

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW010919\
 Data File : VW008140.D
 Acq On : 10 Jan 2019 09:25
 Operator : SY/AP
 Sample : K1006-11
 Misc : 6.78G/5ML/MSVOA W/SOIL
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 JC-03-010819-F

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\82W010819S.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	2.367	68	76	84	rBV2	5316	11290	3.20%	0.199%
2	2.818	143	150	157	rBV	6573	15726	4.45%	0.277%
3	4.141	358	367	376	rBV	3600	11424	3.23%	0.201%
4	4.916	491	494	508	rVB8	2145	7011	1.98%	0.124%
5	5.434	571	579	585	rBV6	1559	4667	1.32%	0.082%
6	7.159	852	862	864	rBV5	2055	5484	1.55%	0.097%
7	7.885	973	981	986	rBV	47331	111914	31.68%	1.972%
8	7.952	986	992	1000	rVB	78987	174952	49.52%	3.083%
9	8.305	1041	1050	1061	rVB	51654	115153	32.59%	2.029%
10	8.842	1130	1138	1152	rBV	128231	257053	72.76%	4.530%
11	9.342	1212	1220	1224	rBV6	2877	6047	1.71%	0.107%
12	10.323	1373	1381	1388	rBV	205301	353289	100.00%	6.226%
13	10.384	1388	1391	1399	rVB	9801	18122	5.13%	0.319%
14	10.658	1432	1436	1441	rBV6	1997	3536	1.00%	0.062%
15	10.762	1448	1453	1459	rVB6	1946	3796	1.07%	0.067%
16	11.177	1517	1521	1525	rVV3	3614	6147	1.74%	0.108%
17	11.232	1526	1530	1536	rVB3	3156	5428	1.54%	0.096%
18	11.433	1561	1563	1571	rVB6	3412	5893	1.67%	0.104%
19	11.512	1571	1576	1584	rBV4	2753	5098	1.44%	0.090%
20	11.634	1589	1596	1605	rBV	157488	267728	75.78%	4.718%
21	11.738	1608	1613	1616	rBV5	2726	4822	1.36%	0.085%
22	11.847	1622	1631	1634	rBV5	6981	16376	4.64%	0.289%
23	12.170	1673	1684	1691	rBV3	10849	28156	7.97%	0.496%
24	12.250	1693	1697	1702	rVB5	2069	3775	1.07%	0.067%
25	12.347	1707	1713	1714	rBV5	2703	4071	1.15%	0.072%
26	12.621	1752	1758	1767	rVB	107166	174542	49.40%	3.076%
27	12.707	1767	1772	1776	rBV7	1904	4021	1.14%	0.071%
28	12.792	1781	1786	1791	rVB5	4794	10815	3.06%	0.191%
29	12.871	1791	1799	1801	rBV3	11169	20399	5.77%	0.359%
30	12.938	1801	1810	1816	rVB3	30159	74757	21.16%	1.317%
31	12.993	1816	1819	1824	rVB4	5625	7996	2.26%	0.141%
32	13.048	1824	1828	1830	rBV3	3270	5225	1.48%	0.092%
33	13.109	1830	1838	1844	rBV	13795	32522	9.21%	0.573%
34	13.170	1844	1848	1854	rVB4	12270	21533	6.10%	0.379%

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35	13.256	1854	1862	1866	rBV	18387	30829	8.73%	0.543%
36	13.347	1875	1877	1880	rBV4	2978	3583	1.01%	0.063%
37	13.384	1880	1883	1887	rVB3	7196	9302	2.63%	0.164%
38	13.451	1887	1894	1897	rBV5	23077	44746	12.67%	0.789%
39	13.493	1897	1901	1903	rBV3	10348	12687	3.59%	0.224%
40	13.560	1907	1912	1921	rVB2	118789	227642	64.44%	4.012%
41	13.634	1921	1924	1928	rVB5	11408	13555	3.84%	0.239%
42	13.701	1929	1935	1936	rBV5	5939	10442	2.96%	0.184%
43	13.725	1936	1939	1940	rBV3	8506	9068	2.57%	0.160%
44	13.755	1940	1944	1947	rVV2	34872	48579	13.75%	0.856%
45	13.798	1947	1951	1959	rVB4	45998	92651	26.23%	1.633%
46	13.877	1959	1964	1970	rBV	105654	168081	47.58%	2.962%
47	13.926	1970	1972	1973	rBV	4485	4130	1.17%	0.073%
48	13.957	1973	1977	1984	rVB	19436	30606	8.66%	0.539%
49	14.018	1984	1987	1988	rBV	12567	13708	3.88%	0.242%
50	14.042	1988	1991	1995	rVB3	21047	31473	8.91%	0.555%
51	14.103	1996	2001	2005	rVB2	45366	67144	19.01%	1.183%
52	14.140	2005	2007	2013	rVB3	10159	17323	4.90%	0.305%
53	14.207	2013	2018	2024	rVB3	60052	99176	28.07%	1.748%
54	14.268	2024	2028	2030	rBV3	16863	26895	7.61%	0.474%
55	14.347	2035	2041	2044	rBV4	32033	65363	18.50%	1.152%
56	14.389	2044	2048	2052	rVV3	57925	103870	29.40%	1.830%
57	14.432	2052	2055	2058	rVV3	32407	51792	14.66%	0.913%
58	14.463	2058	2060	2063	rVB	30535	33138	9.38%	0.584%
59	14.505	2063	2067	2071	rBV2	48448	64876	18.36%	1.143%
60	14.548	2071	2074	2080	rVB4	35603	54975	15.56%	0.969%
61	14.615	2081	2085	2087	rBV3	18980	29890	8.46%	0.527%
62	14.646	2087	2090	2094	rVB	28839	41537	11.76%	0.732%
63	14.694	2094	2098	2100	rBV	38279	59216	16.76%	1.044%
64	14.731	2100	2104	2110	rVB2	136406	213957	60.56%	3.771%
65	14.810	2110	2117	2122	rBV3	120739	304565	86.21%	5.367%
66	14.853	2122	2124	2132	rVB5	56758	107028	30.29%	1.886%
67	14.963	2138	2142	2148	rVB	68114	117744	33.33%	2.075%
68	15.072	2155	2160	2164	rBV3	50057	83010	23.50%	1.463%
69	15.188	2171	2179	2187	rBV3	60234	158552	44.88%	2.794%
70	15.261	2187	2191	2198	rVB7	32091	69308	19.62%	1.221%
71	15.334	2198	2203	2213	rBV5	56234	124608	35.27%	2.196%

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Integration Parameters: RTEINT.P

Integrator: RTE
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72	15.432	2214	2219	2223	rBV	45667	73993	20.94%	1.304%
73	15.481	2223	2227	2230	rBV4	36869	65276	18.48%	1.150%
74	15.566	2236	2241	2250	rVB	121383	233060	65.97%	4.107%
75	15.670	2250	2258	2263	rBV5	53077	136490	38.63%	2.405%
76	15.743	2267	2270	2275	rVV6	25099	42573	12.05%	0.750%
77	15.834	2279	2285	2287	rVV3	34717	75926	21.49%	1.338%
78	15.877	2287	2292	2301	rVB	71374	145945	41.31%	2.572%
79	16.035	2312	2318	2325	rVB5	26668	65089	18.42%	1.147%
80	16.188	2334	2343	2348	rBV5	51358	134807	38.16%	2.376%
81	16.298	2358	2361	2367	rVV	15943	24654	6.98%	0.434%
82	16.383	2370	2375	2381	rVV3	54627	108960	30.84%	1.920%
83	16.450	2382	2386	2390	rVV5	19566	37655	10.66%	0.664%
84	16.511	2391	2396	2401	rVB	49698	86182	24.39%	1.519%

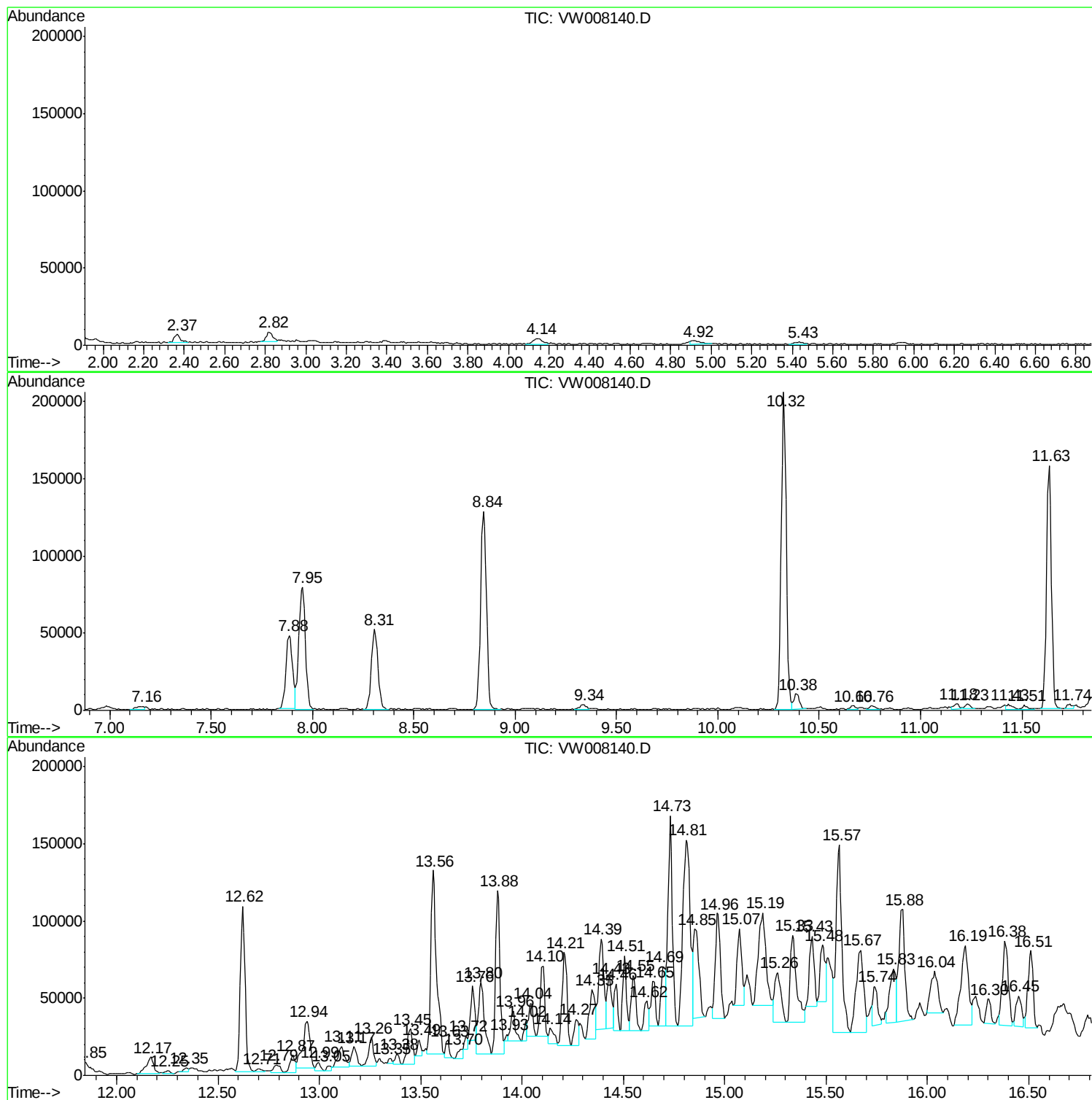
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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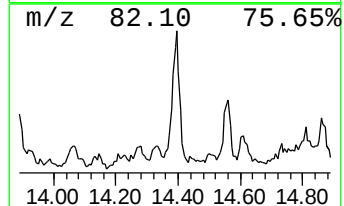
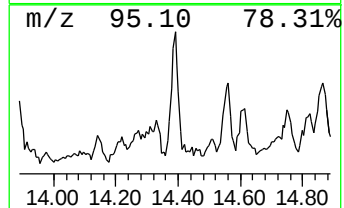
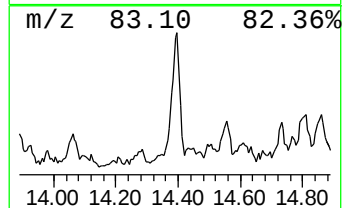
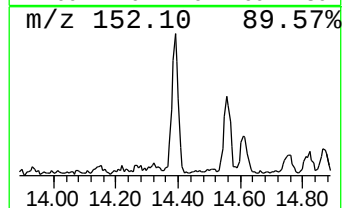
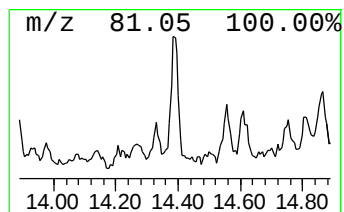
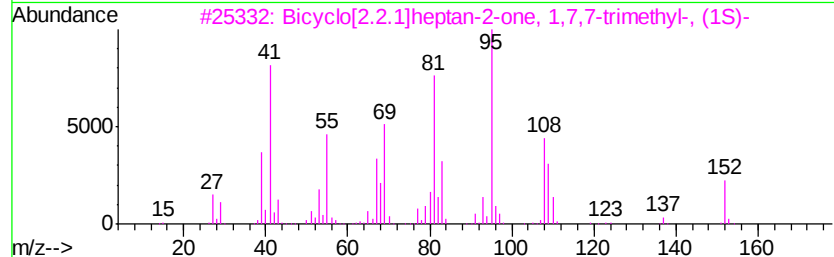
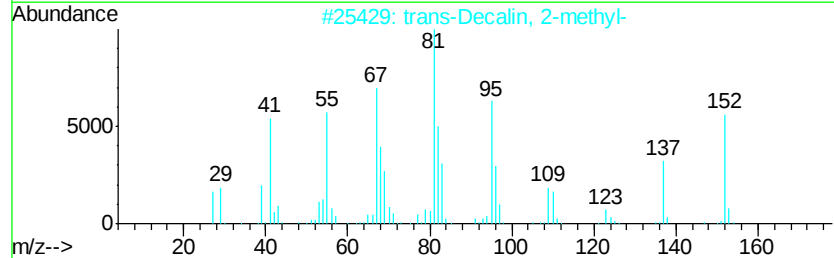
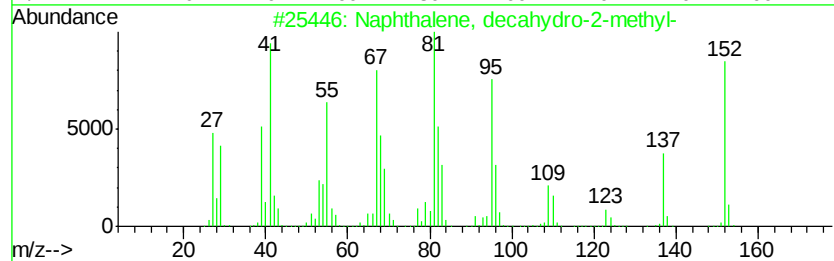
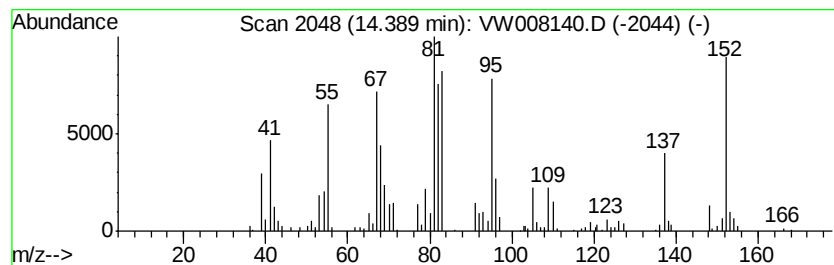
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 1 Naphthalene, decahydro-2-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	22.81 ug/l	103870	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, decahydro-2-methyl-	152 C11H20	002958-76-1 97
2		trans-Decalin, 2-methyl-	152 C11H20	1000152-47-3 74
3		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152 C10H16O	000464-48-2 64
4		Decalin, anti-1-methyl-, cis-	152 C11H20	1000158-89-0 64
5		Pulegone	152 C10H16O	000089-82-7 49



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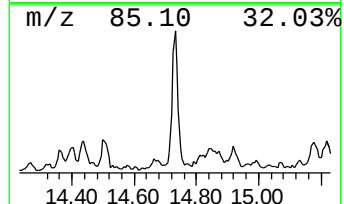
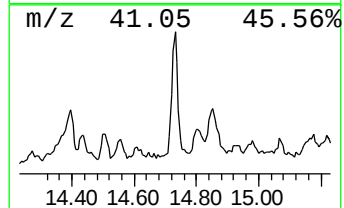
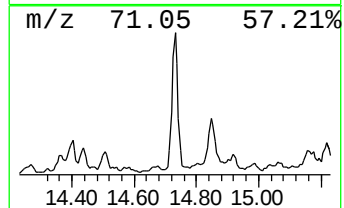
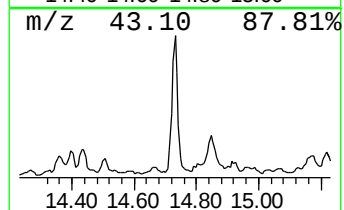
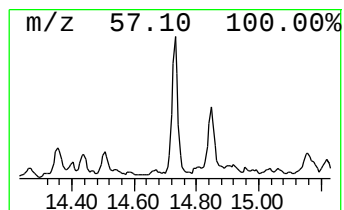
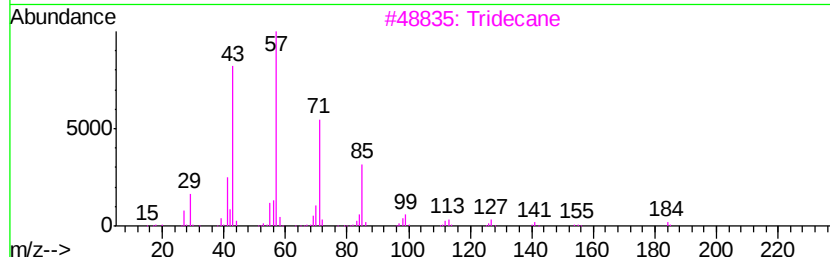
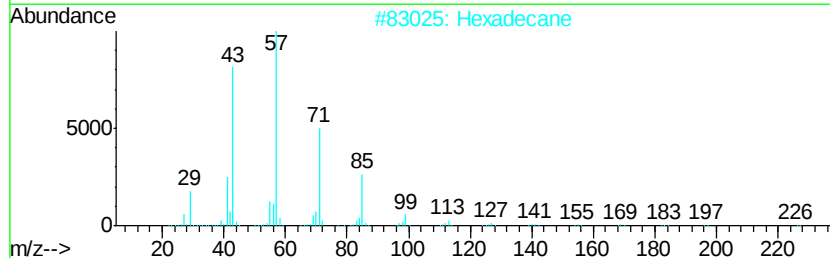
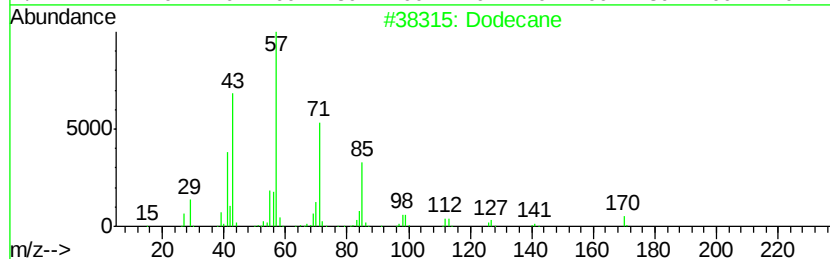
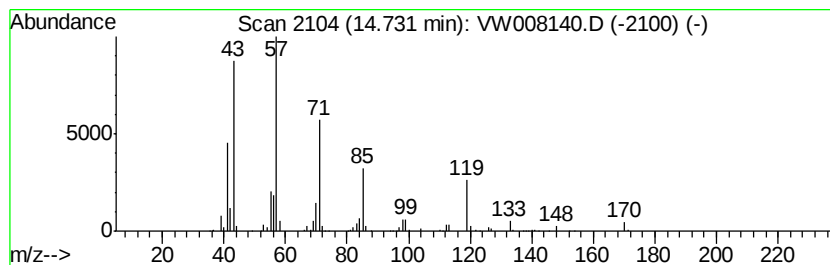
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.73	46.99 ug/l	213957	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dodecane	170	C12H26	000112-40-3	81
2		Hexadecane	226	C16H34	000544-76-3	64
3		Tridecane	184	C13H28	000629-50-5	64
4		Decane, 6-ethyl-2-methyl-	184	C13H28	062108-21-8	59
5		Undecane	156	C11H24	001120-21-4	53



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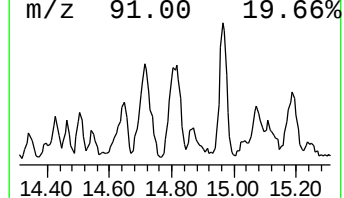
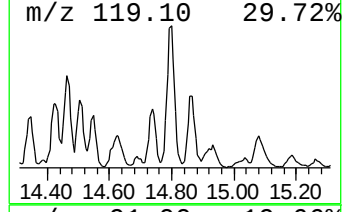
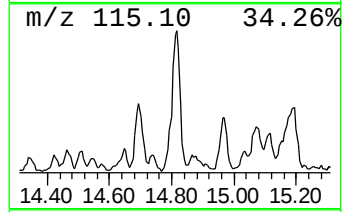
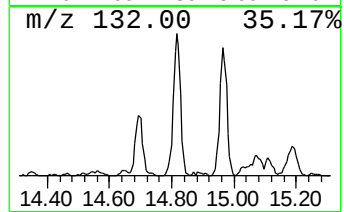
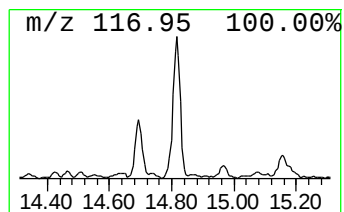
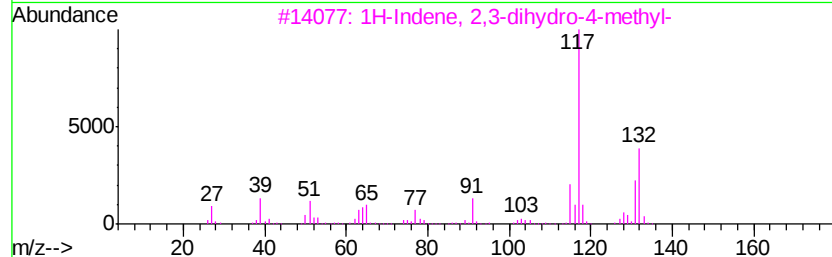
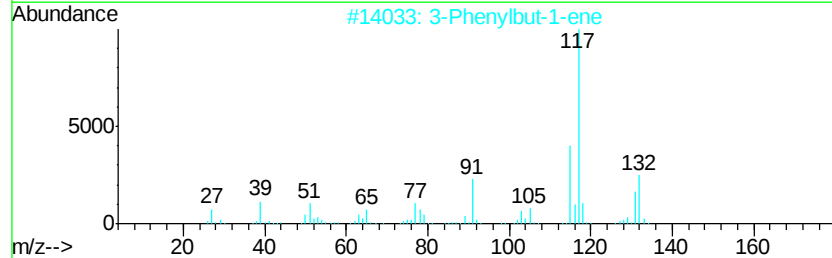
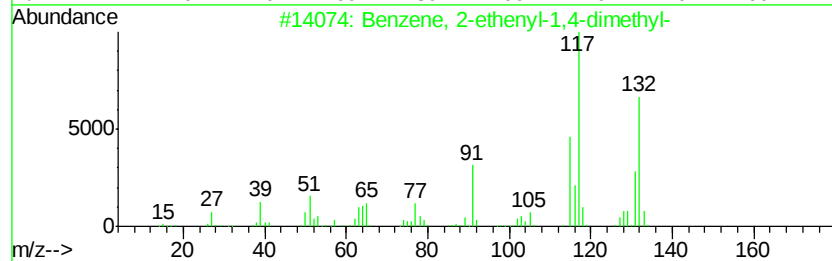
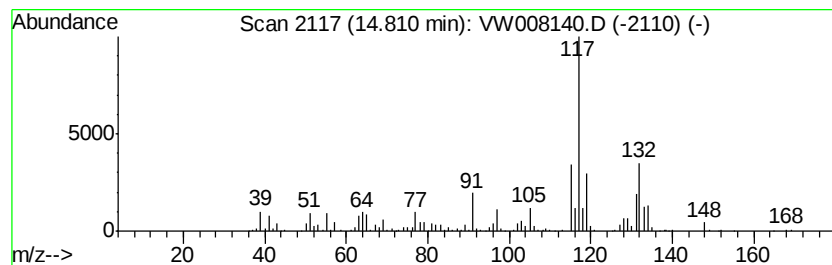
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	66.90 ug/l	304565	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	70
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70



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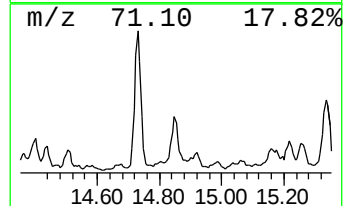
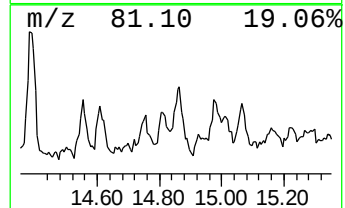
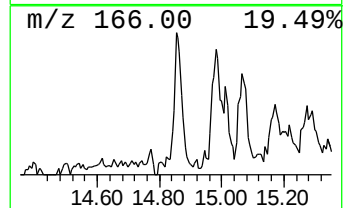
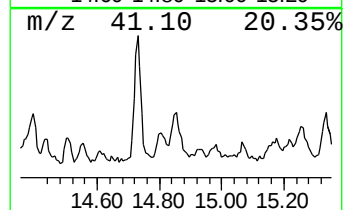
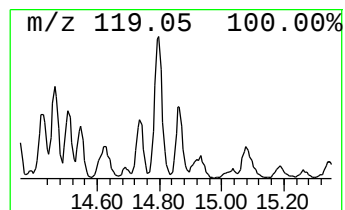
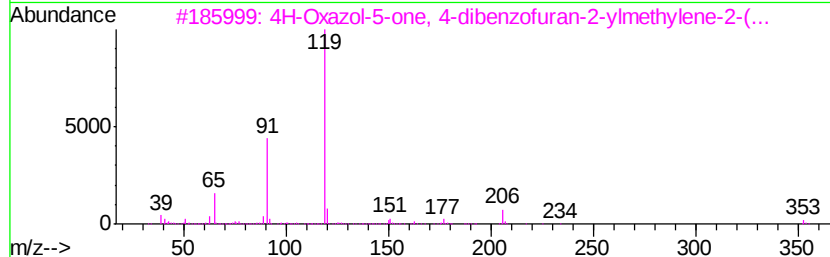
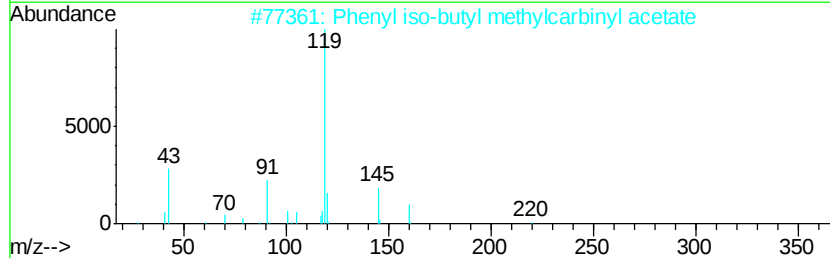
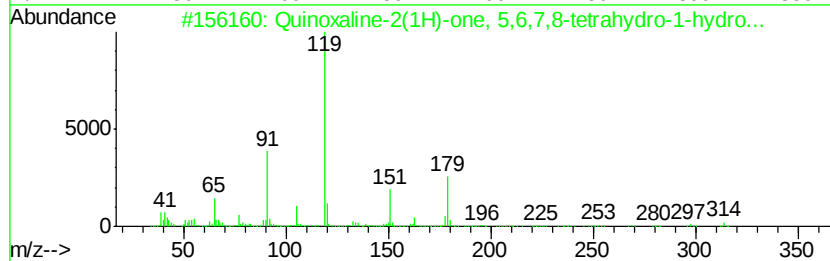
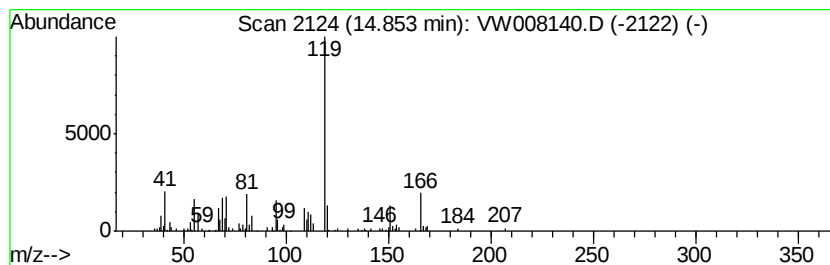
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Quant Title : SW846 8260

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

Peak Number 4 unknown14.85 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	23.51 ug/l	107028	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Quinoxaline-2(1H)-one, 5,6,7,8-t...	314	C17H18N2O4	1000295-63-8	38
2		Phenyl iso-butyl methylcarbinyl ...	220	C14H20O2	1000285-48-0	25
3		4H-Oxazol-5-one, 4-dibenzofuran-...	353	C23H15N03	1000302-18-6	25
4		5-Chlorovaleric acid, 3-tridecyl...	318	C18H35ClO2	1000299-91-3	25
5		Acetamide, N-[1-benzylideno-2-(2...	306	C19H18N2O2	314281-76-0	25



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA W/SOIL
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

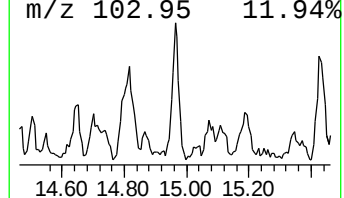
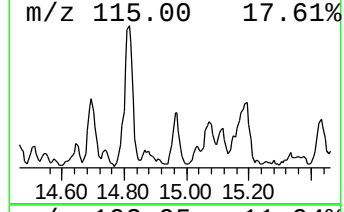
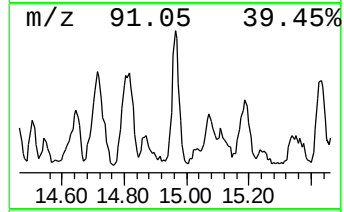
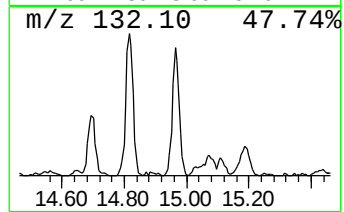
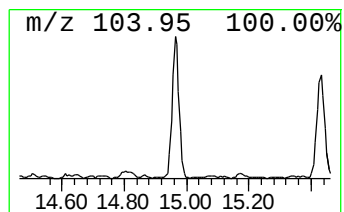
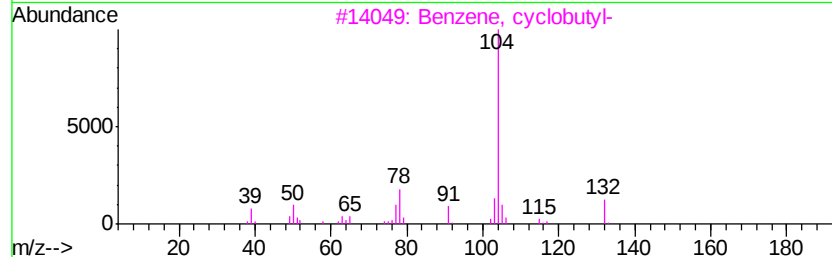
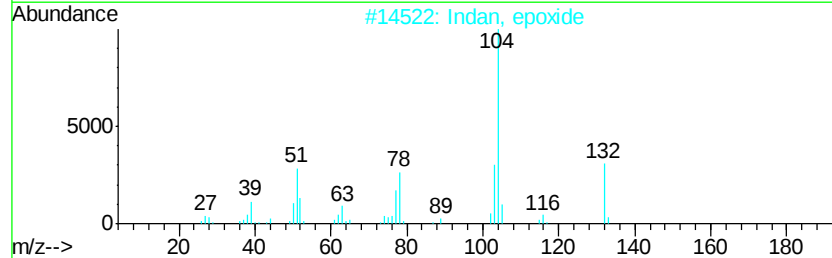
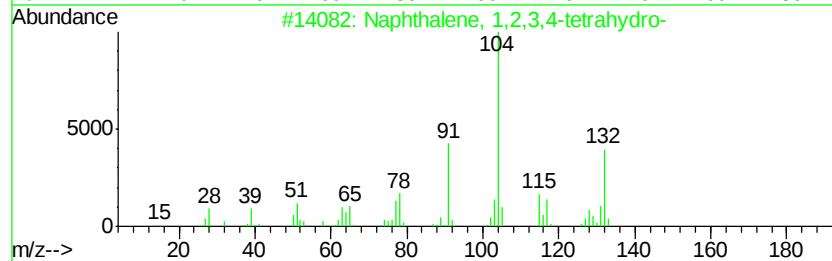
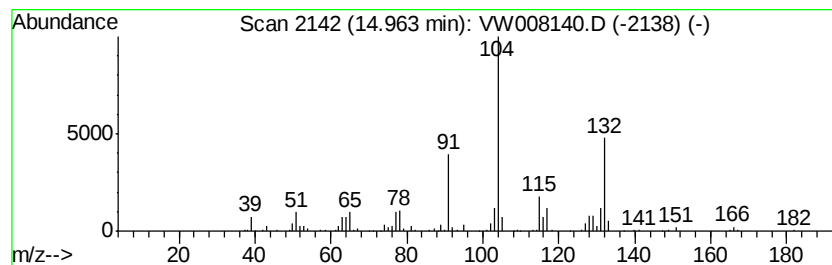
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	25.86 ug/l	117744	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2		Indan, epoxide	132	C9H8O	000768-22-9	62
3		Benzene, cyclobutyl-	132	C10H12	004392-30-7	59
4		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	58
5		2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA W/SOIL
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

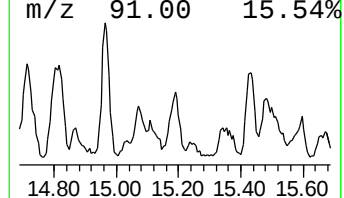
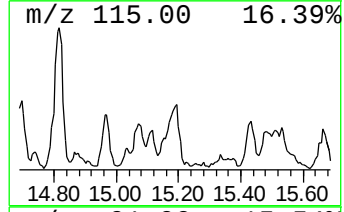
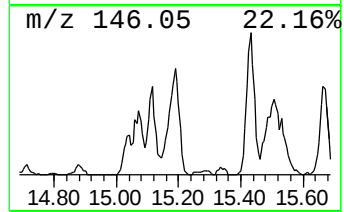
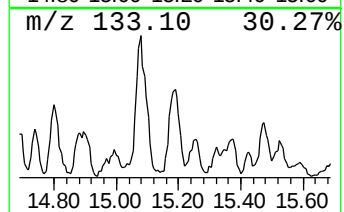
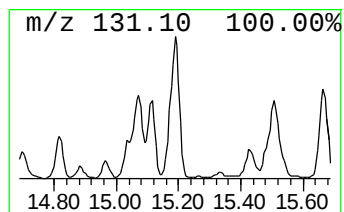
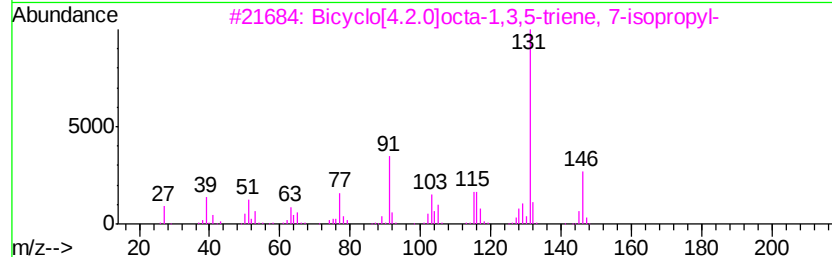
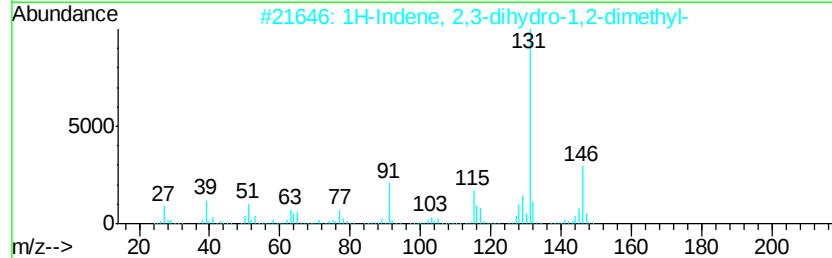
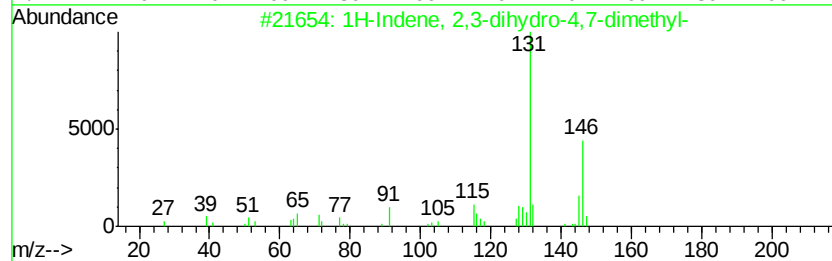
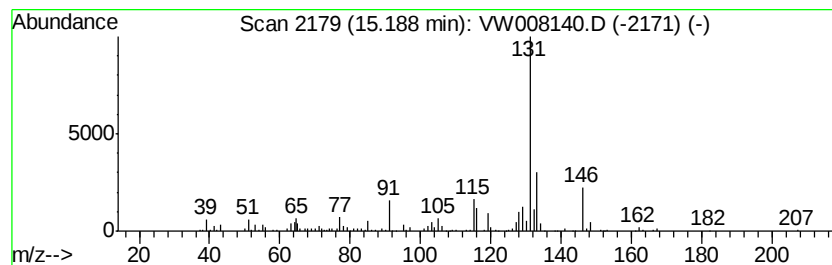
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	34.82 ug/l	158552	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	81
4		Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	81
5		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA W/SOIL
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

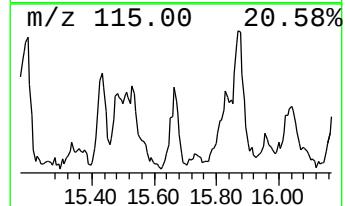
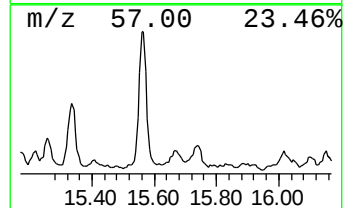
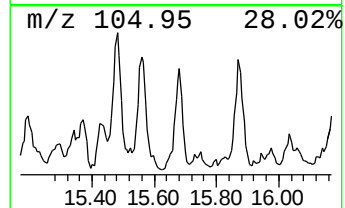
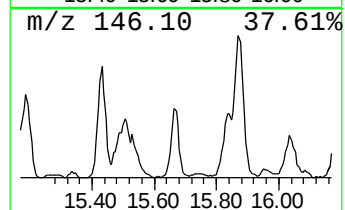
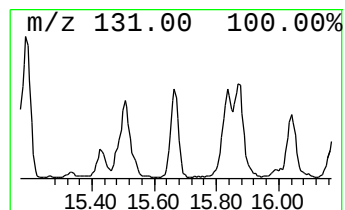
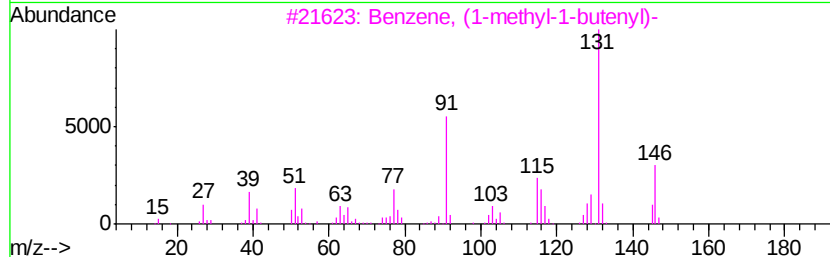
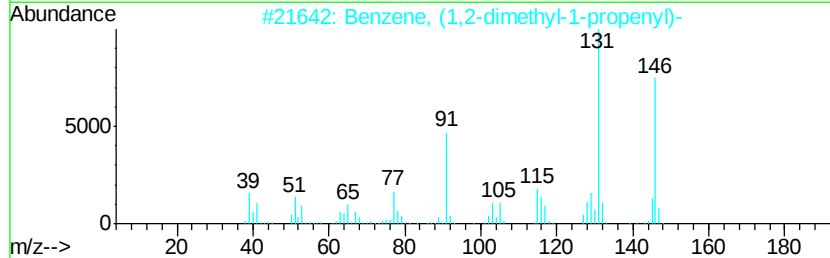
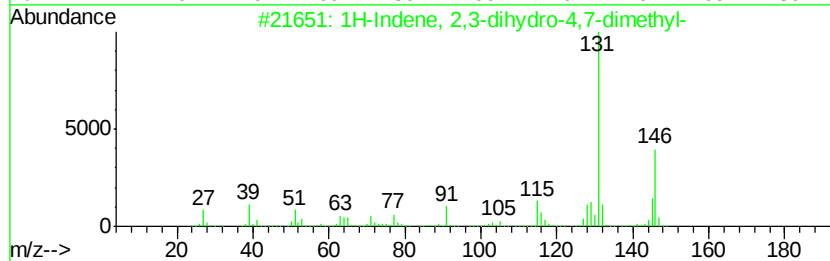
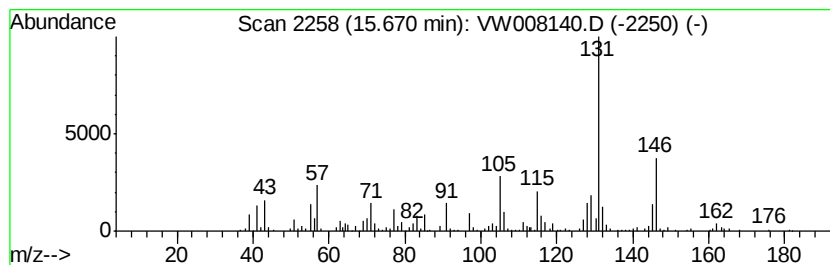
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.67	29.98 ug/l	136490	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	90
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14		000769-57-3	83
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14		053172-84-2	74
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14		004175-53-5	70
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14		001075-22-5	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA W/SOIL
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

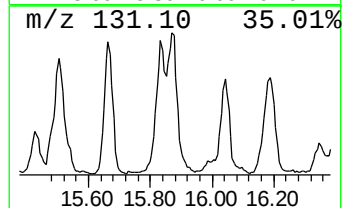
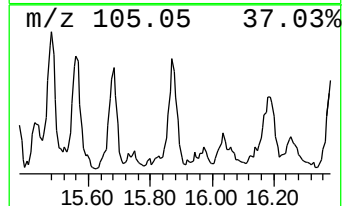
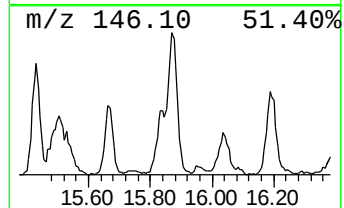
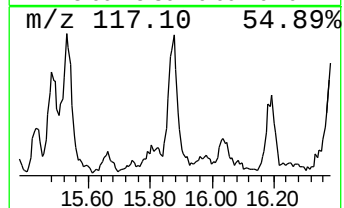
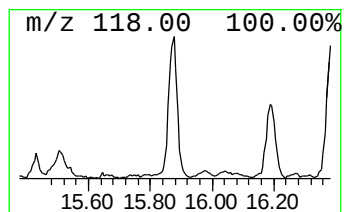
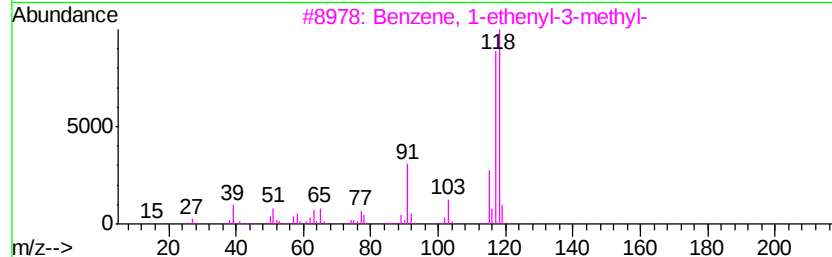
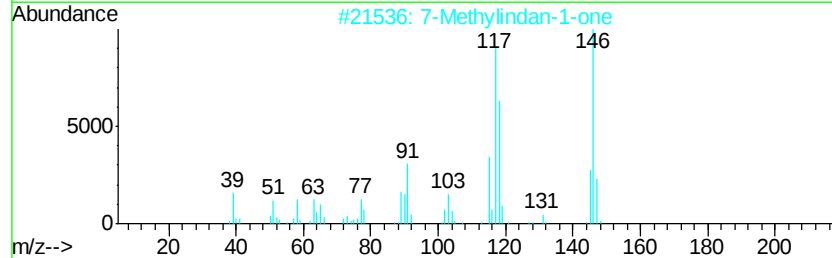
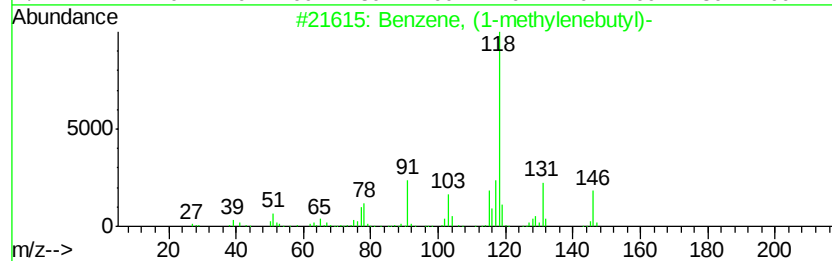
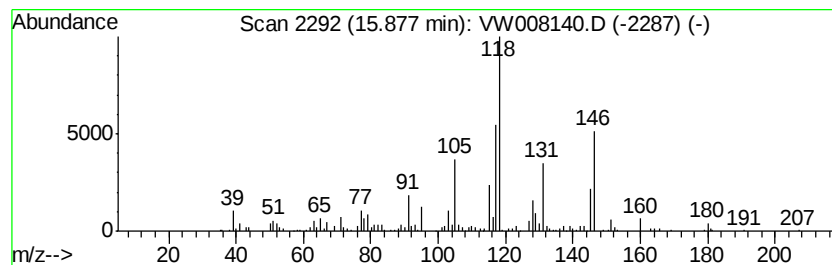
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methylenebutyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	32.06 ug/l	145945	1,4-Dichlorobenzene-d4	13.56

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	64
2	7-Methylindan-1-one	146	C10H10O	039627-61-7	43
3	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	41
4	Benzeneethanol, .beta.-ethenyl-....	162	C11H14O	040596-14-3	38
5	Propanoic acid, 2-methyl-, 3-phe...	206	C13H18O2	000103-58-2	35



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA W/SOIL
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

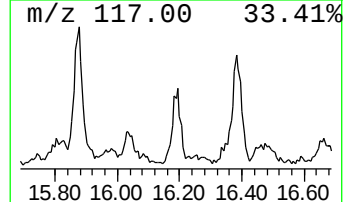
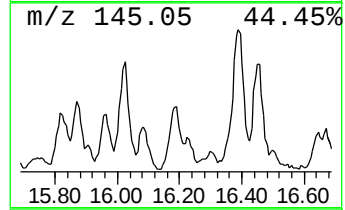
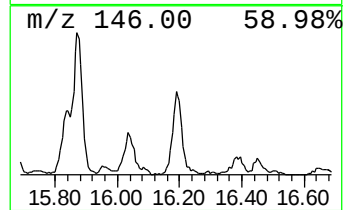
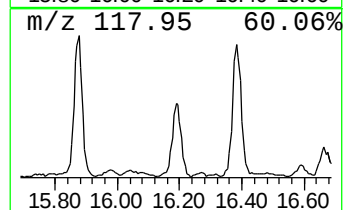
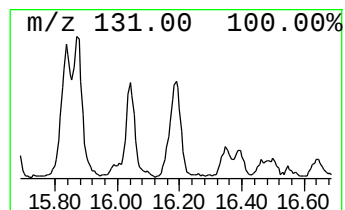
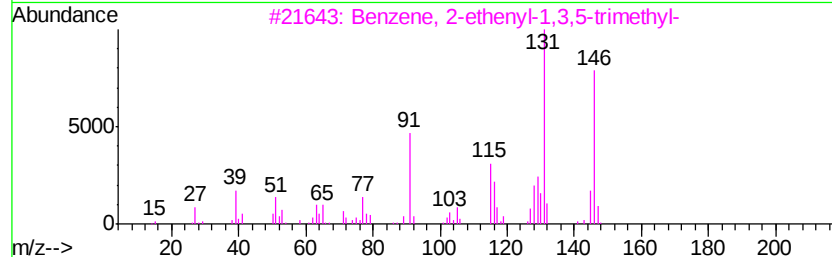
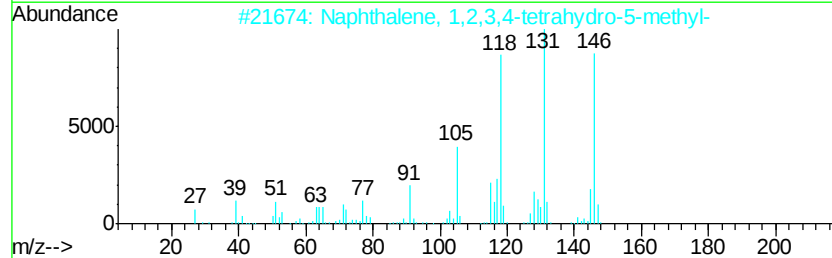
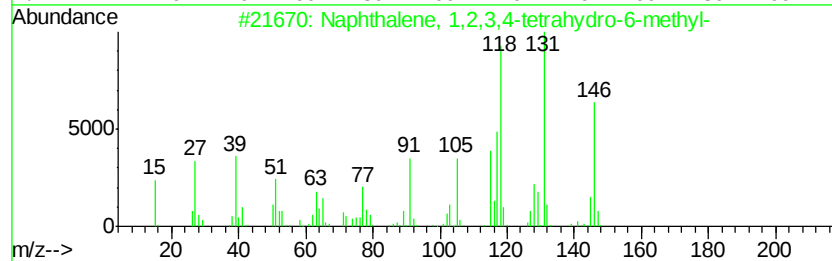
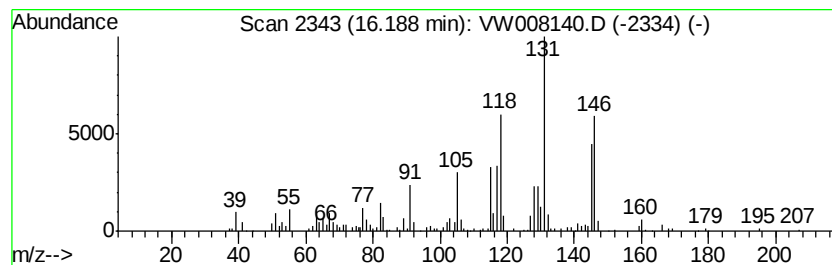
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.19	29.61 ug/l	134807	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	89
2		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	76
3		Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	64
4		1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	60
5		1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA W/SOIL
ALS Vial : 52 Sample Multiplier: 1

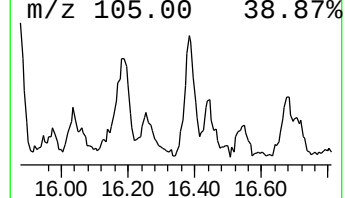
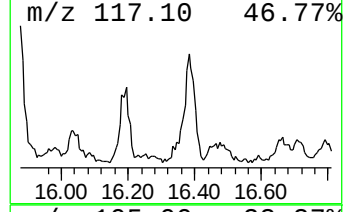
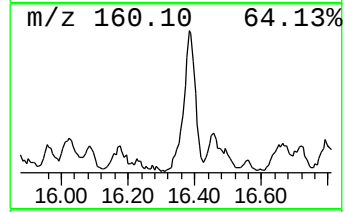
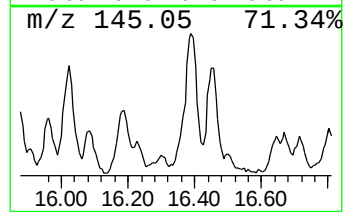
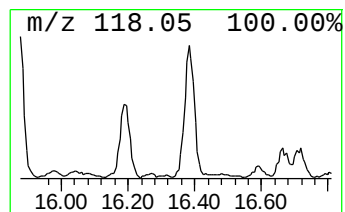
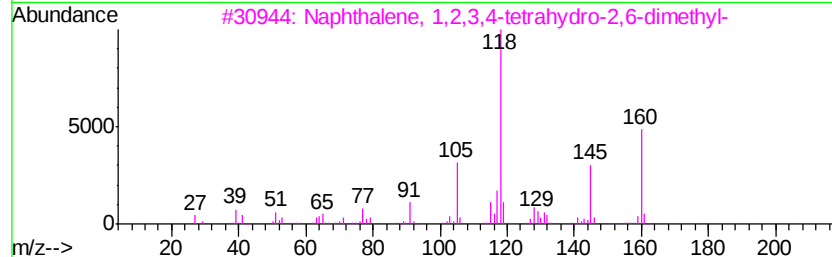
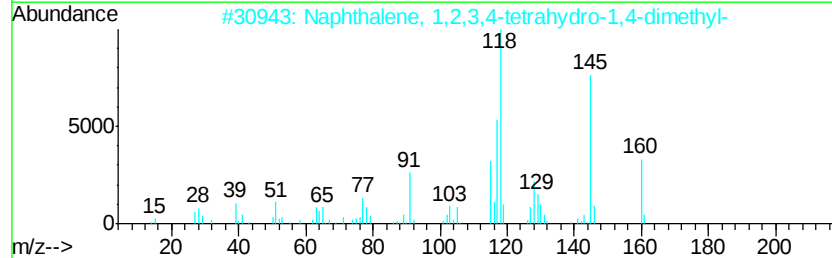
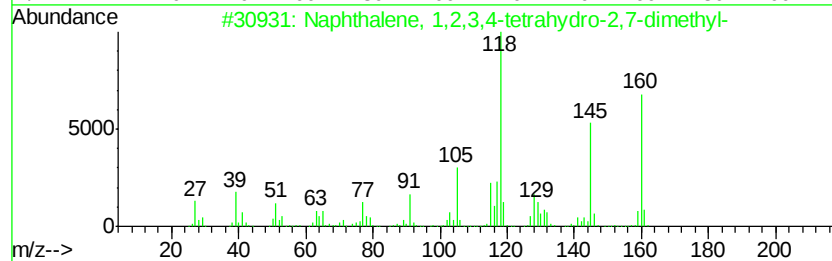
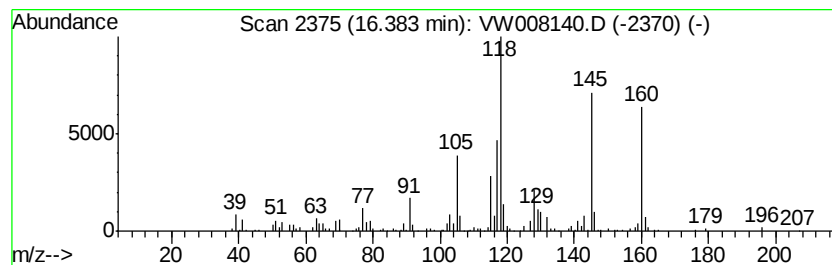
Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.38	23.93 ug/l	108960	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	013065-07-1	90
2	Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6	86
3	Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	007524-63-2	64
4	Benzene, 1-(1-methylethenyl)-3-(...	160 C12H16	001129-29-9	50
5	1(2H)-Naphthalenone, 3,4-dihydro...	160 C11H12O	014944-23-1	49



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW010919\
Data File : VW008140.D
Acq On : 10 Jan 2019 09:25
Operator : SY/AP
Sample : K1006-11
Misc : 6.78G/5ML/MSVOA_W/SOIL
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
JC-03-010819-F

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W010819S.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
Naphthalene, deca...	14.39	22.8	ug/l	103870	4	13.56	227642	50.0
Dodecane	14.73	47.0	ug/l	213957	4	13.56	227642	50.0
Benzene, 2-etheny...	14.81	66.9	ug/l	304565	4	13.56	227642	50.0
unknown14.85	14.85	23.5	ug/l	107028	4	13.56	227642	50.0
Naphthalene, 1,2,...	14.96	25.9	ug/l	117744	4	13.56	227642	50.0
1H-Indene, 2,3-di...	15.19	34.8	ug/l	158552	4	13.56	227642	50.0
Benzene, (1,2-dim...	15.67	30.0	ug/l	136490	4	13.56	227642	50.0
Benzene, (1-methy...	15.88	32.1	ug/l	145945	4	13.56	227642	50.0
Naphthalene, 1,2,...	16.19	29.6	ug/l	134807	4	13.56	227642	50.0
Naphthalene, 1,2,...	16.38	23.9	ug/l	108960	4	13.56	227642	50.0