

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
 Data File : VN075878.D
 Acq On : 19 Dec 2022 17:58
 Operator : JC\MD
 Sample : N6102-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3

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\82N121422W.M
 Title : SW846 8260

Signal : TIC: VN075878.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.524	74	80	86	rBV	40984	70501	3.24%	0.178%
2	2.836	126	133	140	rVB3	10826	26087	1.20%	0.066%
3	3.259	195	205	217	rVB2	265845	613332	28.21%	1.552%
4	3.653	262	272	285	rBV2	109896	285550	13.13%	0.723%
5	4.442	394	406	419	rBV4	20006	65839	3.03%	0.167%
6	5.147	514	526	538	rBV8	8464	35376	1.63%	0.090%
7	5.294	538	551	552	rBV3	45724	117154	5.39%	0.296%
8	5.353	554	561	578	rVB4	127887	486839	22.39%	1.232%
9	5.824	628	641	654	rBV	131746	442466	20.35%	1.120%
10	6.141	684	695	696	rBV3	15591	32719	1.50%	0.083%
11	6.324	712	726	738	rVB3	59797	177411	8.16%	0.449%
12	6.606	762	774	780	rBV2	14275	46192	2.12%	0.117%
13	6.806	798	808	815	rBV5	9756	30261	1.39%	0.077%
14	7.088	847	856	865	rBV7	12890	38065	1.75%	0.096%
15	7.230	872	880	888	rBV2	14791	41711	1.92%	0.106%
16	7.335	888	898	913	rVB	308557	877617	40.37%	2.221%
17	7.735	958	966	973	rBV2	11116	26525	1.22%	0.067%
18	7.941	993	1001	1007	rBV6	10733	26693	1.23%	0.068%
19	8.018	1007	1014	1021	rVB2	20306	48136	2.21%	0.122%
20	8.177	1030	1041	1045	rBV	246845	611773	28.14%	1.548%
21	8.235	1045	1051	1063	rVV2	572490	1810330	83.27%	4.582%
22	8.335	1063	1068	1079	rVB3	70731	174713	8.04%	0.442%
23	8.453	1079	1088	1094	rBV	91980	205150	9.44%	0.519%
24	8.518	1094	1099	1103	rVV3	31753	67656	3.11%	0.171%
25	8.582	1103	1110	1125	rVB	242474	555911	25.57%	1.407%
26	8.724	1125	1134	1141	rBV5	95874	250053	11.50%	0.633%
27	8.794	1141	1146	1152	rVV2	92523	205771	9.46%	0.521%
28	8.859	1152	1157	1166	rVB	105106	239221	11.00%	0.605%
29	8.971	1166	1176	1188	rVB5	26674	75262	3.46%	0.190%
30	9.106	1188	1199	1214	rBV	693409	1484697	68.29%	3.757%
31	9.388	1242	1247	1259	rVB2	13361	32832	1.51%	0.083%
32	9.606	1267	1284	1301	rBV	455375	1147771	52.79%	2.905%
33	9.782	1307	1314	1320	rBV2	56395	106890	4.92%	0.271%
34	9.871	1320	1329	1335	rVB3	25252	54251	2.50%	0.137%
35	10.012	1349	1353	1365	rVB4	35383	77323	3.56%	0.196%
36	10.247	1389	1393	1400	rVB6	13778	28359	1.30%	0.072%

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37	10.335	1400	1408	1415	rBV3	26496	67625	3.11%	0.171%
38	10.459	1420	1429	1433	rBV2	34011	69268	3.19%	0.175%
39	10.571	1440	1448	1455	rBV	1051880	1979064	91.03%	5.009%
40	10.747	1472	1478	1486	rVB5	40202	98811	4.54%	0.250%

41	10.912	1500	1506	1513	rBV2	53635	96707	4.45%	0.245%
42	11.006	1513	1522	1529	rVB6	47365	129612	5.96%	0.328%
43	11.088	1529	1536	1543	rVB7	11397	28579	1.31%	0.072%
44	11.259	1558	1565	1567	rBV5	13457	27431	1.26%	0.069%
45	11.300	1567	1572	1577	rVB	55470	92567	4.26%	0.234%

46	11.394	1583	1588	1589	rBV3	16538	24738	1.14%	0.063%
47	11.423	1589	1593	1598	rVV	44165	79732	3.67%	0.202%
48	11.476	1598	1602	1608	rVB3	25623	45075	2.07%	0.114%
49	11.682	1631	1637	1646	rBV6	16586	39022	1.79%	0.099%
50	11.870	1661	1669	1679	rBV	850375	1572966	72.35%	3.981%

51	11.965	1679	1685	1695	rVB	128799	231285	10.64%	0.585%
52	12.076	1695	1704	1717	rBV	798988	1598615	73.53%	4.046%
53	12.400	1749	1759	1778	rVB	551118	979951	45.07%	2.480%
54	12.582	1780	1790	1796	rBV6	33279	72864	3.35%	0.184%
55	12.700	1801	1810	1816	rVB	163508	273674	12.59%	0.693%

56	12.853	1829	1836	1845	rBV	674756	1138125	52.35%	2.880%
57	13.041	1860	1868	1873	rVV	284288	481230	22.13%	1.218%
58	13.100	1873	1878	1881	rVV	504683	839334	38.61%	2.124%
59	13.135	1881	1884	1888	rVV	327061	521165	23.97%	1.319%
60	13.176	1888	1891	1899	rVB	216434	326273	15.01%	0.826%

61	13.341	1911	1919	1927	rBV	745989	1252961	57.63%	3.171%
62	13.488	1936	1944	1955	rVB	1324997	2174135	100.00%	5.502%
63	13.617	1957	1966	1971	rBV2	53688	123731	5.69%	0.313%
64	13.676	1971	1976	1981	rVB	104074	160772	7.39%	0.407%
65	13.729	1981	1985	1990	rBV	47332	70890	3.26%	0.179%

66	13.794	1990	1996	2008	rVV2	831869	2012078	92.55%	5.092%
67	13.888	2008	2012	2017	rVB	15862	26897	1.24%	0.068%
68	13.953	2017	2023	2025	rBV	167481	251368	11.56%	0.636%
69	13.988	2025	2029	2033	rVV2	353427	676348	31.11%	1.712%
70	14.035	2033	2037	2046	rVB4	225693	504011	23.18%	1.276%

71	14.117	2046	2051	2057	rVB	56396	90471	4.16%	0.229%
72	14.188	2057	2063	2068	rBV	157053	248945	11.45%	0.630%
73	14.247	2068	2073	2076	rBV	310820	514462	23.66%	1.302%
74	14.276	2076	2078	2083	rVV	193380	255372	11.75%	0.646%
75	14.335	2083	2088	2094	rVB	630083	984618	45.29%	2.492%

76	14.400	2094	2099	2102	rBV	76559	133098	6.12%	0.337%
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77	14.447	2102	2107	2115	rVV2	640390	1086200	49.96%	2.749%
78	14.511	2115	2118	2123	rVB5	21433	34969	1.61%	0.088%
79	14.576	2123	2129	2135	rVB2	107543	190000	8.74%	0.481%
80	14.659	2135	2143	2147	rBV	479366	823483	37.88%	2.084%
81	14.694	2147	2149	2154	rVV	239455	338785	15.58%	0.857%
82	14.741	2154	2157	2160	rVV3	76798	98574	4.53%	0.249%
83	14.776	2160	2163	2167	rVB2	21654	30017	1.38%	0.076%
84	14.882	2172	2181	2184	rBV3	75258	126472	5.82%	0.320%
85	14.929	2184	2189	2195	rBV	354009	585494	26.93%	1.482%
86	15.053	2201	2210	2216	rBV3	653287	1582086	72.77%	4.004%
87	15.106	2216	2219	2229	rVB4	50532	126614	5.82%	0.320%
88	15.211	2231	2237	2244	rBV3	107359	210092	9.66%	0.532%
89	15.323	2244	2256	2260	rVV4	218498	603403	27.75%	1.527%
90	15.364	2260	2263	2268	rVV	125609	221936	10.21%	0.562%
91	15.441	2268	2276	2290	rVB3	236949	613937	28.24%	1.554%
92	15.641	2303	2310	2318	rVB	569203	1103092	50.74%	2.792%
93	15.753	2323	2329	2334	rVV3	80225	164622	7.57%	0.417%
94	15.811	2334	2339	2350	rVB3	54025	129289	5.95%	0.327%
95	15.947	2354	2362	2371	rBV	71145	146836	6.75%	0.372%
96	16.135	2383	2394	2410	rVV3	49287	178627	8.22%	0.452%
97	16.258	2410	2415	2421	rVV3	16212	32676	1.50%	0.083%
98	16.353	2422	2431	2443	rVB2	39736	103351	4.75%	0.262%
99	16.505	2447	2457	2465	rBV7	12395	35684	1.64%	0.090%
100	16.717	2481	2493	2498	rBV4	15061	36815	1.69%	0.093%

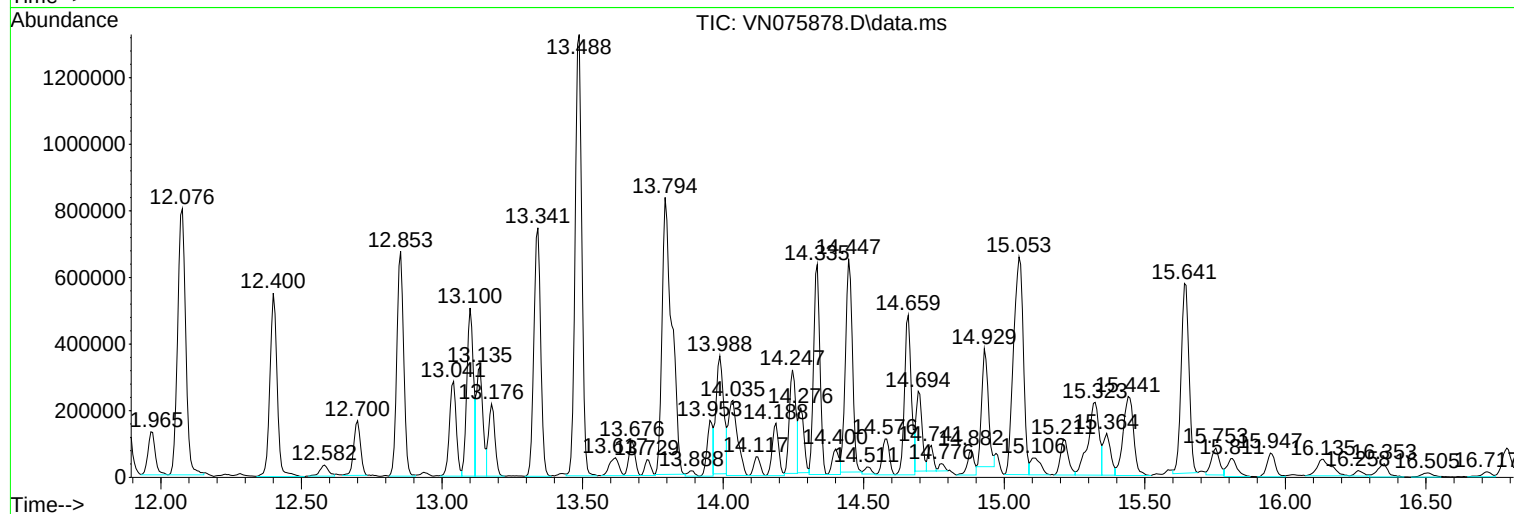
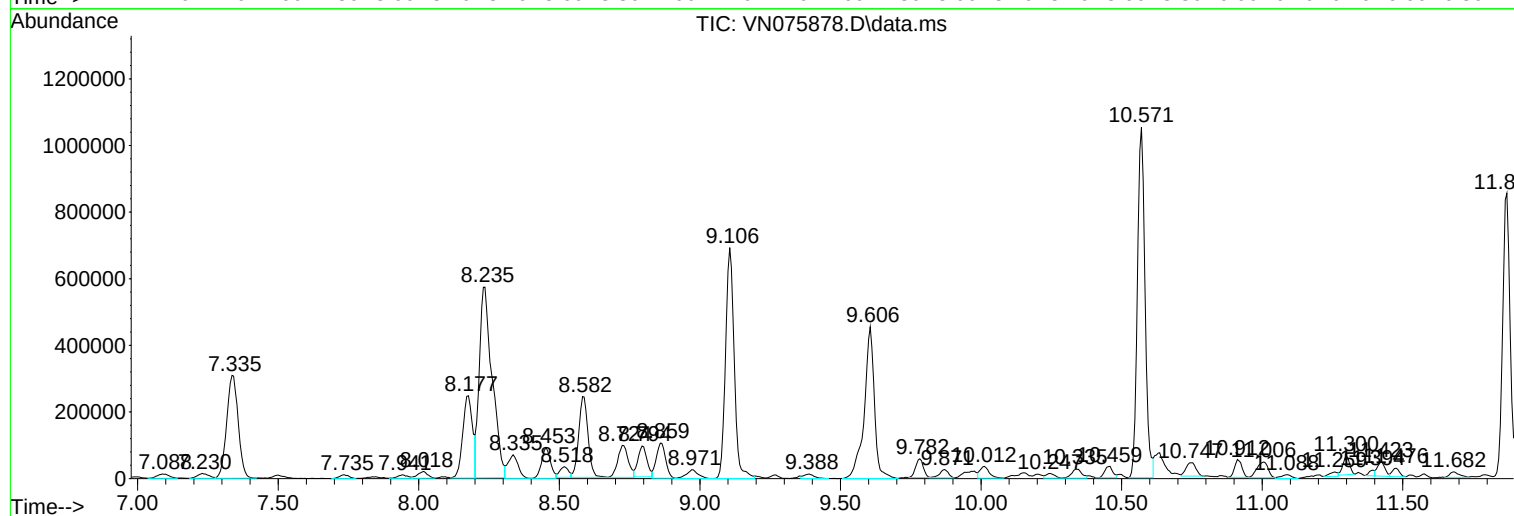
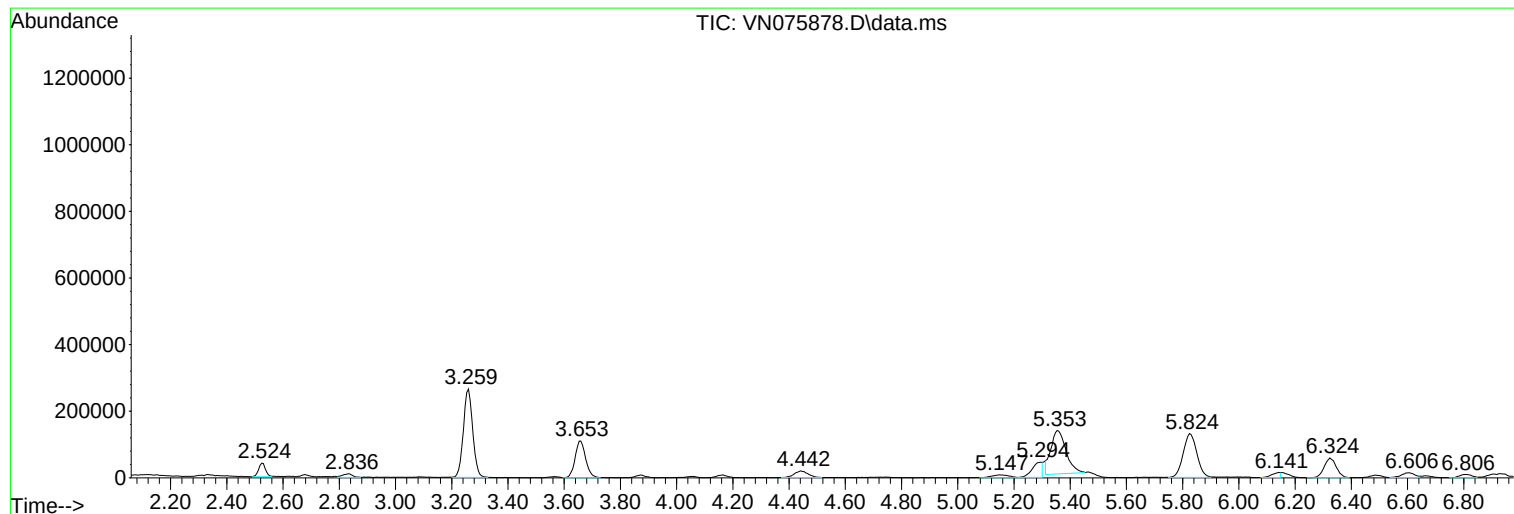
Sum of corrected areas: 39513361

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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



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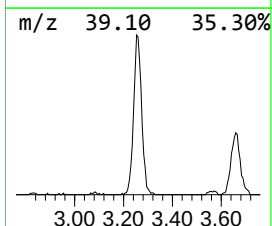
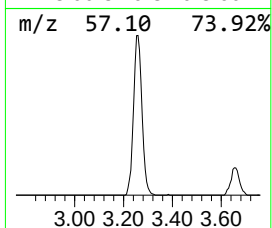
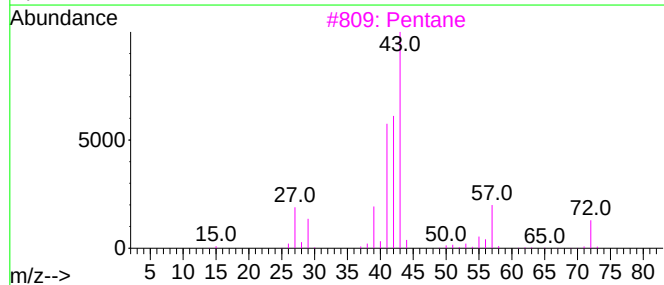
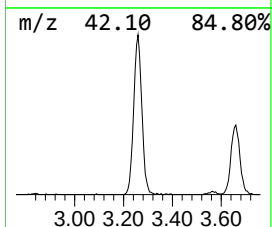
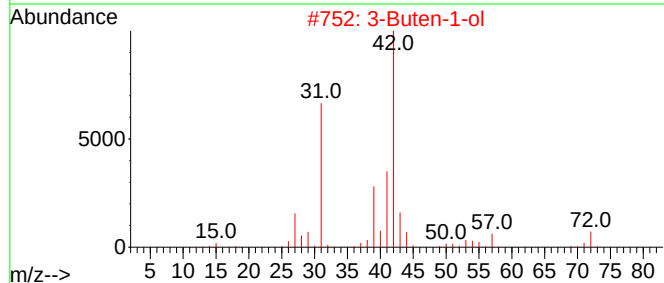
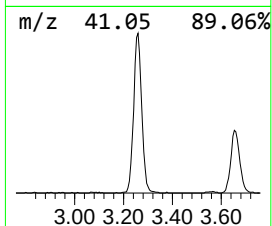
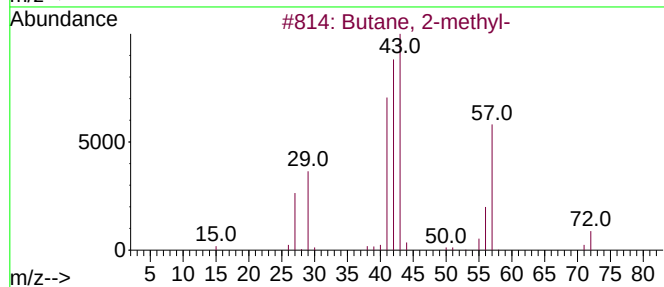
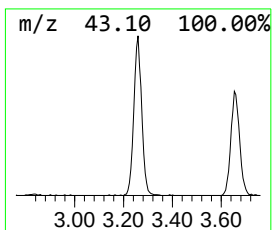
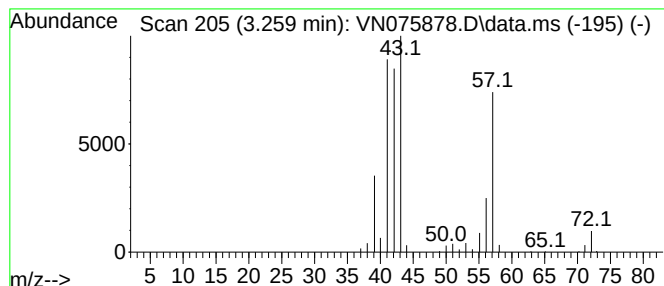
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.259	16.94 ug/l	613332	Pentafluorobenzene	8.230

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	86
2			3-Buten-1-ol	72	C4H8O	000627-27-0	37
3			Pentane	72	C5H12	000109-66-0	28
4			Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	27
5			Propane, 2-methyl-2-nitro-	103	C4H9NO2	000594-70-7	10



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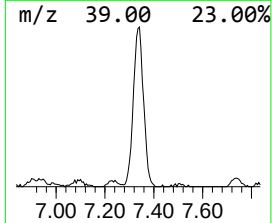
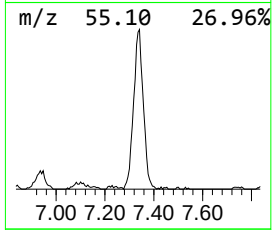
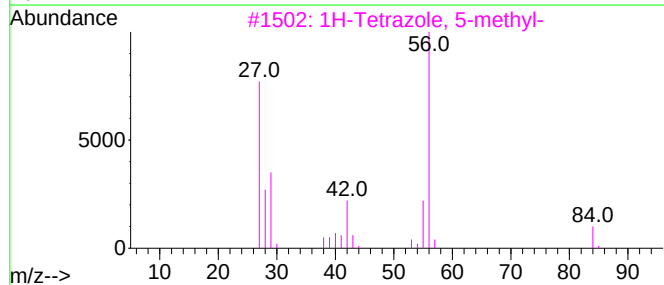
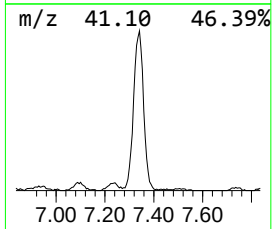
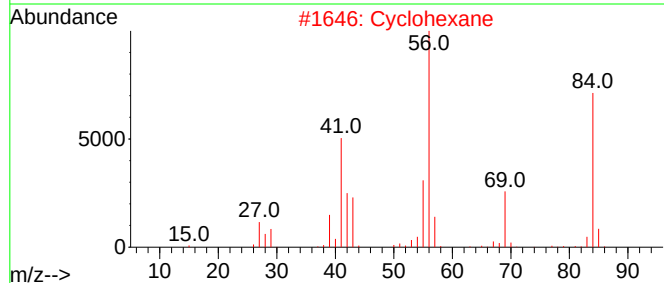
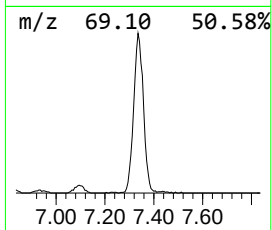
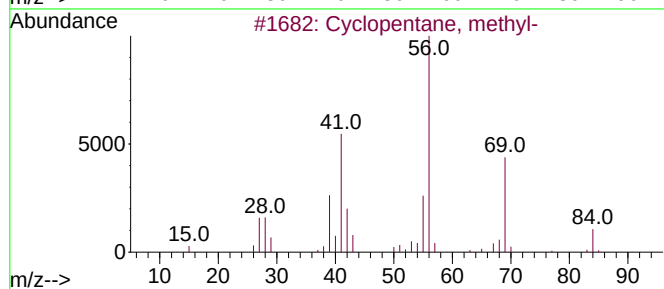
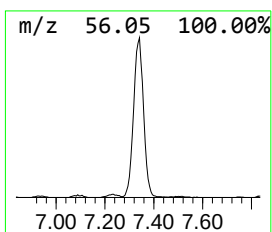
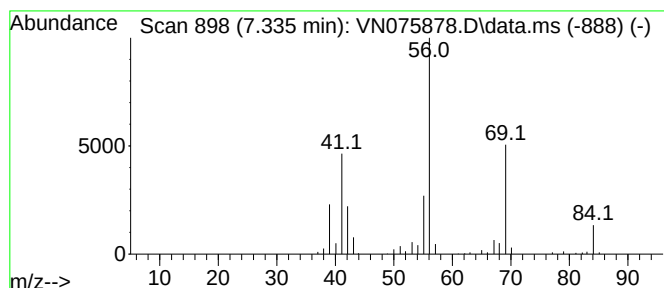
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.335	24.24 ug/l	877617	Pentafluorobenzene	8.230

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclohexane	84	C6H12	000110-82-7	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		Cyclobutane, butyl-	112	C8H16	013152-44-8	50



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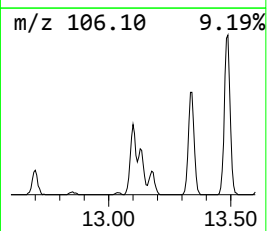
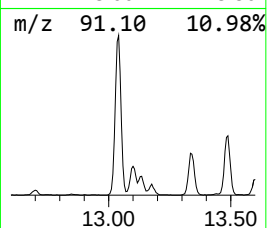
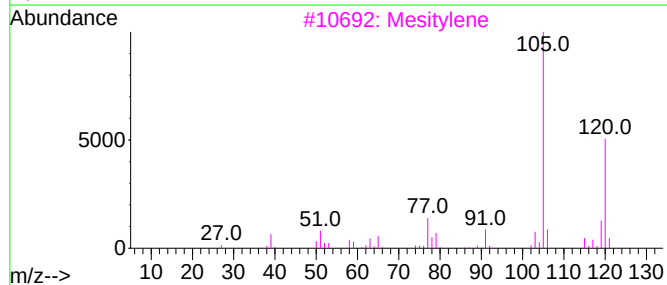
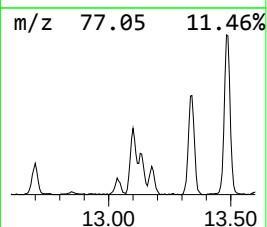
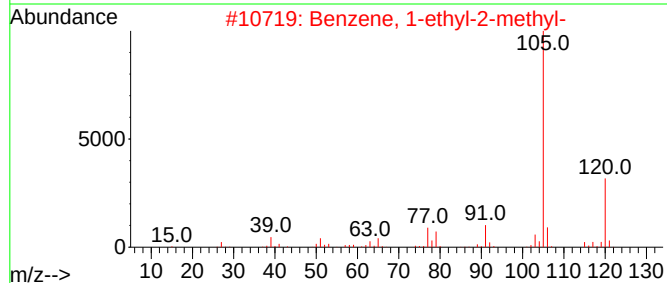
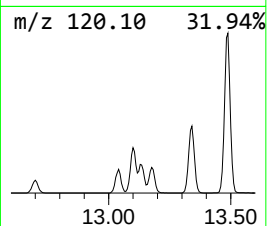
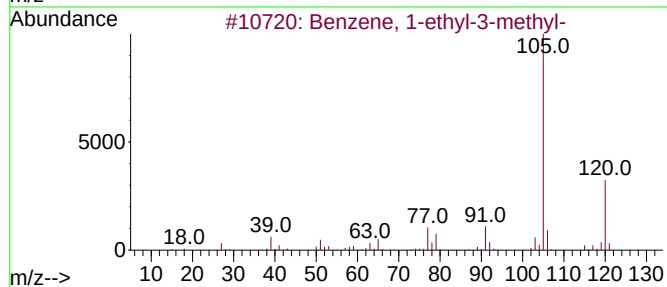
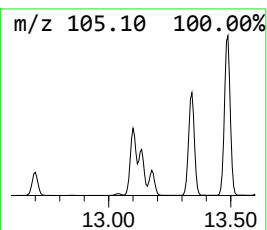
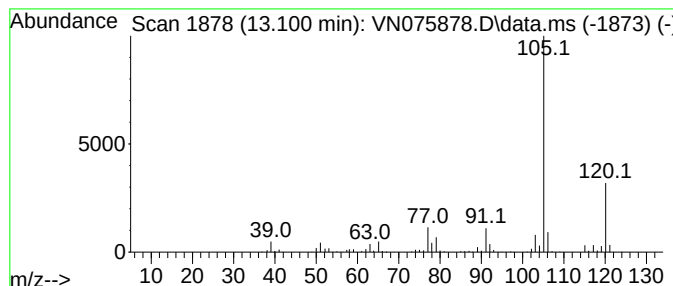
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.100	20.86 ug/l	839334	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

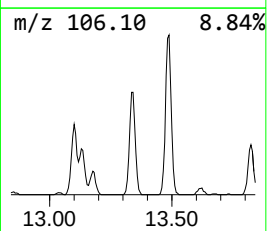
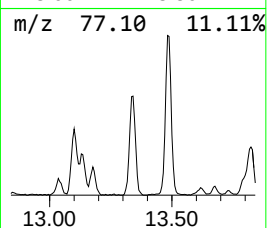
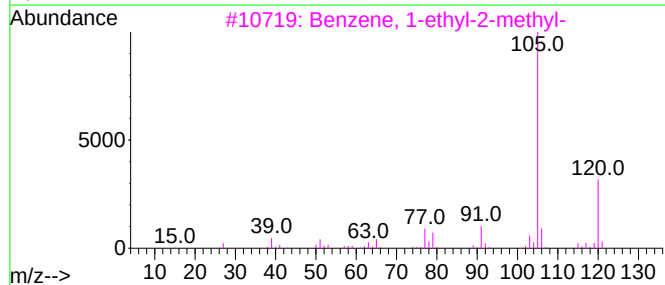
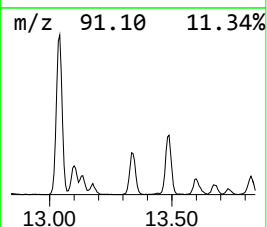
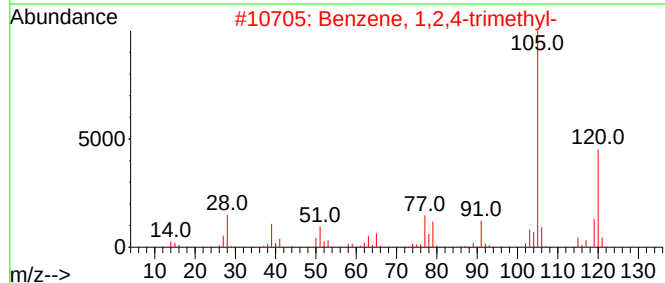
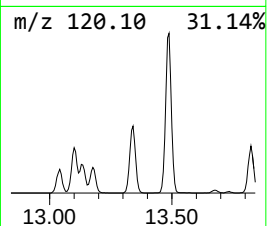
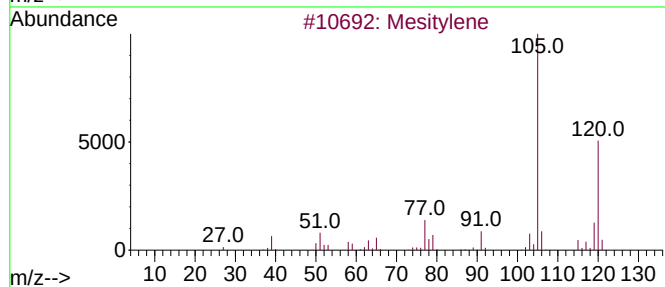
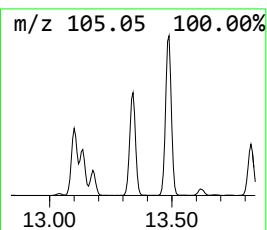
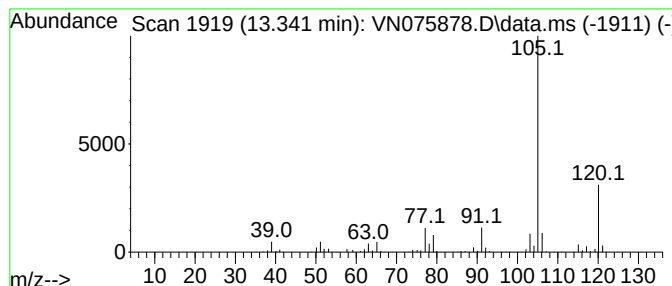
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.341	31.14 ug/l	1252960	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	90
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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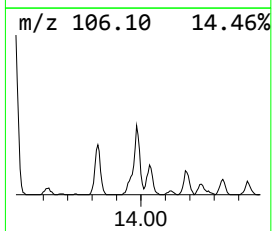
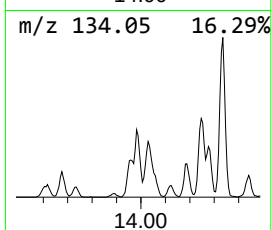
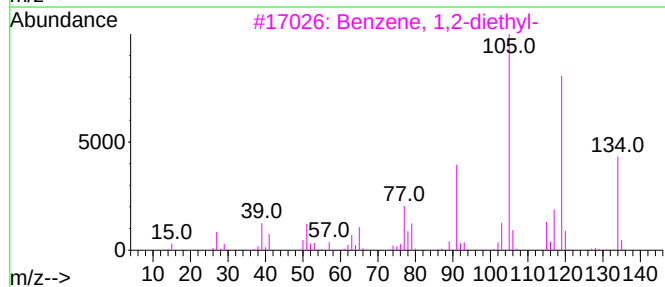
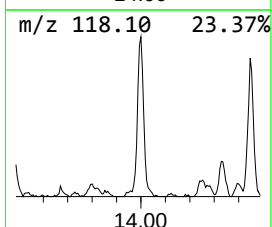
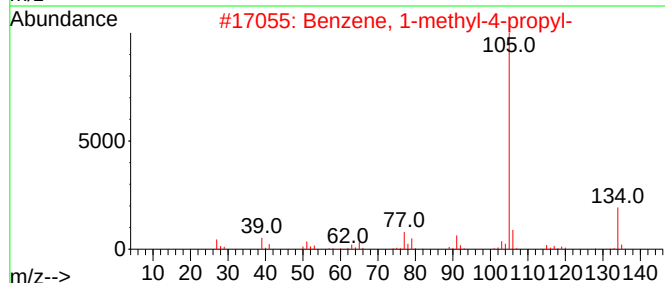
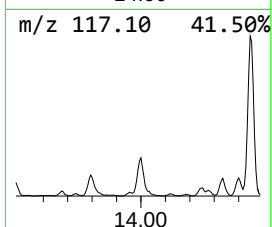
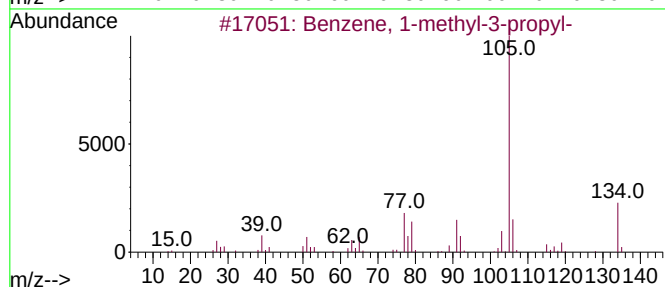
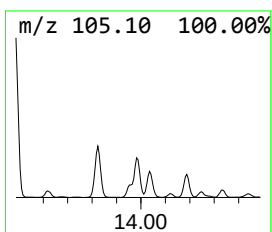
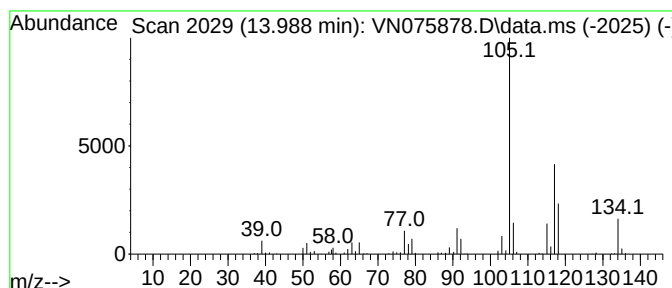
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Quant Title : SW846 8260

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

Peak Number 5 unknown13.988 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.988	16.81 ug/l	676348	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	35
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	35
4			2-Indanol	134	C9H10O	004254-29-9	35
5			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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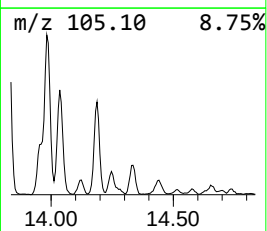
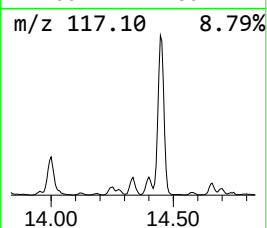
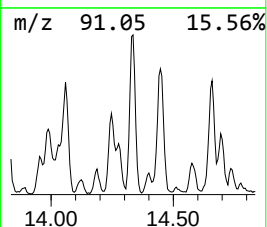
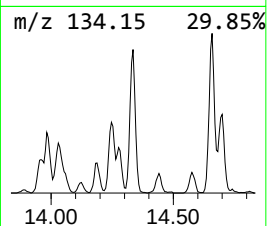
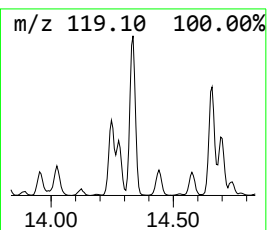
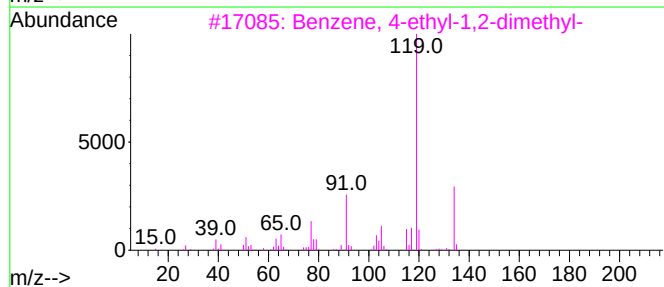
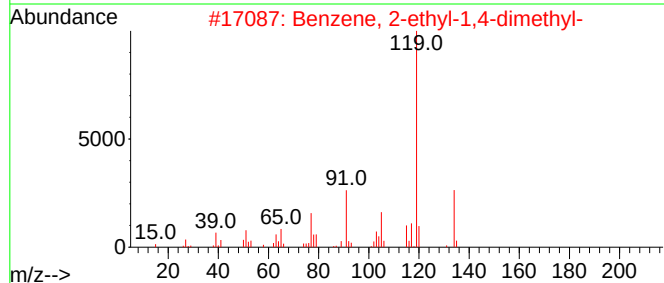
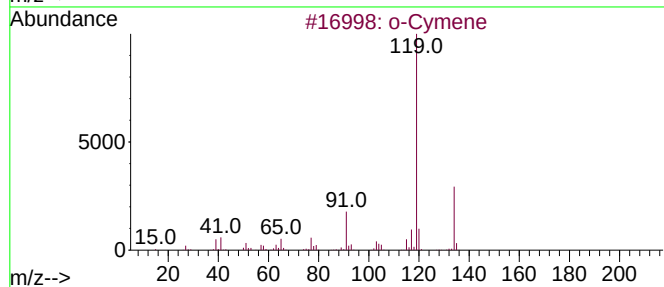
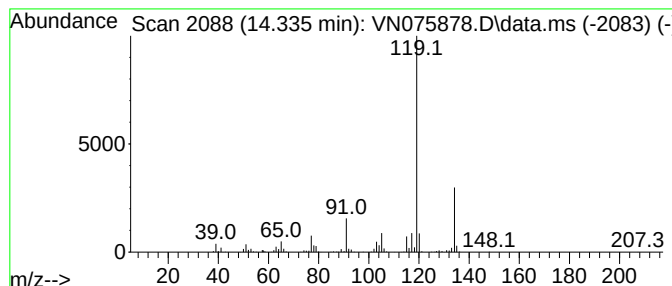
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Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.335	24.47 ug/l	984618	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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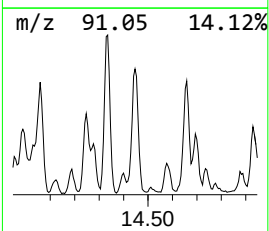
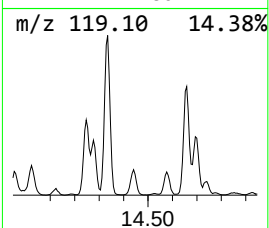
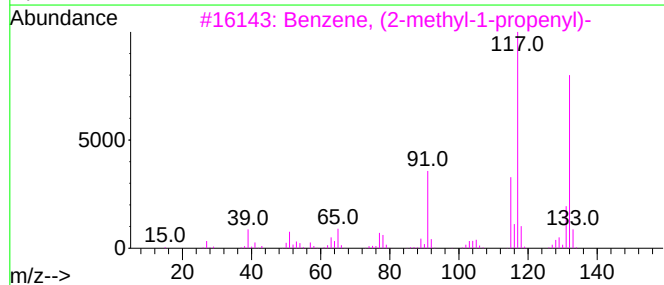
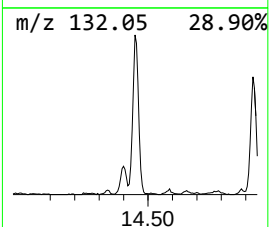
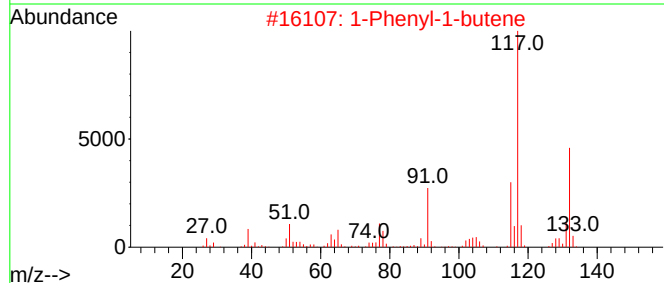
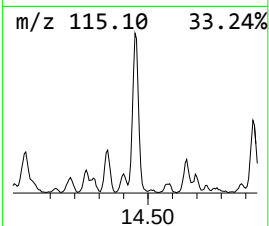
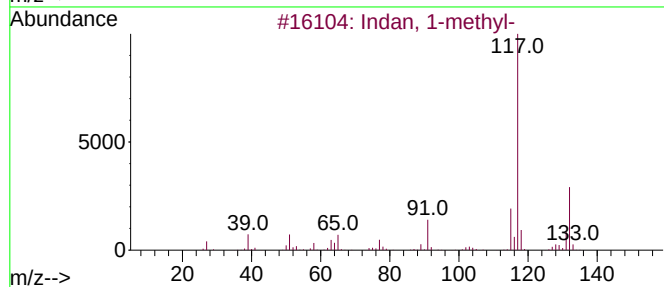
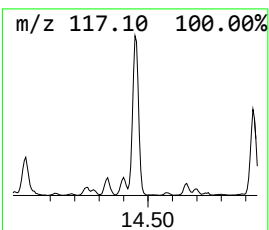
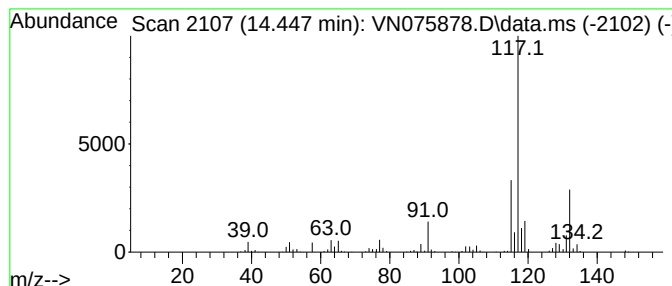
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.447	26.99 ug/l	1086200	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
3			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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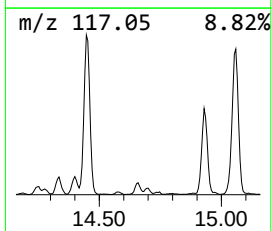
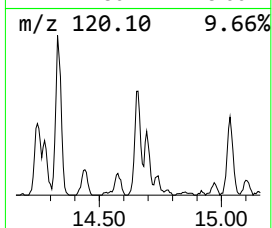
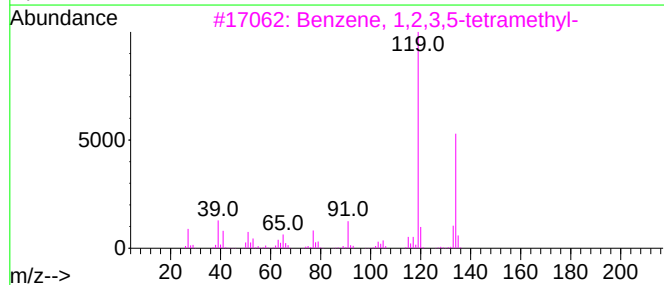
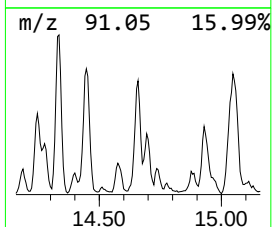
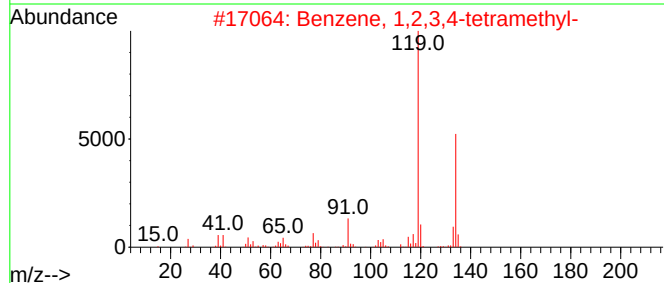
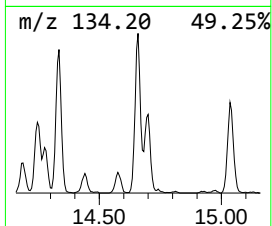
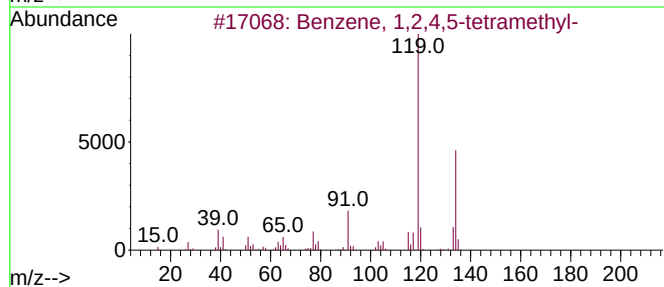
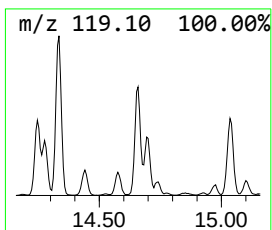
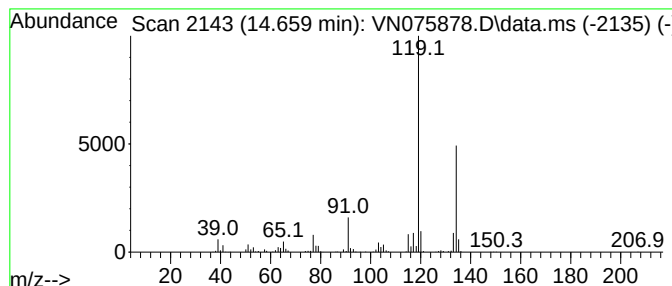
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.659	20.46 ug/l	823483	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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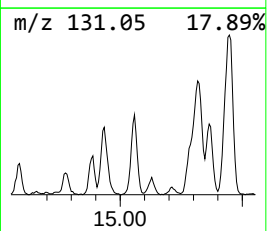
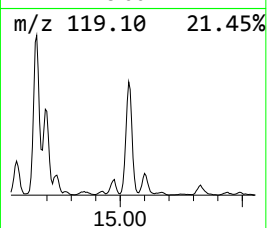
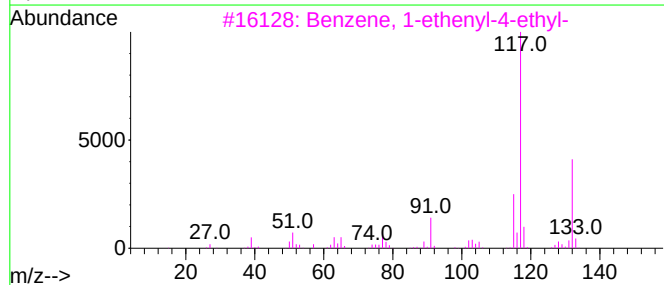
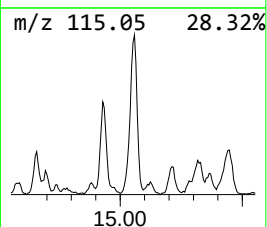
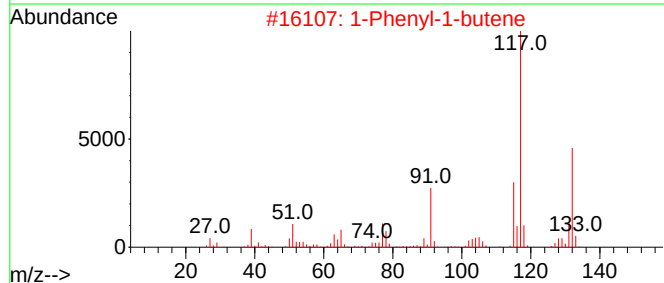
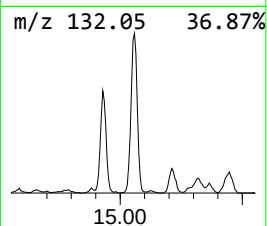
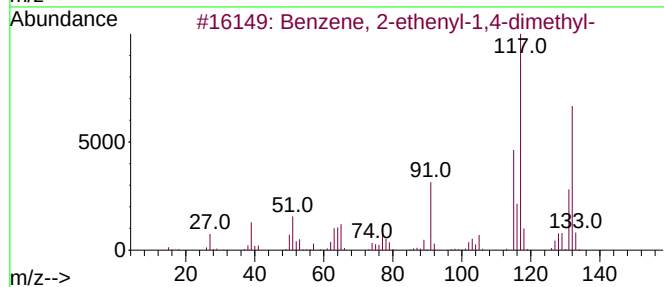
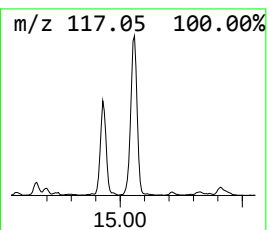
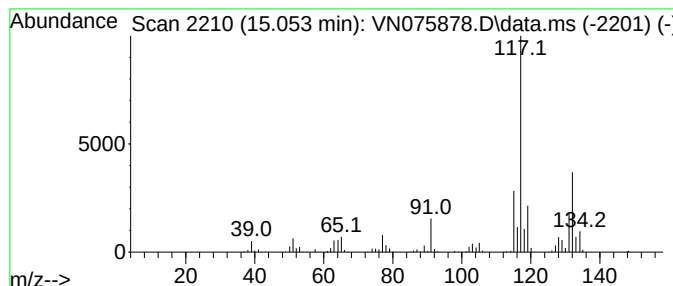
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.053	39.31 ug/l	1582090	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	95
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	91
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
4			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

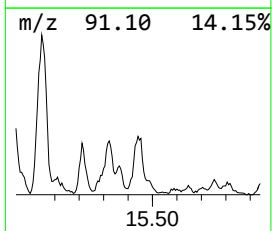
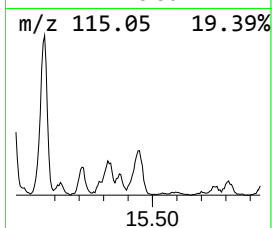
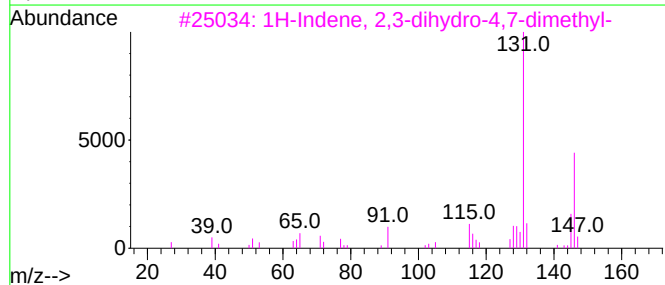
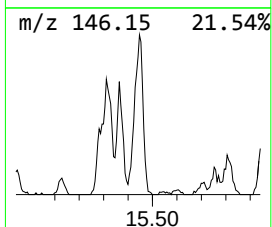
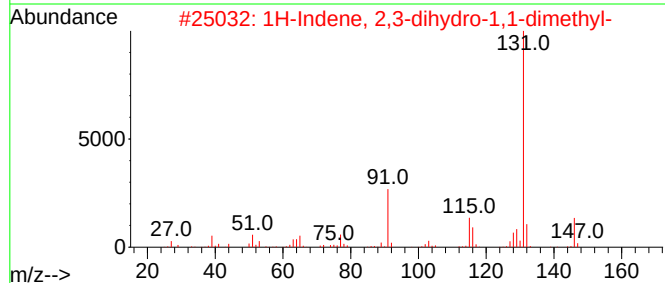
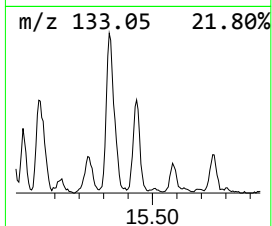
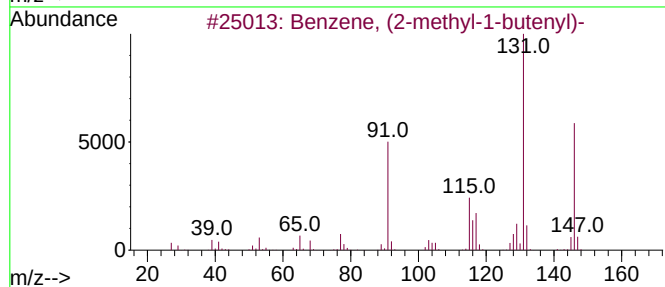
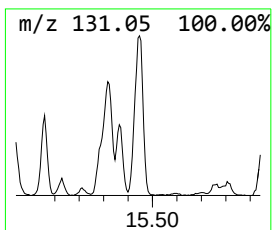
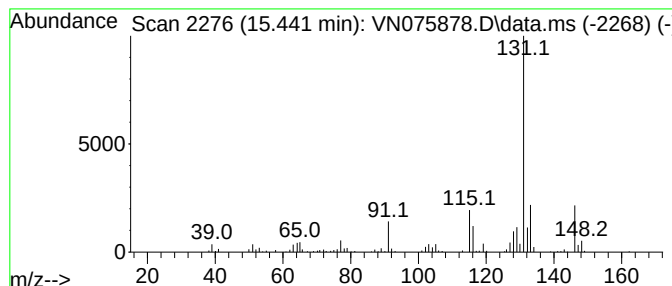
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.441	15.26 ug/l	613937	1,4-Dichlorobenzene-d4	13.794

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
2			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	87
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	87
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121922\
Data File : VN075878.D
Acq On : 19 Dec 2022 17:58
Operator : JC\MD
Sample : N6102-04
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methyl-	3.259	16.9	ug/l	613332	1	8.230	1810330	50.0
Cyclopentane, m...	7.335	24.2	ug/l	877617	1	8.230	1810330	50.0
Benzene, 1-ethy...	13.100	20.9	ug/l	839334	4	13.794	2012080	50.0
Benzene, 1,2,4-...	13.341	31.1	ug/l	1252960	4	13.794	2012080	50.0
unknown13.988	13.988	16.8	ug/l	676348	4	13.794	2012080	50.0
o-Cymene	14.335	24.5	ug/l	984618	4	13.794	2012080	50.0
Indan, 1-methyl-	14.447	27.0	ug/l	1086200	4	13.794	2012080	50.0
Benzene, 1,2,4,...	14.659	20.5	ug/l	823483	4	13.794	2012080	50.0
Benzene, 2-ethe...	15.053	39.3	ug/l	1582090	4	13.794	2012080	50.0
Benzene, (2-met...	15.441	15.3	ug/l	613937	4	13.794	2012080	50.0