

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121720\
 Data File : VX020104.D
 Acq On : 17 Dec 2020 07:04
 Operator : JC/MD
 Sample : L5116-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

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\82X120420W.M

Title : Sw846 8260

Signal : TIC: VX020104.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.587	74	82	88	rBV5	6574	16562	1.05%	0.149%
2	5.464	703	718	731	rBV	178848	509143	32.42%	4.575%
3	5.629	733	745	762	rVB	310485	877939	55.90%	7.889%
4	6.025	800	810	824	rBV	203857	536065	34.13%	4.817%
5	6.830	931	942	953	rBV	487777	1151759	73.33%	10.350%
6	8.696	1235	1248	1255	rBV	999036	1570612	100.00%	14.114%
7	8.763	1255	1259	1266	rVB	9950	16523	1.05%	0.148%
8	10.092	1471	1477	1490	rBV	974286	1361313	86.67%	12.233%
9	10.232	1494	1500	1507	rBV	233629	303931	19.35%	2.731%
10	10.342	1512	1518	1524	rBV	109838	143377	9.13%	1.288%
11	10.683	1569	1574	1584	rVB	29483	39710	2.53%	0.357%
12	11.000	1621	1626	1633	rBV	25309	30929	1.97%	0.278%
13	11.122	1640	1646	1658	rBV	782894	1013379	64.52%	9.106%
14	11.341	1675	1682	1689	rVB	72706	87636	5.58%	0.788%
15	11.415	1689	1694	1696	rBV	57630	76965	4.90%	0.692%
16	11.439	1696	1698	1702	rVV	61972	66393	4.23%	0.597%
17	11.488	1702	1706	1715	rVB	153155	183540	11.69%	1.649%
18	11.652	1727	1733	1739	rBV	80914	102579	6.53%	0.922%
19	11.793	1750	1756	1762	rBV	403027	517399	32.94%	4.649%
20	12.061	1794	1800	1806	rBV	1029543	1263641	80.46%	11.355%
21	12.128	1806	1811	1815	rVB	33249	43794	2.79%	0.394%
22	12.280	1829	1836	1844	rBV2	207728	295337	18.80%	2.654%
23	12.353	1844	1848	1857	rVV2	76281	111913	7.13%	1.006%
24	12.445	1857	1863	1868	rVV2	11496	15870	1.01%	0.143%
25	12.567	1878	1883	1886	rBV	41128	58916	3.75%	0.529%
26	12.591	1886	1887	1892	rVV	29866	24242	1.54%	0.218%
27	12.652	1892	1897	1900	rVV	75193	90603	5.77%	0.814%
28	12.683	1900	1902	1906	rVB	17301	18584	1.18%	0.167%
29	12.738	1906	1911	1919	rBV	51356	68506	4.36%	0.616%
30	12.957	1940	1947	1951	rBV	41655	53158	3.38%	0.478%
31	13.000	1951	1954	1961	rVB	64826	75042	4.78%	0.674%
32	13.207	1979	1988	1993	rBV	46468	58495	3.72%	0.526%
33	13.329	2003	2008	2014	rVB2	99430	130130	8.29%	1.169%
34	13.463	2025	2030	2035	rVB	33626	41160	2.62%	0.370%
35	13.628	2052	2057	2063	rVV5	7452	15742	1.00%	0.141%
36	13.817	2082	2088	2095	rBV	106637	134660	8.57%	1.210%
37	14.676	2225	2229	2236	rVB	16692	22549	1.44%	0.203%

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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 >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Title : Sw846 8260

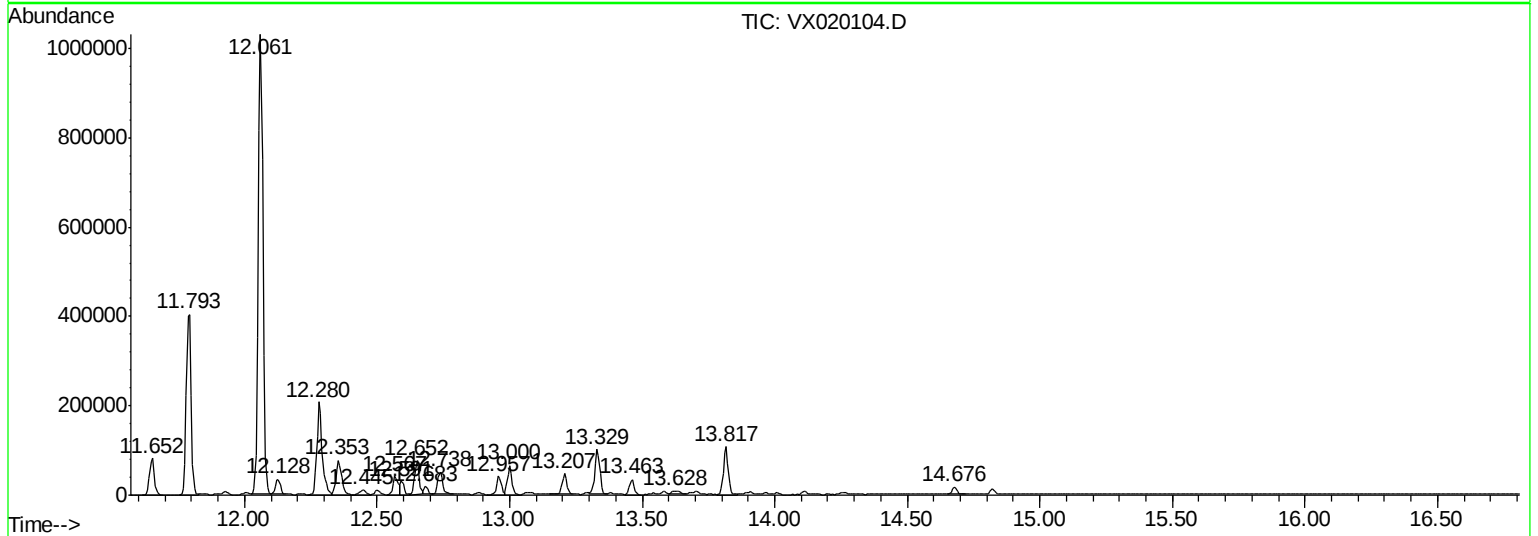
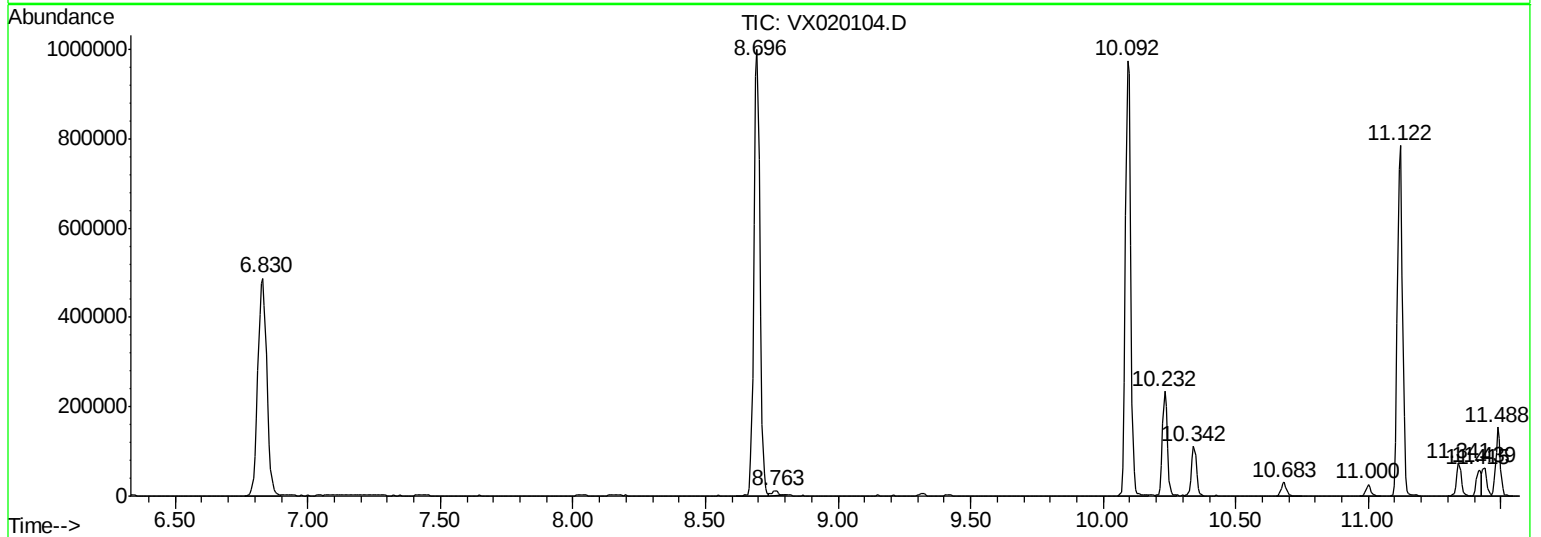
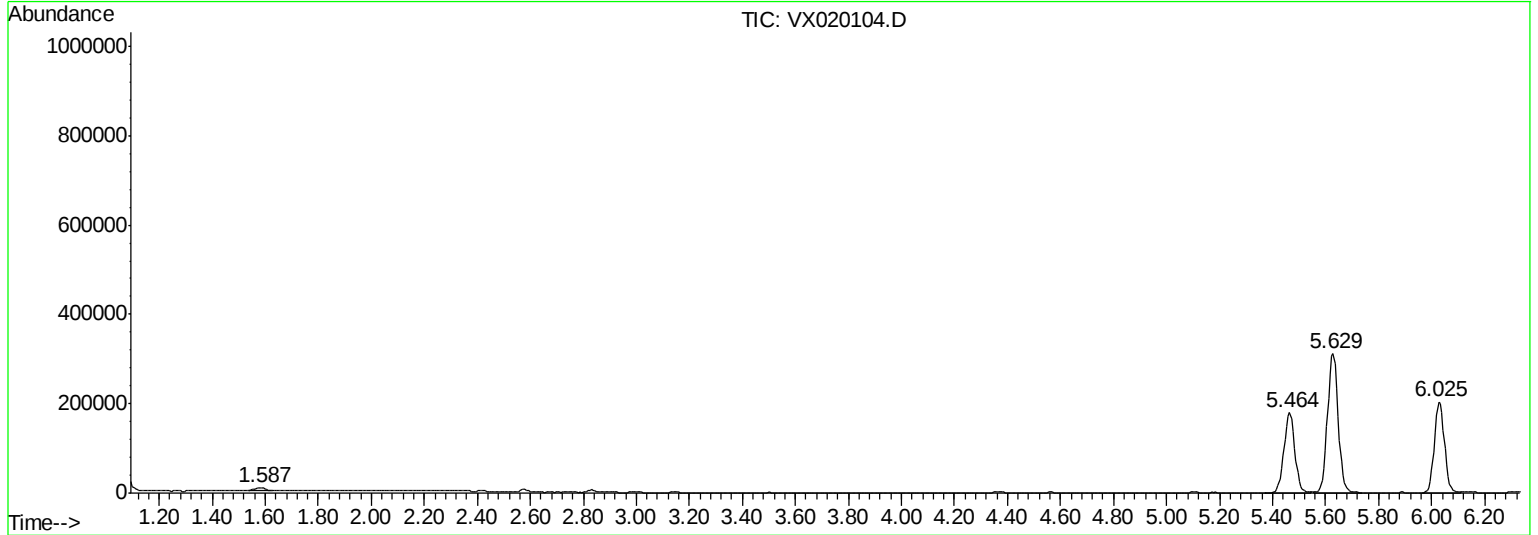
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X120420W.M
Quant Title : SW846 8260

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



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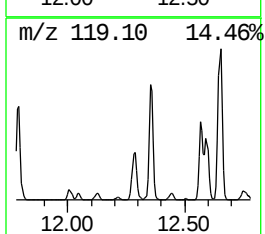
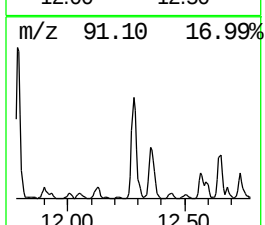
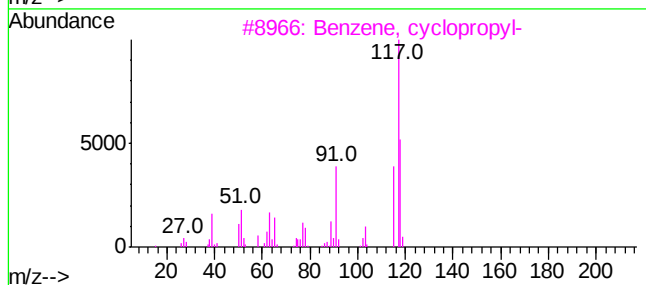
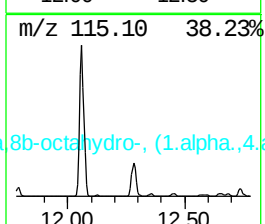
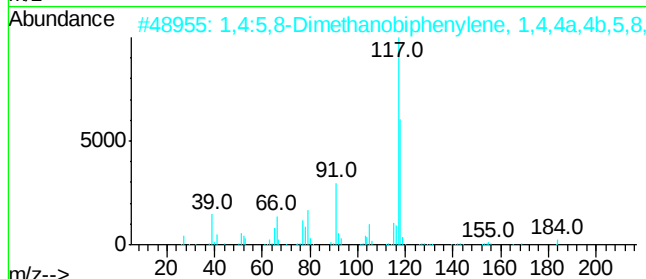
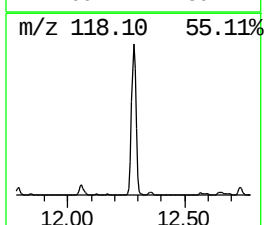
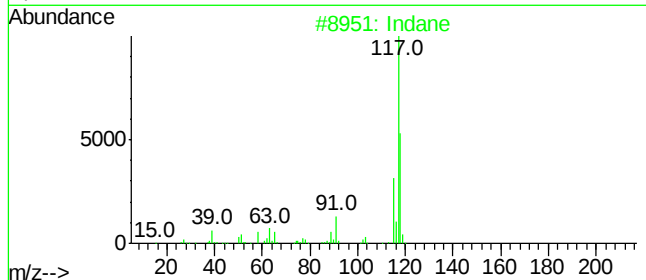
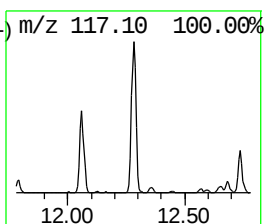
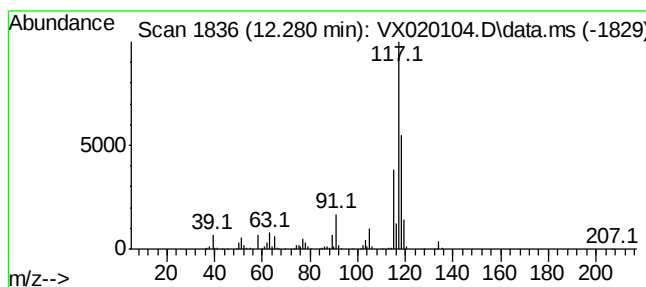
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.280	11.69 ug/l	295337	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	70
2	1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	64
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	58
5	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	58



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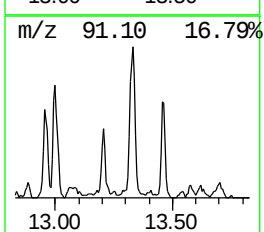
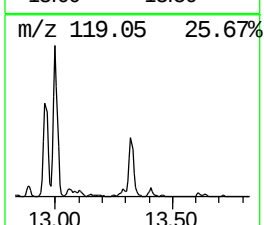
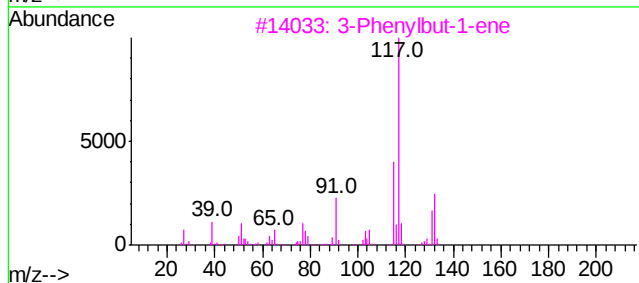
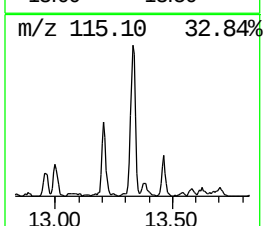
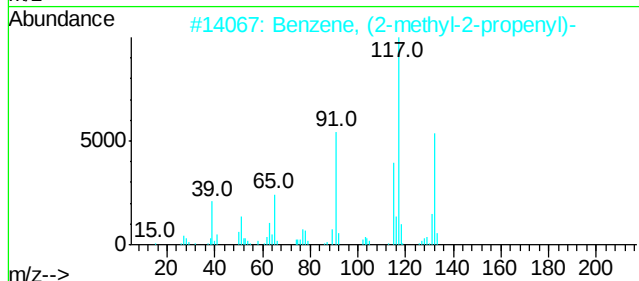
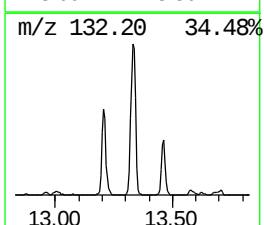
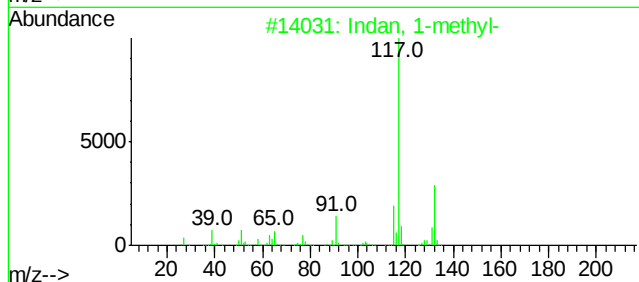
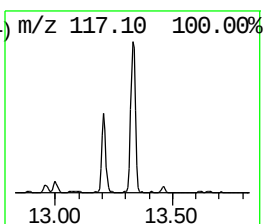
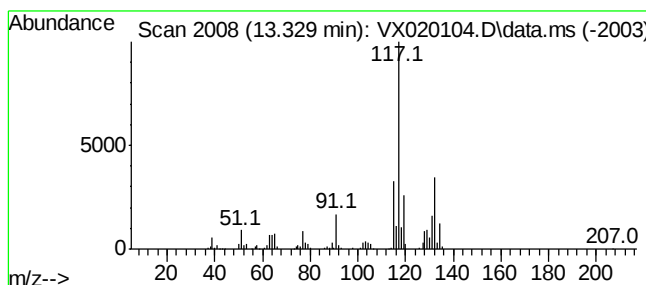
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Peak Number 2 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.330	5.15 ug/l	130130	1,4-Dichlorobenzene-d4	12.061

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	81
2	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	81
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	12.280	11.7	ug/l	295337	4	12.061	1263640	50.0
Indan, 1-methyl-	13.330	5.2	ug/l	130130	4	12.061	1263640	50.0