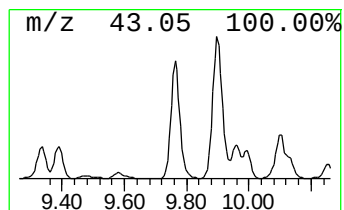
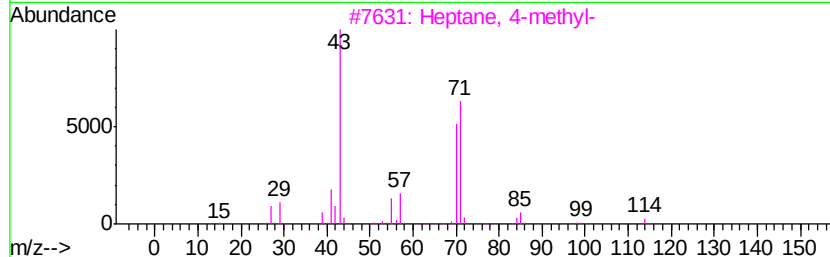
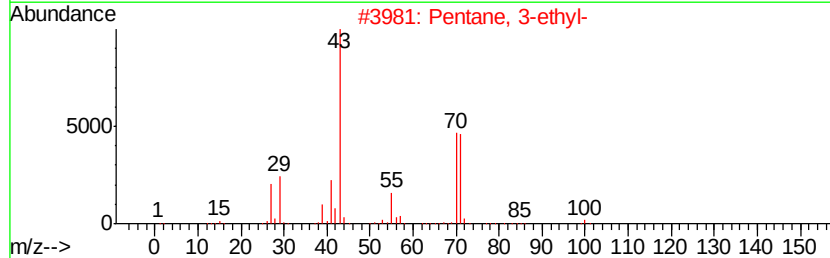
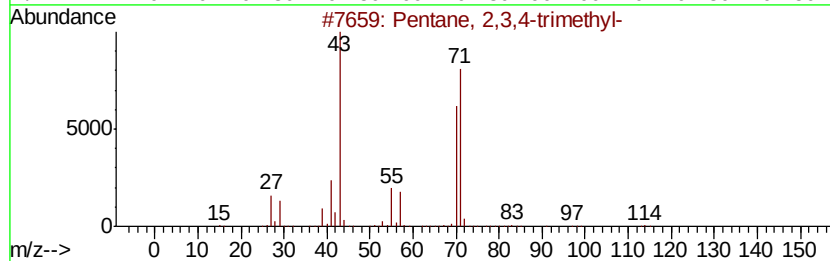
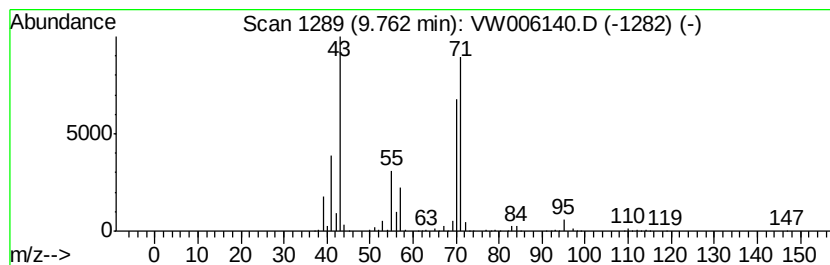
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	41.67 ug/l	1038540	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	86
2		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	72
4		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	64
5		Heptane, 5-ethyl-2-methyl-	142	C10H22	013475-78-0	59



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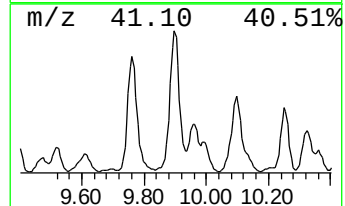
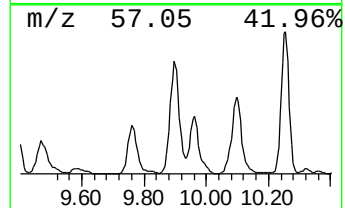
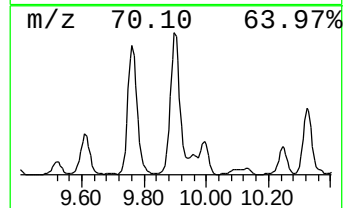
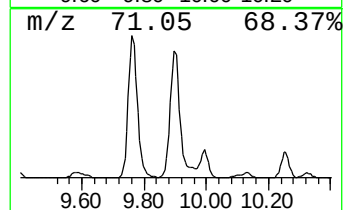
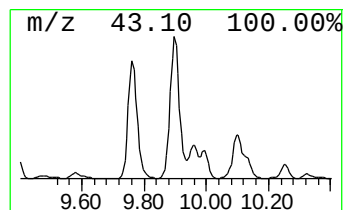
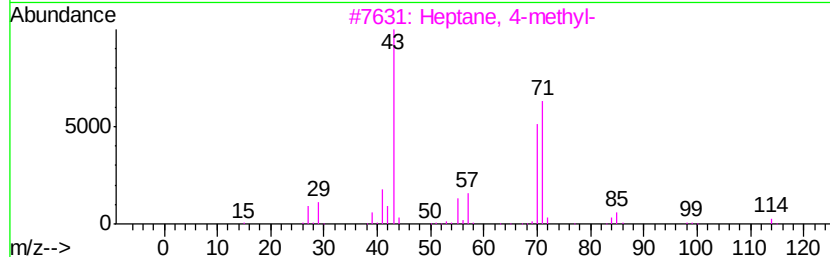
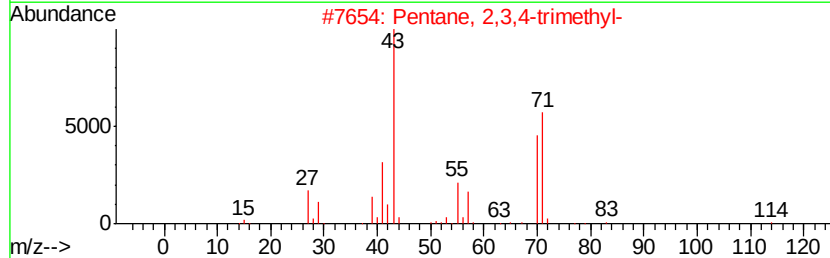
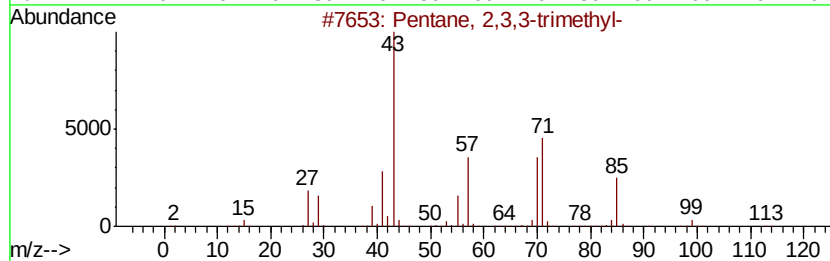
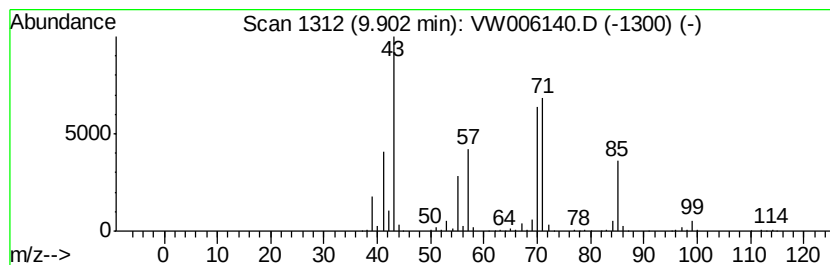
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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.
9.90	56.40 ug/l	1405840	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	81
3		Heptane, 4-methyl-	114	C8H18	000589-53-7	72
4		Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5		Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
Misc : 5.76G/5ML/MSVOA W/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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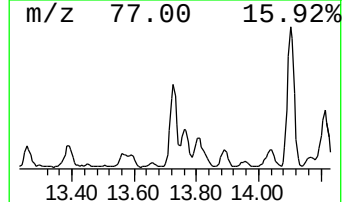
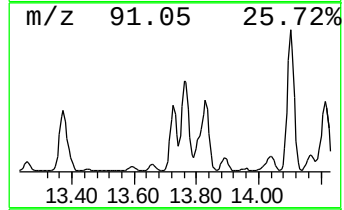
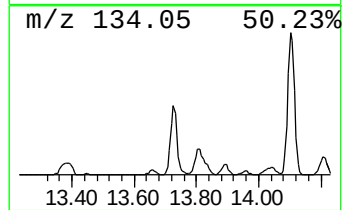
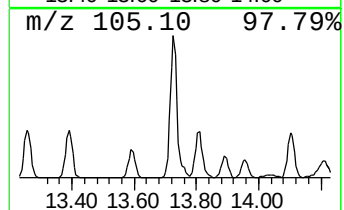
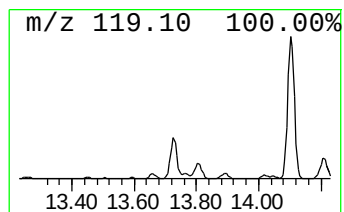
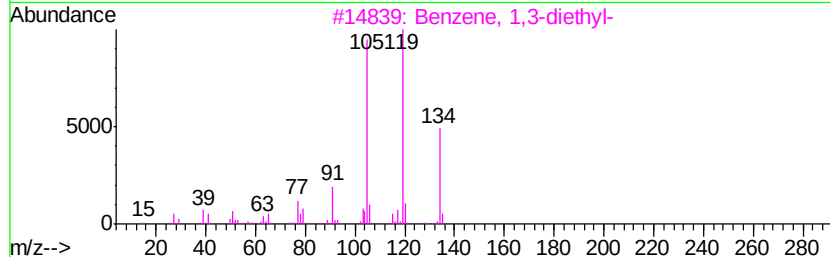
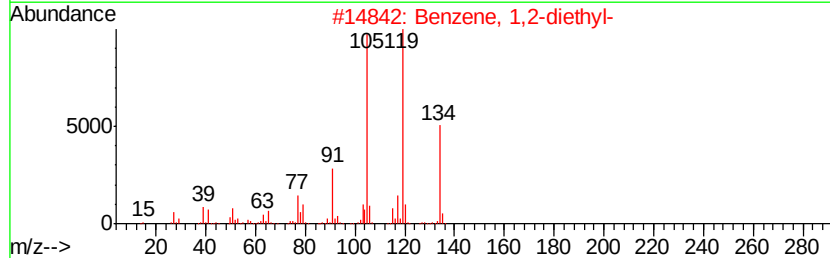
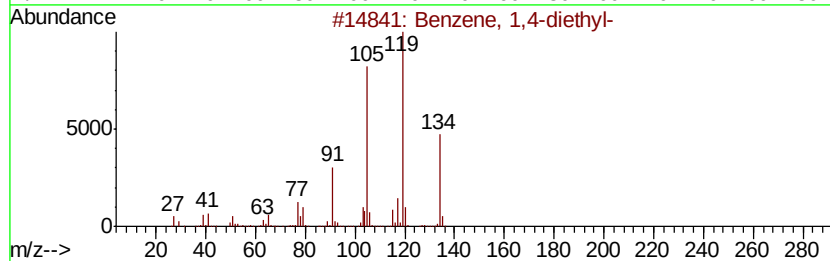
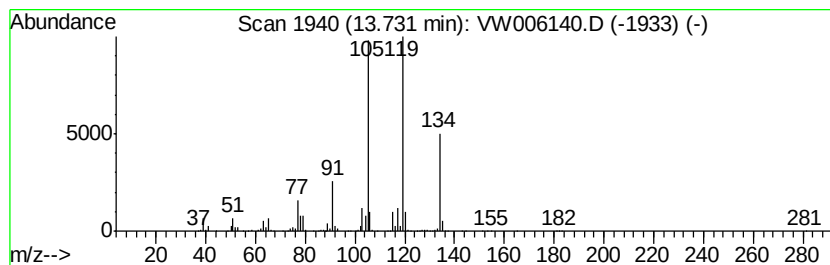
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	54.63 ug/l	792742	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
Misc : 5.76G/5ML/MSVOA W/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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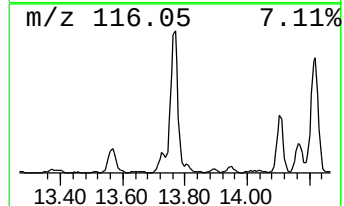
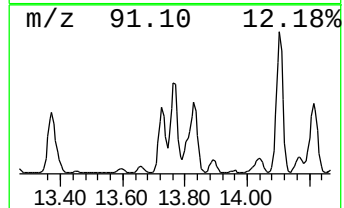
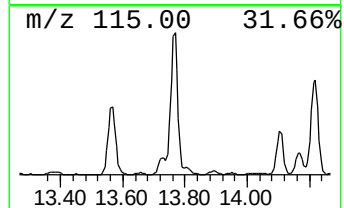
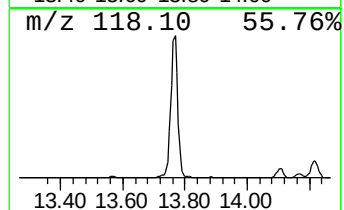
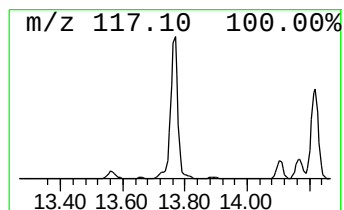
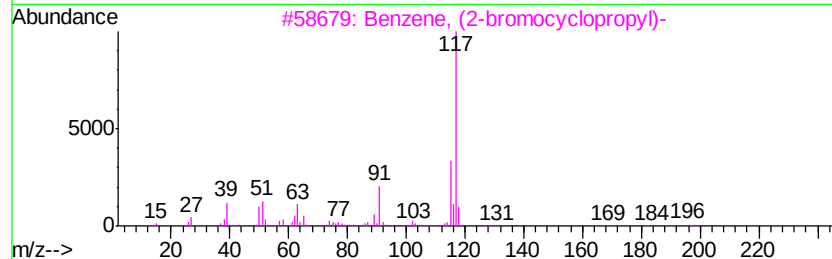
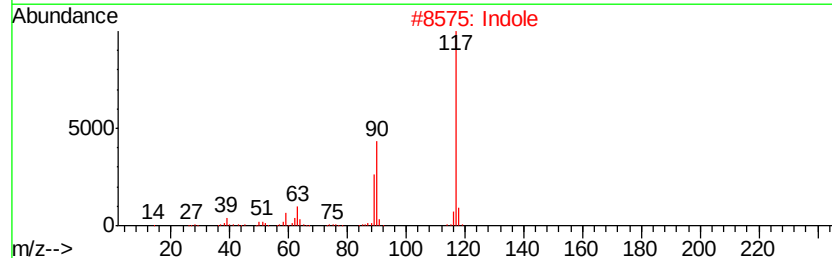
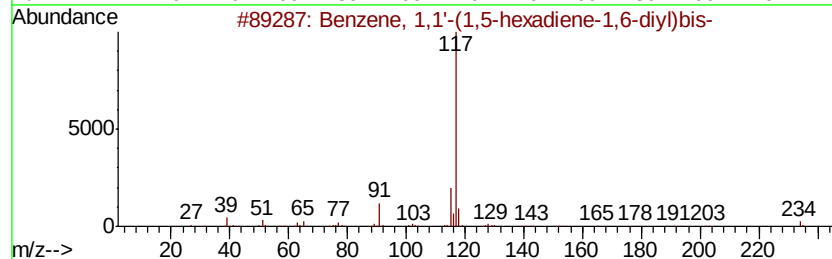
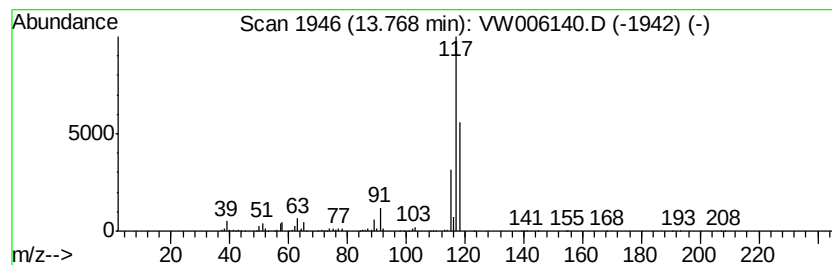
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	87.00 ug/l	1262320	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	53
2		Indole	117	C8H7N	000120-72-9	46
3		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
4		m-Aminophenylacetylene	117	C8H7N	054060-30-9	43
5		Benzyl nitrile	117	C8H7N	000140-29-4	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
Misc : 5.76G/5ML/MSVOA W/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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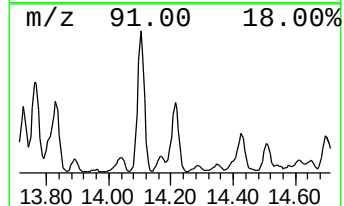
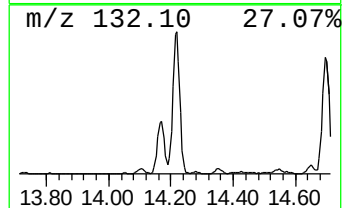
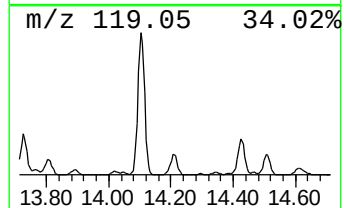
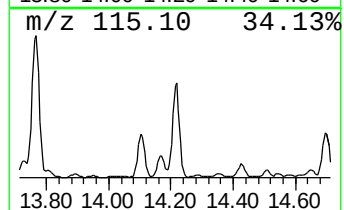
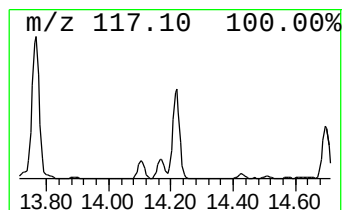
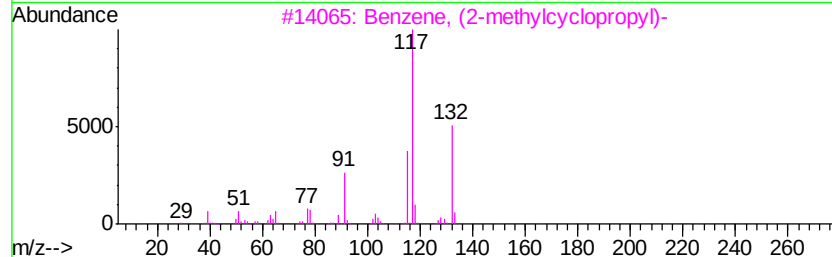
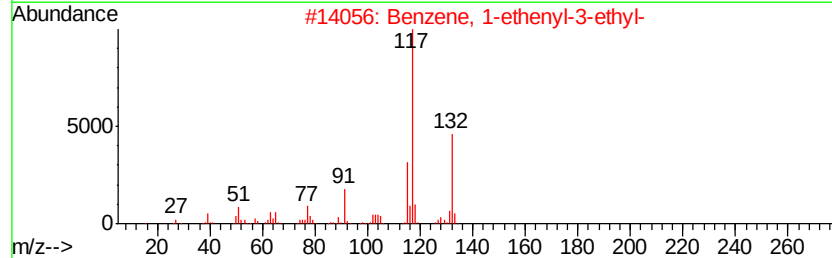
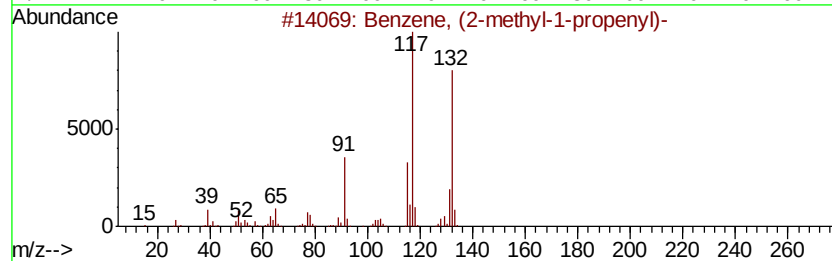
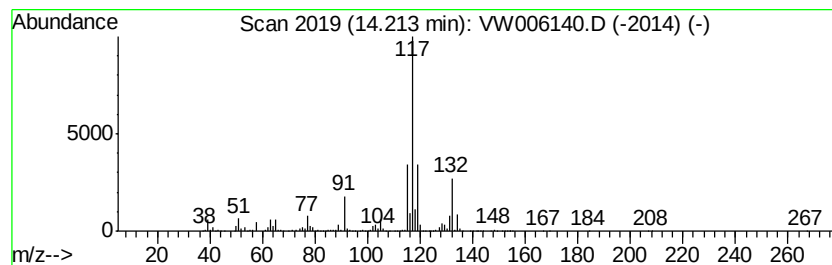
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (2-methyl-1-propen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	69.86 ug/l	1013720	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	76
4		Indan, 1-methyl-	132	C10H12	000767-58-8	76
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
Misc : 5.76G/5ML/MSVOA W/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
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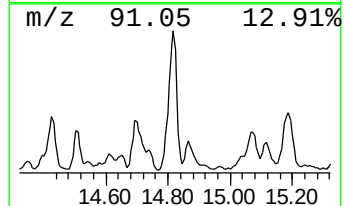
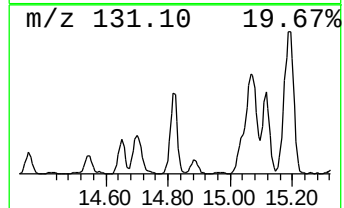
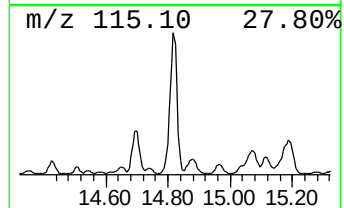
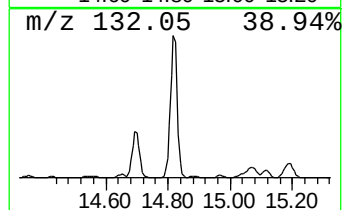
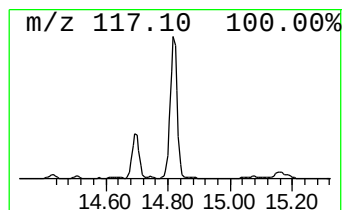
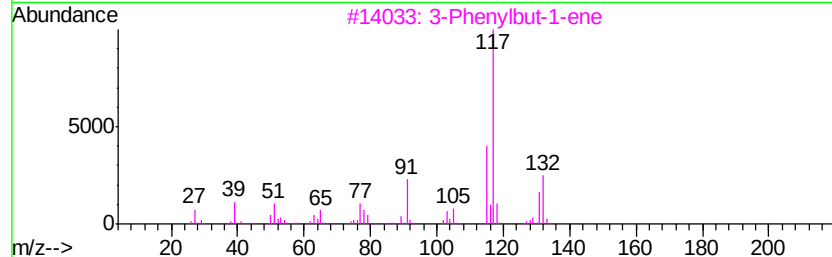
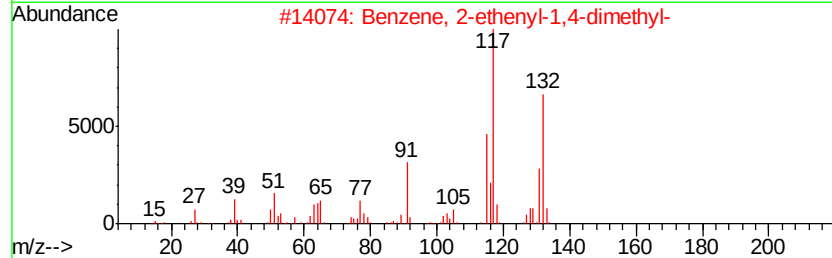
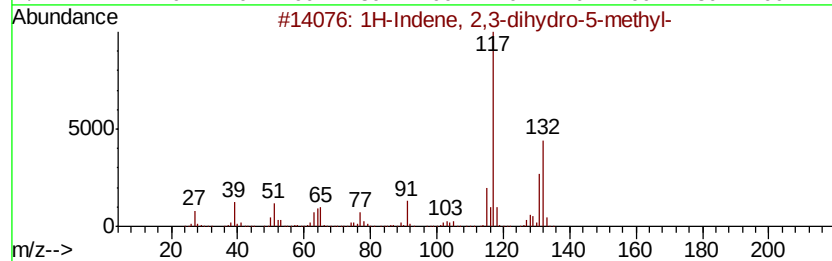
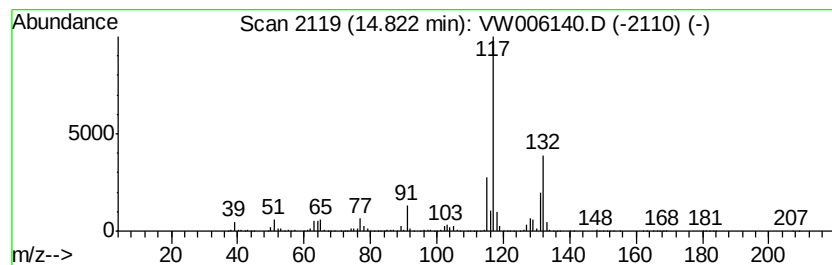
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	136.04 ug/l	1973910	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93	
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92	
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91	
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91	
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
Misc : 5.76G/5ML/MSVOA W/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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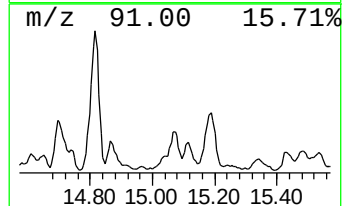
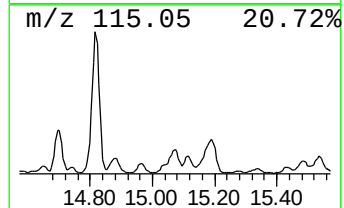
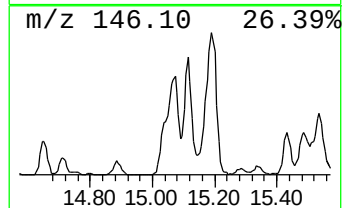
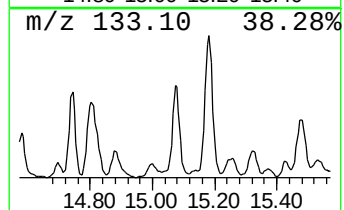
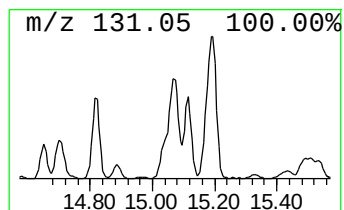
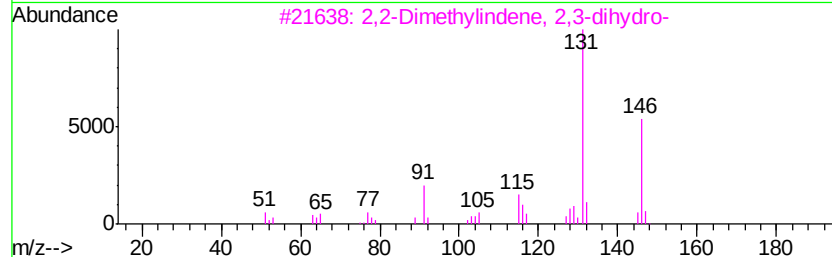
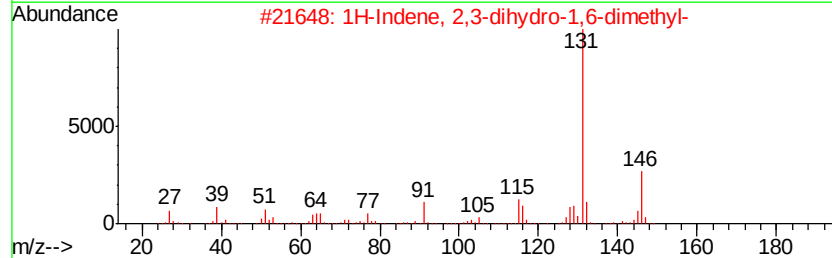
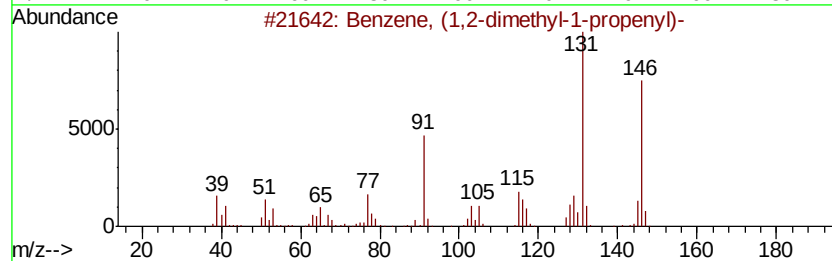
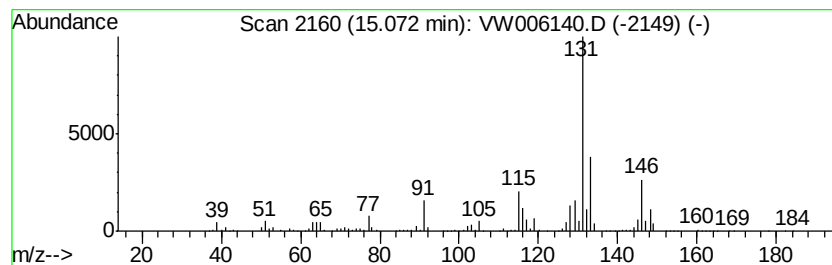
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	44.44 ug/l	644778	1,4-Dichlorobenzene-d4	13.57

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
2		1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	92
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	76
4		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	76
5		1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
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Sample : J5515-20
Misc : 5.76G/5ML/MSVOA W/SOIL
ALS Vial : 22 Sample Multiplier: 1

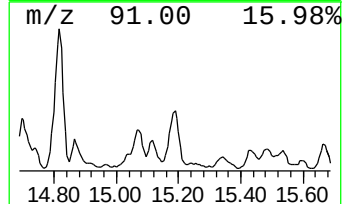
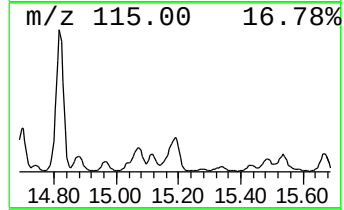
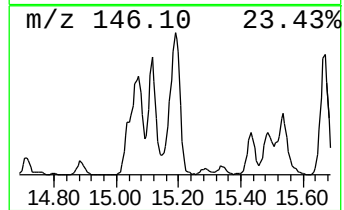
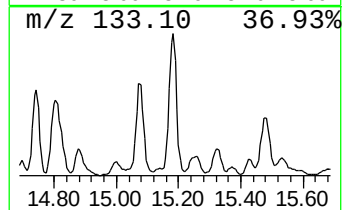
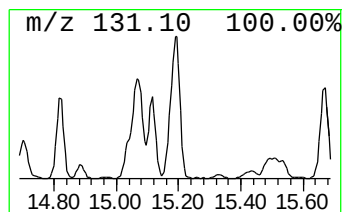
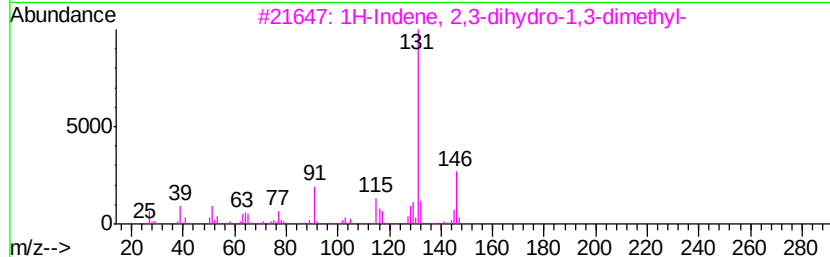
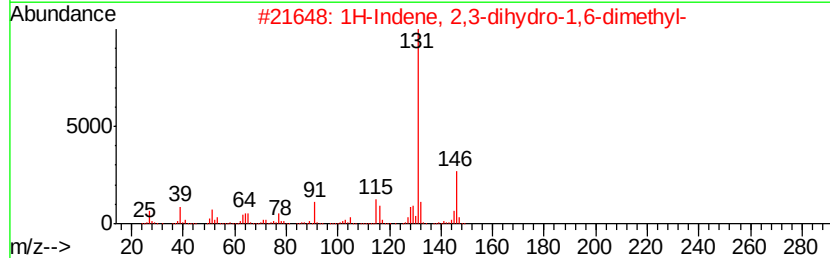
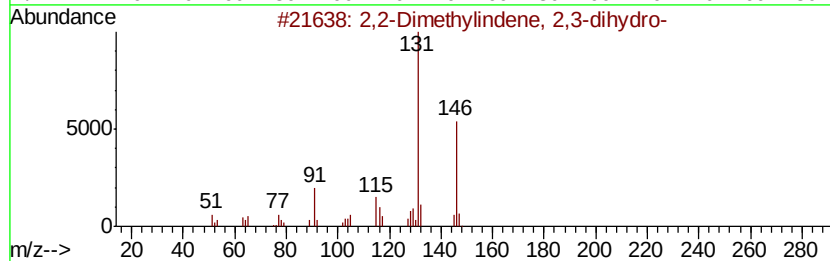
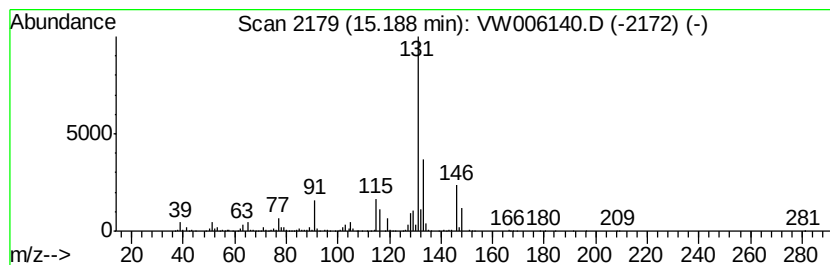
Instrument :
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ClientSampleId :
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Quant Title : SW846 8260

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

Peak Number 9 2,2-Dimethylindene, 2,3-dih... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	56.09 ug/l	813929	1,4-Dichlorobenzene-d4	13.57
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7	81
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	81
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	81
4	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	74
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
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ALS Vial : 22 Sample Multiplier: 1

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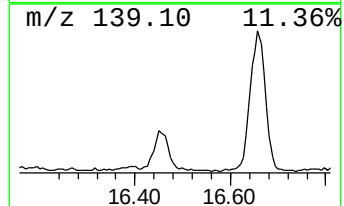
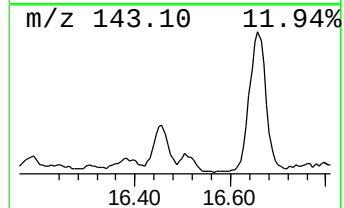
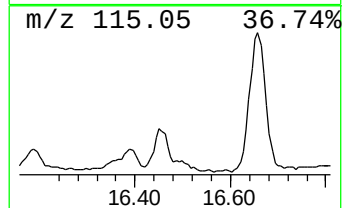
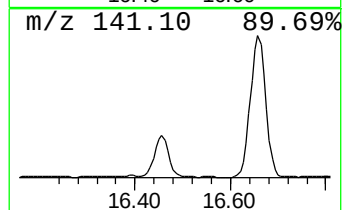
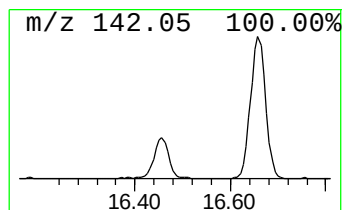
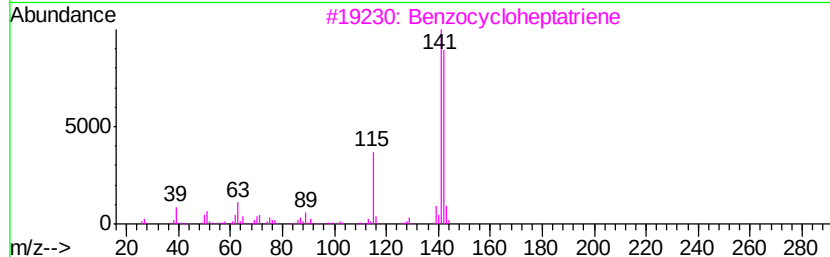
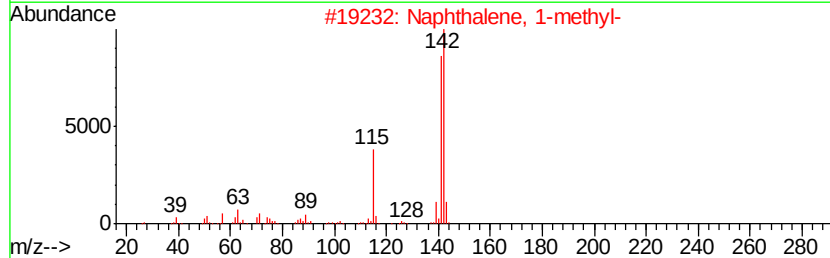
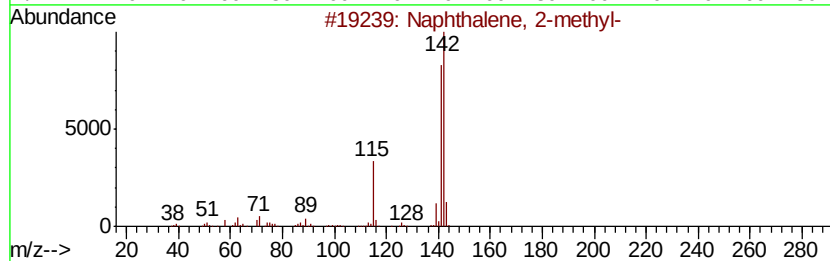
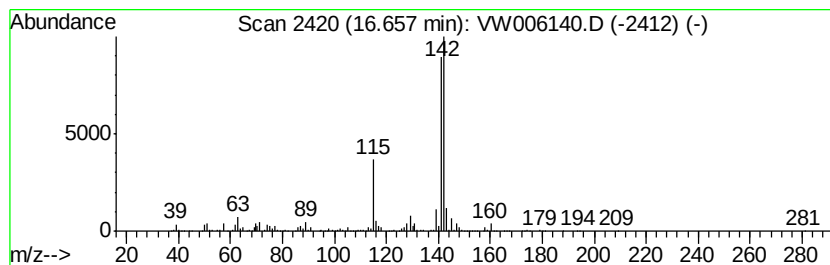
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.66	60.60 ug/l	879235	1,4-Dichlorobenzene-d4	13.57

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	90
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5		1-Naphthaleneacetamide	185	C12H11NO	000086-86-2	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW101718\
Data File : VW006140.D
Acq On : 17 Oct 2018 21:11
Operator : VA/AP
Sample : J5515-20
Misc : 5.76G/5ML/MSVOA_W/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
AOC1-02A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.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
Pentane, 2,2,4-tr...	8.48	69.8	ug/l	1740330	2	8.85	1246270	50.0
Pentane, 2,3,4-tr...	9.76	41.7	ug/l	1038540	2	8.85	1246270	50.0
Pentane, 2,3,3-tr...	9.90	56.4	ug/l	1405840	2	8.85	1246270	50.0
Benzene, 1,4-diet...	13.73	54.6	ug/l	792742	4	13.57	725497	50.0
Benzene, 1,1'-(1,...	13.77	87.0	ug/l	1262320	4	13.57	725497	50.0
Benzene, (2-methy...	14.21	69.9	ug/l	1013720	4	13.57	725497	50.0
1H-Indene, 2,3-di...	14.82	136.0	ug/l	1973910	4	13.57	725497	50.0
Benzene, (1,2-dim...	15.07	44.4	ug/l	644778	4	13.57	725497	50.0
2,2-Dimethylinden...	15.19	56.1	ug/l	813929	4	13.57	725497	50.0
Naphthalene, 2-me...	16.66	60.6	ug/l	879235	4	13.57	725497	50.0