

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW042020\
 Data File : VW015229.D
 Acq On : 20 Apr 2020 18:00
 Operator : SY/VA
 Sample : L2330-02
 Misc : 5.24G/5ML/MSVOA W/SOIL
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WB-9-2

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\82W042020S.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	2.154	34	41	48	rBV2	11646	24985	3.71%	0.273%
2	2.330	65	70	80	rBV2	11167	20699	3.07%	0.226%
3	2.635	115	120	129	rBV9	5294	14351	2.13%	0.157%
4	4.129	356	365	379	rVB	7419	21301	3.16%	0.233%
5	4.391	398	408	421	rBV2	6686	18540	2.75%	0.202%
6	4.928	485	496	507	rBV2	6277	22446	3.33%	0.245%
7	5.165	528	535	547	rVB3	2634	9643	1.43%	0.105%
8	5.952	655	664	674	rBV6	3089	9444	1.40%	0.103%
9	6.903	811	820	829	rBV2	4709	15518	2.30%	0.169%
10	7.135	849	858	876	rVB2	19369	64354	9.55%	0.703%
11	7.890	972	982	986	rBV	80874	190786	28.32%	2.083%
12	7.951	986	992	1001	rVV	203671	455054	67.55%	4.968%
13	8.311	1034	1051	1062	rVV3	79440	210071	31.18%	2.293%
14	8.476	1072	1078	1089	rVB	7432	19873	2.95%	0.217%
15	8.701	1108	1115	1123	rBV3	3904	9272	1.38%	0.101%
16	8.842	1131	1138	1149	rVB	248906	495016	73.48%	5.404%
17	9.031	1162	1169	1171	rBV	7490	13504	2.00%	0.147%
18	9.073	1171	1176	1185	rVB2	12065	28700	4.26%	0.313%
19	9.335	1206	1219	1224	rBV4	8734	24977	3.71%	0.273%
20	9.756	1282	1288	1296	rVB2	7258	14922	2.22%	0.163%
21	9.896	1300	1311	1316	rBV3	7297	17601	2.61%	0.192%
22	10.097	1335	1344	1348	rBV7	3455	8082	1.20%	0.088%
23	10.323	1373	1381	1388	rVV	375276	673664	100.00%	7.355%
24	10.396	1388	1393	1396	rVV2	13574	30121	4.47%	0.329%
25	10.427	1396	1398	1404	rVV2	14300	20707	3.07%	0.226%
26	10.512	1404	1412	1418	rVB5	5855	15218	2.26%	0.166%
27	10.701	1440	1443	1448	rVV	11490	18223	2.71%	0.199%
28	10.829	1458	1464	1474	rVB	38655	75308	11.18%	0.822%
29	10.939	1476	1482	1487	rBV2	5934	12744	1.89%	0.139%
30	11.061	1490	1502	1507	rBV3	13634	33964	5.04%	0.371%
31	11.176	1517	1521	1525	rVV2	4819	7472	1.11%	0.082%
32	11.225	1525	1529	1535	rVB2	8252	13598	2.02%	0.148%
33	11.341	1544	1548	1552	rBV3	5565	9277	1.38%	0.101%
34	11.439	1559	1564	1571	rVB	17007	31190	4.63%	0.341%

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35	11.512	1572	1576	1587	rVB3	7715	18633	2.77%	0.203%
36	11.628	1588	1595	1605	rBV	306380	524822	77.91%	5.730%
37	11.731	1605	1612	1621	rBV	84021	139196	20.66%	1.520%
38	11.835	1622	1629	1641	rBV	58033	133538	19.82%	1.458%
39	12.054	1659	1665	1672	rVB	38883	67631	10.04%	0.738%
40	12.164	1672	1683	1691	rBV	76228	154763	22.97%	1.690%
41	12.249	1691	1697	1701	rVV2	9837	16682	2.48%	0.182%
42	12.341	1707	1712	1714	rBV4	11218	19905	2.95%	0.217%
43	12.463	1726	1732	1738	rBV	190042	311029	46.17%	3.396%
44	12.615	1750	1757	1766	rVB2	262911	463044	68.74%	5.055%
45	12.780	1778	1784	1792	rVB3	16388	44662	6.63%	0.488%
46	12.865	1792	1798	1801	rBV	65071	111677	16.58%	1.219%
47	12.896	1801	1803	1805	rVV2	25864	34528	5.13%	0.377%
48	12.938	1805	1810	1816	rVB3	95369	177818	26.40%	1.941%
49	13.103	1830	1837	1843	rBV	34126	78101	11.59%	0.853%
50	13.164	1843	1847	1855	rVB3	25454	40884	6.07%	0.446%
51	13.249	1856	1861	1865	rBV	96776	142805	21.20%	1.559%
52	13.298	1865	1869	1874	rVB5	15945	25352	3.76%	0.277%
53	13.377	1879	1882	1884	rBV3	10752	13580	2.02%	0.148%
54	13.444	1889	1893	1896	rBV2	25401	40166	5.96%	0.439%
55	13.560	1906	1912	1920	rVB2	252686	440008	65.32%	4.804%
56	13.621	1920	1922	1930	rVB5	15685	24818	3.68%	0.271%
57	13.713	1930	1937	1939	rBV6	16423	40792	6.06%	0.445%
58	13.755	1939	1944	1948	rBV2	77028	133209	19.77%	1.454%
59	13.871	1958	1963	1968	rBV2	140435	231047	34.30%	2.522%
60	13.944	1971	1975	1978	rVB2	30791	53895	8.00%	0.588%
61	14.011	1983	1986	1987	rBV	17193	18554	2.75%	0.203%
62	14.091	1996	1999	2004	rVB	54346	81757	12.14%	0.893%
63	14.133	2004	2006	2012	rVB3	20240	31273	4.64%	0.341%
64	14.206	2012	2018	2023	rBV3	55765	104862	15.57%	1.145%
65	14.261	2023	2027	2034	rBV7	16503	40028	5.94%	0.437%
66	14.334	2034	2039	2043	rBV5	23255	49928	7.41%	0.545%
67	14.383	2044	2047	2050	rVV2	53447	82847	12.30%	0.904%
68	14.420	2050	2053	2056	rVV2	48809	75886	11.26%	0.828%
69	14.456	2056	2059	2063	rVV	51024	67836	10.07%	0.741%
70	14.499	2063	2066	2069	rVV2	35494	43883	6.51%	0.479%
71	14.542	2069	2073	2081	rVB7	38330	79731	11.84%	0.870%

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72	14.688	2093	2097	2099	rBV2	24628	36887	5.48%	0.403%
73	14.725	2099	2103	2109	rVB2	107873	158731	23.56%	1.733%
74	14.798	2109	2115	2120	rBV3	71922	183587	27.25%	2.004%
75	14.956	2136	2141	2148	rVB3	61983	112224	16.66%	1.225%
76	15.066	2154	2159	2167	rBV3	55371	123127	18.28%	1.344%
77	15.170	2171	2176	2184	rVB3	57025	139340	20.68%	1.521%
78	15.255	2186	2190	2196	rVB5	15266	35705	5.30%	0.390%
79	15.328	2197	2202	2203	rBV3	37429	56764	8.43%	0.620%
80	15.365	2203	2208	2215	rVB	363405	623557	92.56%	6.808%
81	15.554	2235	2239	2248	rVB3	61564	120179	17.84%	1.312%
82	15.657	2248	2256	2262	rBV3	31588	85964	12.76%	0.939%
83	15.950	2300	2304	2308	rVB4	16139	26367	3.91%	0.288%
84	16.169	2332	2340	2345	rBV6	24389	70061	10.40%	0.765%
85	16.285	2356	2359	2364	rVB4	16094	26397	3.92%	0.288%
86	16.371	2369	2373	2379	rVV4	30208	70693	10.49%	0.772%
87	16.438	2379	2384	2390	rVV	66670	149811	22.24%	1.636%
88	16.493	2390	2393	2399	rVB3	29433	50051	7.43%	0.546%
89	16.645	2410	2418	2430	rVB	112822	280271	41.60%	3.060%

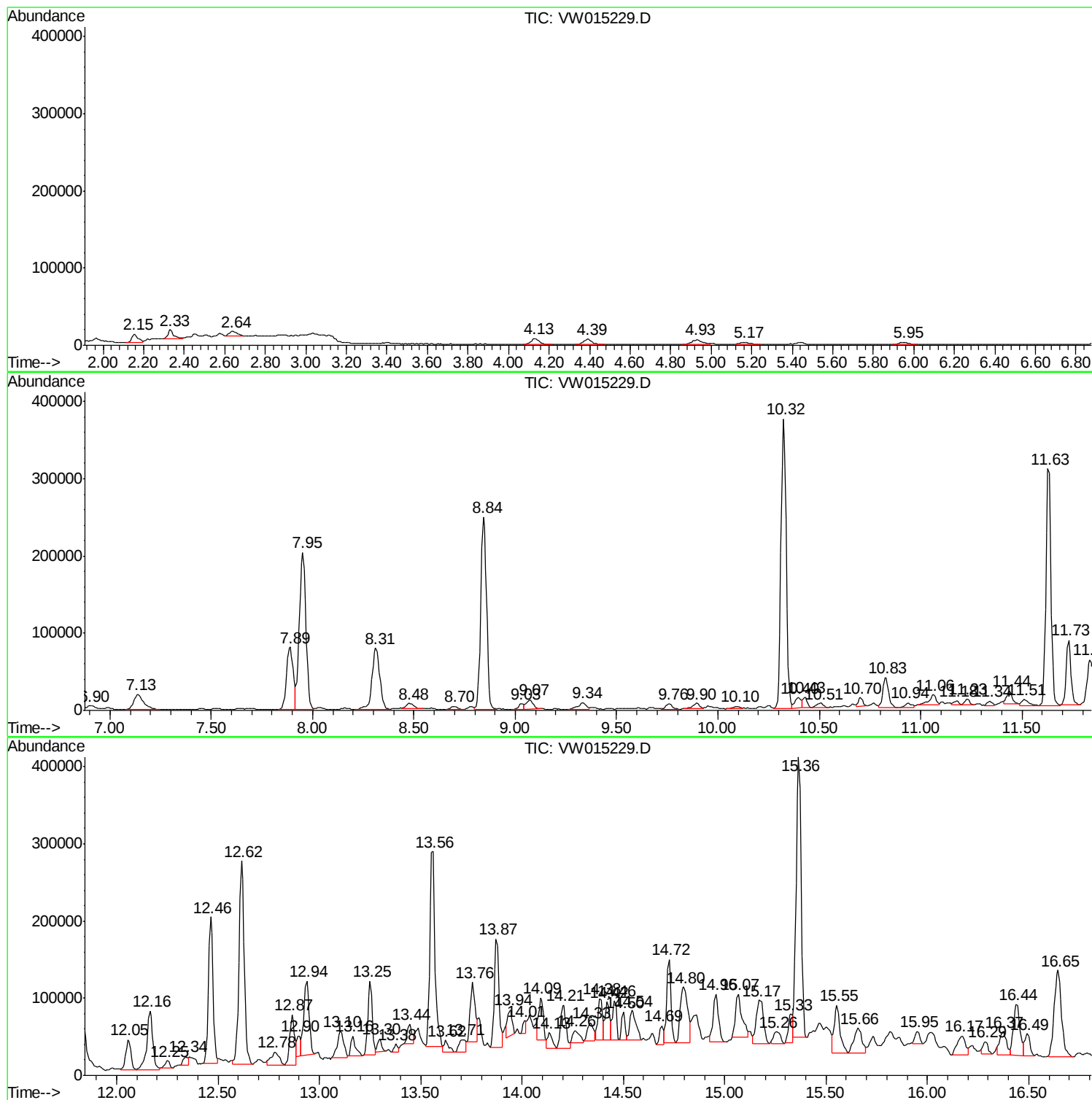
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W042020S.M
Quant Title : SW846 8260

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



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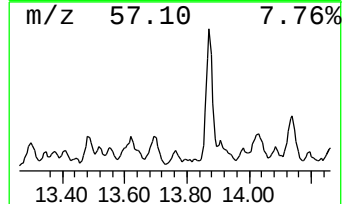
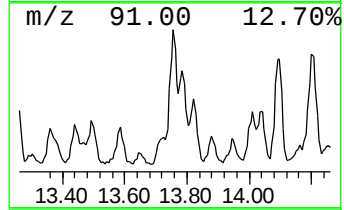
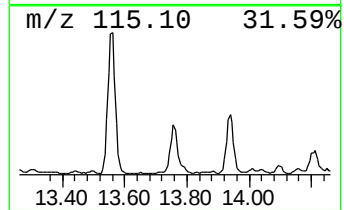
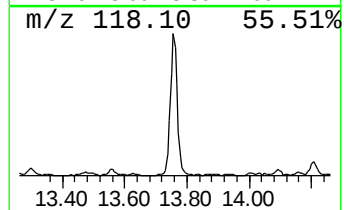
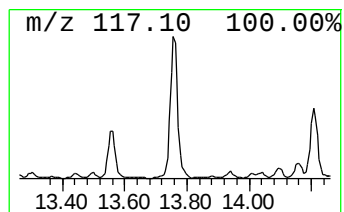
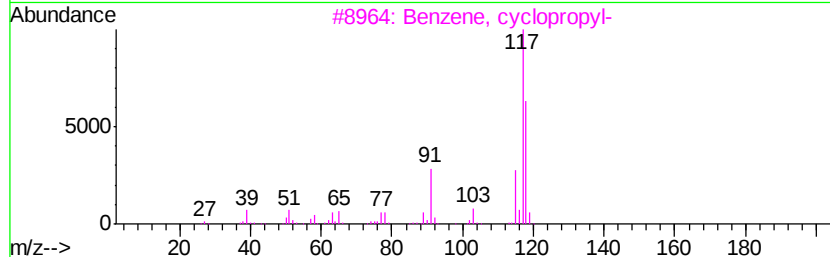
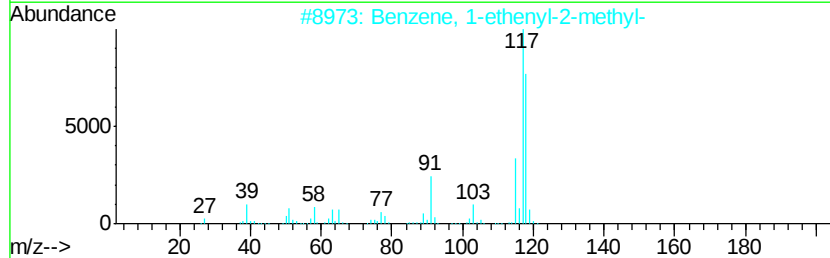
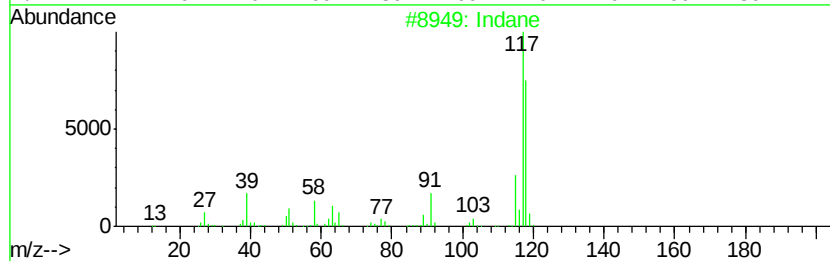
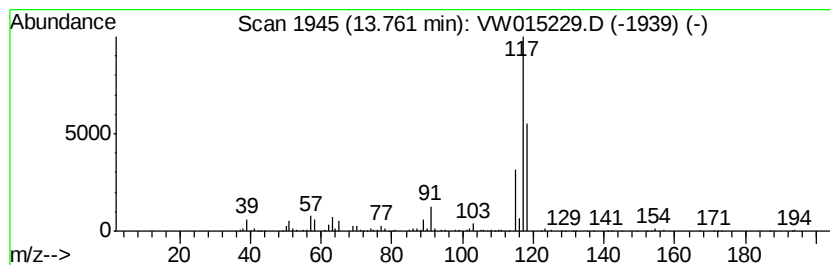
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W042020S.M
Quant Title : SW846 8260

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

Peak Number 1 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.76	15.14 ug/l	133209	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	81
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76
3		Benzene, cyclopropyl-	118	C9H10	000873-49-4	68
4		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	64
5		Benzene, 2-propenyl-	118	C9H10	000300-57-2	58



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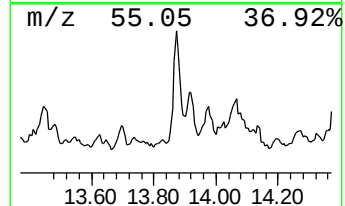
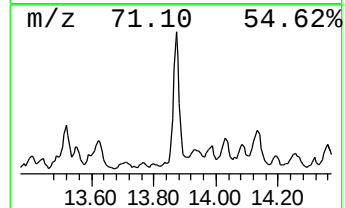
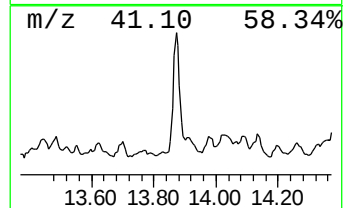
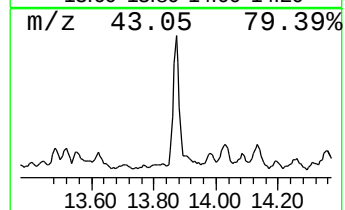
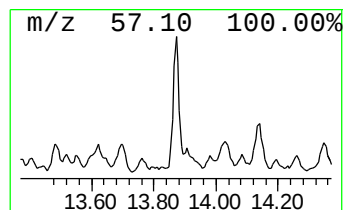
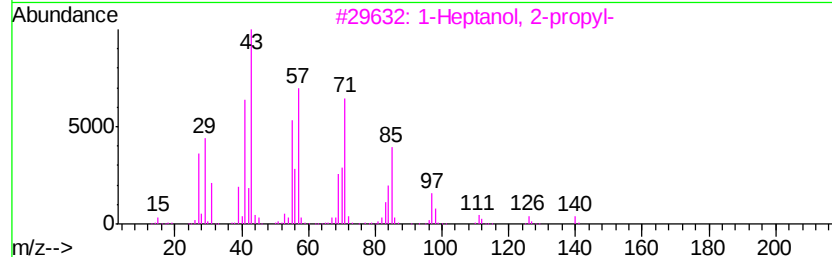
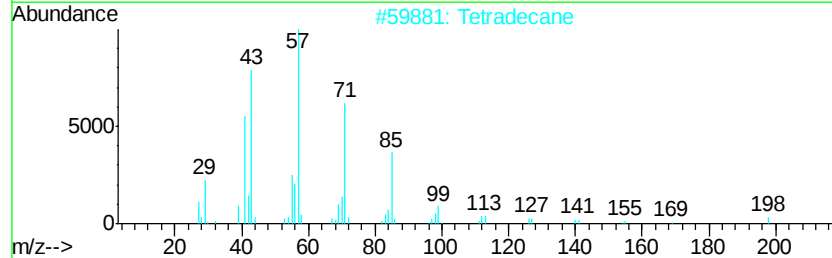
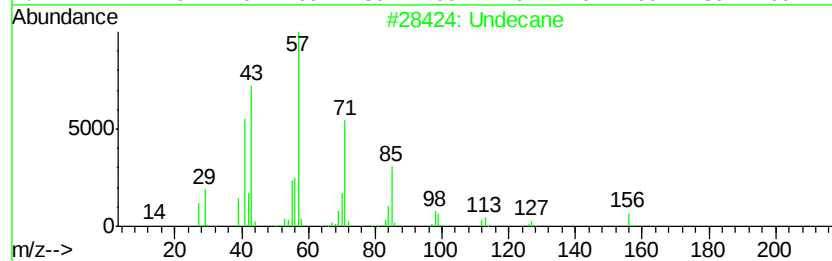
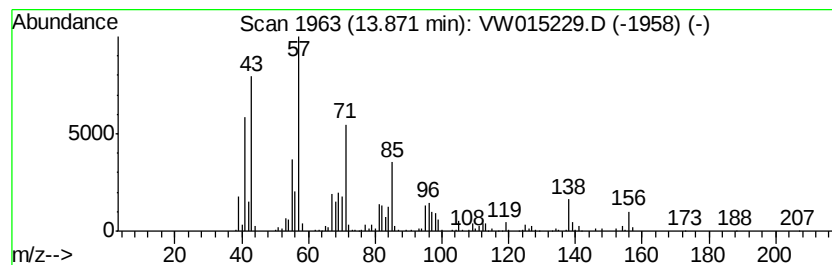
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	26.25 ug/l	231047	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	96
2	Tetradecane		198	C14H30	000629-59-4	52
3	1-Heptanol, 2-propyl-		158	C10H22O	010042-59-8	49
4	Decane, 2,4,6-trimethyl-		184	C13H28	062108-27-4	49
5	Undecane, 4,6-dimethyl-		184	C13H28	017312-82-2	47



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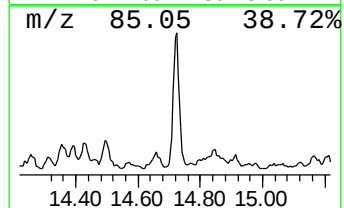
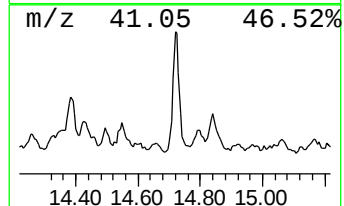
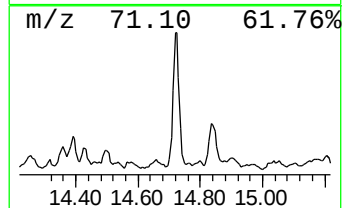
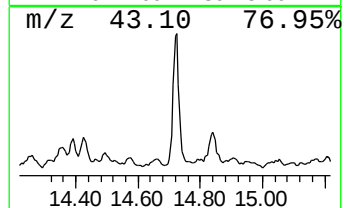
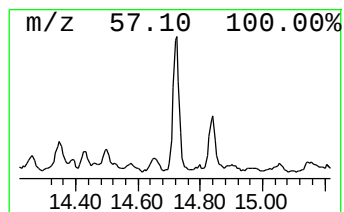
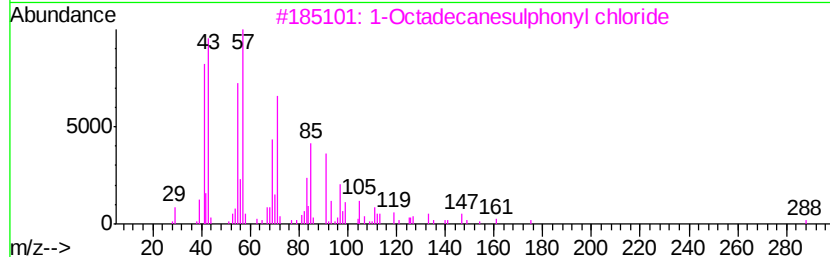
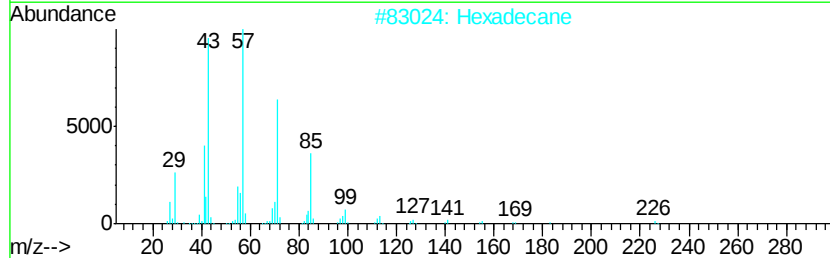
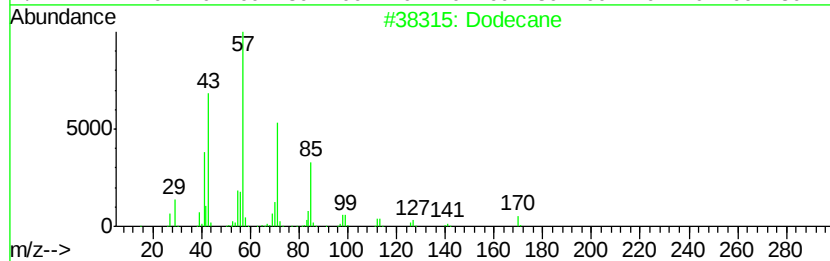
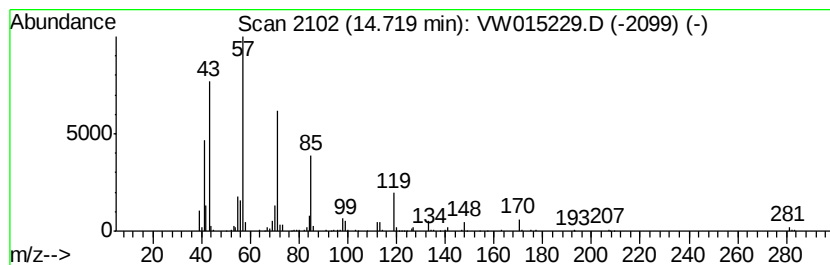
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Quant Title : SW846 8260

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

Peak Number 3 Dodecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.72	18.04 ug/l	158731	1,4-Dichlorobenzene-d4	13.56

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane		170	C12H26	000112-40-3	96
2	Hexadecane		226	C16H34	000544-76-3	58
3	1-Octadecanesulphonvl chloride		352	C18H37ClO2S	1000342-70-4	58
4	Tridecane		184	C13H28	000629-50-5	58
5	Bacchotricuneatin c		342	C20H22O5	066563-30-2	58



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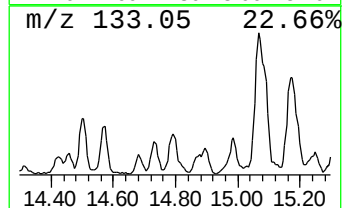
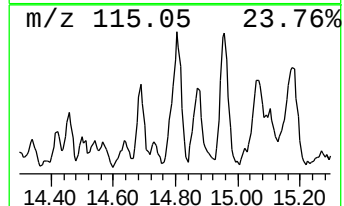
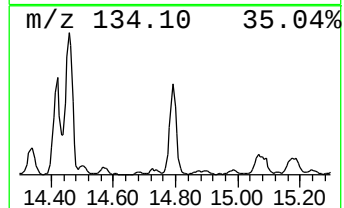
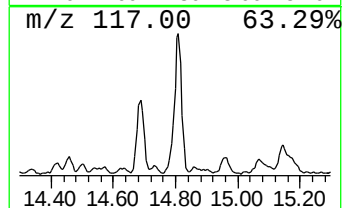
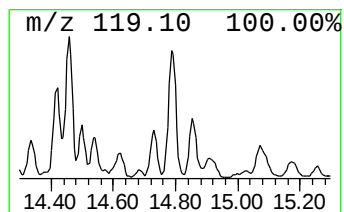
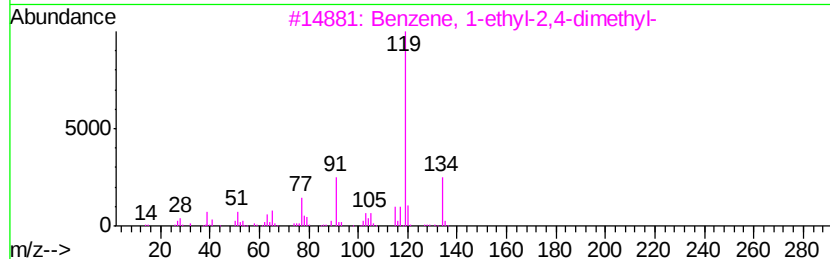
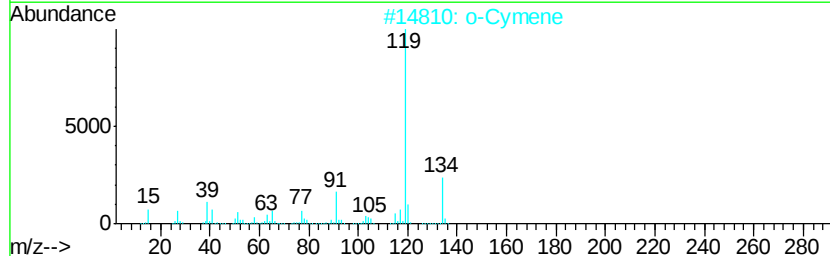
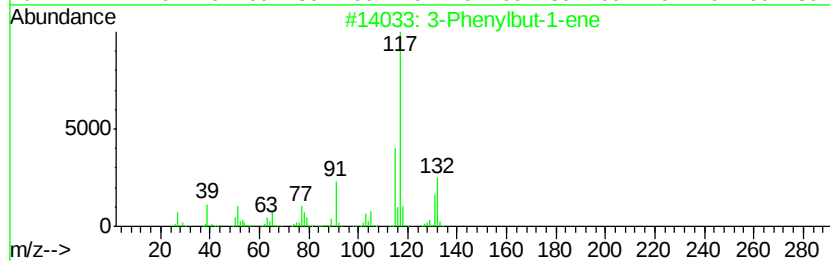
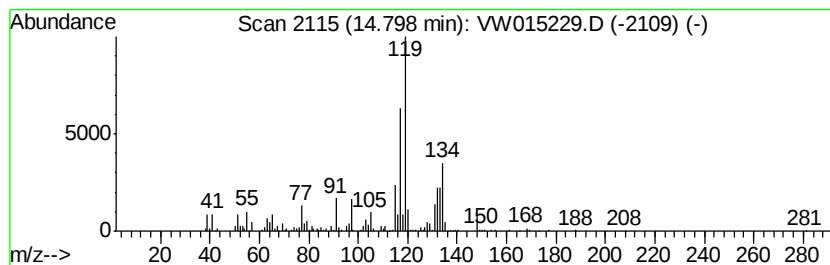
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Quant Title : SW846 8260

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

Peak Number 4 3-Phenylbut-1-ene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	20.86 ug/l	183587	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	83
2	o-Cymene		134	C10H14	000527-84-4	50
3	Benzene, 1-ethyl-2,4-dimethyl-		134	C10H14	000874-41-9	50
4	Benzene, 2-ethyl-1,3-dimethyl-		134	C10H14	002870-04-4	50
5	Benzene, 1-ethyl-2,3-dimethyl-		134	C10H14	000933-98-2	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW042020\
Data File : VW015229.D
Acq On : 20 Apr 2020 18:00
Operator : SY/VA
Sample : L2330-02
Misc : 5.24G/5ML/MSVOA W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
WB-9-2

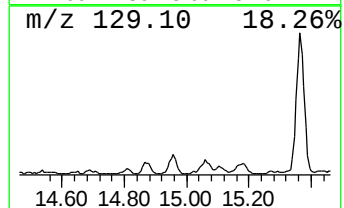
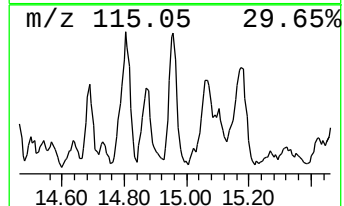
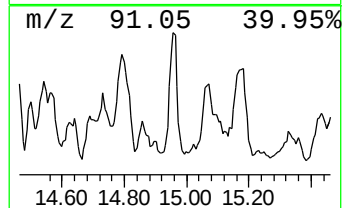
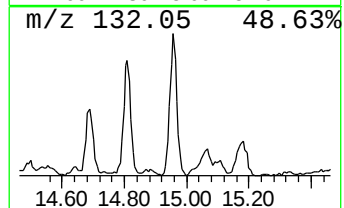
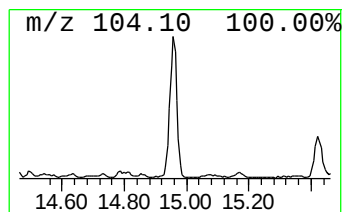
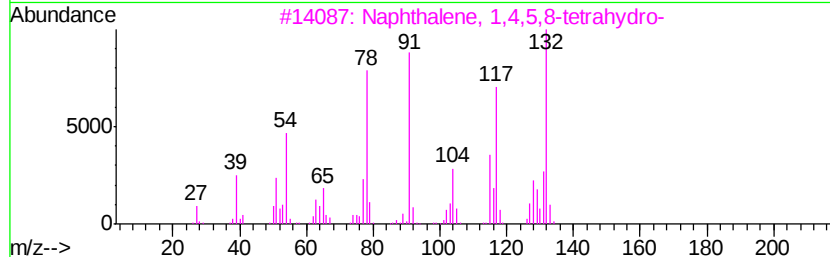
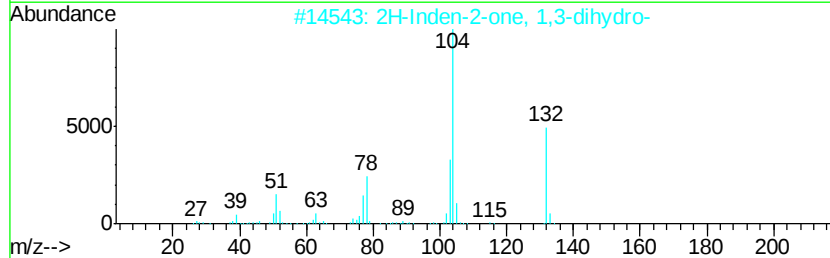
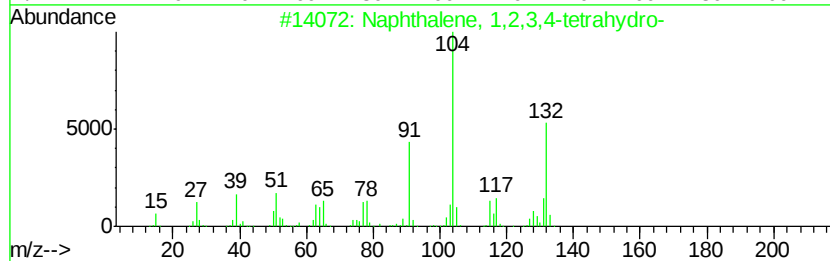
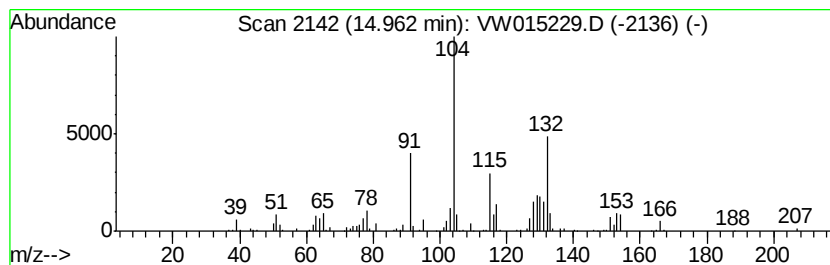
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	12.75 ug/l	112224	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	43
3		Naphthalene, 1,4,5,8-tetrahydro-	132	C10H12	000493-04-9	25
4		Bromoacetic acid, 2-phenylethyl ...	242	C10H11BrO2	003785-33-9	22
5		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW042020\
Data File : VW015229.D
Acq On : 20 Apr 2020 18:00
Operator : SY/VA
Sample : L2330-02
Misc : 5.24G/5ML/MSVOA W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
WB-9-2

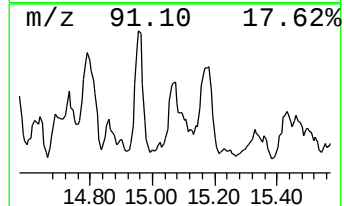
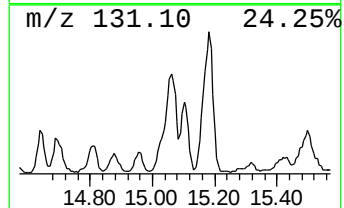
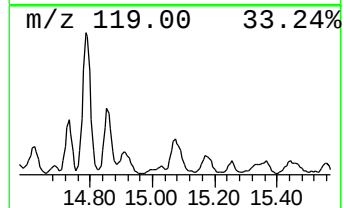
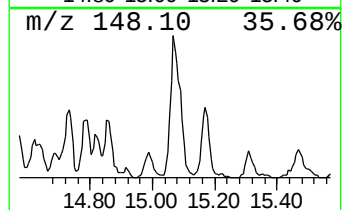
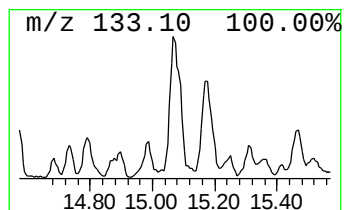
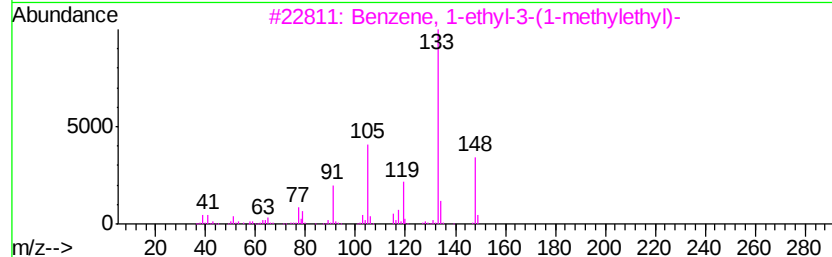
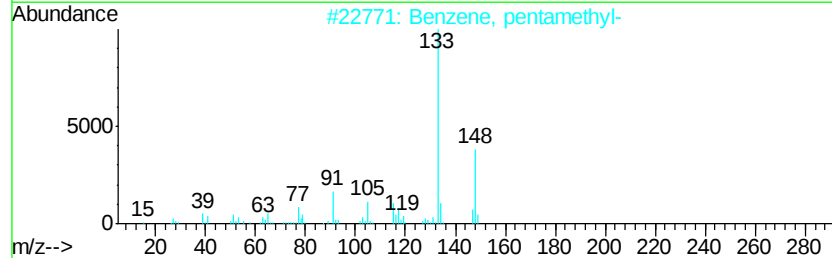
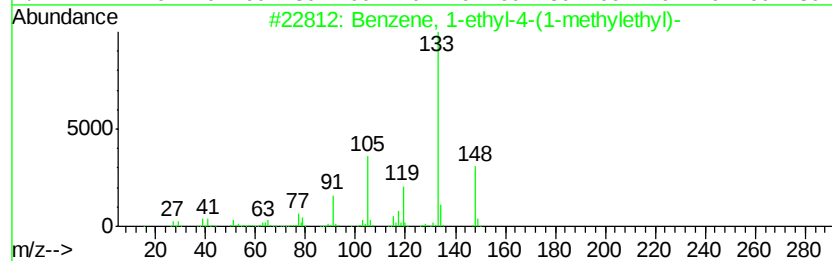
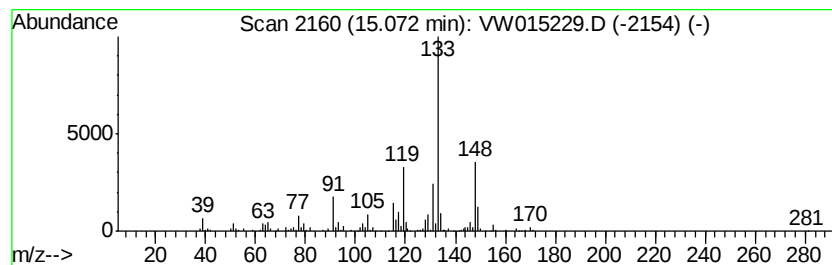
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	13.99 ug/l	123127	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	74
2		Benzene, pentamethyl-	148	C11H16	000700-12-9	64
3		Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	64
4		Phenol, 2-(2-methyl-2-propenyl)-	148	C10H12O	020944-88-1	53
5		4-tert-Butyltoluene	148	C11H16	000098-51-1	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW042020\
Data File : VW015229.D
Acq On : 20 Apr 2020 18:00
Operator : SY/VA
Sample : L2330-02
Misc : 5.24G/5ML/MSVOA W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
WB-9-2

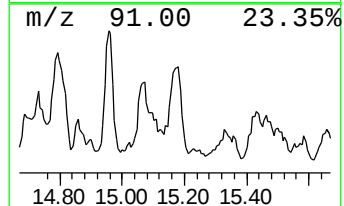
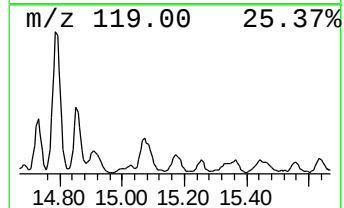
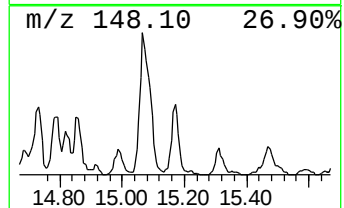
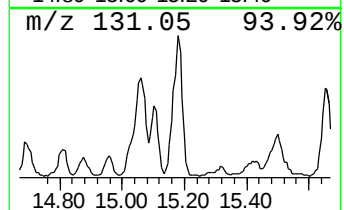
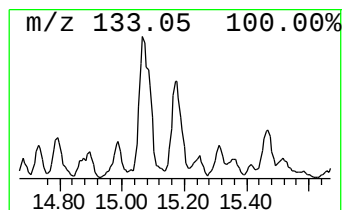
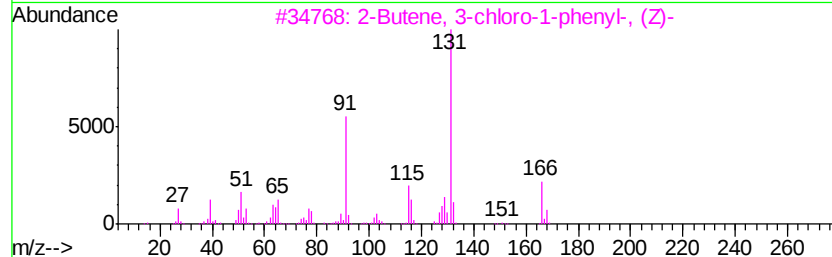
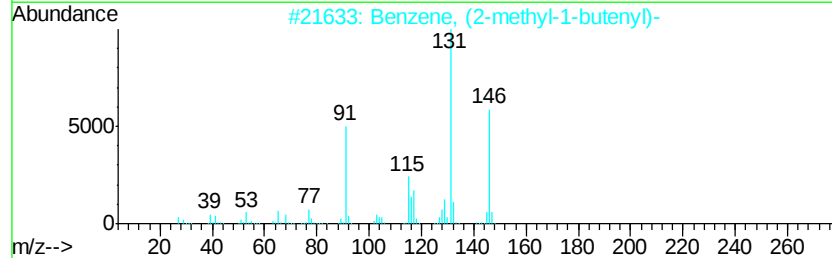
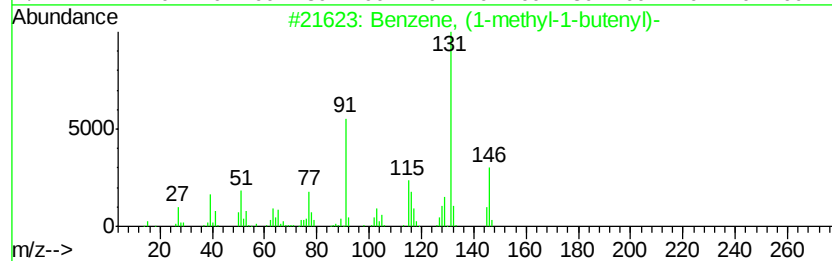
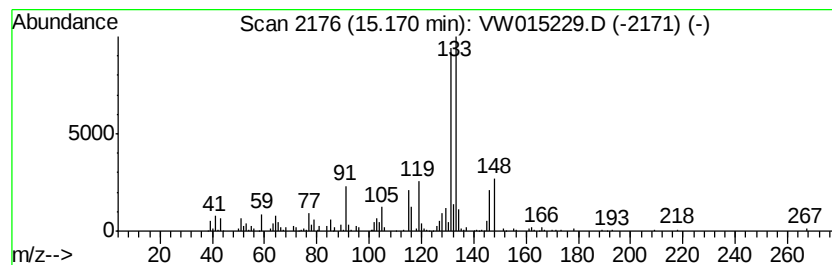
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W042020S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, (1-methyl-1-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	15.83 ug/l	139340	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	86
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
3		2-Butene, 3-chloro-1-phenyl-, (Z)-	166	C10H11Cl	016608-68-7	52
4		Benzene, (2-chloro-2-butenyl)-	166	C10H11Cl	054411-12-0	52
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW042020\
Data File : VW015229.D
Acq On : 20 Apr 2020 18:00
Operator : SY/VA
Sample : L2330-02
Misc : 5.24G/5ML/MSVOA W/SOIL
ALS Vial : 16 Sample Multiplier: 1

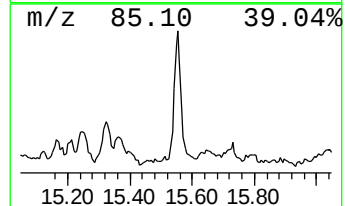
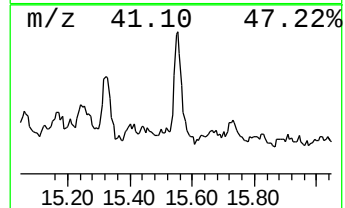
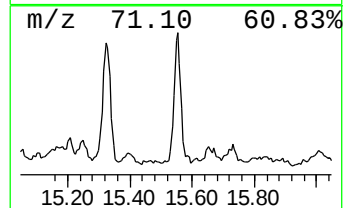
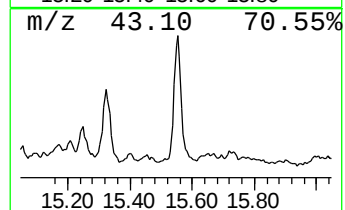
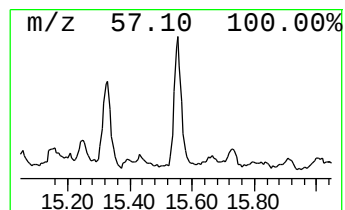
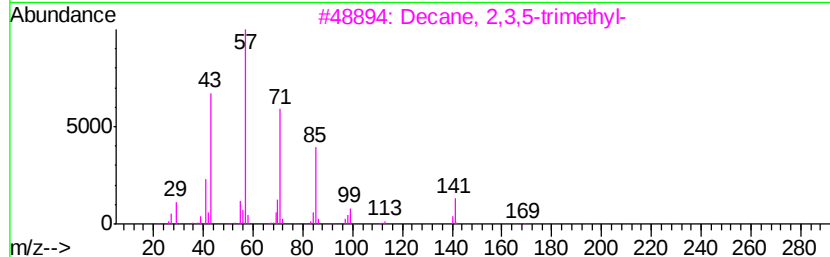
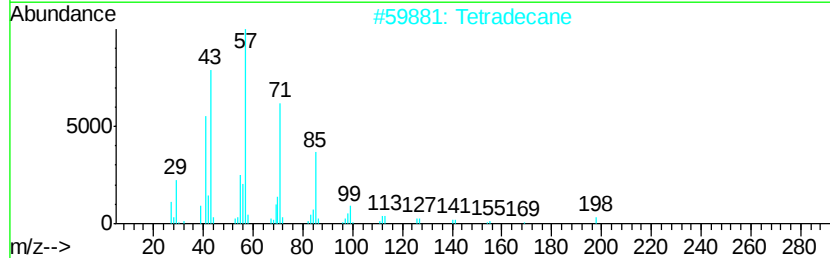
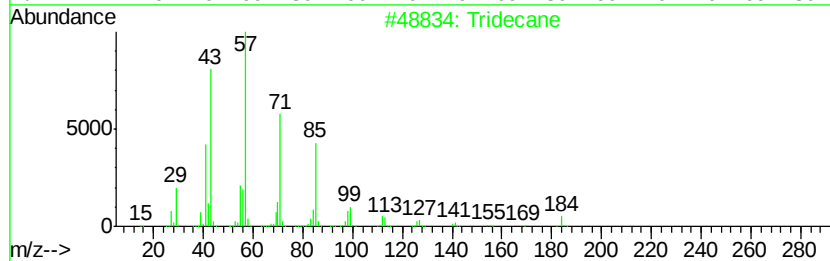
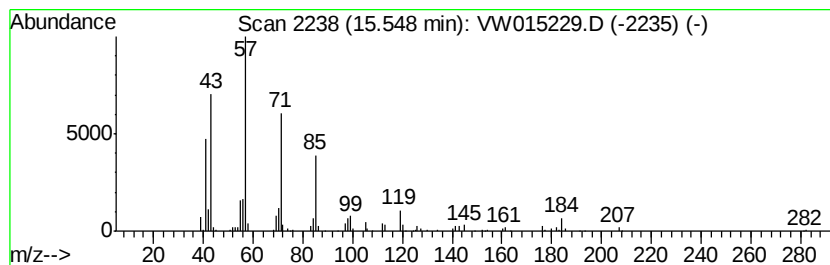
Instrument :
MSVOA_W
ClientSampleId :
WB-9-2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W042020S.M
Quant Title : SW846 8260

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

Peak Number 8 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	13.66 ug/l	120179	1,4-Dichlorobenzene-d4	13.56
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Tridecane	184 C13H28	000629-50-5	98
2	Tetradecane	198 C14H30	000629-59-4	74
3	Decane, 2,3,5-trimethyl-	184 C13H28	062238-11-3	64
4	Decane, 6-ethyl-2-methyl-	184 C13H28	062108-21-8	59
5	Bacchotricuneatin c	342 C20H22O5	066563-30-2	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA W\DATA\VW042020\
Data File : VW015229.D
Acq On : 20 Apr 2020 18:00
Operator : SY/VA
Sample : L2330-02
Misc : 5.24G/5ML/MSVOA W/SOIL
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
WB-9-2

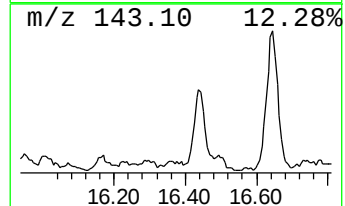
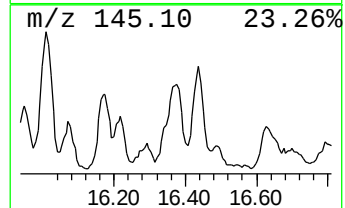
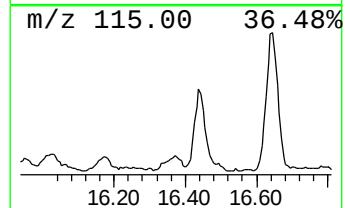
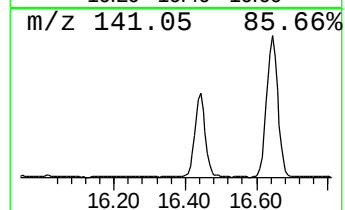
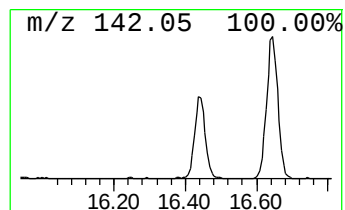
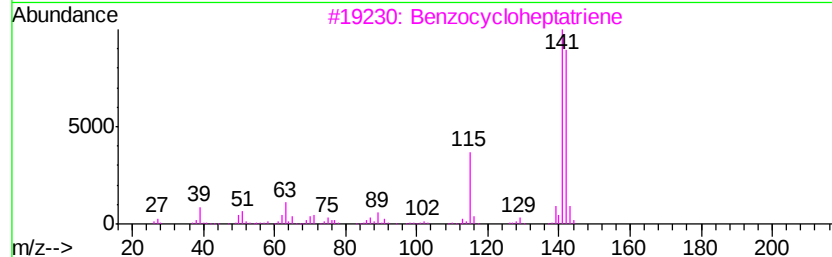
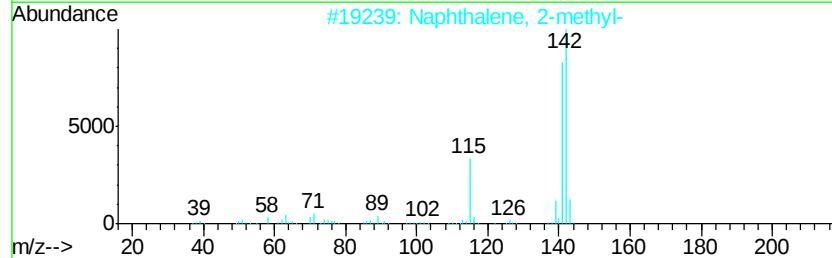
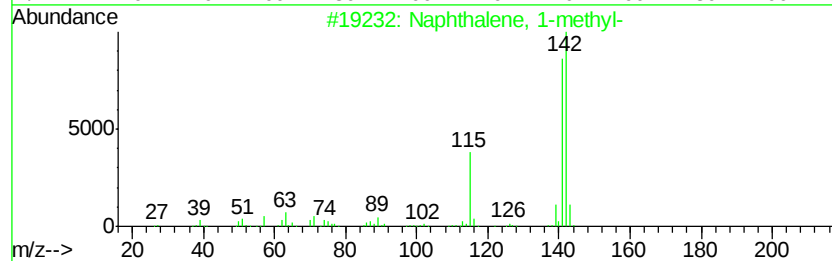
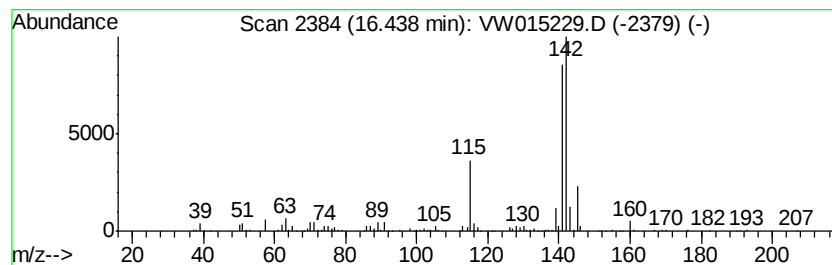
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W042020S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.44	17.02 ug/l	149811	1,4-Dichlorobenzene-d4	13.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	90
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_W\DATA\VW042020\
 Data File : VW015229.D
 Acq On : 20 Apr 2020 18:00
 Operator : SY/VA
 Sample : L2330-02
 Misc : 5.24G/5ML/MSVOA_W/SOIL
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 WB-9-2

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W042020S.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
Indane	13.76	15.1	ug/l	133209	4	13.56	440008	50.0
Undecane	13.87	26.3	ug/l	231047	4	13.56	440008	50.0
Dodecane	14.72	18.0	ug/l	158731	4	13.56	440008	50.0
3-Phenylbut-1-ene	14.80	20.9	ug/l	183587	4	13.56	440008	50.0
Naphthalene, 1,2,...	14.96	12.8	ug/l	112224	4	13.56	440008	50.0
Benzene, 1-ethyl-...	15.07	14.0	ug/l	123127	4	13.56	440008	50.0
Benzene, (1-methy...	15.17	15.8	ug/l	139340	4	13.56	440008	50.0
Tridecane	15.55	13.7	ug/l	120179	4	13.56	440008	50.0
Naphthalene, 1-me...	16.44	17.0	ug/l	149811	4	13.56	440008	50.0