

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN110122\
 Data File : VN075104.D
 Acq On : 01 Nov 2022 17:58
 Operator : JC\MD
 Sample : N5417-09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 AUD-1906

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_N\methods\82N102622W.M
 Title : SW846 8260

Signal : TIC: VN075104.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.301	35	42	51	rBV2	34152	60941	5.69%	1.128%
2	2.818	124	130	132	rBV2	11836	18396	1.72%	0.340%
3	4.448	400	407	414	rBV2	4458	11890	1.11%	0.220%
4	4.736	449	456	462	rBV2	4331	11684	1.09%	0.216%
5	7.441	907	916	918	rBV	6148	11890	1.11%	0.220%
6	7.494	919	925	935	rVB	21927	57971	5.41%	1.073%
7	8.171	1030	1040	1045	rBV	91819	219798	20.51%	4.068%
8	8.230	1045	1050	1059	rVB	156086	330233	30.81%	6.111%
9	8.583	1101	1110	1120	rBV	98984	224719	20.96%	4.159%
10	9.106	1188	1199	1208	rBV	226285	465969	43.47%	8.623%
11	10.571	1440	1448	1453	rBV	355275	653587	60.97%	12.096%
12	10.635	1453	1459	1476	rVB	582903	1071910	100.00%	19.837%
13	11.865	1660	1668	1679	rBV	310773	563726	52.59%	10.433%
14	12.071	1696	1703	1708	rBV2	9619	17445	1.63%	0.323%
15	12.400	1754	1759	1764	rBV3	11710	19079	1.78%	0.353%
16	12.853	1826	1836	1843	rVB	232850	416781	38.88%	7.713%
17	13.100	1872	1878	1882	rBV2	16949	30602	2.85%	0.566%
18	13.176	1887	1891	1897	rVB2	11375	18447	1.72%	0.341%
19	13.335	1912	1918	1925	rBV	13670	22627	2.11%	0.419%
20	13.482	1938	1943	1951	rVB2	25621	45494	4.24%	0.842%
21	13.794	1988	1996	2005	rBV2	283975	532633	49.69%	9.857%
22	13.994	2024	2030	2047	rVV5	26405	99831	9.31%	1.848%
23	14.182	2058	2062	2068	rVB3	8206	15034	1.40%	0.278%
24	14.247	2068	2073	2075	rBV3	9538	14247	1.33%	0.264%
25	14.329	2082	2087	2094	rVB3	12399	19673	1.84%	0.364%
26	14.447	2101	2107	2113	rVB2	17411	27581	2.57%	0.510%
27	14.570	2121	2128	2134	rVB5	7529	16954	1.58%	0.314%
28	14.659	2138	2143	2145	rVV	8579	14909	1.39%	0.276%
29	14.694	2145	2149	2155	rVB	15171	22071	2.06%	0.408%
30	14.929	2184	2189	2193	rBV2	18665	29370	2.74%	0.544%
31	15.053	2201	2210	2216	rBV4	38531	91617	8.55%	1.696%
32	15.212	2229	2237	2245	rBV2	37407	71075	6.63%	1.315%
33	15.317	2250	2255	2259	rBV6	7881	14899	1.39%	0.276%
34	15.435	2269	2275	2286	rVB5	11090	30934	2.89%	0.572%
35	15.823	2337	2341	2352	rVB8	8754	22130	2.06%	0.410%
36	15.947	2357	2362	2368	rVB	6610	12934	1.21%	0.239%

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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

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Title : SW846 8260

37	16.164	2393	2399	2409	rVB3	20922	48150	4.49%	0.891%
38	16.341	2419	2429	2437	rBV7	5467	15336	1.43%	0.284%
39	16.500	2449	2456	2465	rBV3	13752	30970	2.89%	0.573%

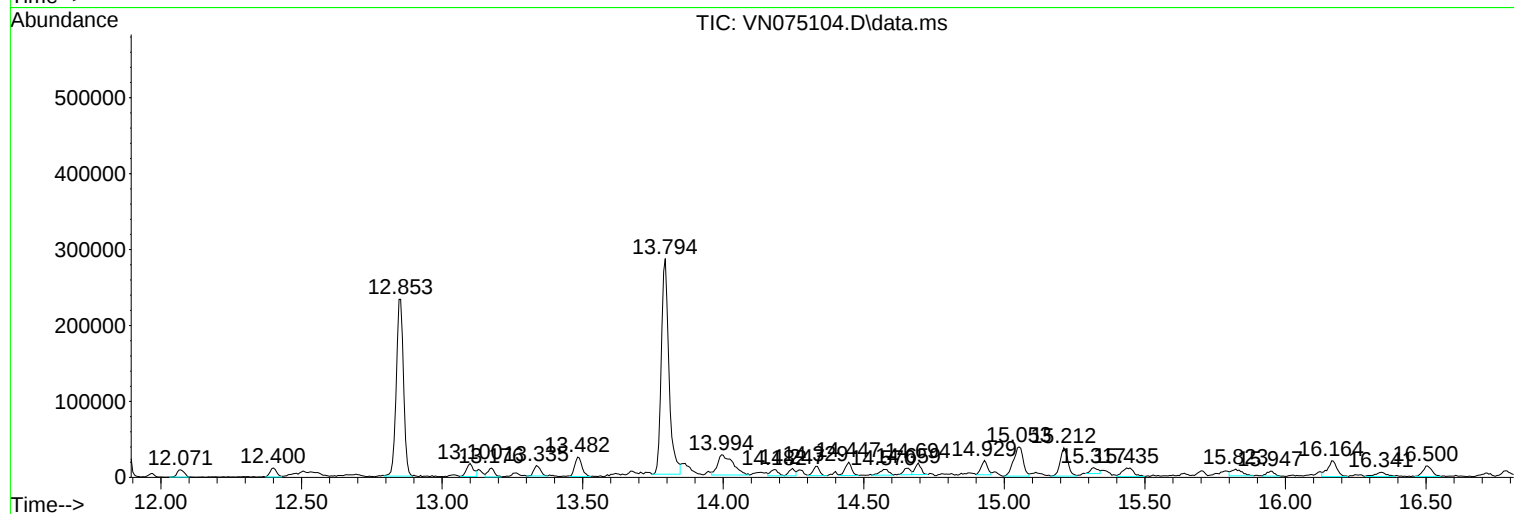
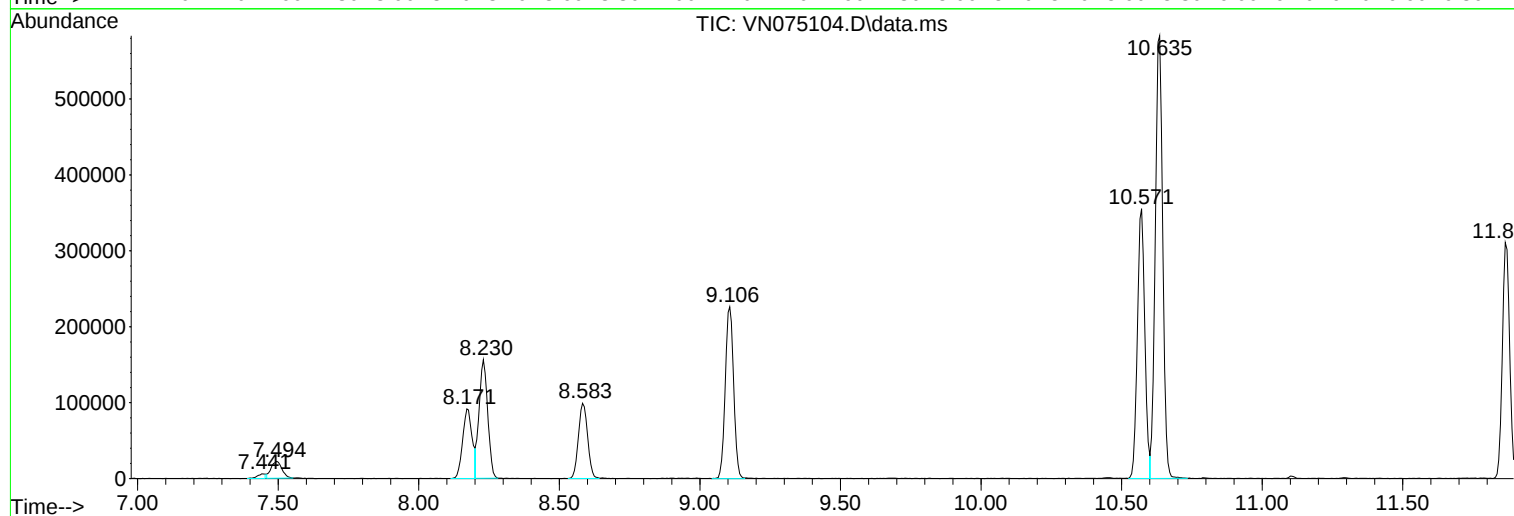
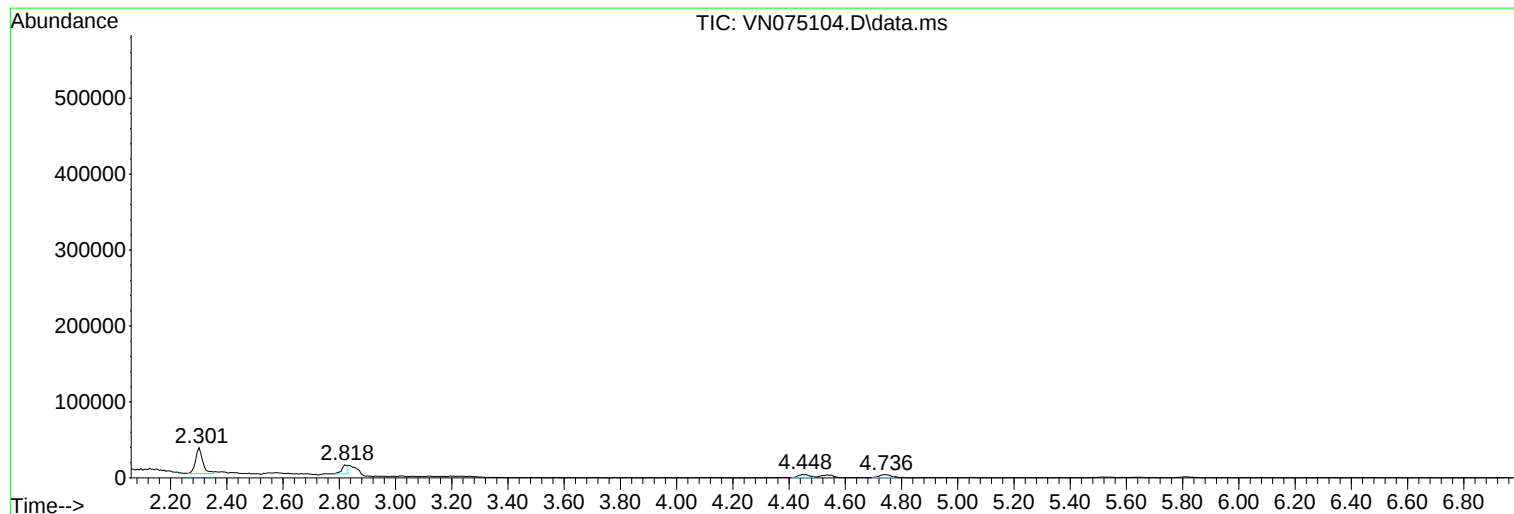
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
Quant Title : SW846 8260

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



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Instrument :
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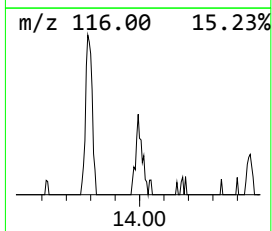
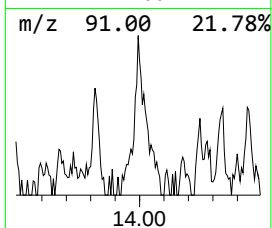
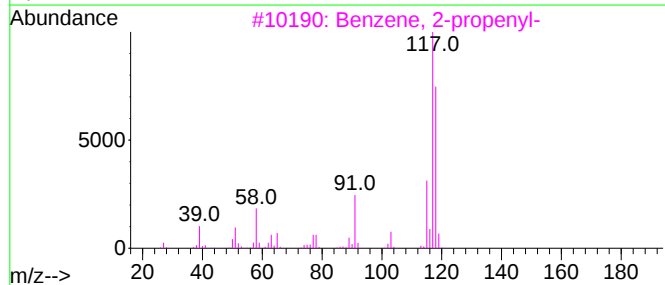
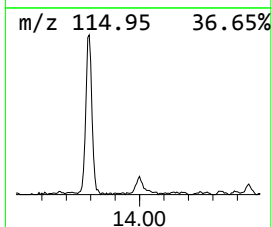
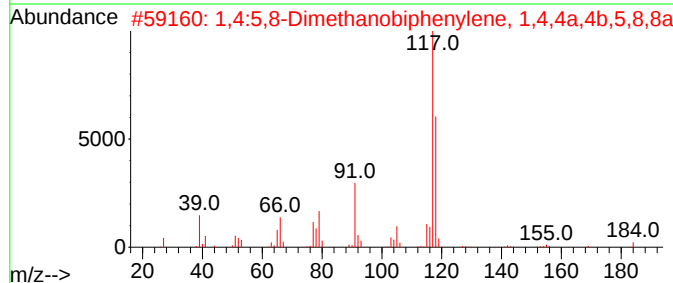
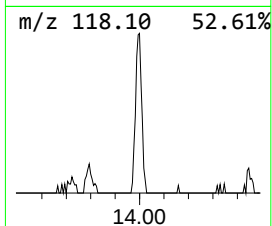
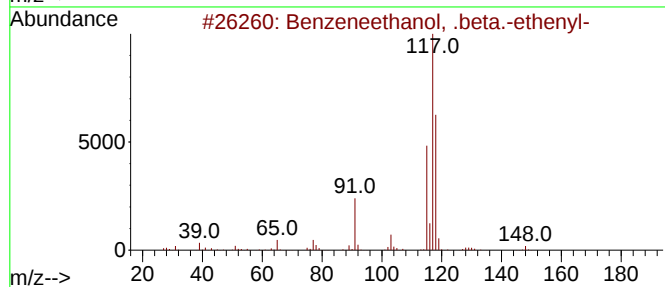
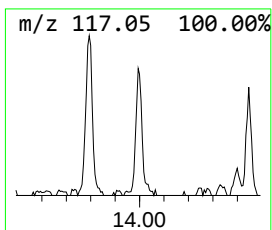
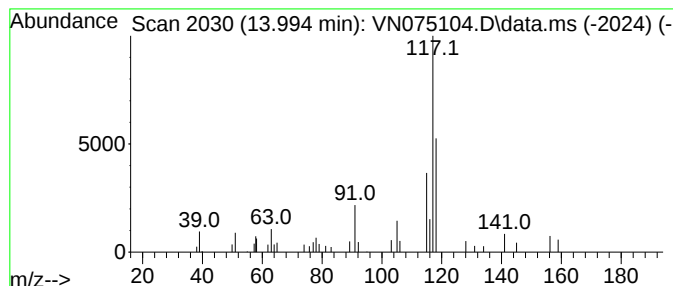
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N102622W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzeneethanol, .beta.-ethe... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.994	9.37 ug/l	99831	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneethanol, .beta.-ethenyl-	148	C10H12O	006052-63-7	64
2			1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	64
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	62
4			Indane	118	C9H10	000496-11-7	62
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	58



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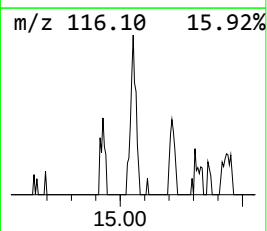
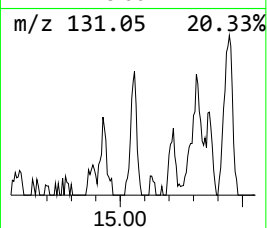
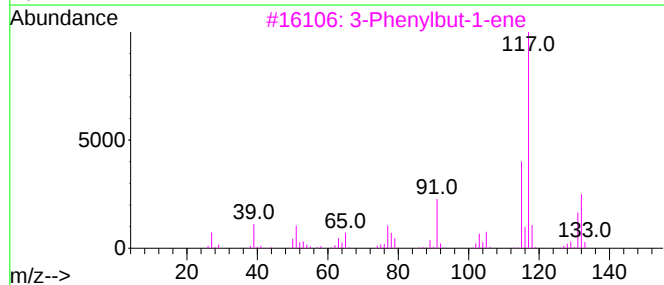
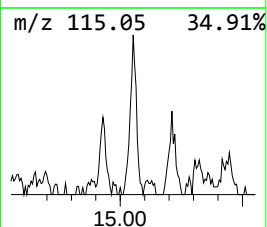
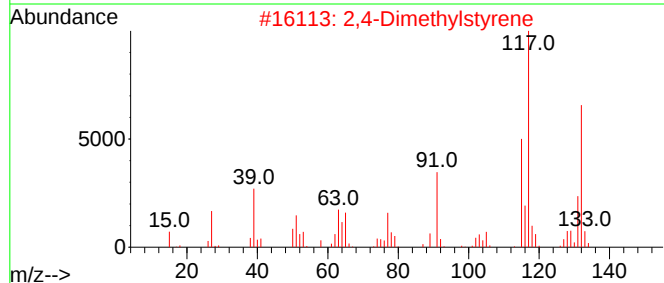
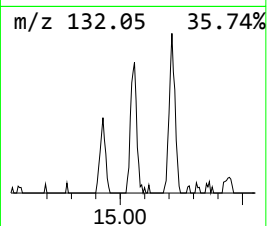
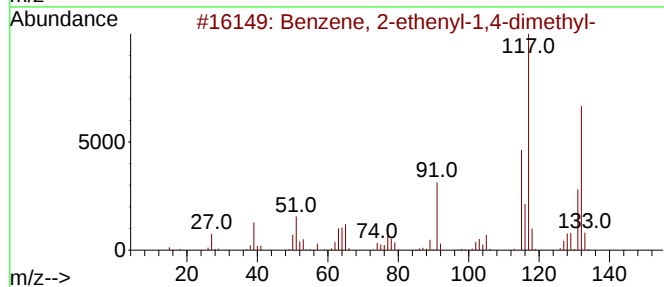
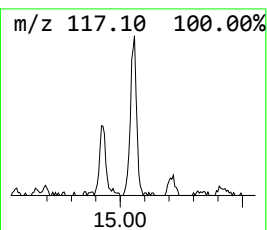
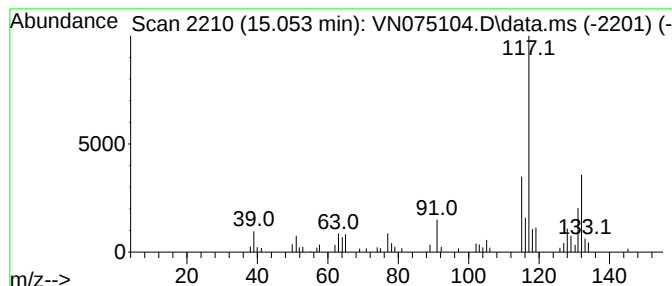
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	8.60 ug/l	91617	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	81



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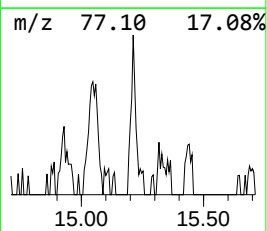
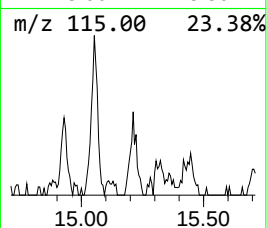
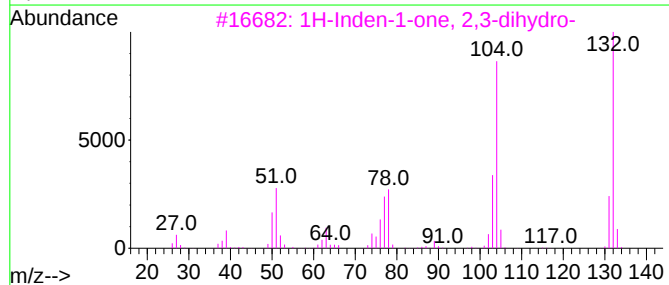
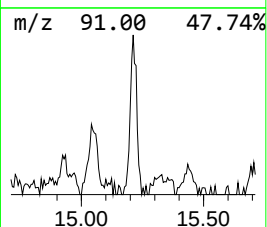
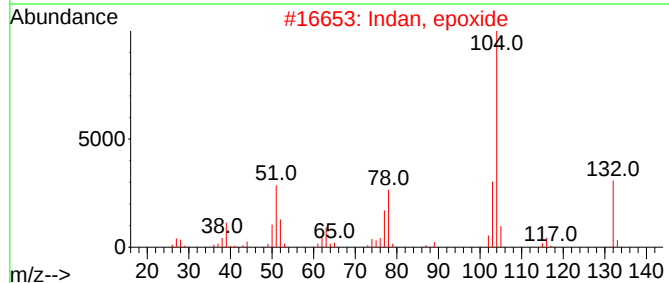
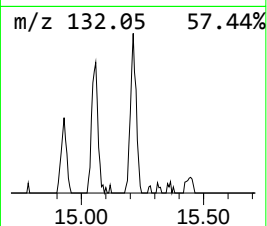
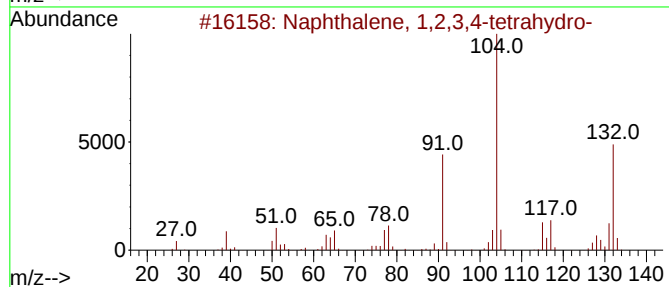
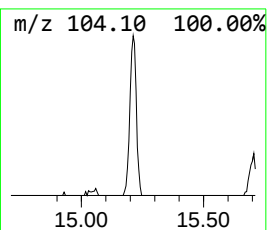
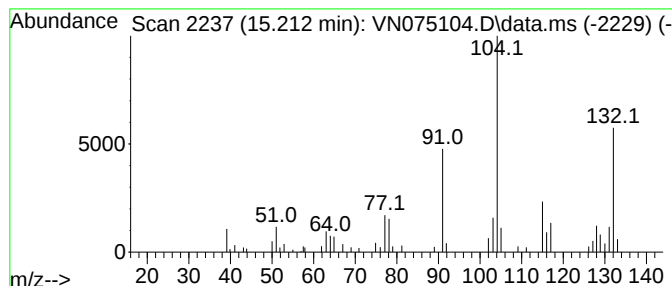
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.212	6.67 ug/l	71075	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	58
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
5			5-Aminoindole	132	C8H8N2	005192-03-0	50



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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
Benzeneethanol,...	13.994	9.4	ug/l	99831	4	13.794	532633	50.0
Benzene, 2-ethe...	15.053	8.6	ug/l	91617	4	13.794	532633	50.0
Naphthalene, 1,...	15.212	6.7	ug/l	71075	4	13.794	532633	50.0