

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091820\
 Data File : VX018493.D
 Acq On : 18 Sep 2020 15:32
 Operator : JC/SP
 Sample : L3939-07
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PAV-B44-090820-GW

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\82X091720W.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	2.087	158	164	176	rBV	285668	465708	17.28%	2.456%
2	2.422	214	219	225	rVB	23604	40171	1.49%	0.212%
3	2.581	236	245	252	rBV	54263	108481	4.03%	0.572%
4	4.641	575	583	594	rBV2	13142	45873	1.70%	0.242%
5	5.184	660	672	682	rBV2	133485	410342	15.23%	2.164%
6	5.464	707	718	730	rBV	328126	939730	34.87%	4.956%
7	5.629	733	745	760	rVV	557180	1585444	58.83%	8.362%
8	6.037	797	812	829	rBV	356796	957458	35.53%	5.050%
9	6.830	932	942	953	rBV	819661	1977953	73.39%	10.432%
10	8.695	1240	1248	1255	rBV	1762347	2695162	100.00%	14.214%
11	8.762	1255	1259	1264	rVB	43633	68269	2.53%	0.360%
12	10.097	1472	1478	1487	rBV	1810425	2435734	90.37%	12.846%
13	11.122	1641	1646	1653	rBV	1616820	1988965	73.80%	10.490%
14	11.518	1707	1711	1714	rVB2	36095	45127	1.67%	0.238%
15	11.750	1747	1749	1752	rVB4	29505	29711	1.10%	0.157%
16	12.061	1795	1800	1806	rBV	2162769	2619510	97.19%	13.815%
17	12.256	1828	1832	1839	rBV	130533	185787	6.89%	0.980%
18	12.463	1861	1866	1870	rBV	108253	129988	4.82%	0.686%
19	13.304	2000	2004	2007	rBV	126279	165640	6.15%	0.874%
20	13.426	2020	2024	2025	rBV4	46230	60017	2.23%	0.317%
21	13.810	2084	2087	2095	rVV8	18516	38199	1.42%	0.201%
22	13.877	2095	2098	2105	rVB8	31437	58065	2.15%	0.306%
23	14.078	2127	2131	2134	rBV	80370	87925	3.26%	0.464%
24	14.652	2222	2225	2227	rBV3	47412	64954	2.41%	0.343%
25	14.676	2227	2229	2233	rVB	66299	77415	2.87%	0.408%
26	14.792	2245	2248	2251	rBV	66165	75693	2.81%	0.399%
27	14.822	2251	2253	2257	rVV3	42743	46925	1.74%	0.247%
28	14.889	2261	2264	2270	rVB7	26909	39088	1.45%	0.206%
29	15.231	2316	2320	2321	rBV3	29076	33685	1.25%	0.178%
30	15.255	2321	2324	2327	rVB3	44170	59118	2.19%	0.312%
31	15.420	2347	2351	2354	rBV2	58090	87526	3.25%	0.462%
32	15.529	2364	2369	2376	rVB2	152056	269177	9.99%	1.420%
33	15.664	2387	2391	2394	rBV	118955	183111	6.79%	0.966%
34	15.700	2394	2397	2405	rVB2	91763	159207	5.91%	0.840%

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35	15.895	2425	2429	2442	rVB9	36393	93703	3.48%	0.494%
36	16.145	2465	2470	2476	rBV2	43865	72180	2.68%	0.381%
37	16.249	2484	2487	2493	rVB5	32400	61393	2.28%	0.324%
38	16.322	2493	2499	2502	rBV8	17312	42090	1.56%	0.222%
39	16.432	2509	2517	2523	rBV4	68255	178452	6.62%	0.941%
40	16.535	2530	2534	2540	rVB5	17142	27358	1.02%	0.144%
41	16.621	2545	2548	2550	rVV4	26441	34148	1.27%	0.180%
42	16.657	2550	2554	2559	rVB3	58937	107463	3.99%	0.567%
43	16.724	2559	2565	2569	rBV2	60522	109104	4.05%	0.575%

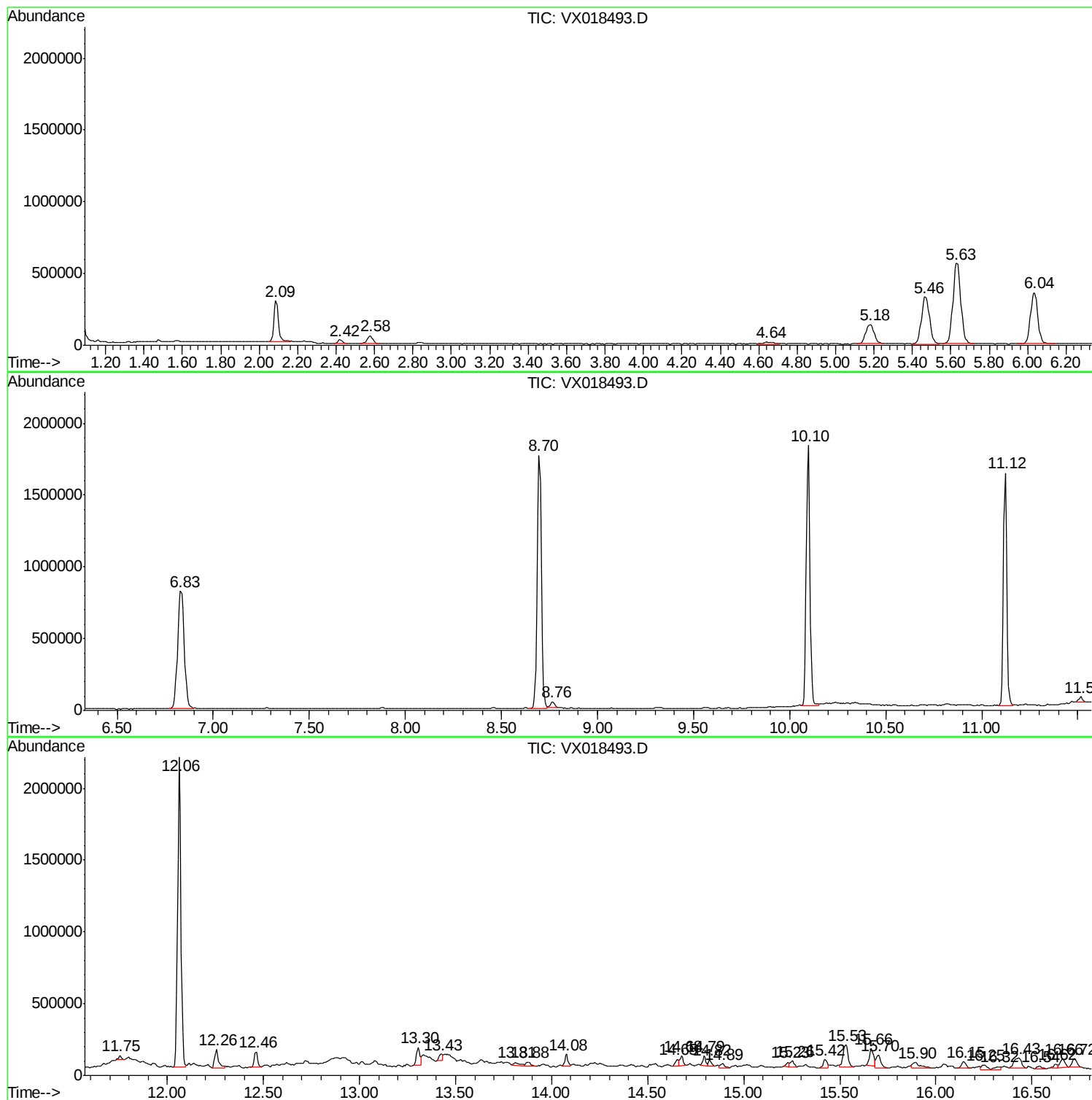
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091720W.M
Quant Title : SW846 8260

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



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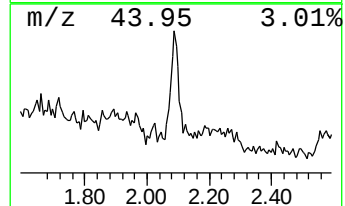
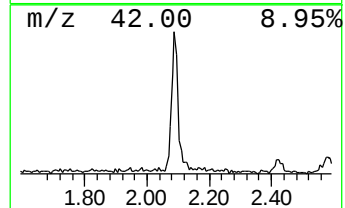
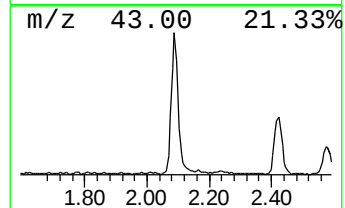
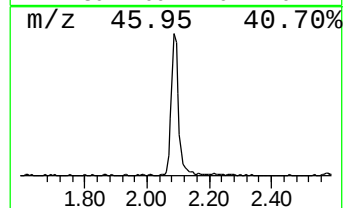
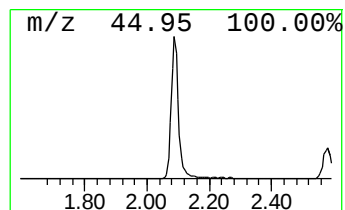
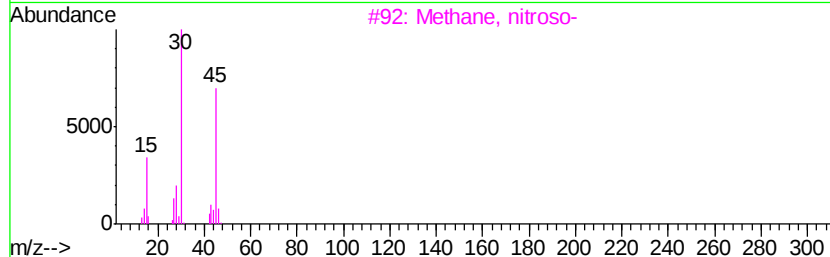
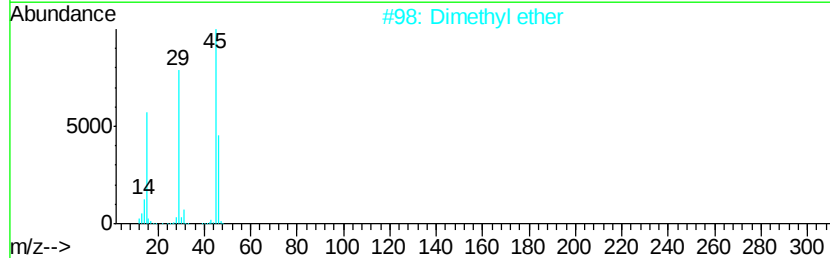
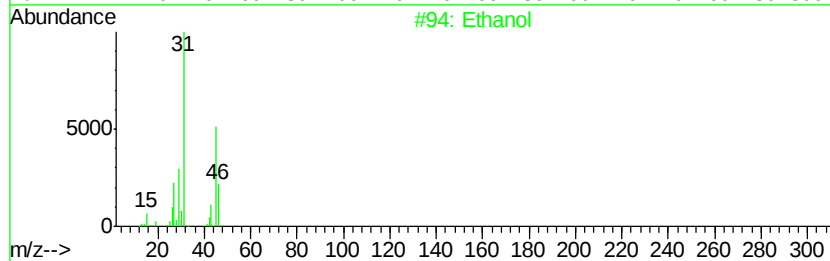
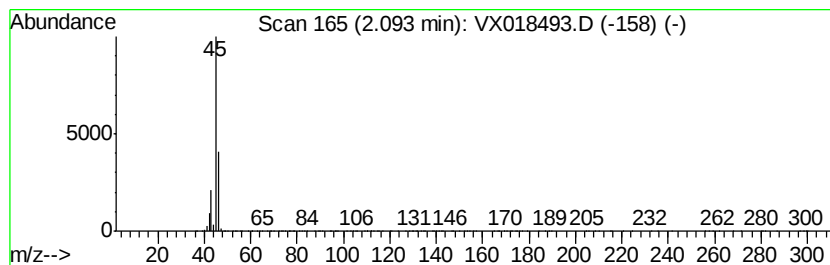
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Peak Number 1 Ethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.09	14.69 ug/l	465708	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol	46	C2H6O	000064-17-5	64
2		Dimethyl ether	46	C2H6O	000115-10-6	9
3		Methane, nitroso-	45	CH3NO	000865-40-7	4
4		Formic acid	46	CH2O2	000064-18-6	4
5		L-Lactic acid	90	C3H6O3	000079-33-4	2



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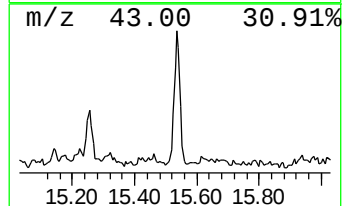
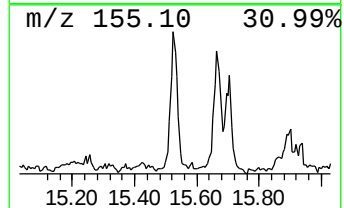
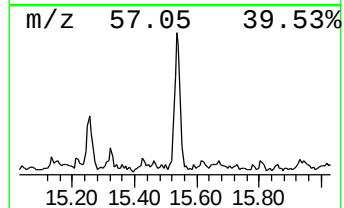
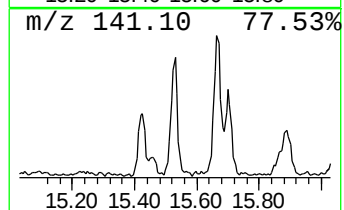
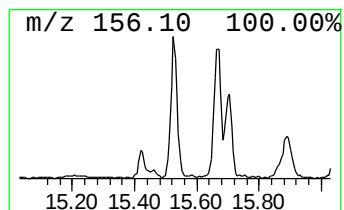
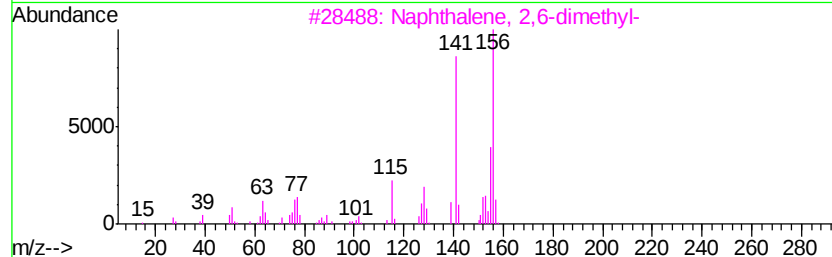
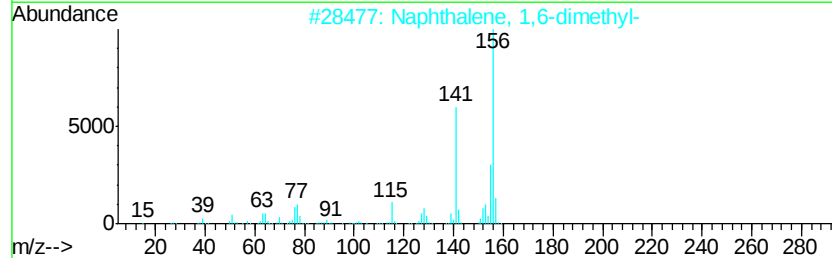
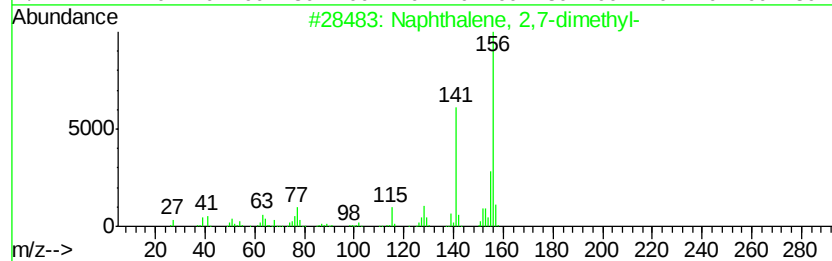
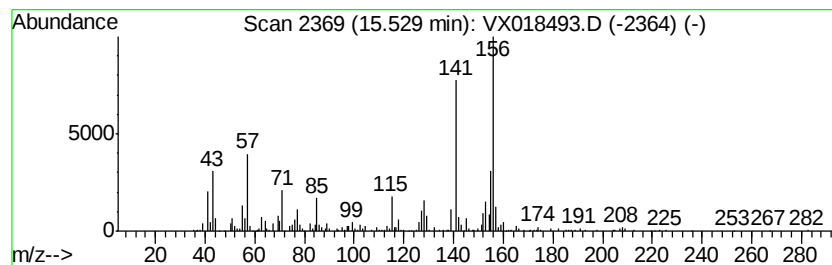
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Peak Number 2 Naphthalene, 2,7-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	5.14 ug/l	269177	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
2		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
3		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	95
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	93
5		Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Ethanol	2.09	14.7	ug/l	465708	1	5.63	1585440	50.0
Naphthalene, 2,7-...	15.53	5.1	ug/l	269177	4	12.06	2619510	50.0