

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092222\
 Data File : VX031580.D
 Acq On : 22 Sep 2022 19:04
 Operator : JC/MD
 Sample : N4614-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10R

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 >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
 Title : SW846 8260

Signal : TIC: VX031580.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.618	77	87	88	rBV3	3876	6872	1.08%	0.203%
2	2.867	286	292	301	rVB4	3126	6831	1.08%	0.201%
3	2.983	301	311	324	rBV3	6569	20381	3.21%	0.601%
4	3.117	324	333	342	rVB	24254	56207	8.86%	1.658%
5	4.306	519	528	537	rBV2	4224	11240	1.77%	0.332%
6	5.391	692	706	716	rBV	73725	215999	34.07%	6.371%
7	5.556	724	733	749	rVB	69513	197834	31.20%	5.835%
8	5.958	788	799	806	rBV2	81818	212978	33.59%	6.282%
9	6.044	806	813	825	rVB	45894	125150	19.74%	3.691%
10	6.763	921	931	942	rBV2	108668	254106	40.08%	7.495%
11	8.647	1234	1240	1249	rBV	392053	634044	100.00%	18.701%
12	8.958	1284	1291	1299	rVB2	6789	11009	1.74%	0.325%
13	9.049	1299	1306	1311	rBV	5288	8269	1.30%	0.244%
14	9.458	1367	1373	1379	rBV3	7306	11569	1.82%	0.341%
15	9.945	1448	1453	1460	rBV	9035	13960	2.20%	0.412%
16	10.055	1465	1471	1479	rVV	238053	317818	50.13%	9.374%
17	10.195	1489	1494	1501	rBV	26949	33966	5.36%	1.002%
18	10.305	1507	1512	1516	rBV	15129	19797	3.12%	0.584%
19	10.963	1615	1620	1626	rBV	23005	27610	4.35%	0.814%
20	11.079	1634	1639	1652	rBV	326857	424980	67.03%	12.535%
21	11.305	1671	1676	1682	rBV	34143	42296	6.67%	1.248%
22	11.402	1685	1692	1696	rBV2	5912	9067	1.43%	0.267%
23	11.756	1745	1750	1754	rBV	28125	35228	5.56%	1.039%
24	12.024	1788	1794	1801	rBV	243037	300264	47.36%	8.856%
25	12.085	1801	1804	1811	rVB	10968	14145	2.23%	0.417%
26	12.244	1825	1830	1837	rVB	90448	111246	17.55%	3.281%
27	12.317	1837	1842	1851	rVV2	5780	9110	1.44%	0.269%
28	12.610	1886	1890	1893	rBV	10611	13195	2.08%	0.389%
29	12.646	1893	1896	1900	rVB	9872	11220	1.77%	0.331%
30	12.701	1900	1905	1912	rBV	35684	48196	7.60%	1.422%
31	12.920	1935	1941	1945	rVV	17036	21341	3.37%	0.629%
32	12.963	1945	1948	1955	rVB2	7471	9416	1.49%	0.278%
33	13.140	1967	1977	1979	rBV5	4740	6624	1.04%	0.195%
34	13.170	1979	1982	1990	rVB	8867	10800	1.70%	0.319%
35	13.292	1994	2002	2008	rBV2	52994	68637	10.83%	2.024%
36	13.420	2020	2023	2028	rBV3	6183	7291	1.15%	0.215%

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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

37	13.585	2047	2050	2055	rVV4	5718	8434	1.33%	0.249%
38	13.664	2061	2063	2069	rVB2	7318	8858	1.40%	0.261%
39	13.774	2075	2081	2090	rVB	24548	32232	5.08%	0.951%
40	14.780	2240	2246	2252	rBV	9631	12132	1.91%	0.358%

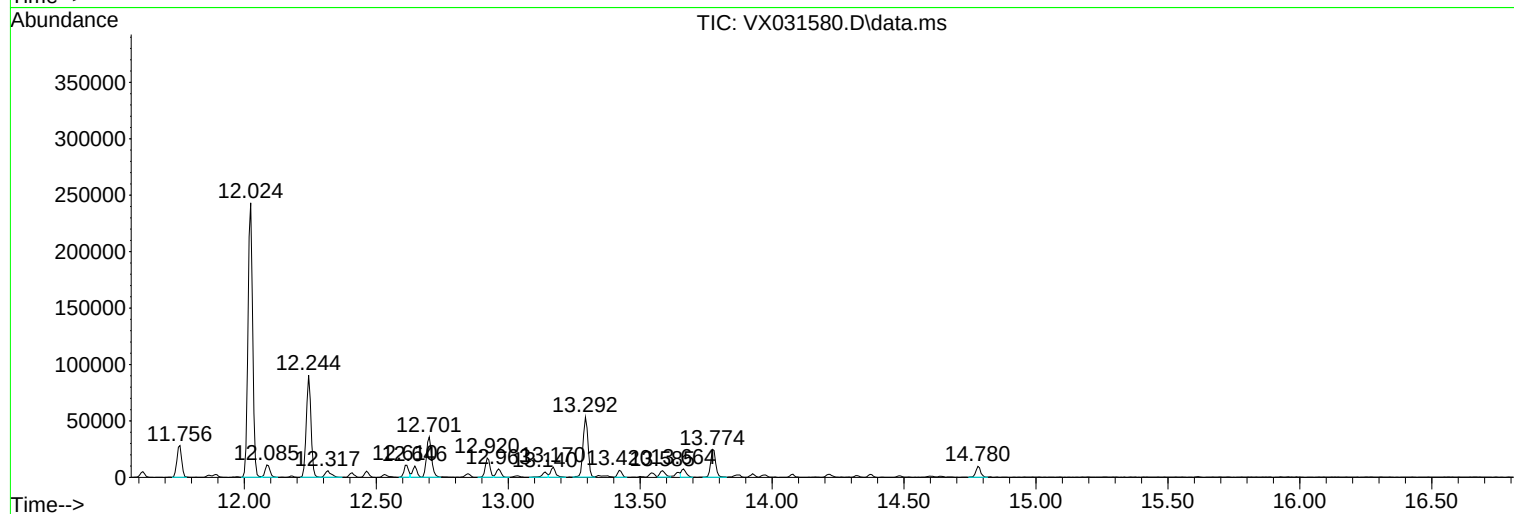
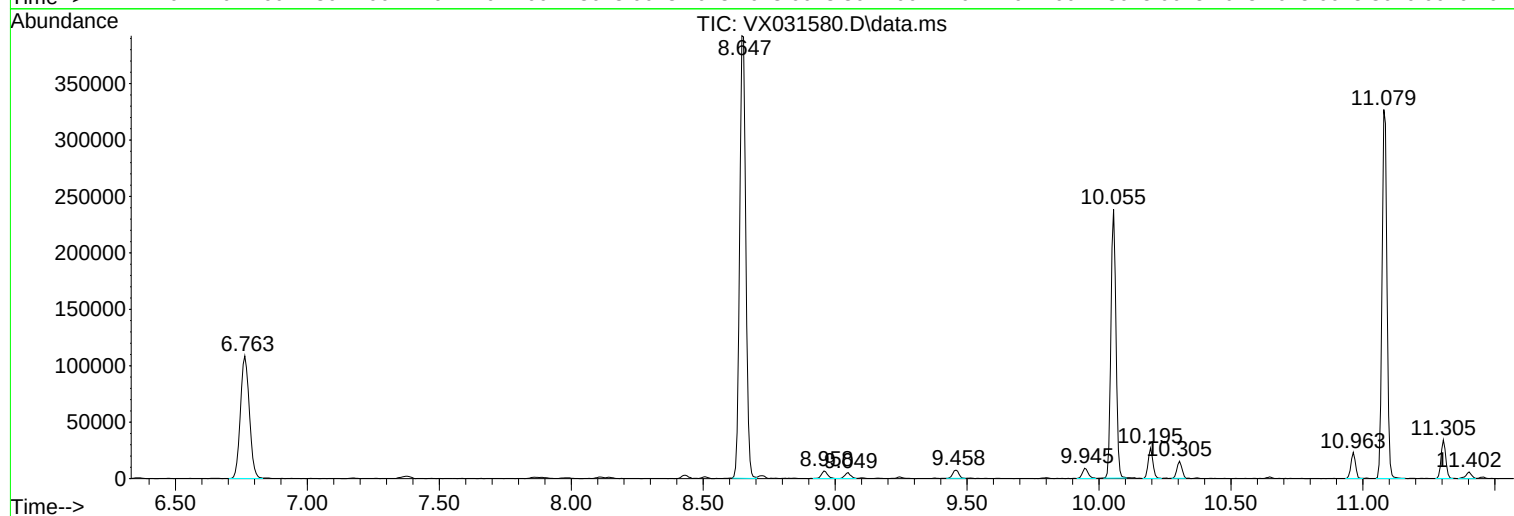
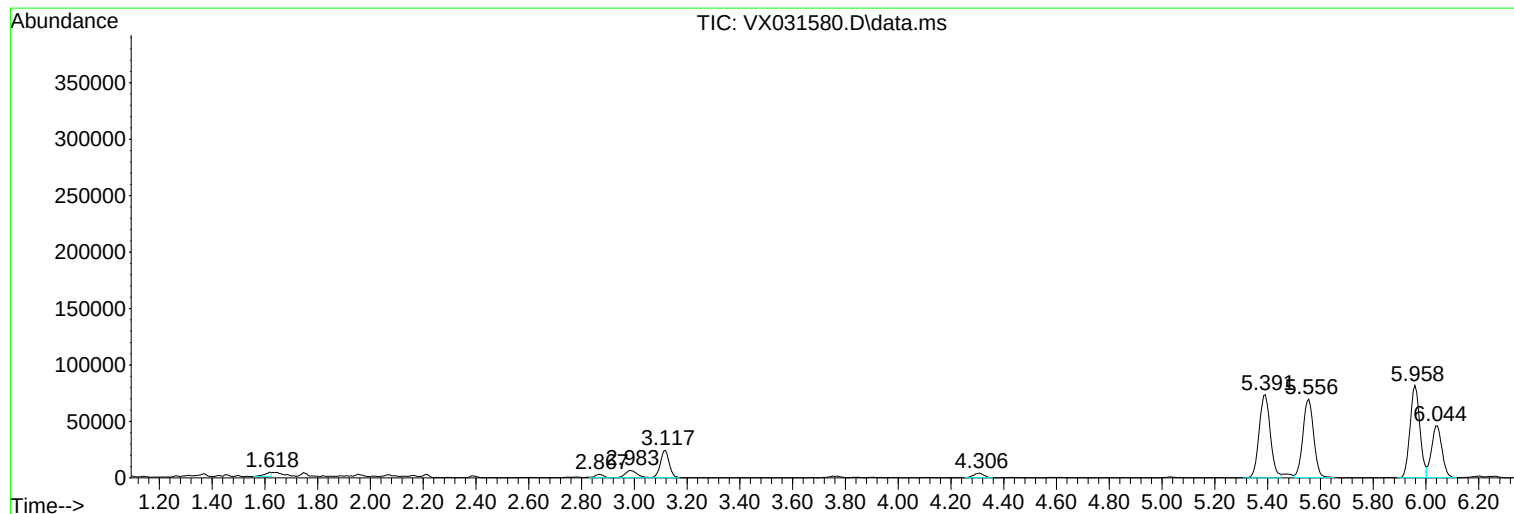
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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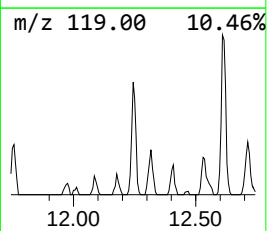
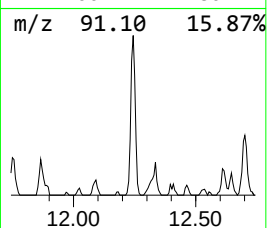
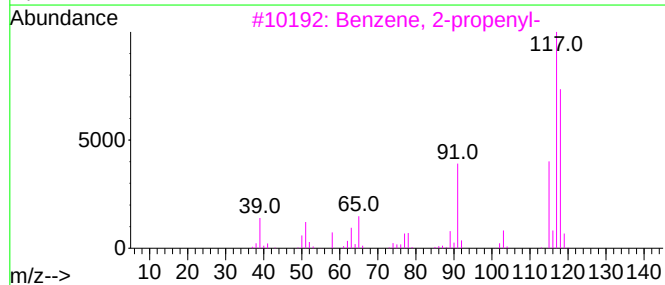
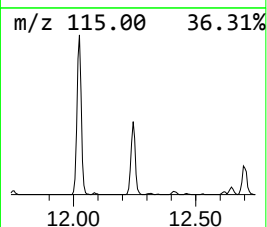
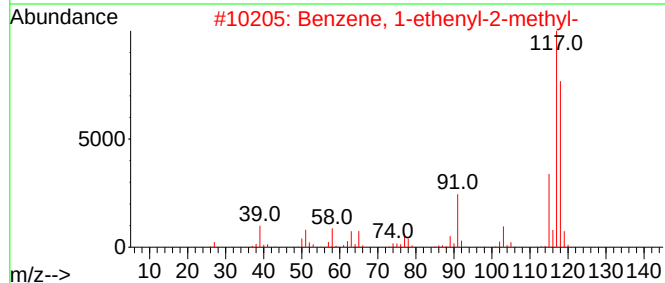
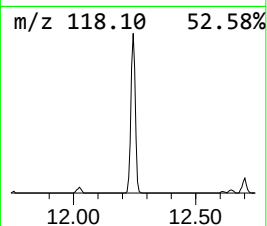
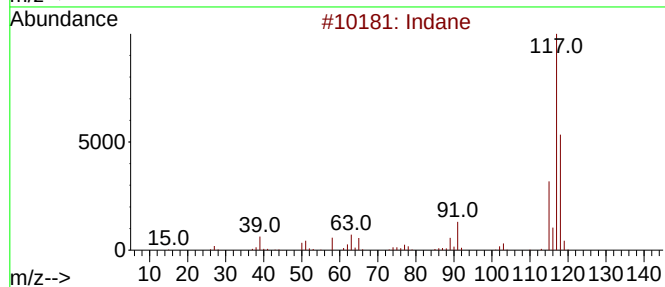
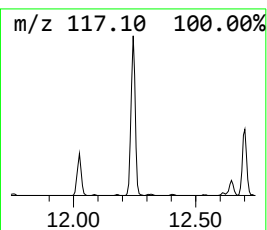
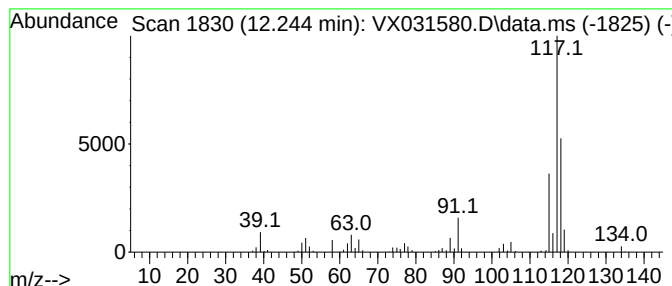
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	18.52 ug/l	111246	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
4			Delta-cyclene	118	C9H10	007785-10-6	68
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	58



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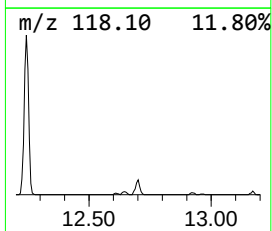
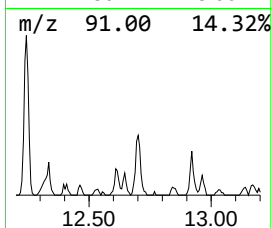
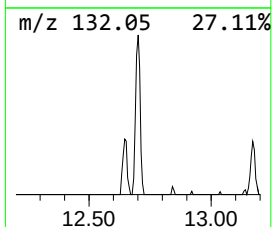
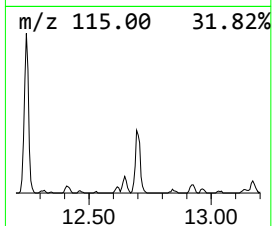
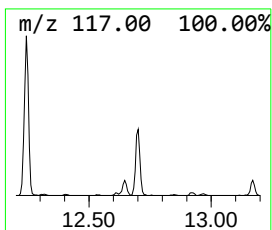
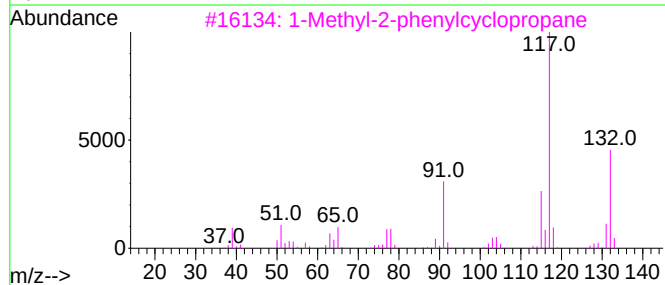
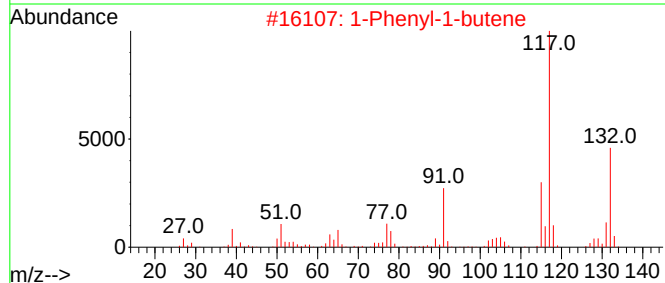
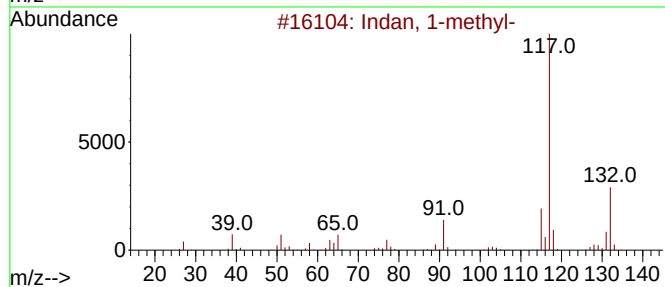
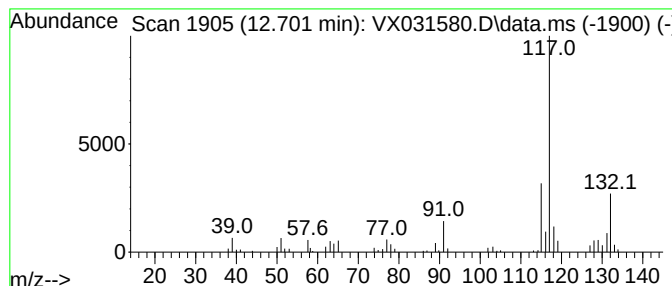
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Peak Number 2 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	8.03 ug/l	48196	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	86
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86



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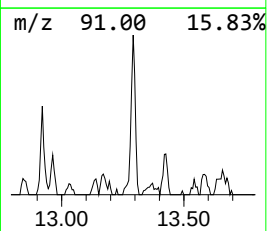
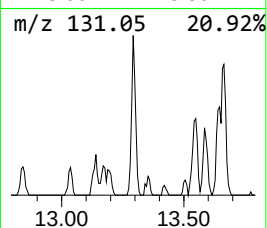
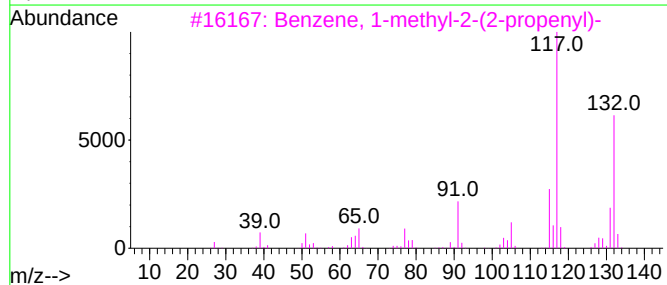
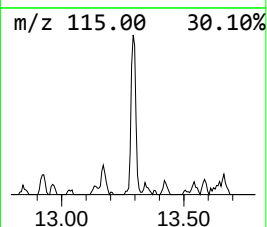
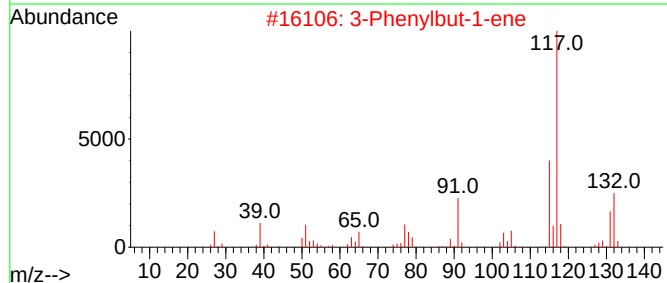
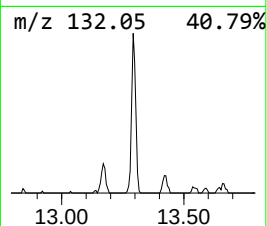
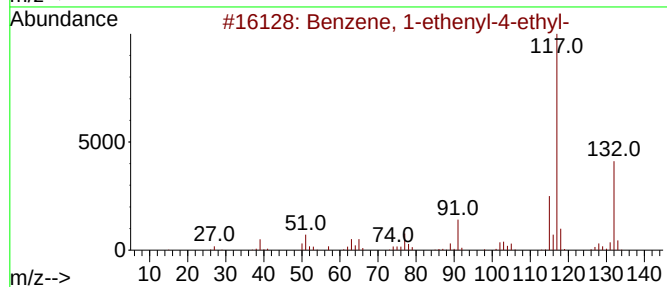
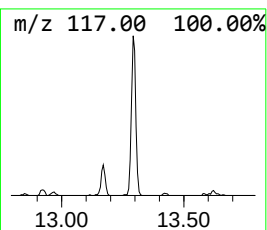
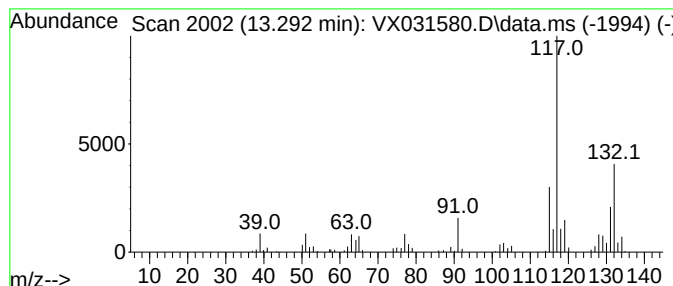
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	11.43 ug/l	68637	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87



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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	12.244	18.5	ug/l	111246	4	12.024	300264	50.0
Indan, 1-methyl-	12.701	8.0	ug/l	48196	4	12.024	300264	50.0
Benzene, 1-ethe...	13.292	11.4	ug/l	68637	4	12.024	300264	50.0