

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092122\
 Data File : VX031532.D
 Acq On : 21 Sep 2022 18:14
 Operator : JC/MD
 Sample : N4614-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-23R

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: VX031532.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	92	rBV8	4408	14741	1.96%	0.427%
2	1.746	104	108	114	rBV4	6107	8185	1.09%	0.237%
3	2.337	199	205	211	rBV2	5357	10480	1.39%	0.304%
4	2.782	267	278	287	rBV2	8857	20450	2.72%	0.592%
5	5.391	692	706	719	rBV	90469	259038	34.40%	7.503%
6	5.550	721	732	743	rBV	84330	241762	32.11%	7.002%
7	5.958	787	799	808	rBV	98859	258869	34.38%	7.498%
8	6.044	808	813	827	rVB	10844	27629	3.67%	0.800%
9	6.251	831	847	859	rBV6	13191	44926	5.97%	1.301%
10	6.763	922	931	940	rBV	130558	307151	40.79%	8.896%
11	7.982	1123	1131	1137	rVB5	3202	7641	1.01%	0.221%
12	8.104	1144	1151	1163	rBV6	5759	15203	2.02%	0.440%
13	8.507	1213	1217	1225	rVB2	7808	13854	1.84%	0.401%
14	8.653	1228	1241	1250	rBV	473480	753029	100.00%	21.811%
15	8.842	1266	1272	1282	rBV2	7535	13592	1.80%	0.394%
16	8.958	1286	1291	1299	rVB2	10436	18118	2.41%	0.525%
17	9.049	1299	1306	1312	rBV2	10540	16871	2.24%	0.489%
18	10.055	1461	1471	1477	rBV	282603	377415	50.12%	10.932%
19	10.195	1490	1494	1499	rVB	7542	9226	1.23%	0.267%
20	10.964	1614	1620	1626	rBV	11017	14108	1.87%	0.409%
21	11.079	1634	1639	1649	rBV	380236	497052	66.01%	14.397%
22	11.305	1672	1676	1681	rVB	9235	11773	1.56%	0.341%
23	12.024	1788	1794	1801	rBV	295118	353208	46.90%	10.230%
24	12.250	1825	1831	1837	rBV	15687	20186	2.68%	0.585%
25	12.317	1837	1842	1848	rVB2	7286	10729	1.42%	0.311%
26	12.646	1893	1896	1901	rVB2	11250	13079	1.74%	0.379%
27	13.170	1978	1982	1985	rBV	10176	12390	1.65%	0.359%
28	13.292	1998	2002	2007	rBV2	21334	26357	3.50%	0.763%
29	13.542	2040	2043	2046	rVV2	7572	9376	1.25%	0.272%
30	13.585	2046	2050	2053	rVV3	5890	8318	1.10%	0.241%
31	13.664	2053	2063	2071	rVB2	17416	35590	4.73%	1.031%
32	13.780	2074	2082	2090	rVB3	7645	12783	1.70%	0.370%
33	14.786	2242	2247	2252	rBV2	6067	9412	1.25%	0.273%

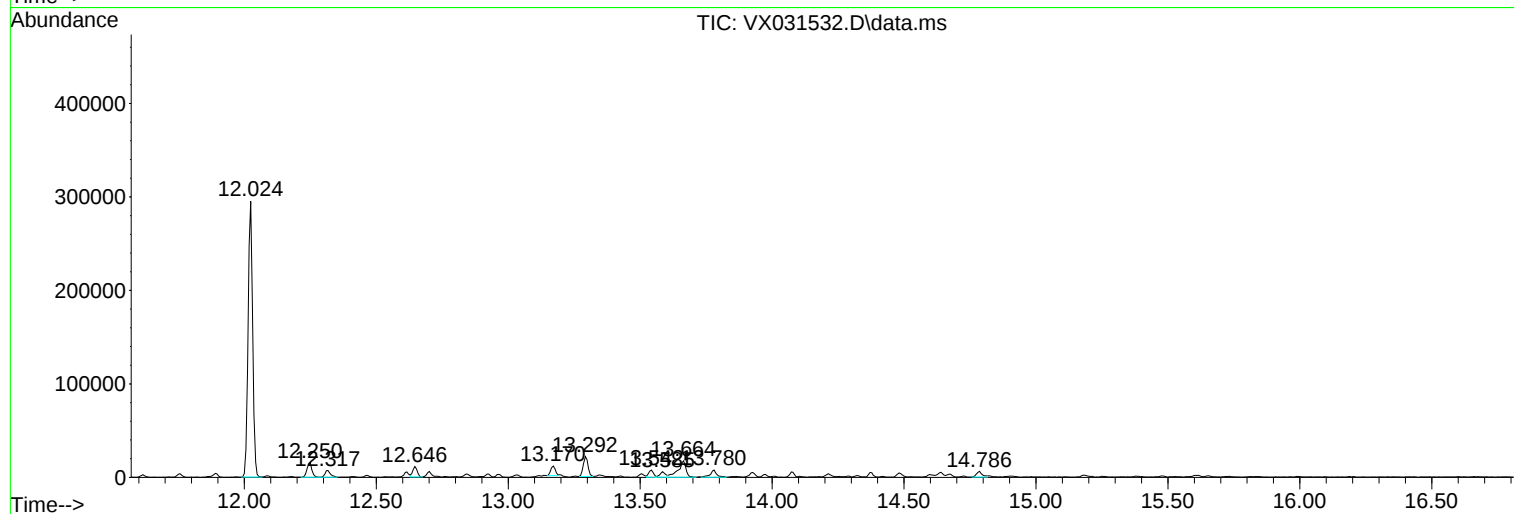
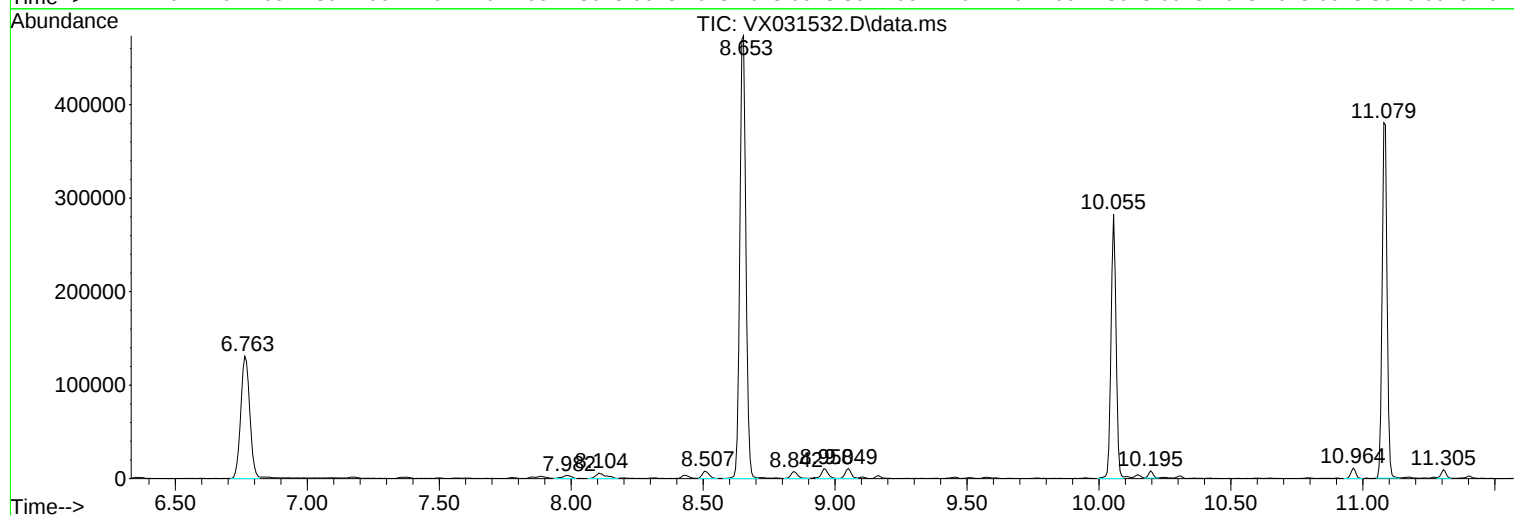
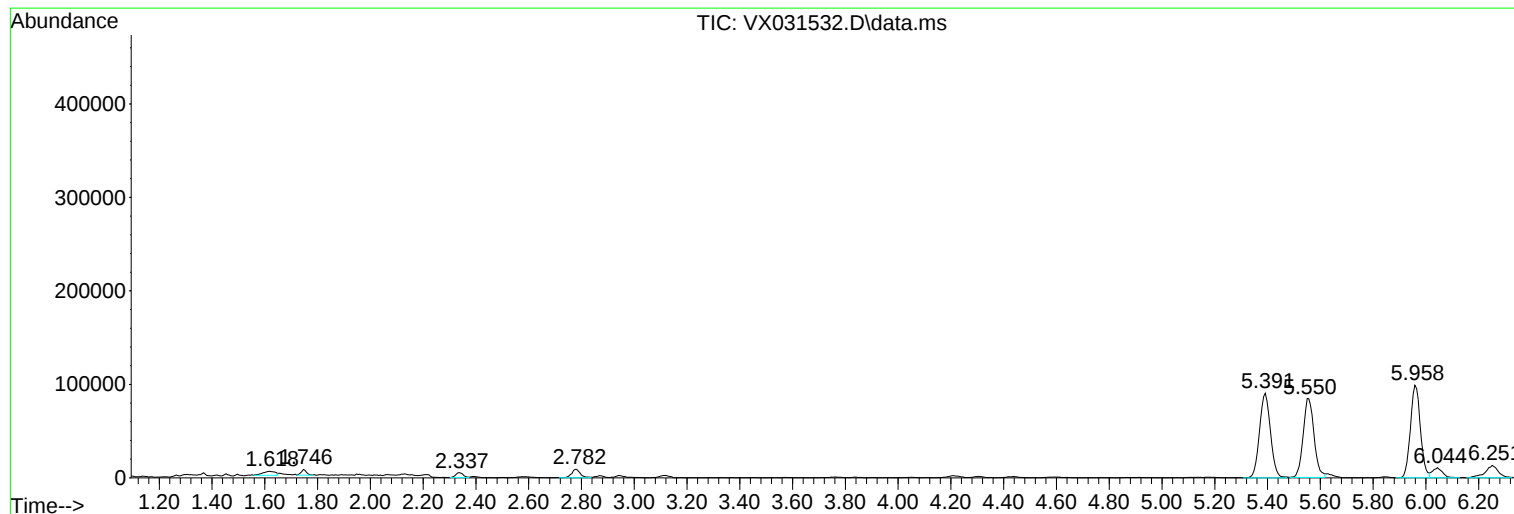
Sum of corrected areas: 3452541

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Instrument :
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ClientSampleId :
MW-23R

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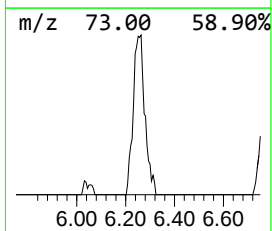
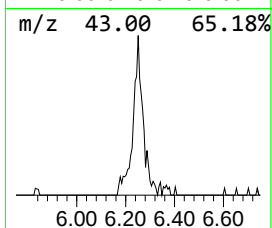
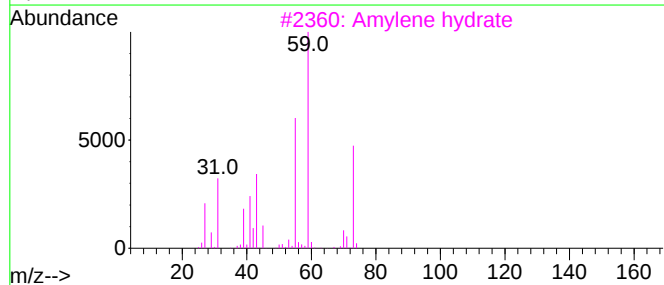
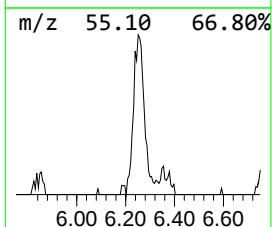
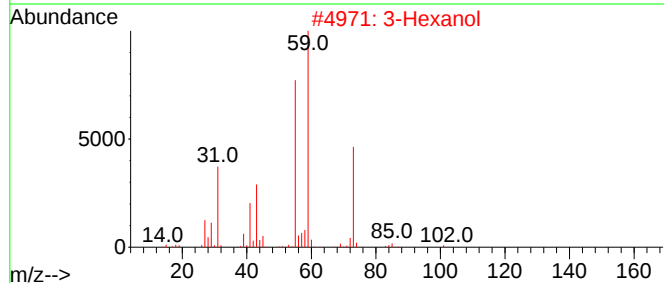
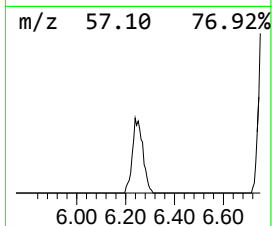
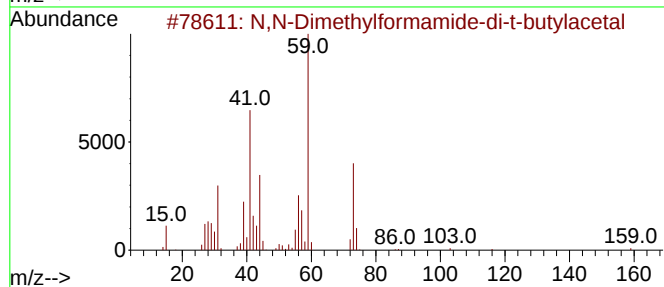
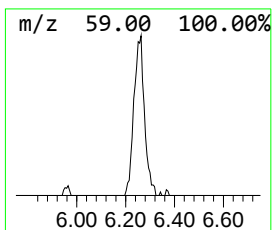
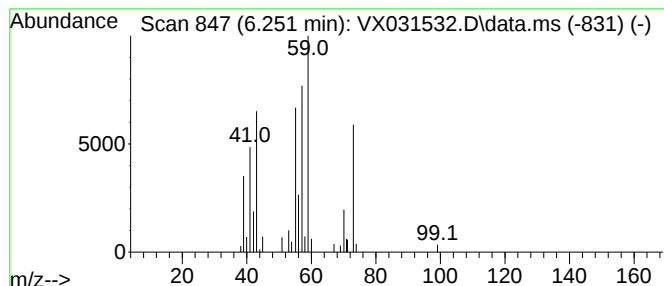
Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091922W.M
Quant Title : SW846 8260

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

Peak Number 1 N,N-Dimethylformamide-di-t-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.251	7.31 ug/l	44926	1,4-Difluorobenzene	6.763	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	N,N-Dimethylformamide-di-t-butyl...	203	C11H25NO2	036805-97-7	50
2	3-Hexanol	102	C6H14O	000623-37-0	47
3	Amylene hydrate	88	C5H12O	000075-85-4	42
4	Silane, trimethyl-	74	C3H10Si	000993-07-7	38
5	(3,3-Dimethyloxiranyl)methanol	102	C5H10O2	1010306-71-7	38



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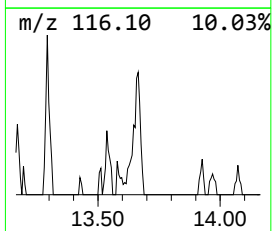
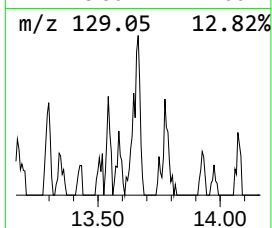
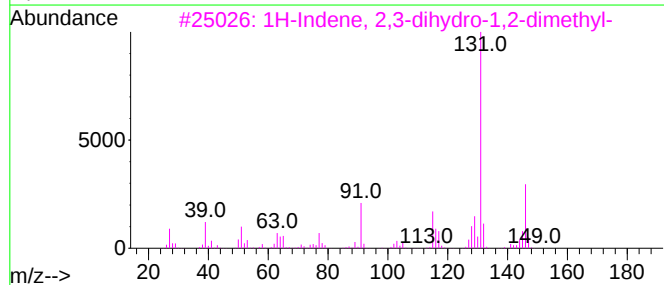
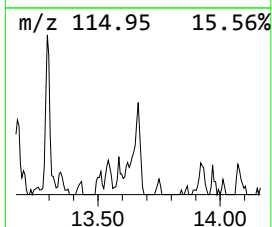
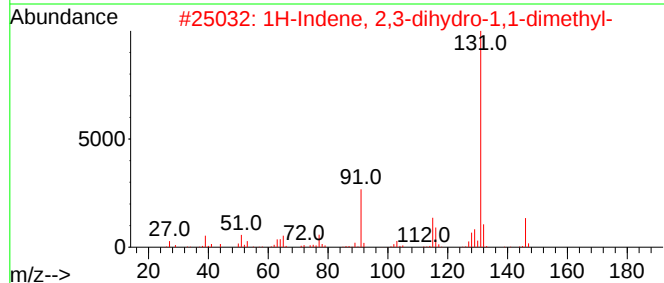
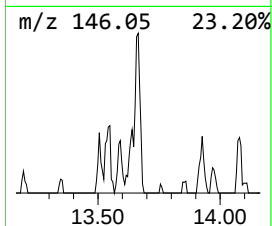
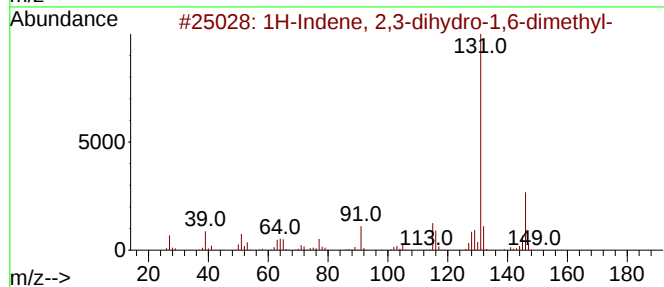
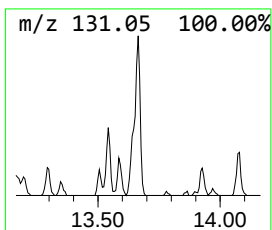
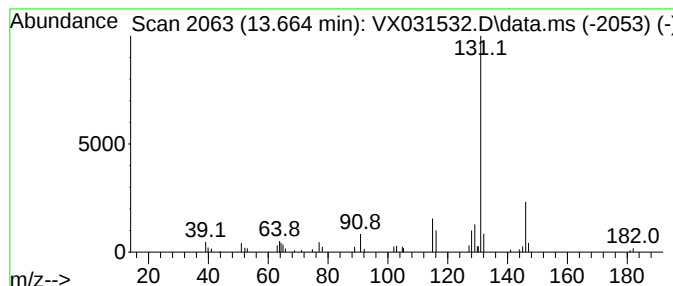
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 2 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	5.04 ug/l	35590	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91		
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	91		
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91		
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90		
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
N,N-Dimethylfor...	6.251	7.3	ug/l	44926	2	6.763	307151	50.0
1H-Indene, 2,3-...	13.664	5.0	ug/l	35590	4	12.024	353208	50.0