

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX021119\
 Data File : VX007477.D
 Acq On : 11 Feb 2019 19:19
 Operator : JC/SP
 Sample : K1449-03
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-6-OW

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_X\METHOD\82X020119W.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	1.281	17	21	27	rVB2	28859	42040	1.67%	0.256%
2	1.611	66	75	79	rBV2	29375	56376	2.23%	0.343%
3	3.171	325	331	347	rVB4	19926	69248	2.74%	0.421%
4	5.262	665	674	678	rBV5	10331	25839	1.02%	0.157%
5	5.500	699	713	725	rBV	314570	920566	36.47%	5.598%
6	5.665	729	740	758	rVB	521563	1476185	58.48%	8.977%
7	5.933	773	784	789	rBV	8560	28448	1.13%	0.173%
8	6.067	795	806	817	rBV	309182	796909	31.57%	4.846%
9	6.463	861	871	878	rVB5	19530	55799	2.21%	0.339%
10	6.860	926	936	946	rBV	770871	1812018	71.79%	11.020%
11	7.256	994	1001	1009	rVB4	19464	41347	1.64%	0.251%
12	7.445	1023	1032	1044	rVB8	13116	43435	1.72%	0.264%
13	8.664	1225	1232	1234	rBV3	17024	28737	1.14%	0.175%
14	8.713	1234	1240	1251	rVB	1636240	2524050	100.00%	15.350%
15	9.036	1289	1293	1299	rVB3	16765	30290	1.20%	0.184%
16	10.109	1462	1469	1476	rBV	1603156	2201714	87.23%	13.390%
17	11.018	1613	1618	1622	rBV	65556	85270	3.38%	0.519%
18	11.133	1632	1637	1643	rBV	1347899	1766245	69.98%	10.741%
19	11.359	1669	1674	1682	rVB	86394	106981	4.24%	0.651%
20	11.450	1682	1689	1694	rVV	62807	87637	3.47%	0.533%
21	11.505	1694	1698	1703	rVB	33021	41711	1.65%	0.254%
22	11.664	1715	1724	1733	rBV	95219	136608	5.41%	0.831%
23	11.804	1743	1747	1754	rVB	272990	333515	13.21%	2.028%
24	12.078	1786	1792	1798	rVV	1737095	2169833	85.97%	13.196%
25	12.139	1798	1802	1808	rVB	53648	67319	2.67%	0.409%
26	12.298	1820	1828	1835	rBV	99218	140619	5.57%	0.855%
27	12.371	1835	1840	1846	rVB4	36662	62519	2.48%	0.380%
28	12.584	1871	1875	1878	rBV	42418	57679	2.29%	0.351%
29	12.664	1883	1888	1891	rBV	74454	93659	3.71%	0.570%
30	12.700	1891	1894	1898	rVV	44435	60315	2.39%	0.367%
31	12.755	1898	1903	1910	rVB2	216666	294345	11.66%	1.790%
32	12.901	1920	1927	1931	rBV3	29006	41178	1.63%	0.250%
33	12.975	1933	1939	1942	rBV	38264	46743	1.85%	0.284%
34	13.017	1942	1946	1952	rVB	51139	62011	2.46%	0.377%

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Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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35	13.090	1952	1958	1963	rBV2	18813	30157	1.19%	0.183%
36	13.224	1976	1980	1989	rVB	113732	159008	6.30%	0.967%
37	13.346	1995	2000	2006	rVB2	205259	267538	10.60%	1.627%
38	13.596	2038	2041	2045	rVB	36843	47792	1.89%	0.291%
39	13.645	2045	2049	2052	rBV3	23151	31986	1.27%	0.195%
40	13.718	2058	2061	2067	rVB2	58055	74345	2.95%	0.452%
41	13.834	2077	2080	2089	rVB4	10487	25367	1.01%	0.154%

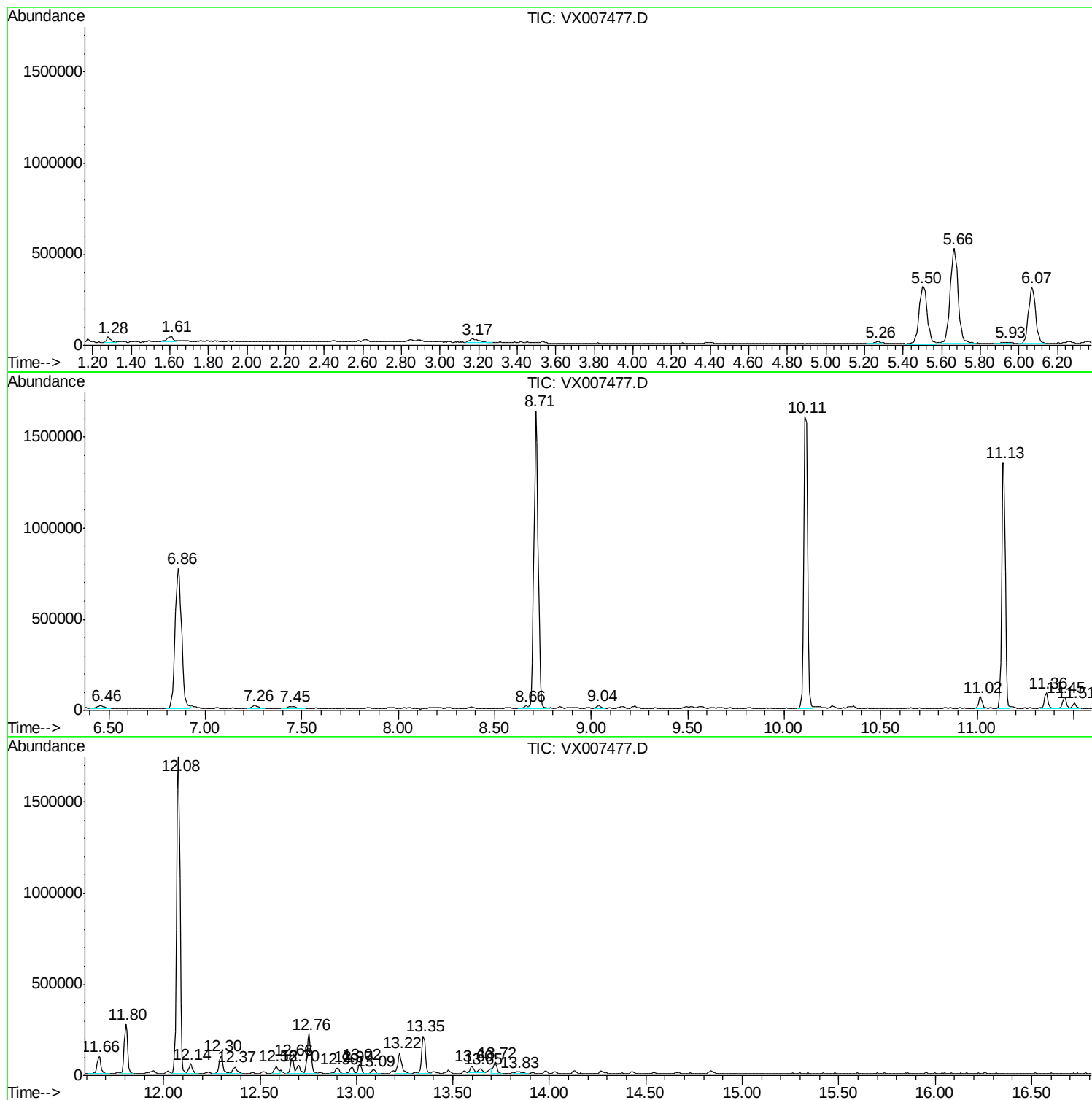
Sum of corrected areas: 16443381

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X020119W.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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ClientSampled :
B-6-OW

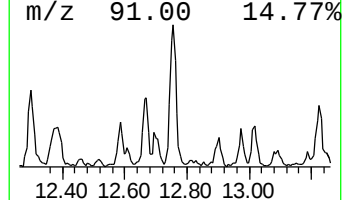
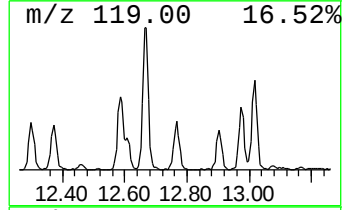
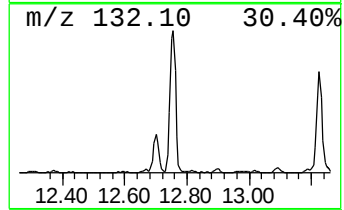
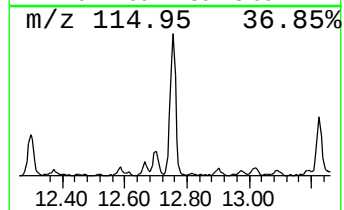
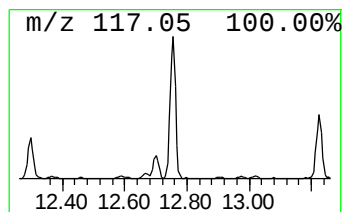
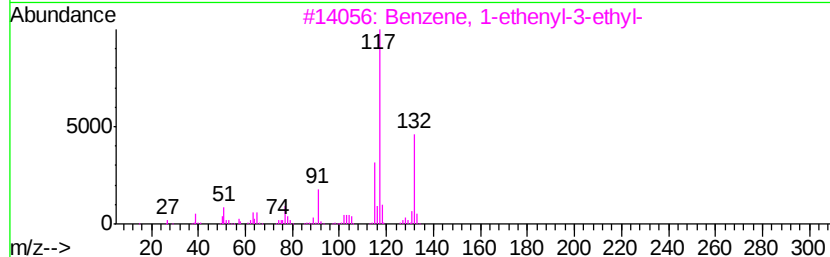
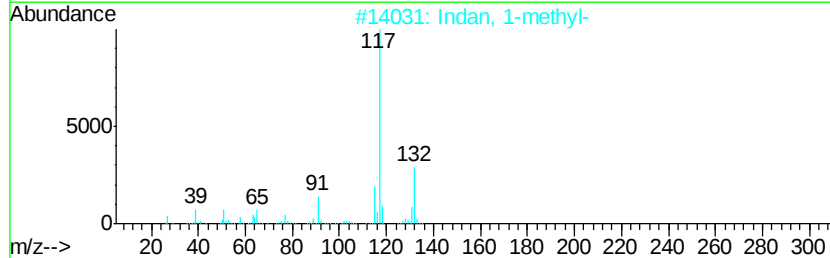
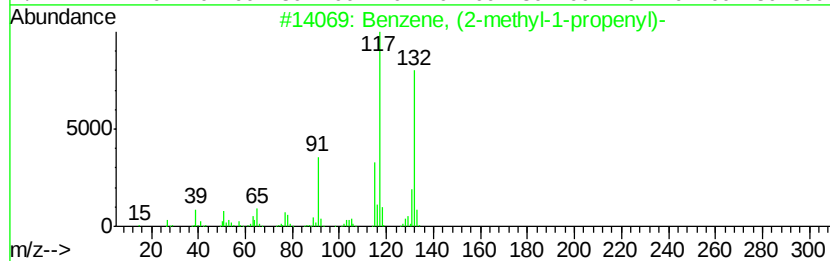
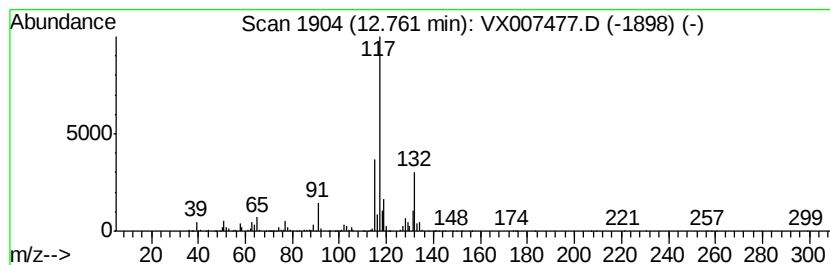
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Peak Number 1 Benzene, (2-methyl-1-propen... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	6.78 ug/l	294345	1,4-Dichlorobenzene-d4	12.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
2		Indan, 1-methyl-	132	C10H12	000767-58-8	87
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	80



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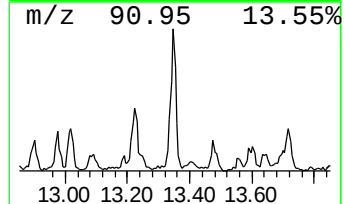
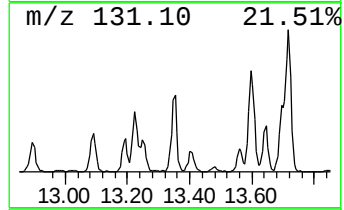
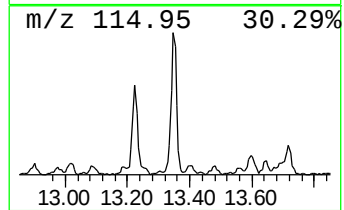
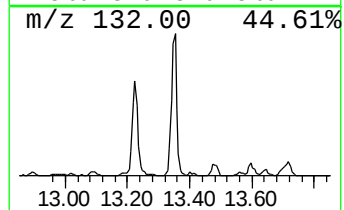
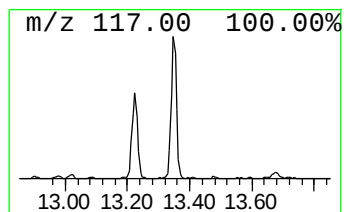
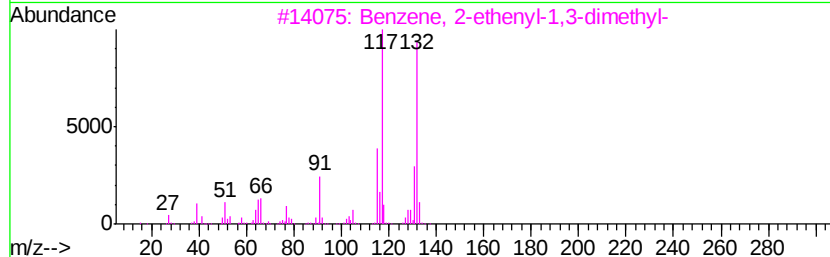
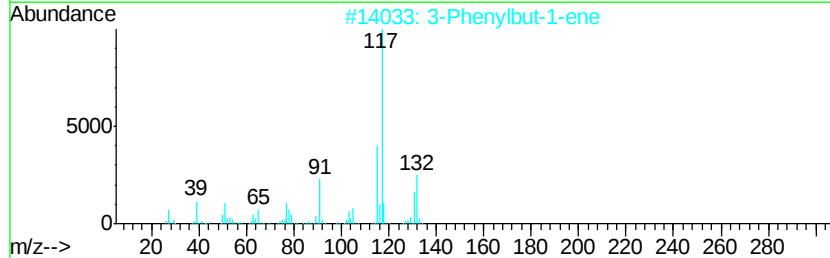
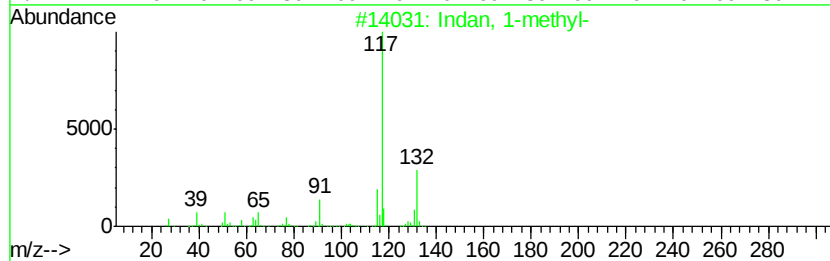
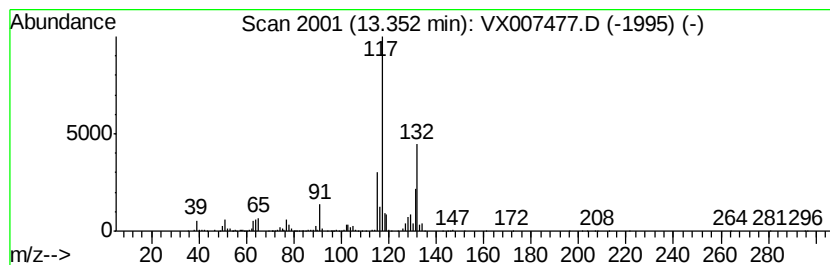
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Peak Number 2 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	6.16 ug/l	267538	1,4-Dichlorobenzene-d4	12.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-			132	C10H12	000767-58-8	90
2	3-Phenylbut-1-ene			132	C10H12	000934-10-1	90
3	Benzene, 2-ethenyl-1,3-dimethyl-			132	C10H12	002039-90-9	87
4	Benzene, 1-ethenyl-3-ethyl-			132	C10H12	007525-62-4	87
5	Benzene, 1-ethenyl-4-ethyl-			132	C10H12	003454-07-7	87



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, (2-methy...	12.76	6.8	ug/l	294345	4	12.08	2169830	50.0
Indan, 1-methyl-	13.35	6.2	ug/l	267538	4	12.08	2169830	50.0