

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW1111318\
 Data File : VW006857.D
 Acq On : 13 Nov 2018 16:28
 Operator : SY/AP
 Sample : J5860-24
 Misc : 4.87G/5ML/MSVOA W/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EP2-SB-111(14-16)

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\82W111218S.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	7.890	971	982	986	rBV	33002	76459	1.65%	0.271%
2	7.951	986	992	1001	rVB	80249	189312	4.10%	0.671%
3	8.311	1043	1051	1063	rVB	27019	60190	1.30%	0.213%
4	8.847	1127	1139	1151	rBV	111919	234478	5.07%	0.831%
5	9.335	1205	1219	1225	rBV2	49074	117606	2.54%	0.417%
6	9.762	1282	1289	1299	rVB	40669	80598	1.74%	0.286%
7	9.896	1307	1311	1317	rVB2	50545	94415	2.04%	0.334%
8	9.963	1317	1322	1326	rBV2	46202	83459	1.81%	0.296%
9	10.103	1335	1345	1356	rBV	54367	121608	2.63%	0.431%
10	10.256	1364	1370	1375	rVB	57770	100743	2.18%	0.357%
11	10.329	1375	1382	1392	rBV	160481	318116	6.88%	1.127%
12	10.506	1405	1411	1418	rVB2	115813	212479	4.60%	0.753%
13	10.750	1440	1451	1462	rVB6	93856	364583	7.89%	1.292%
14	10.939	1473	1482	1488	rBV2	28595	57430	1.24%	0.203%
15	11.042	1488	1499	1506	rBV	44118	124679	2.70%	0.442%
16	11.341	1544	1548	1552	rBV2	31945	49984	1.08%	0.177%
17	11.414	1552	1560	1571	rVV3	164627	400701	8.67%	1.419%
18	11.518	1571	1577	1587	rVV	142481	252099	5.45%	0.893%
19	11.634	1589	1596	1602	rVV	168637	309754	6.70%	1.097%
20	11.737	1603	1613	1621	rVV	502422	896367	19.39%	3.175%
21	11.847	1621	1631	1640	rVV	639817	1226213	26.53%	4.344%
22	11.920	1640	1643	1647	rVB2	33366	50803	1.10%	0.180%
23	12.073	1655	1668	1674	rBV5	40685	119268	2.58%	0.423%
24	12.140	1674	1679	1685	rBV2	66386	131446	2.84%	0.466%
25	12.255	1694	1698	1702	rVB	59454	90471	1.96%	0.320%
26	12.335	1702	1711	1715	rBV4	55070	145447	3.15%	0.515%
27	12.383	1715	1719	1723	rVB4	31206	60236	1.30%	0.213%
28	12.469	1729	1733	1737	rBV	73804	139054	3.01%	0.493%
29	12.518	1737	1741	1743	rVV4	25060	47729	1.03%	0.169%
30	12.572	1748	1750	1754	rVV2	95066	170989	3.70%	0.606%
31	12.621	1755	1758	1761	rVV	124968	182533	3.95%	0.647%
32	12.652	1761	1763	1768	rVB	111421	144615	3.13%	0.512%
33	12.707	1768	1772	1776	rVB5	31244	51494	1.11%	0.182%
34	12.761	1776	1781	1783	rBV4	28193	46656	1.01%	0.165%

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW111318\
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 Misc : 4.87G/5ML/MSVOA W/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 EP2-SB-111(14-16)

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

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 Title : SW846 8260

35	12.810	1783	1789	1794	rVB	406188	634735	13.73%	2.249%
36	12.902	1798	1804	1808	rBV	767755	1292257	27.96%	4.578%
37	12.950	1808	1812	1817	rVB	717100	1133123	24.51%	4.014%
38	13.109	1829	1838	1845	rBV	194895	402961	8.72%	1.427%
39	13.255	1856	1862	1868	rBV	3014671	4622388	100.00%	16.375%
40	13.371	1874	1881	1886	rBV3	86670	242448	5.25%	0.859%
41	13.493	1897	1901	1908	rBV4	71030	152094	3.29%	0.539%
42	13.591	1913	1917	1930	rVB	621028	1222825	26.45%	4.332%
43	13.725	1933	1939	1942	rBV	218221	364812	7.89%	1.292%
44	13.761	1942	1945	1948	rVV	446161	715595	15.48%	2.535%
45	13.798	1948	1951	1961	rVB4	582151	1298457	28.09%	4.600%
46	13.889	1961	1966	1969	rBV3	60666	104519	2.26%	0.370%
47	13.956	1972	1977	1982	rBV	144942	227548	4.92%	0.806%
48	14.017	1982	1987	1989	rBV	292534	450660	9.75%	1.596%
49	14.048	1989	1992	1996	rVB	309698	436387	9.44%	1.546%
50	14.103	1996	2001	2007	rVB	628567	931323	20.15%	3.299%
51	14.164	2007	2011	2014	rBV	91194	121222	2.62%	0.429%
52	14.212	2014	2019	2025	rVB2	341063	537545	11.63%	1.904%
53	14.279	2025	2030	2035	rBV5	27098	58378	1.26%	0.207%
54	14.346	2035	2041	2046	rBV2	125315	225297	4.87%	0.798%
55	14.426	2046	2054	2057	rBV	312493	531963	11.51%	1.884%
56	14.462	2057	2060	2064	rVB	347641	464931	10.06%	1.647%
57	14.505	2064	2067	2071	rVB	153481	190703	4.13%	0.676%
58	14.548	2071	2074	2080	rVB	93101	131446	2.84%	0.466%
59	14.651	2087	2091	2094	rBV	90357	118049	2.55%	0.418%
60	14.694	2094	2098	2103	rBV2	373930	663493	14.35%	2.350%
61	14.816	2110	2118	2123	rBV2	450992	1050989	22.74%	3.723%
62	14.865	2123	2126	2133	rVB	108266	189998	4.11%	0.673%
63	14.968	2138	2143	2149	rBV2	61651	106543	2.30%	0.377%
64	15.035	2150	2154	2155	rBV	64767	79747	1.73%	0.283%
65	15.072	2155	2160	2164	rBV2	152189	253131	5.48%	0.897%
66	15.188	2172	2179	2186	rVB2	159317	373606	8.08%	1.323%
67	15.377	2198	2210	2216	rBV	462929	888216	19.22%	3.146%
68	15.480	2221	2227	2231	rVV3	87270	178166	3.85%	0.631%
69	15.535	2232	2236	2250	rVB3	62515	181593	3.93%	0.643%
70	15.663	2250	2257	2265	rBV	82863	188506	4.08%	0.668%
71	15.834	2277	2285	2295	rBV5	34213	111083	2.40%	0.394%

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EP2-SB-111(14-16)

Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Title : SW846 8260

72	16.035	2311	2318	2324	rVB5	28396	75507	1.63%	0.267%
73	16.450	2379	2386	2400	rVB	169960	374808	8.11%	1.328%
74	16.657	2412	2420	2432	rVB	63643	149839	3.24%	0.531%

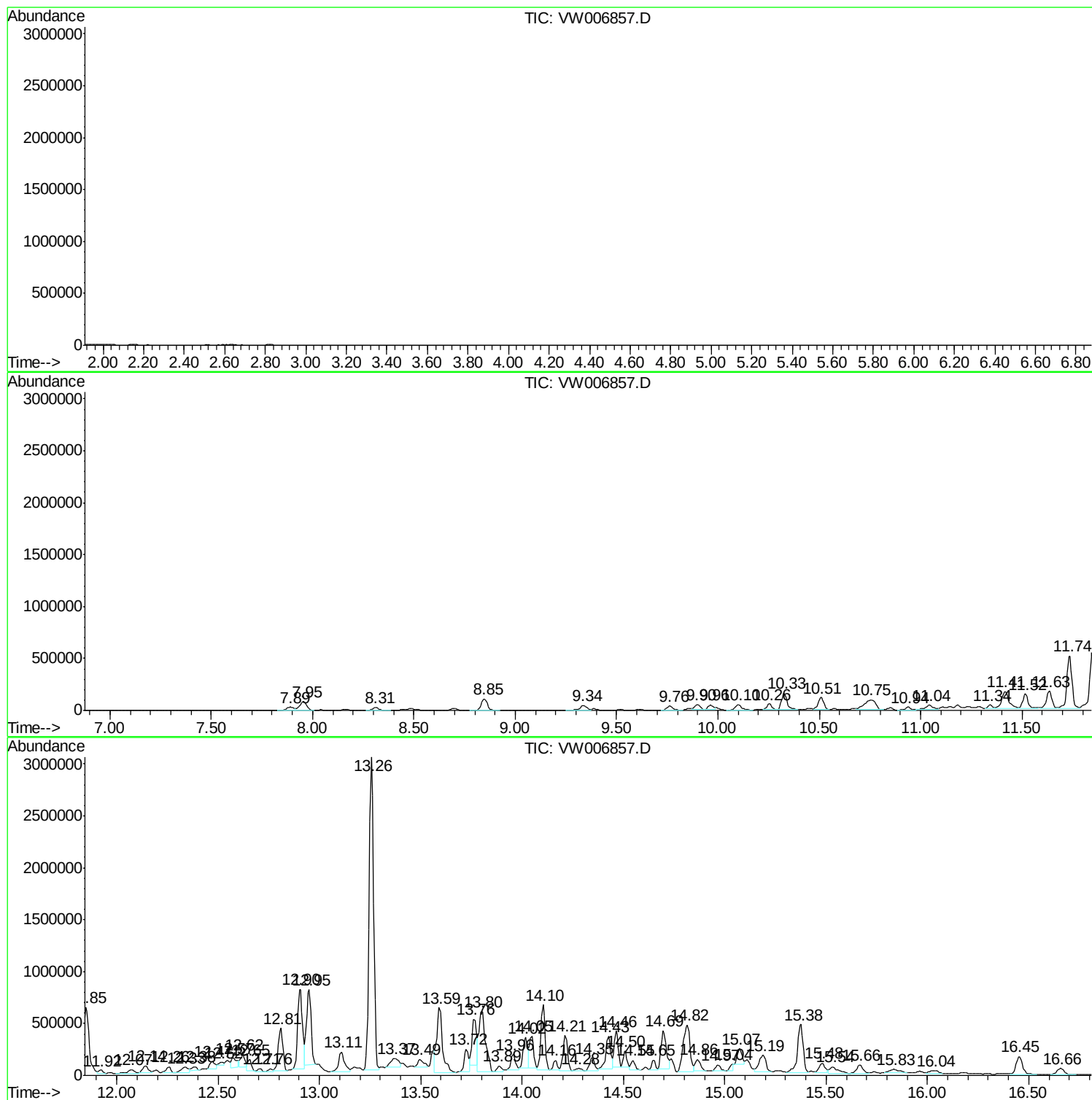
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Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
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Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-111(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006857.D
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Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-111(14-16)

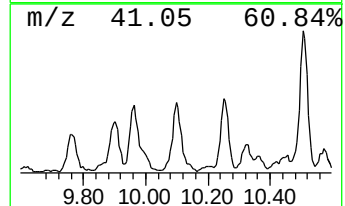
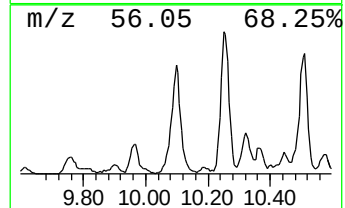
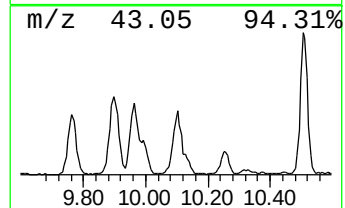
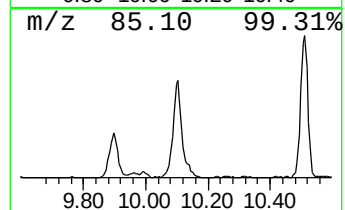
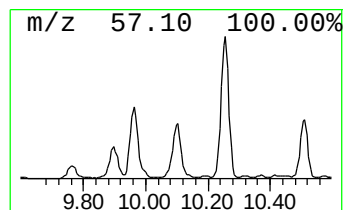
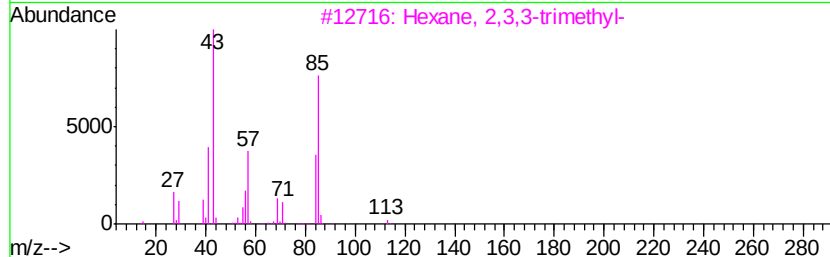
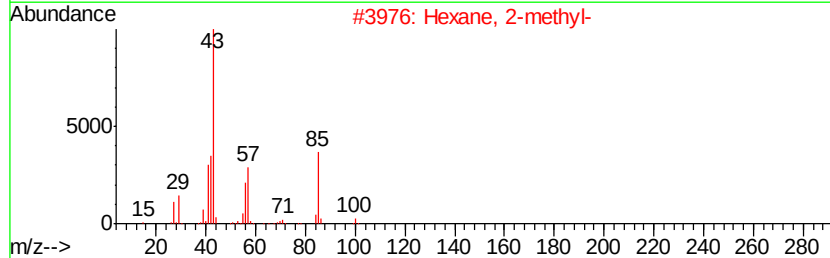
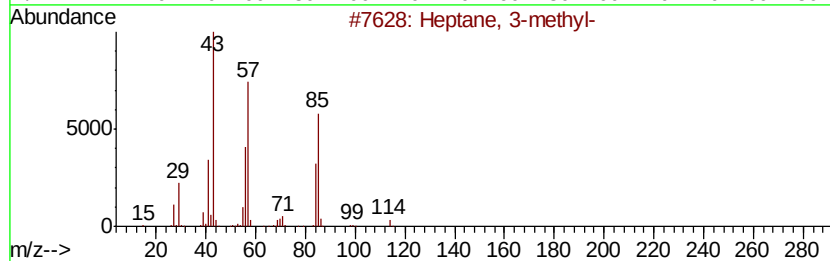
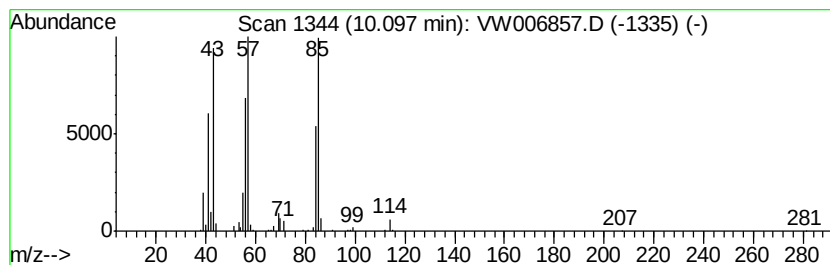
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.10	25.93 ug/l	121608	1,4-Difluorobenzene	8.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2		Hexane, 2-methyl-	100	C7H16	000591-76-4	59
3		Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	59
4		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
5		Pentane, 2,3,3,4-tetramethyl-	128	C9H20	016747-38-9	53



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Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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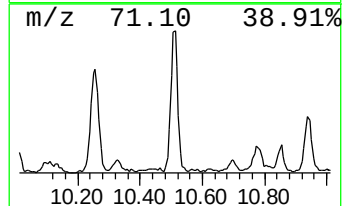
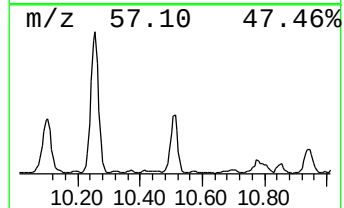
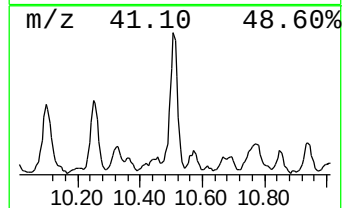
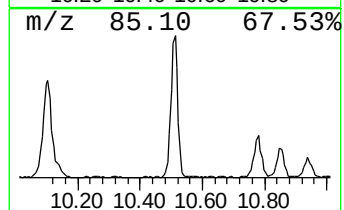
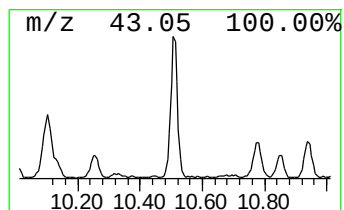
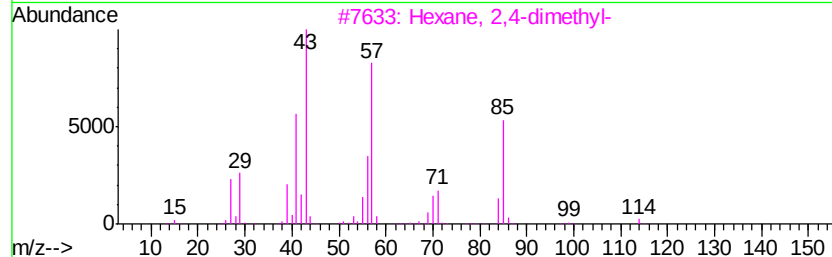
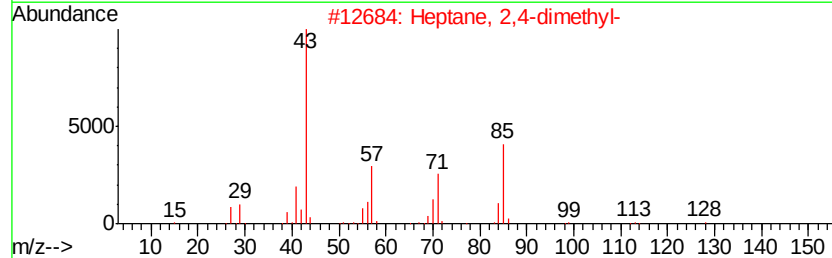
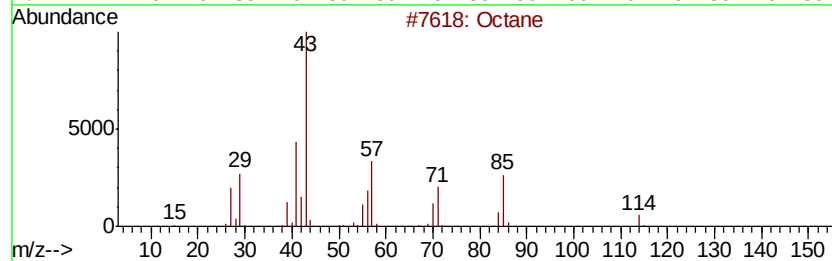
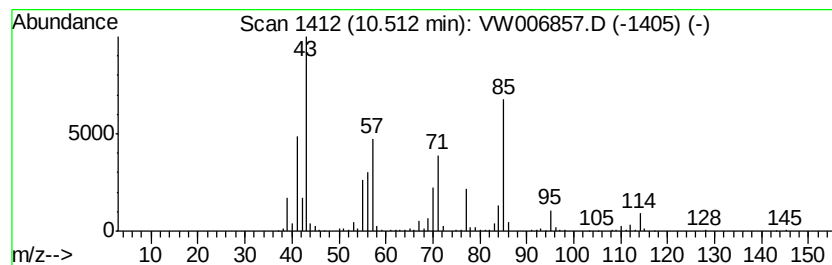
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 2 Octane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.51	34.30 ug/l	212479	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane	114	C8H18	000111-65-9	81
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50
3		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	47
4		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	43
5		Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
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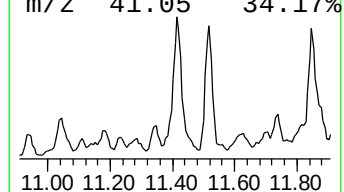
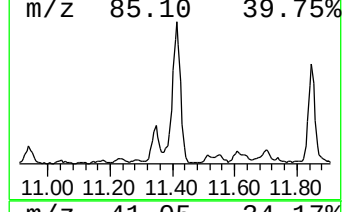
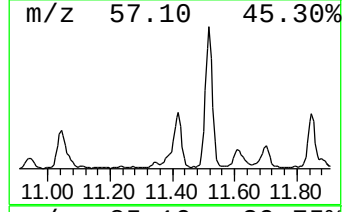
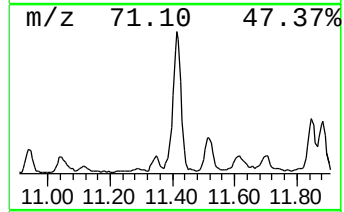
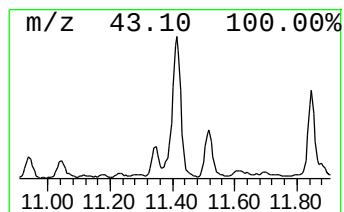
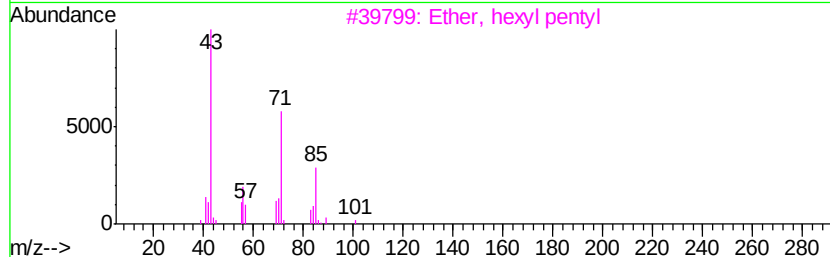
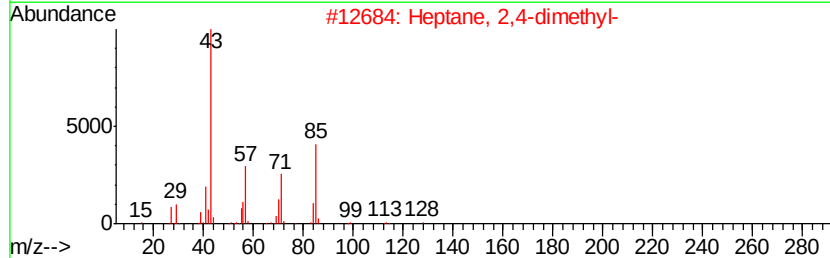
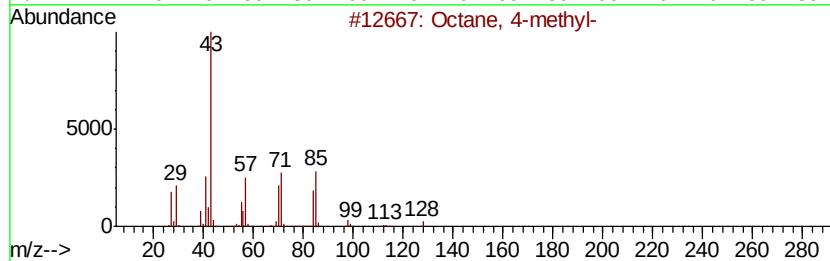
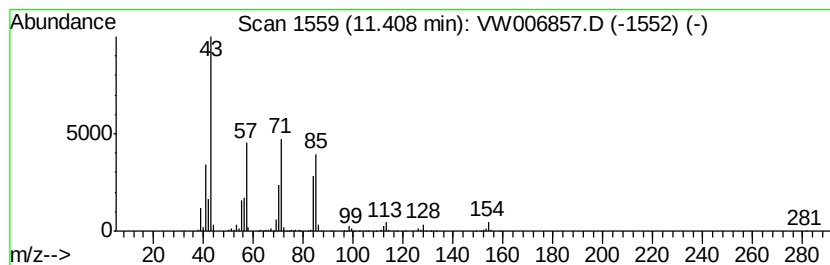
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.41	64.68 ug/l	400701	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 4-methyl-	128	C9H20	002216-34-4	80
2		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	64
3		Ether, hexyl pentyl	172	C11H24O	032357-83-8	53
4		Octane, 2-methyl-	128	C9H20	003221-61-2	49
5		Hexane, 3-ethyl-	114	C8H18	000619-99-8	47



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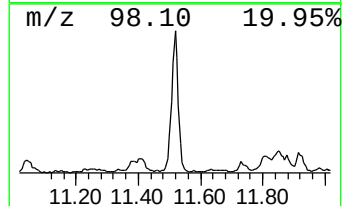
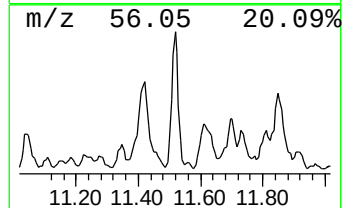
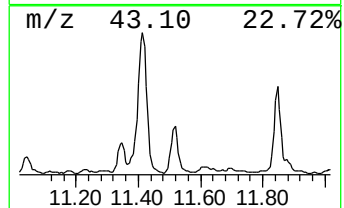
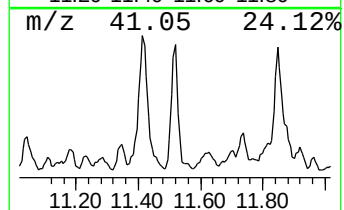
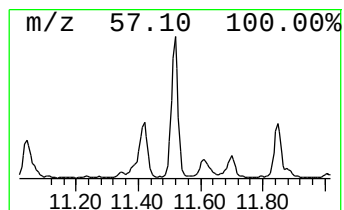
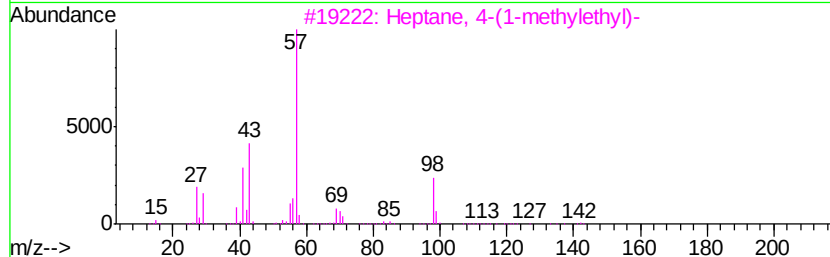
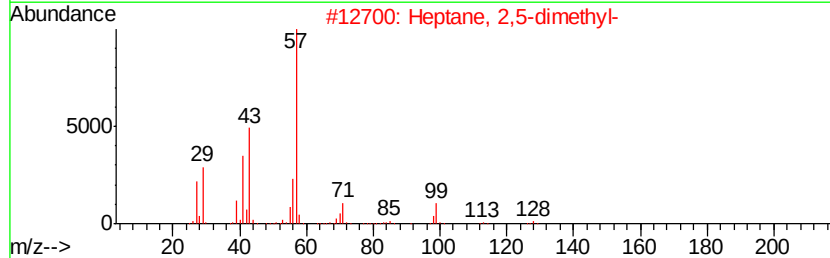
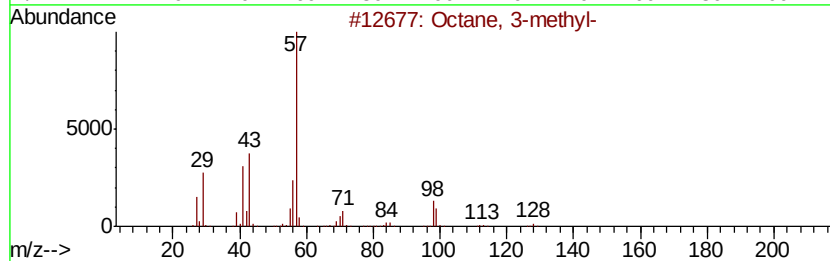
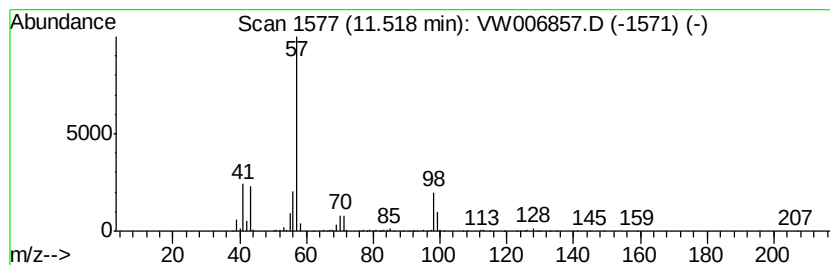
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Quant Title : SW846 8260

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

Peak Number 5 Octane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.52	40.69 ug/l	252099	Chlorobenzene-d5	11.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octane, 3-methyl-	128	C9H20	002216-33-3	91
2		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	81
3		Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	72
4		Undecane, 6-methyl-	170	C12H26	017302-33-9	59
5		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006857.D
Acq On : 13 Nov 2018 16:28
Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-111(14-16)

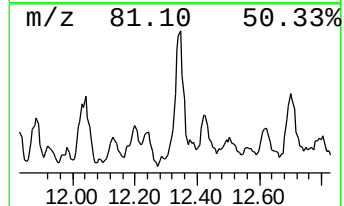
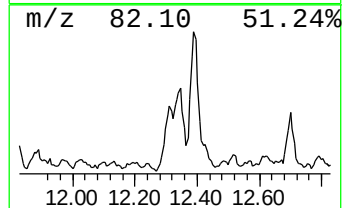
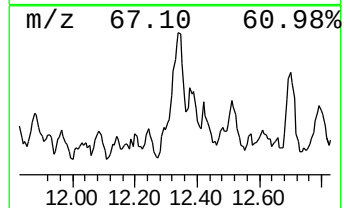
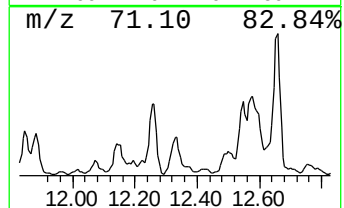
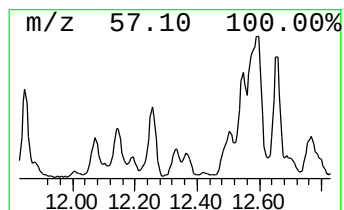
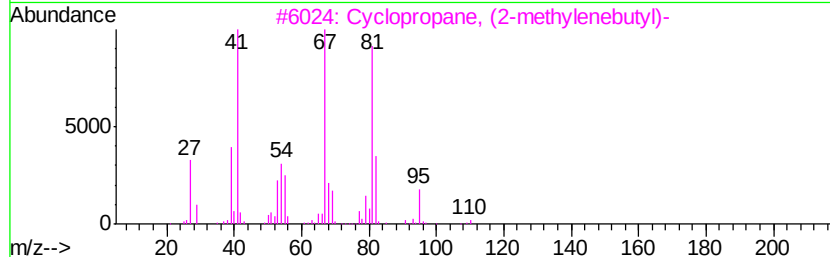
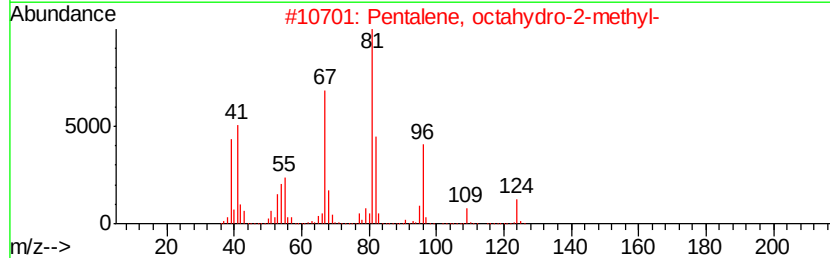
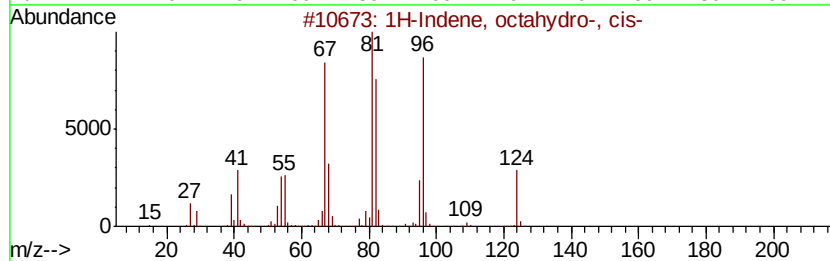
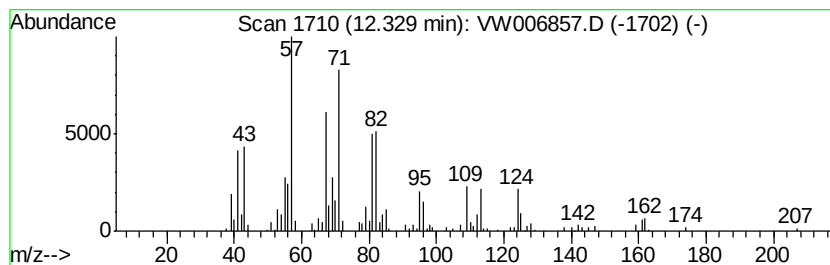
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Quant Title : SW846 8260

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

Peak Number 6 unknown12.33 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	23.48 ug/l	145447	Chlorobenzene-d5	11.63

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	46
2			Pentalene, octahydro-2-methyl-	124	C9H16	003868-64-2	38
3			Cyclopropane, (2-methylenebutyl)-	110	C8H14	074685-56-6	38
4			Butanoic acid, 3-hexenyl ester, ...	170	C10H18O2	016491-36-4	35
5			Cyclohexane, methylene-	96	C7H12	001192-37-6	30



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006857.D
Acq On : 13 Nov 2018 16:28
Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

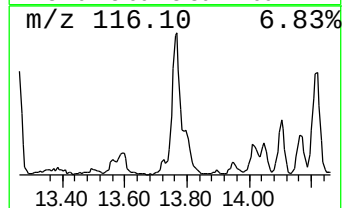
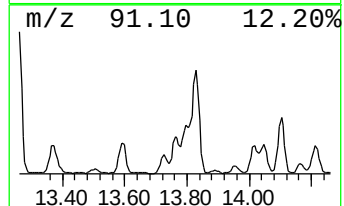
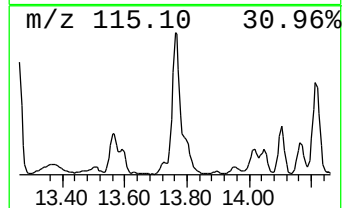
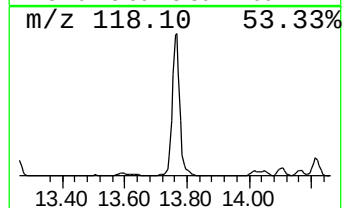
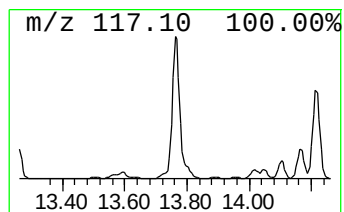
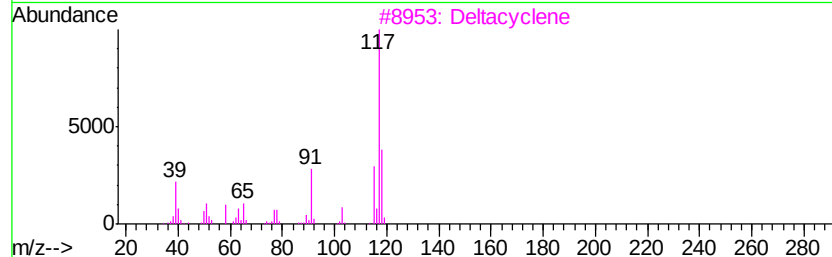
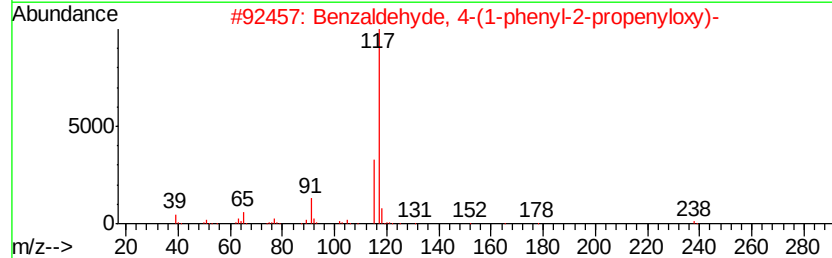
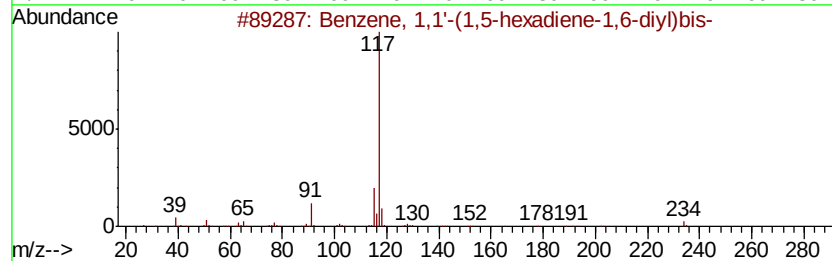
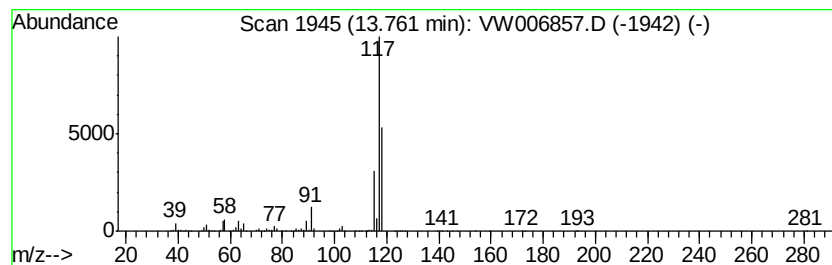
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MSVOA_W
ClientSampleId :
EP2-SB-111(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	29.26 ug/l	715595	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,1'-(1,5-hexadiene-1,6...	234 C18H18	004439-45-6 53
2		Benzaldehyde, 4-(1-phenyl-2-prop...	238 C16H14O2	1000277-56-1 53
3		Deltacyclene	118 C9H10	007785-10-6 50
4		Benzene, (2-bromocyclopropyl)-	196 C9H9Br	036617-02-4 45
5		trans-Cinnamyl bromide	196 C9H9Br	026146-77-0 37



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006857.D
Acq On : 13 Nov 2018 16:28
Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

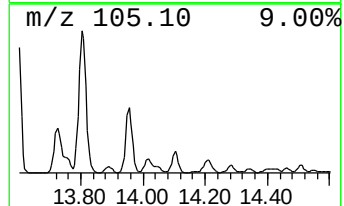
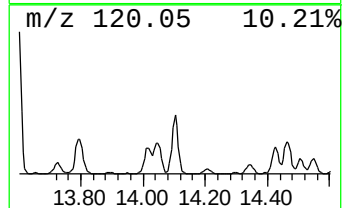
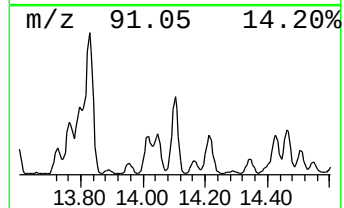
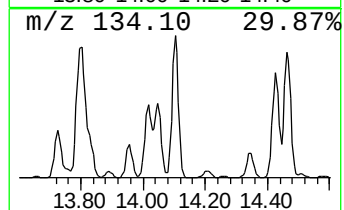
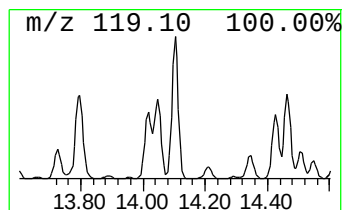
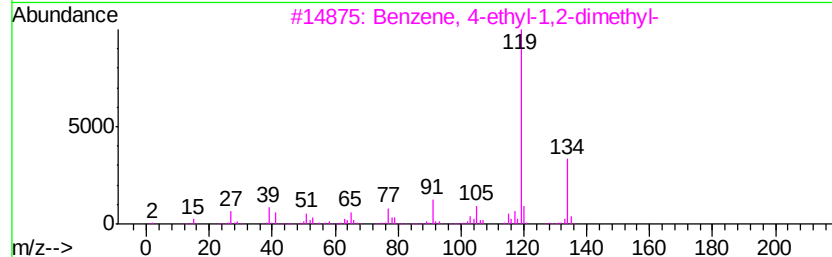
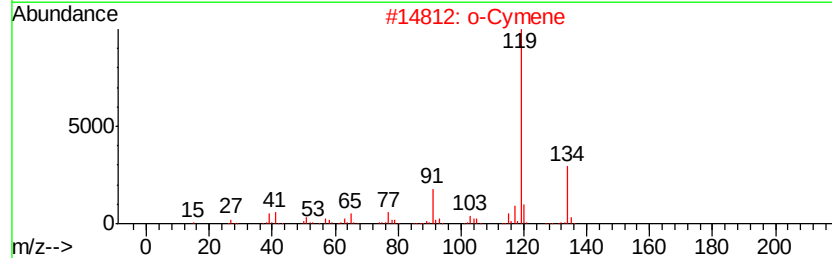
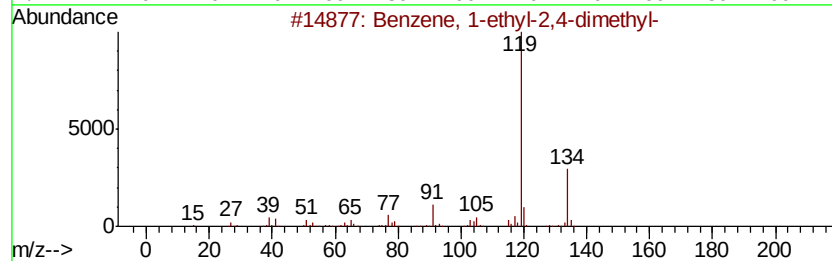
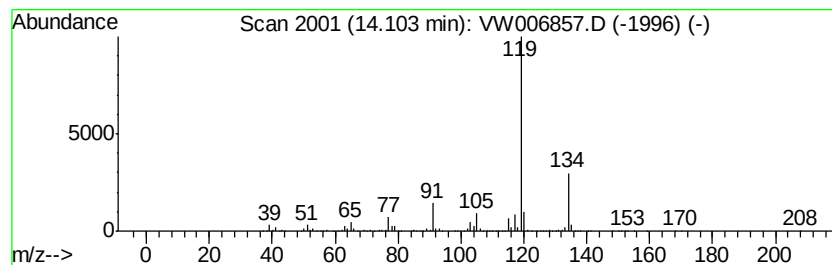
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ClientSampleId :
EP2-SB-111(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.10	38.08 ug/l	931323	1,4-Dichlorobenzene-d4	13.56	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006857.D
Acq On : 13 Nov 2018 16:28
Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

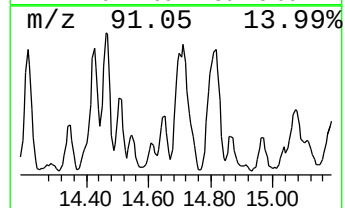
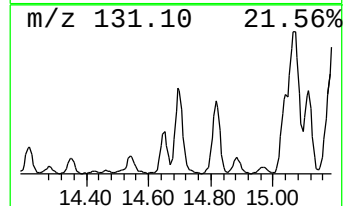
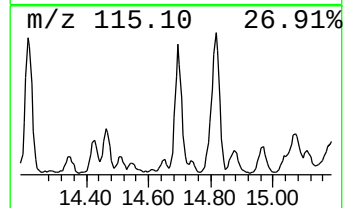
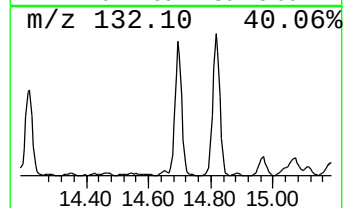
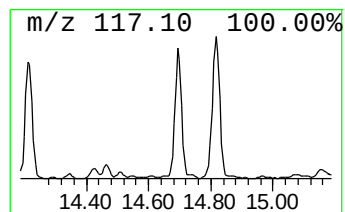
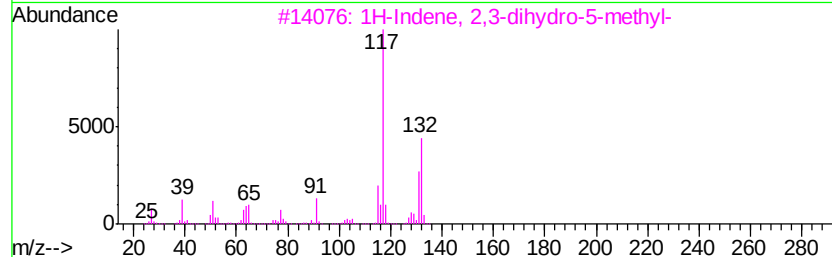
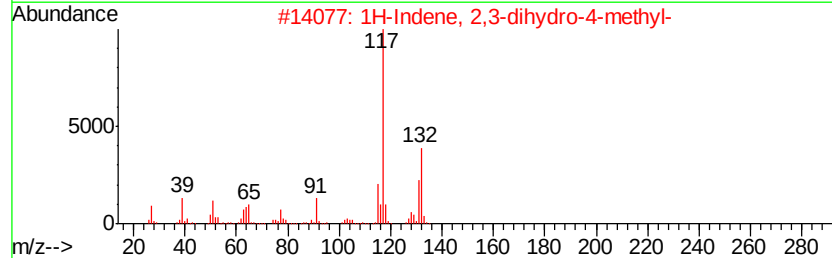
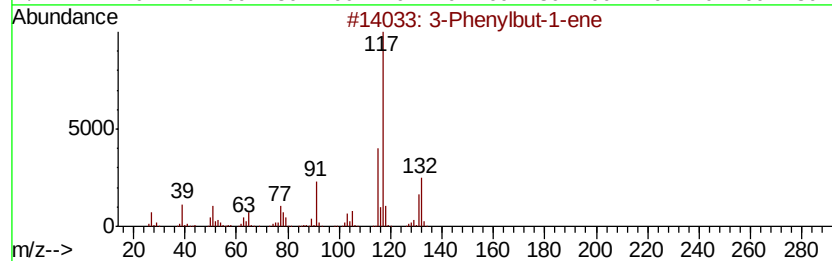
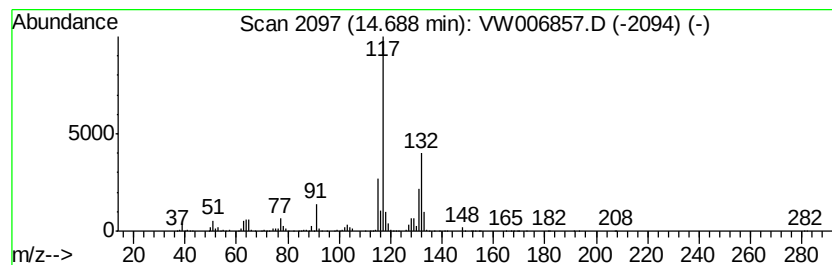
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ClientSampleId :
EP2-SB-111(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 9 3-Phenylbut-1-ene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	27.13 ug/l	663493	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	3-Phenylbut-1-ene	132 C10H12	000934-10-1	91
2	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	91
3	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	87
4	Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7	87
5	Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW111318\
Data File : VW006857.D
Acq On : 13 Nov 2018 16:28
Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EP2-SB-111(14-16)

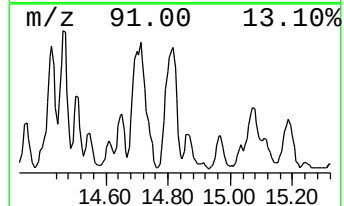
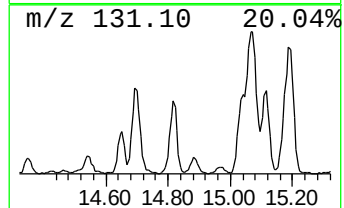
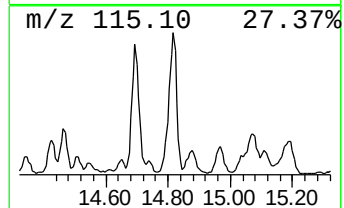
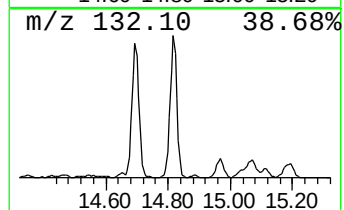
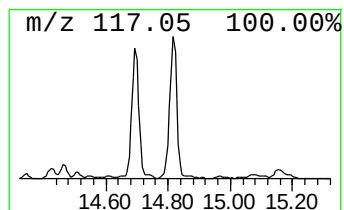
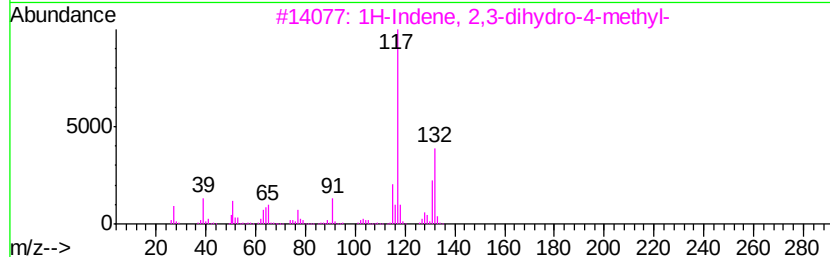
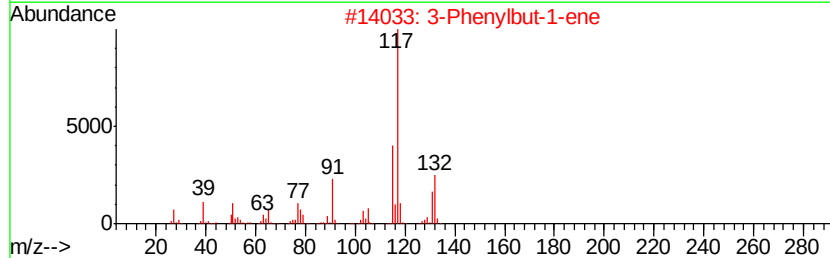
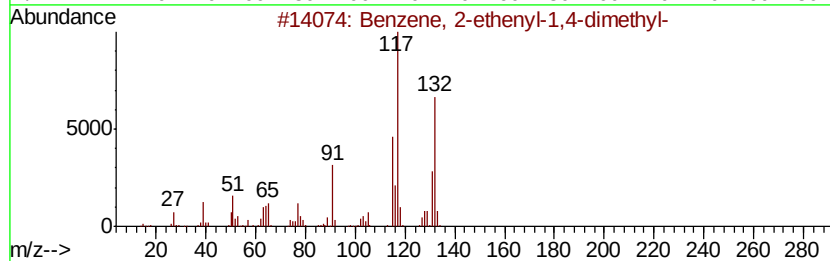
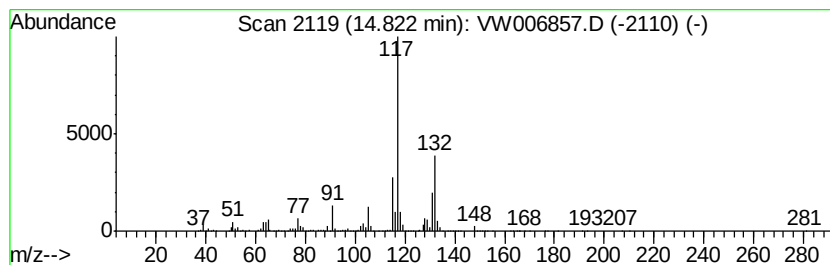
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	42.97 ug/l	1050990	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	92
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW111318\
Data File : VW006857.D
Acq On : 13 Nov 2018 16:28
Operator : SY/AP
Sample : J5860-24
Misc : 4.87G/5ML/MSVOA_W/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
EP2-SB-111(14-16)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W111218S.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
Heptane, 3-methyl-	10.10	25.9	ug/l	121608	2	8.85	234478	50.0
Octane	10.51	34.3	ug/l	212479	3	11.63	309754	50.0
Octane, 4-methyl-	11.41	64.7	ug/l	400701	3	11.63	309754	50.0
Octane, 3-methyl-	11.52	40.7	ug/l	252099	3	11.63	309754	50.0
unknown12.33	12.33	23.5	ug/l	145447	3	11.63	309754	50.0
Benzene, 1,1'-(1,...	13.76	29.3	ug/l	715595	4	13.56	1222830	50.0
Benzene, 1-ethyl-...	14.10	38.1	ug/l	931323	4	13.56	1222830	50.0
3-Phenylbut-1-ene	14.69	27.1	ug/l	663493	4	13.56	1222830	50.0
Benzene, 2-etheny...	14.82	43.0	ug/l	1050990	4	13.56	1222830	50.0