

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX112120\
 Data File : VX019520.D
 Acq On : 21 Nov 2020 02:23
 Operator : JC/MD
 Sample : L4779-05 10X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1555

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\82X112120W.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	1.581	73	81	82	rBV6	9351	20117	1.34%	0.212%
2	2.995	305	313	331	rVB	10722	38197	2.54%	0.402%
3	5.464	706	718	732	rBV	167894	482081	32.02%	5.069%
4	5.629	732	745	760	rVV	309988	879902	58.44%	9.252%
5	6.031	799	811	826	rBV	191847	507856	33.73%	5.340%
6	6.830	931	942	965	rBV	471694	1137630	75.56%	11.962%
7	7.957	1119	1127	1135	rBV2	7858	15817	1.05%	0.166%
8	8.695	1241	1248	1255	rBV	982469	1505590	100.00%	15.830%
9	8.762	1255	1259	1268	rVB	40073	67289	4.47%	0.708%
10	10.097	1472	1478	1490	rBV	909699	1254804	83.34%	13.194%
11	10.341	1514	1518	1524	rVB	15359	22574	1.50%	0.237%
12	10.683	1570	1574	1578	rBV	12247	16514	1.10%	0.174%
13	11.122	1639	1646	1656	rBV	723775	928992	61.70%	9.768%
14	11.420	1691	1695	1697	rBV	13319	17921	1.19%	0.188%
15	11.652	1729	1733	1738	rBV	13470	17477	1.16%	0.184%
16	11.792	1752	1756	1762	rBV	47169	59449	3.95%	0.625%
17	12.060	1795	1800	1807	rBV	901076	1098681	72.97%	11.552%
18	12.127	1807	1811	1819	rVV2	34879	71049	4.72%	0.747%
19	12.256	1829	1832	1834	rBV	12266	15141	1.01%	0.159%
20	12.353	1845	1848	1855	rVB3	16258	26166	1.74%	0.275%
21	12.737	1908	1911	1918	rBV3	11136	26535	1.76%	0.279%
22	12.883	1933	1935	1939	rVB3	15099	17220	1.14%	0.181%
23	12.999	1951	1954	1957	rBV	18340	22446	1.49%	0.236%
24	13.329	2004	2008	2013	rBV2	45248	70346	4.67%	0.740%
25	13.463	2026	2030	2036	rVB	87428	120684	8.02%	1.269%
26	13.548	2041	2044	2047	rBV4	9892	15742	1.05%	0.166%
27	13.621	2053	2056	2061	rBV4	15267	26170	1.74%	0.275%
28	13.889	2097	2100	2106	rVB	36119	45017	2.99%	0.473%
29	13.963	2109	2112	2116	rVB2	27769	36641	2.43%	0.385%
30	14.267	2157	2162	2167	rVB	75225	108626	7.21%	1.142%
31	14.408	2183	2185	2189	rBV2	13205	15567	1.03%	0.164%
32	14.517	2199	2203	2211	rBV2	68415	102399	6.80%	1.077%
33	14.658	2222	2226	2232	rBV2	58404	101586	6.75%	1.068%
34	14.712	2232	2235	2239	rVB	47636	56294	3.74%	0.592%

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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

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Peak separation: 5

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Title : SW846 8260

35	14.859	2257	2259	2261	rBV2	14985	18360	1.22%	0.193%
36	14.883	2261	2263	2267	rVB	32093	37512	2.49%	0.394%
37	15.225	2315	2319	2321	rBV	37439	54575	3.62%	0.574%
38	15.298	2328	2331	2337	rVB3	23648	35444	2.35%	0.373%
39	15.426	2348	2352	2355	rBV5	18685	27973	1.86%	0.294%
40	15.529	2366	2369	2373	rBV3	16592	21597	1.43%	0.227%
41	15.609	2377	2382	2387	rVB	228828	327968	21.78%	3.448%
42	16.139	2466	2469	2476	rVB4	23070	38785	2.58%	0.408%

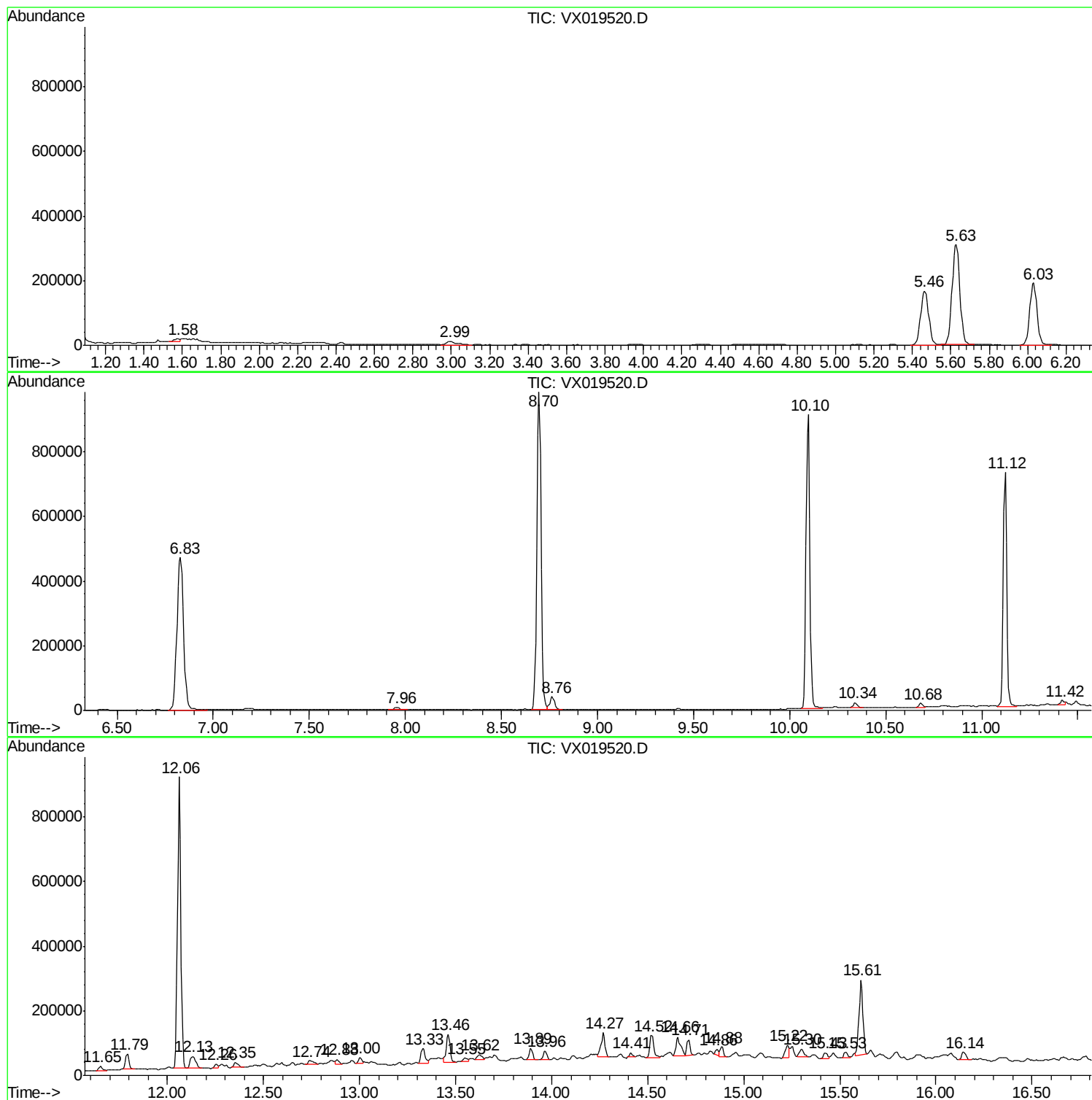
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Quant Title : SW846 8260

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



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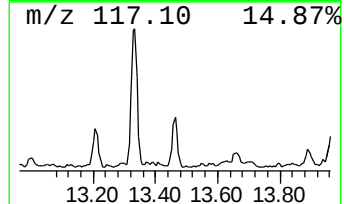
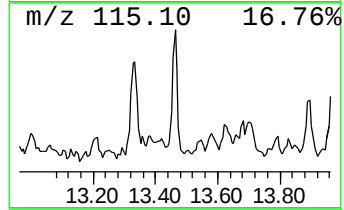
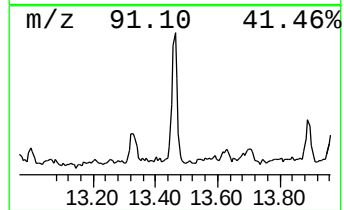
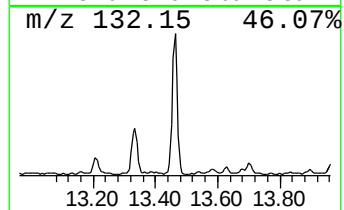
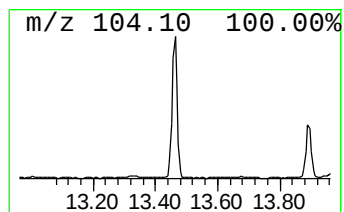
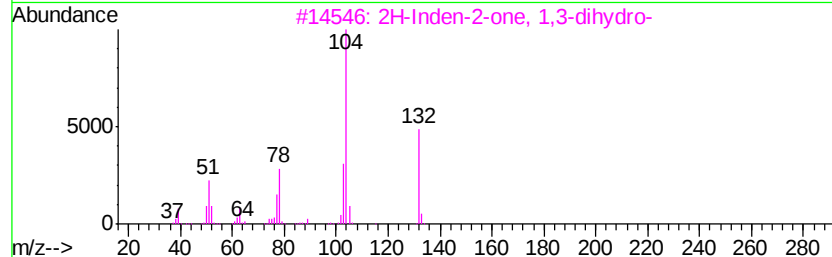
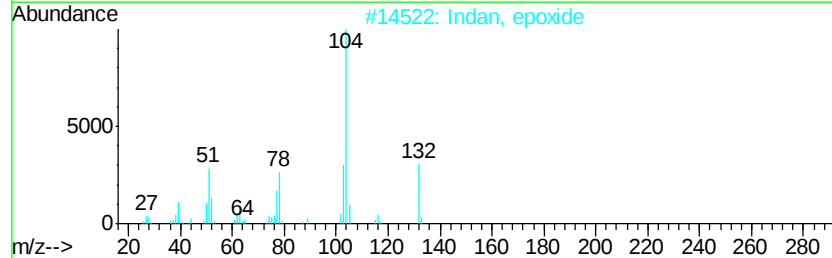
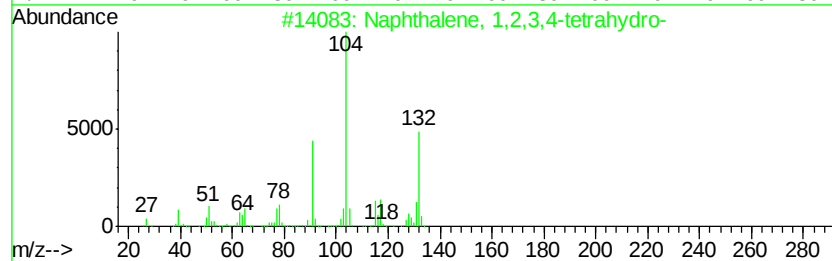
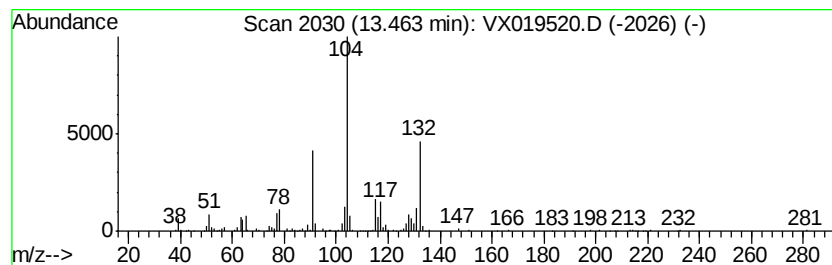
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.46	5.49 ug/l	120684	1,4-Dichlorobenzene-d4	12.06	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2	Indan, epoxide	132	C9H8O	000768-22-9	58
3	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	49
5	Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	43



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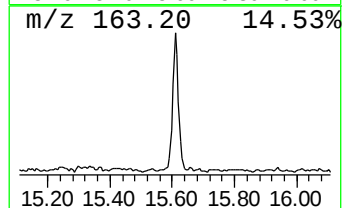
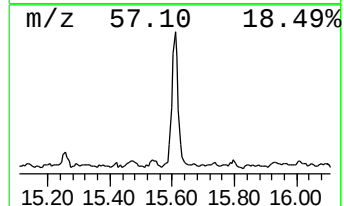
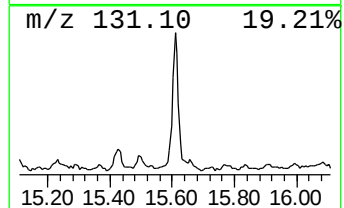
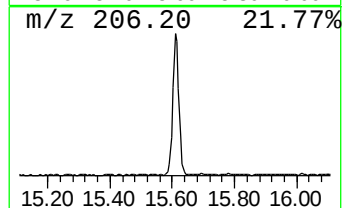
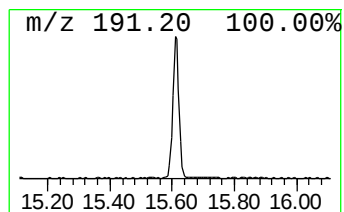
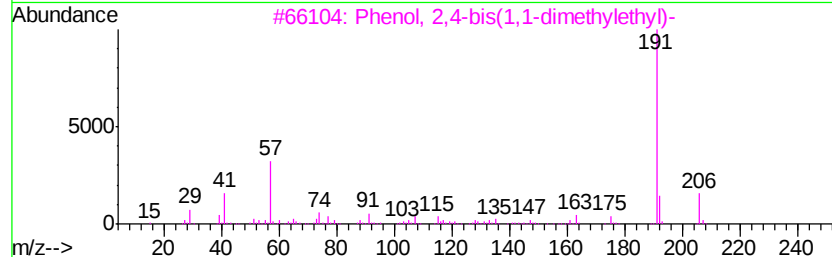
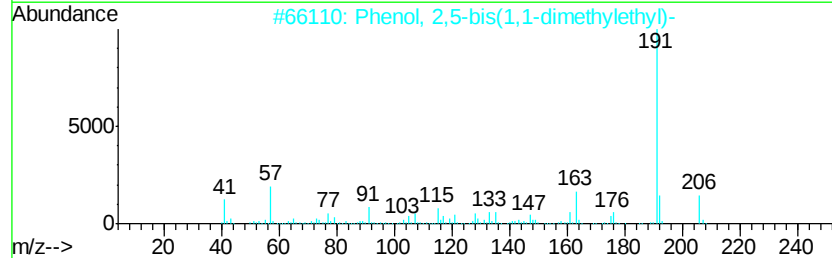
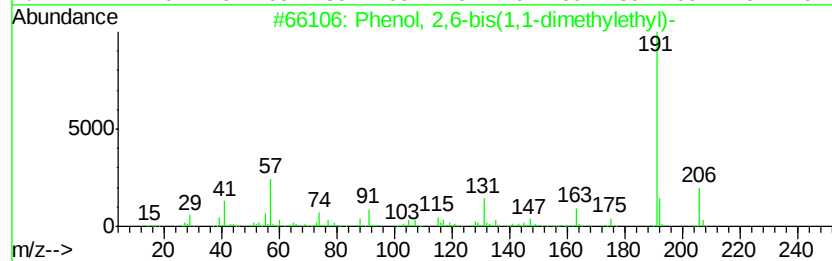
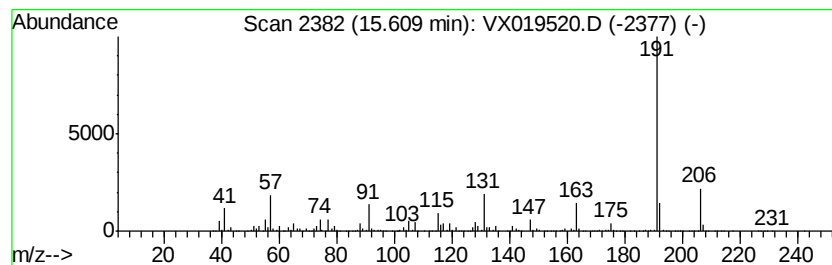
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Peak Number 2 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	14.93 ug/l	327968	1,4-Dichlorobenzene-d4	12.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	95
2	Phenol,	2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	83
3	Phenol,	2,4-bis(1,1-dimethylethyl)-	206	C14H22O	000096-76-4	81
4	Phenol,	3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	70
5	Ethyl	4-t-butylbenzoate	206	C13H18O2	005406-57-5	68



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Naphthalene, 1,2,...	13.46	5.5	ug/l	120684	4	12.06	1098680	50.0
Phenol, 2,6-bis(1...	15.61	14.9	ug/l	327968	4	12.06	1098680	50.0