

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
 Data File : VX043125.D
 Acq On : 24 Sep 2024 16:12
 Operator : JC/MD
 Sample : IBLK
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 IBLK

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\82X092324W.M
 Title : SW846 8260

Signal : TIC: VX043125.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.568	562	571	582	rBV2	4547	14732	2.17%	0.326%
2	5.385	695	705	722	rBV	64597	189263	27.91%	4.187%
3	5.550	722	732	749	rVB	90288	259671	38.30%	5.745%
4	5.958	789	799	807	rBV	73000	193760	28.58%	4.287%
5	6.037	807	812	826	rVB3	17123	48196	7.11%	1.066%
6	6.763	921	931	943	rBV	160553	395306	58.30%	8.745%
7	8.647	1234	1240	1247	rBV	359116	578017	85.25%	12.788%
8	8.720	1247	1252	1266	rVB	33399	64510	9.51%	1.427%
9	10.055	1465	1471	1483	rBV	361617	512987	75.66%	11.349%
10	10.299	1507	1511	1519	rBV	29151	41961	6.19%	0.928%
11	10.646	1562	1568	1581	rVB2	25252	43034	6.35%	0.952%
12	11.079	1634	1639	1649	rBV	325963	419880	61.93%	9.289%
13	11.451	1696	1700	1710	rVB	6825	9542	1.41%	0.211%
14	11.750	1745	1749	1756	rBV2	14941	22380	3.30%	0.495%
15	11.963	1779	1784	1788	rBV2	6494	8920	1.32%	0.197%
16	12.018	1788	1793	1800	rBV	423587	542029	79.94%	11.991%
17	12.091	1800	1805	1809	rVB2	6539	10500	1.55%	0.232%
18	12.414	1853	1858	1864	rBV	80375	107071	15.79%	2.369%
19	12.804	1912	1922	1930	rBV3	14613	34404	5.07%	0.761%
20	12.969	1942	1949	1957	rVB6	7109	17723	2.61%	0.392%
21	13.268	1993	1998	2006	rBV4	12918	21735	3.21%	0.481%
22	13.585	2042	2050	2055	rBV8	3416	7080	1.04%	0.157%
23	13.713	2066	2071	2074	rBV2	15609	24948	3.68%	0.552%
24	13.774	2076	2081	2088	rVV	515590	678033	100.00%	15.000%
25	13.835	2088	2091	2093	rVV2	12082	17925	2.64%	0.397%
26	13.865	2093	2096	2105	rVB	21801	32243	4.76%	0.713%
27	14.511	2198	2202	2209	rBV4	4429	6817	1.01%	0.151%
28	14.633	2218	2222	2232	rVB	45225	63827	9.41%	1.412%
29	14.731	2232	2238	2242	rBV	13302	22024	3.25%	0.487%
30	14.780	2242	2246	2251	rVB	24828	32897	4.85%	0.728%
31	15.206	2311	2316	2320	rVB2	8767	12207	1.80%	0.270%
32	15.274	2322	2327	2332	rBV3	7858	11759	1.73%	0.260%
33	15.481	2356	2361	2368	rVB6	4057	7577	1.12%	0.168%
34	15.682	2390	2394	2406	rVB8	6971	16461	2.43%	0.364%
35	16.560	2532	2538	2542	rBV3	8236	16470	2.43%	0.364%
36	16.615	2542	2547	2553	rVV5	6648	13896	2.05%	0.307%

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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\82X092324W.M
Title : SW846 8260

37	16.682	2553	2558	2565	rVV	9773	20328	3.00%	0.450%
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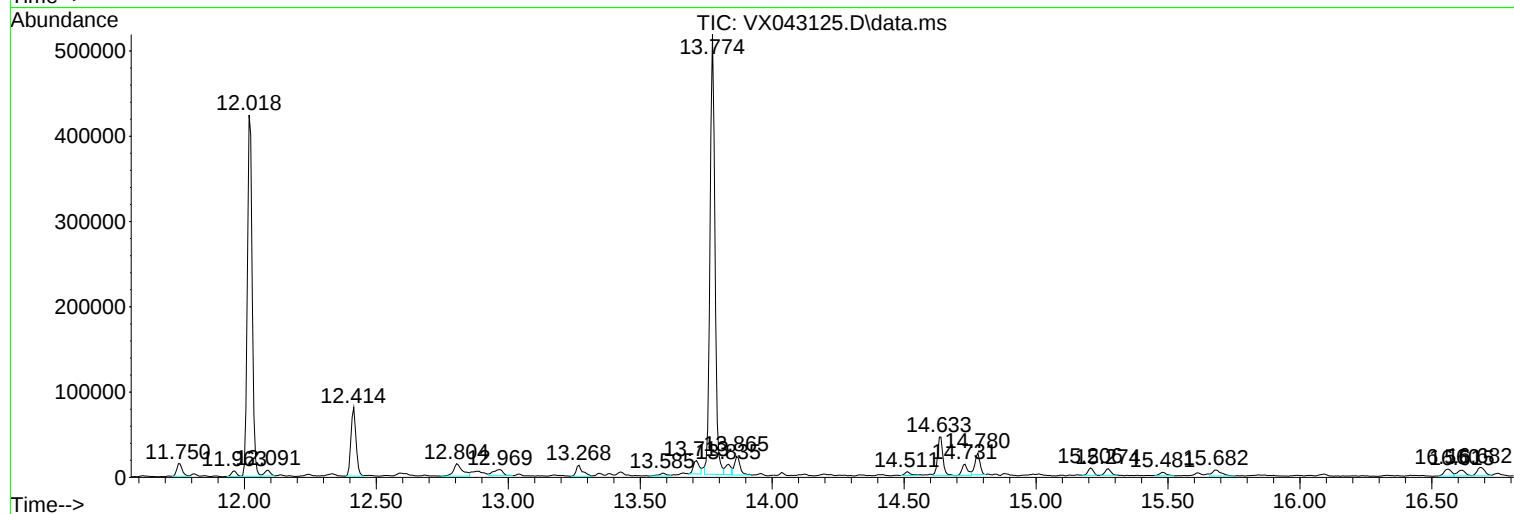
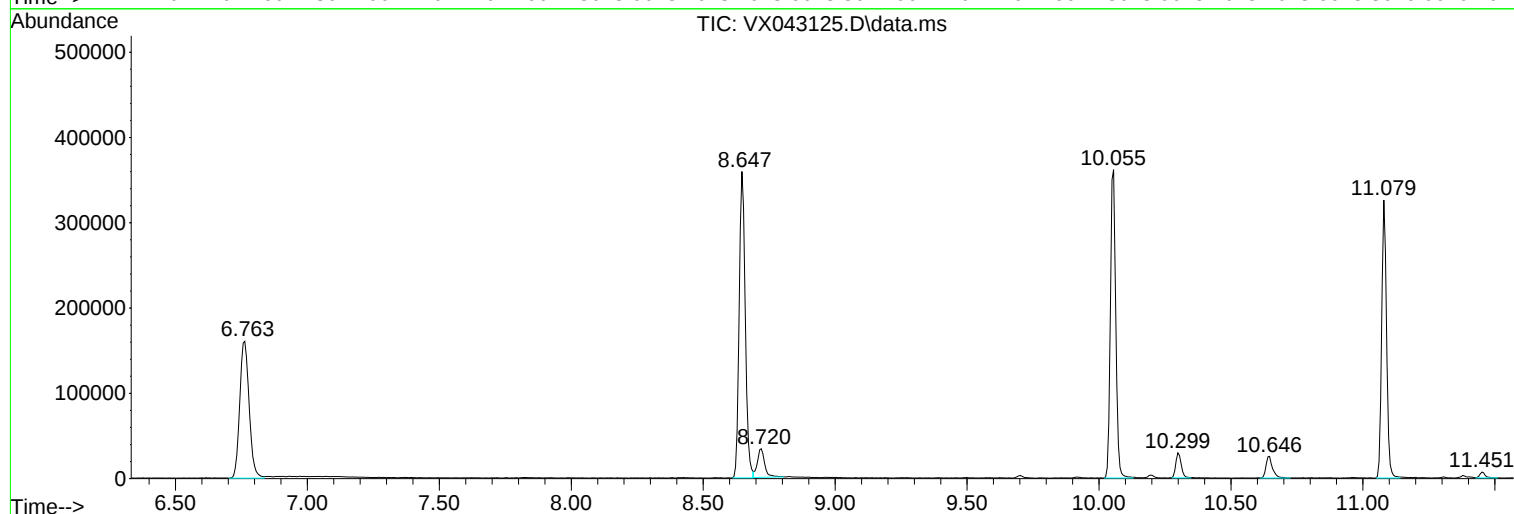
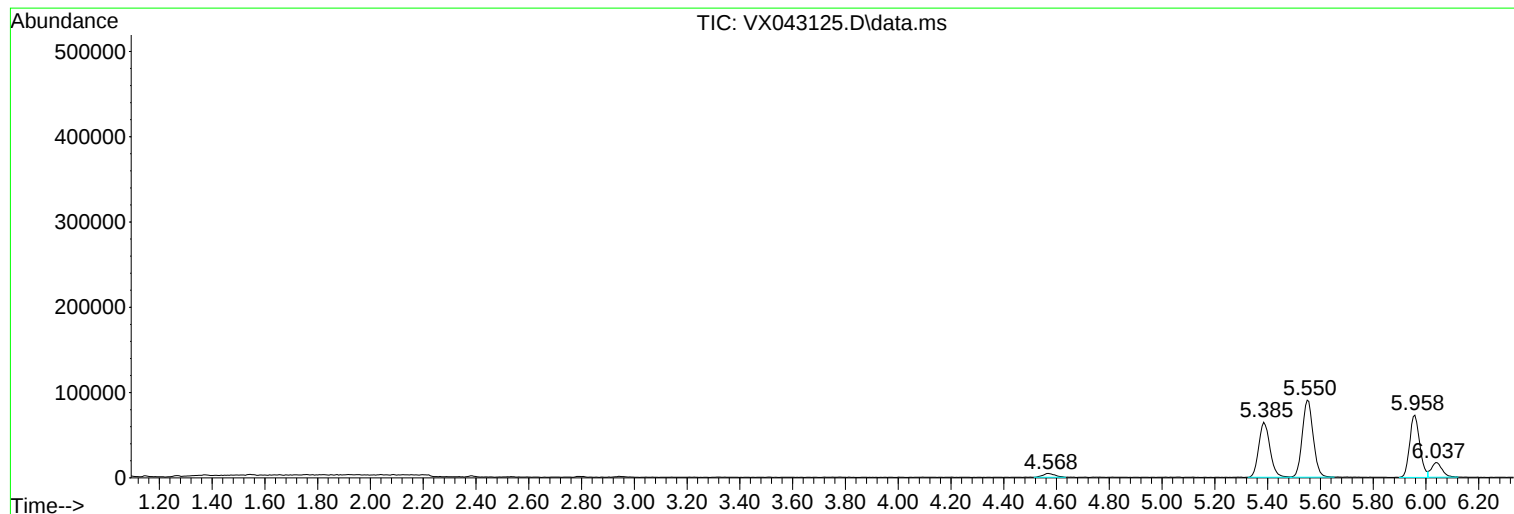
Sum of corrected areas: 4520113

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
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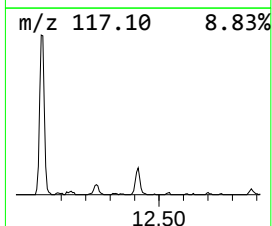
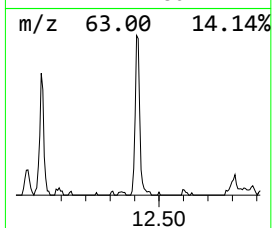
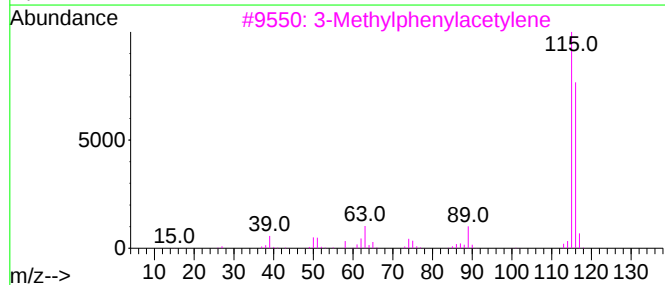
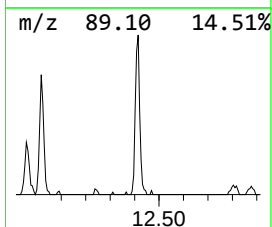
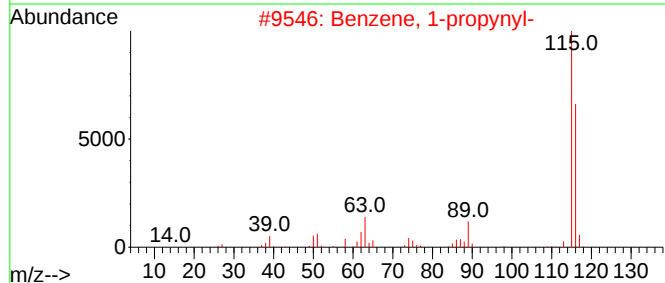
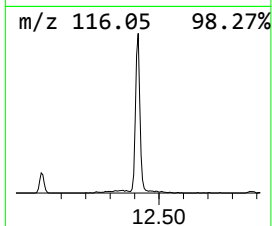
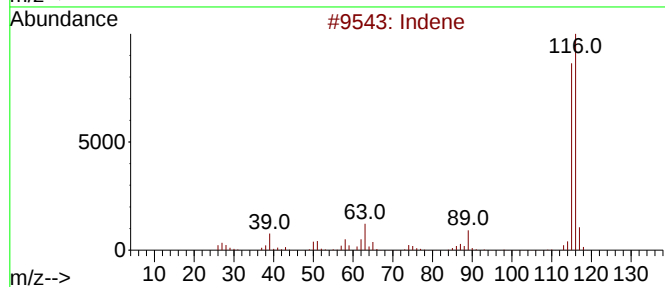
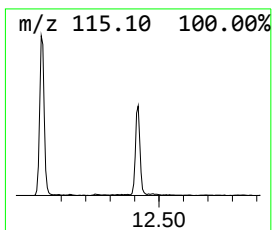
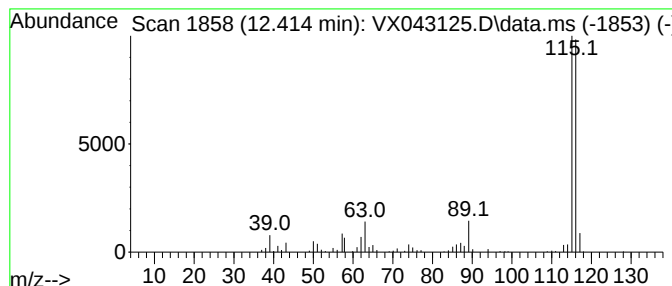
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
Quant Title : SW846 8260

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

Peak Number 1 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.414	9.88 ug/l	107071	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	94
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
3			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4			Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90
5			2-Methylphenylacetylene	116	C9H8	000766-47-2	90



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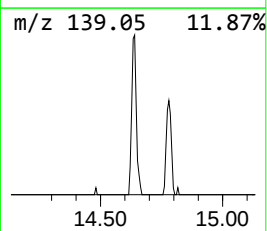
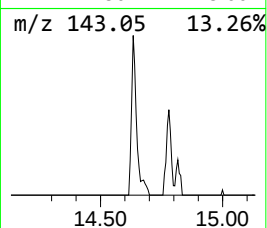
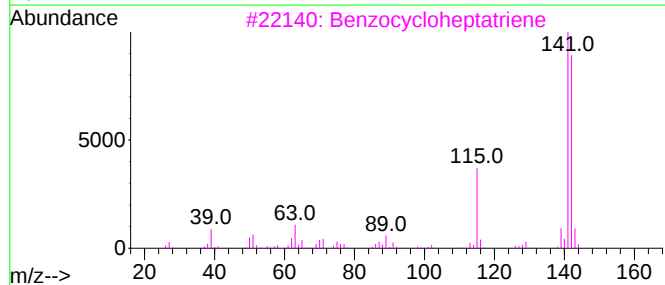
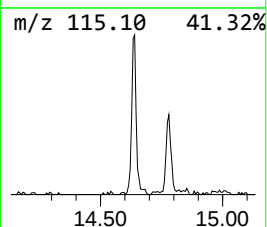
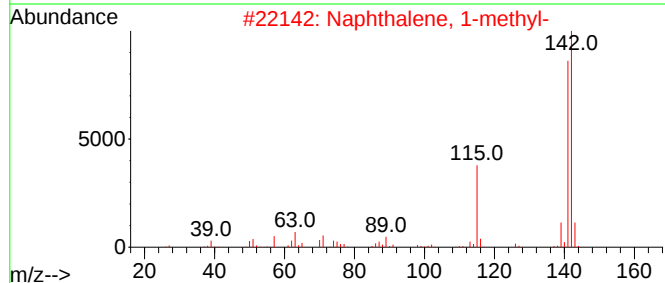
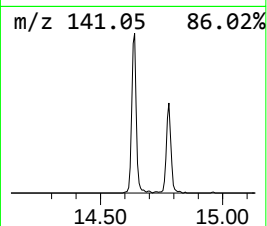
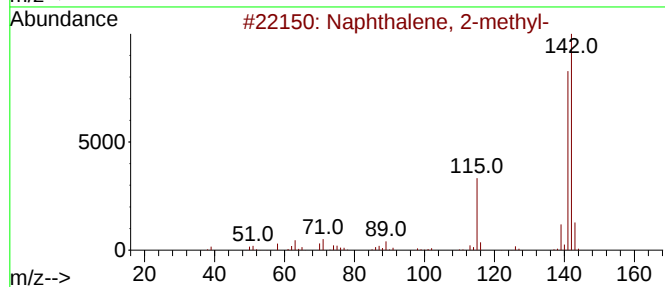
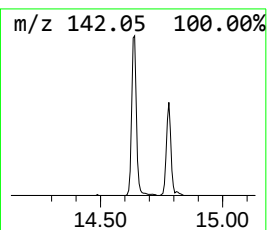
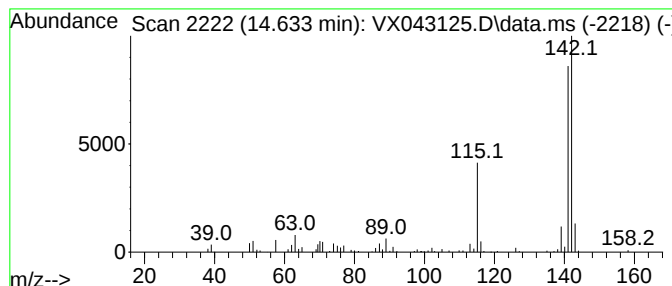
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Peak Number 2 Naphthalene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.633	5.89 ug/l	63827	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indene	12.414	9.9	ug/l	107071	4	12.024	542029	50.0
Naphthalene, 2-...	14.633	5.9	ug/l	63827	4	12.024	542029	50.0