

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
 Data File : VN072772.D
 Acq On : 10 Jun 2022 18:12
 Operator : JC\MD
 Sample : N3293-13
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 40805

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\82N060822W.M
 Title : SW846 8260

Signal : TIC: VN072772.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.287	381	391	404	rBV	49714	151013	1.32%	0.246%
2	8.016	1016	1025	1031	rBV2	244703	602154	5.27%	0.979%
3	8.081	1031	1036	1047	rVB2	267452	610434	5.35%	0.993%
4	8.451	1086	1099	1117	rBV3	650713	1872462	16.40%	3.046%
5	8.963	1178	1186	1199	rBV	399571	831576	7.28%	1.353%
6	10.439	1427	1437	1442	rBV	884952	1667099	14.60%	2.712%
7	10.504	1442	1448	1471	rVB	6296883	11418014	100.00%	18.573%
8	11.739	1651	1658	1668	rBV	576268	1034972	9.06%	1.683%
9	11.845	1668	1676	1685	rVV	1252062	2155565	18.88%	3.506%
10	11.945	1685	1693	1713	rVB	4448060	8320064	72.87%	13.533%
11	12.274	1741	1749	1760	rBV	2956166	4999791	43.79%	8.133%
12	12.574	1794	1800	1807	rBV	81574	130285	1.14%	0.212%
13	12.727	1819	1826	1837	rVB	754027	1260373	11.04%	2.050%
14	12.916	1851	1858	1863	rBV	249200	399724	3.50%	0.650%
15	12.980	1863	1869	1873	rVV	1293840	2272081	19.90%	3.696%
16	13.010	1873	1874	1878	rVV	538373	633431	5.55%	1.030%
17	13.057	1878	1882	1894	rVB	643747	1001332	8.77%	1.629%
18	13.215	1903	1909	1917	rBV	697589	1134061	9.93%	1.845%
19	13.363	1927	1934	1949	rVB	2979658	4762265	41.71%	7.746%
20	13.674	1980	1987	1988	rBV	581831	882138	7.73%	1.435%
21	13.704	1988	1992	1999	rVB	1126651	1899314	16.63%	3.089%
22	13.874	2015	2021	2037	rVB3	859450	2108175	18.46%	3.429%
23	14.062	2047	2053	2058	rBV2	112792	195818	1.71%	0.319%
24	14.127	2058	2064	2067	rVV	248328	431782	3.78%	0.702%
25	14.157	2067	2069	2074	rVV	222582	282759	2.48%	0.460%
26	14.215	2074	2079	2084	rVV	380268	600054	5.26%	0.976%
27	14.274	2084	2089	2093	rVV	77874	133579	1.17%	0.217%
28	14.327	2093	2098	2107	rVB2	275510	465367	4.08%	0.757%
29	14.457	2114	2120	2126	rVB	154134	247184	2.16%	0.402%
30	14.539	2126	2134	2137	rBV	298855	516590	4.52%	0.840%
31	14.574	2137	2140	2145	rVB	443753	646711	5.66%	1.052%
32	14.804	2174	2179	2185	rBV	361901	573732	5.02%	0.933%
33	14.927	2191	2200	2206	rBV3	705505	1486119	13.02%	2.417%
34	15.080	2220	2226	2234	rVV3	134249	252700	2.21%	0.411%
35	15.192	2234	2245	2257	rVV3	153439	546622	4.79%	0.889%
36	15.304	2257	2264	2273	rVB4	136290	328495	2.88%	0.534%

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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_N\methods\82N060822W.M
Title : SW846 8260

37	15.504	2290	2298	2306	rBV	1368310	2548022	22.32%	4.145%
38	15.668	2321	2326	2340	rVB2	59371	140986	1.23%	0.229%
39	15.804	2340	2349	2357	rBV	105734	217148	1.90%	0.353%
40	15.980	2365	2379	2384	rBV2	76855	230786	2.02%	0.375%
41	16.186	2406	2414	2428	rVB5	64613	180372	1.58%	0.293%
42	16.615	2478	2487	2502	rVB	584641	1306767	11.44%	2.126%

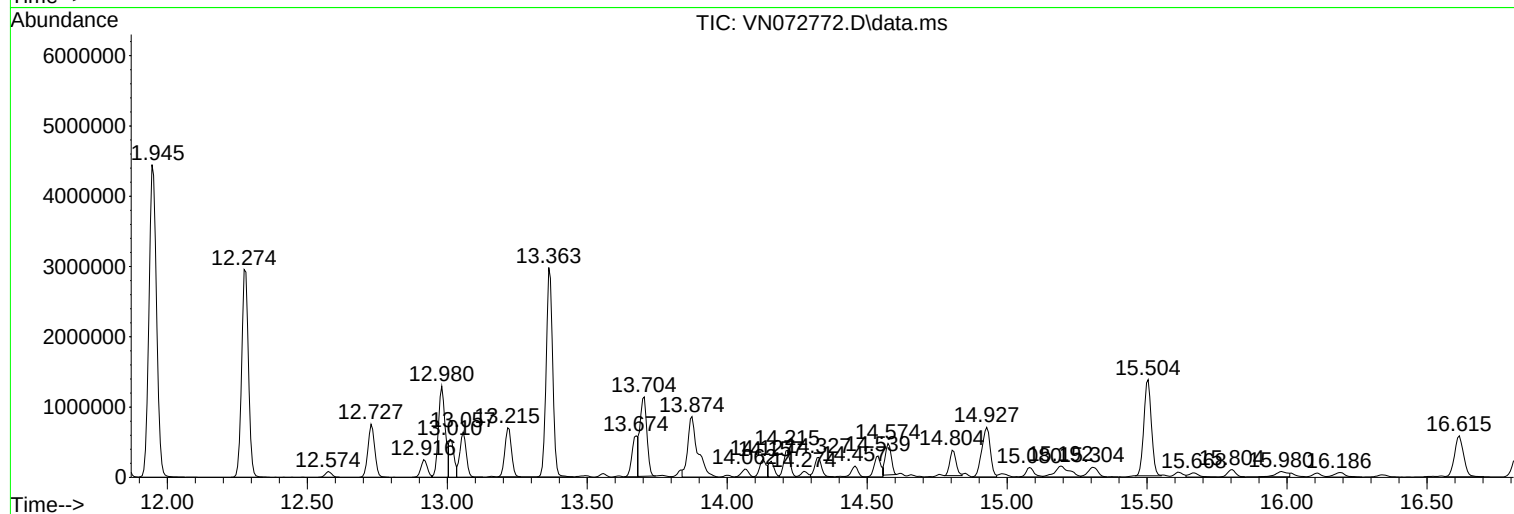
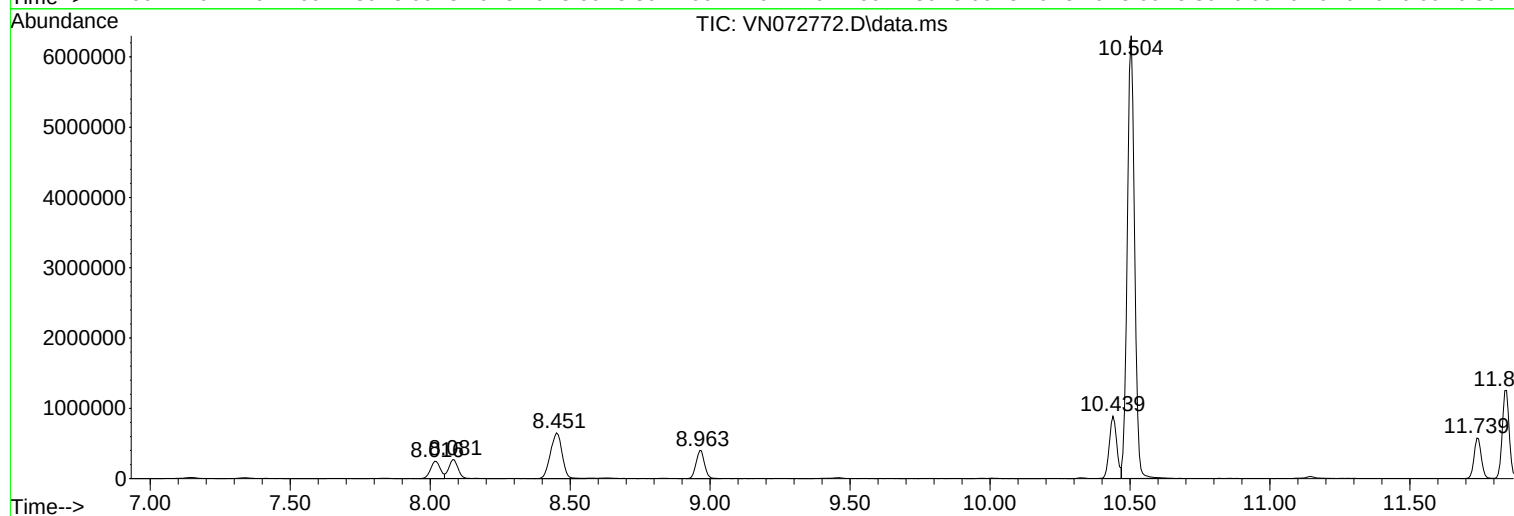
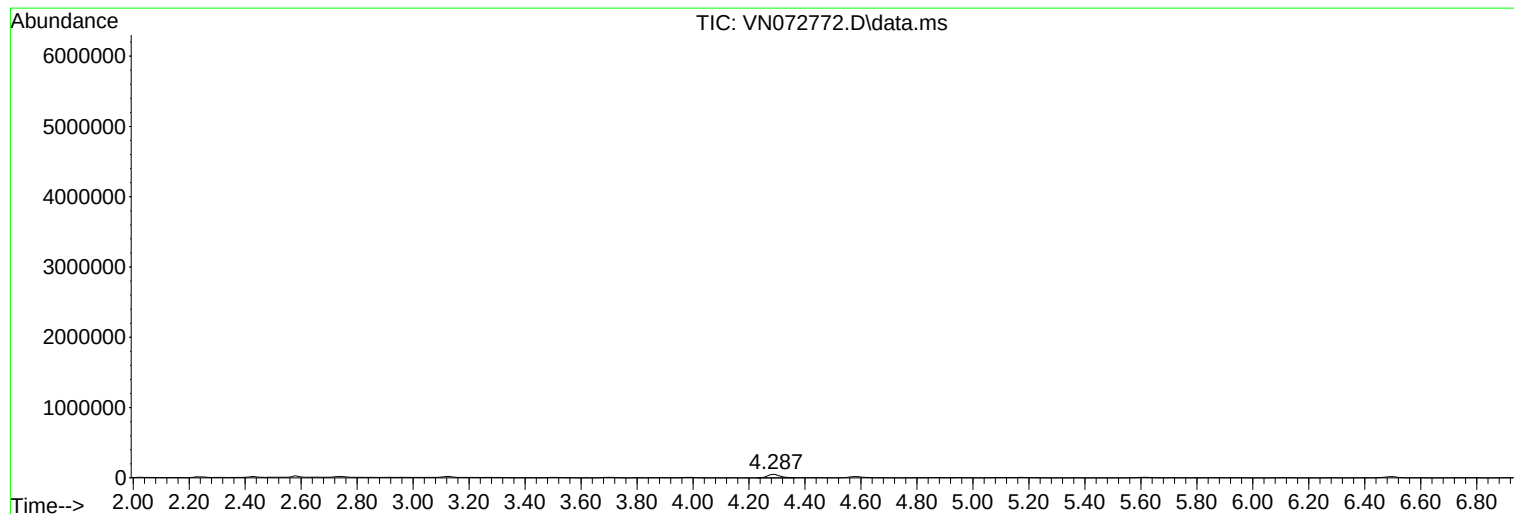
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

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



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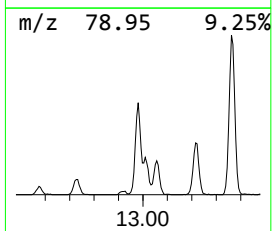
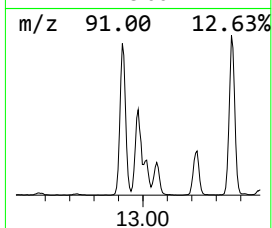
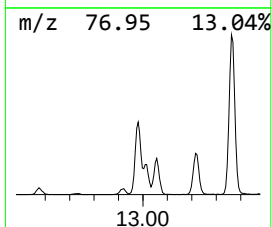
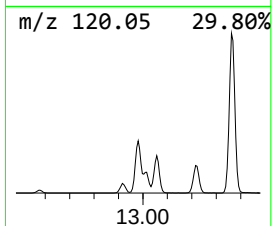
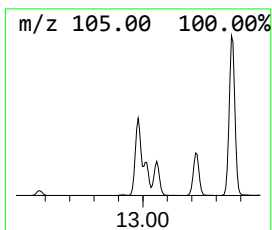
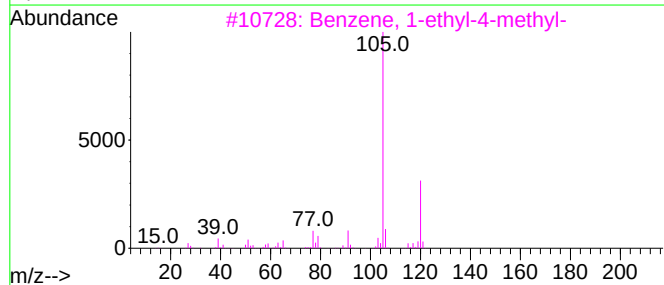
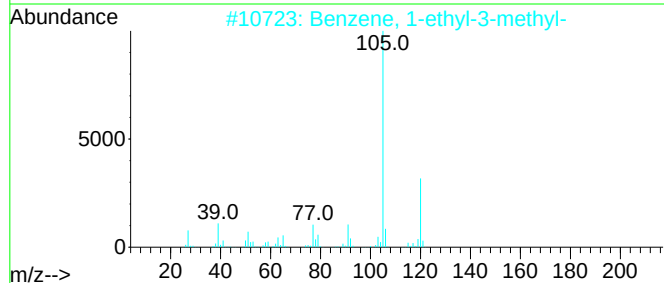
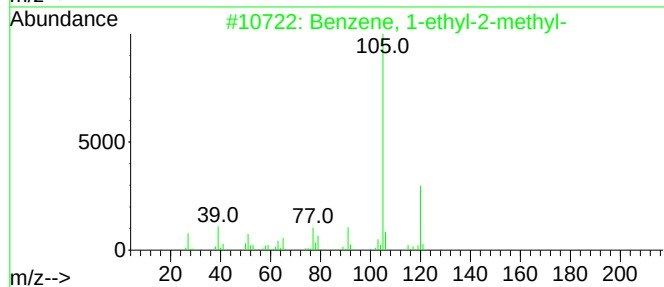
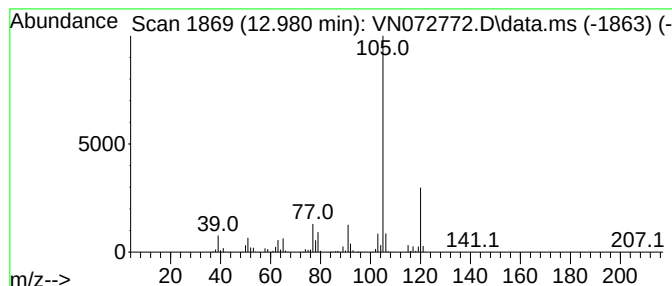
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	128.78 ug/l	2272080	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



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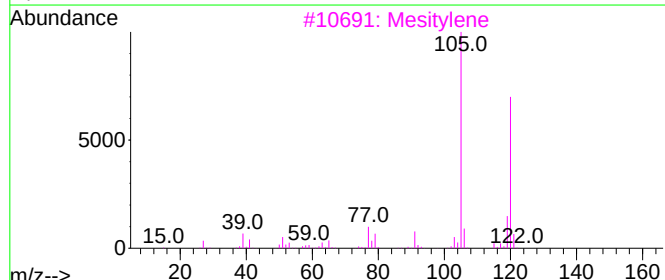
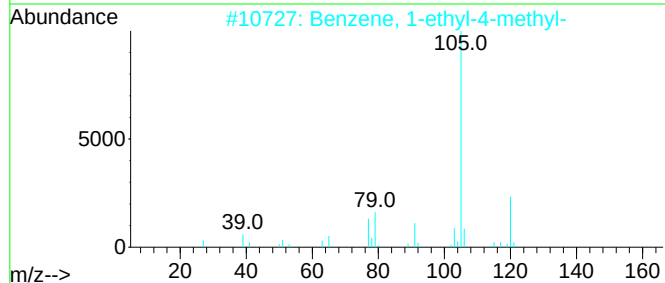
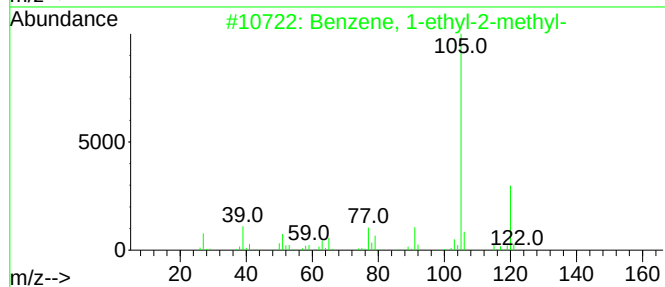
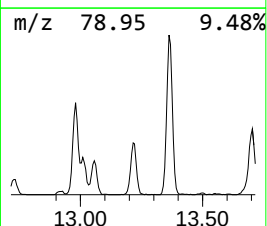
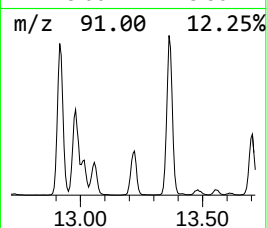
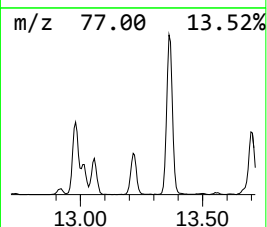
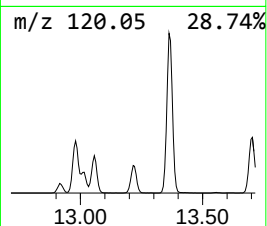
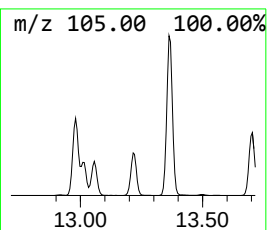
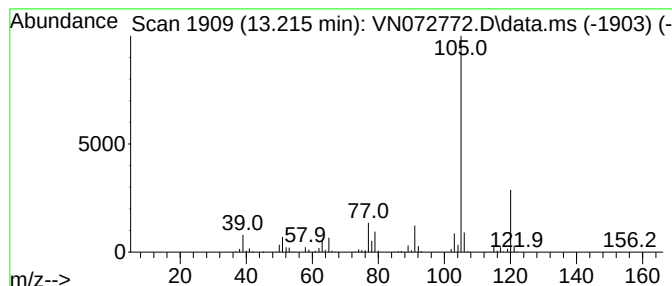
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	64.28 ug/l	1134060	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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Instrument :
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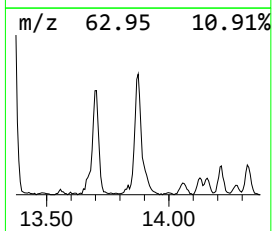
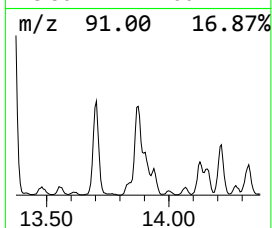
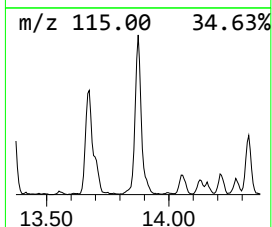
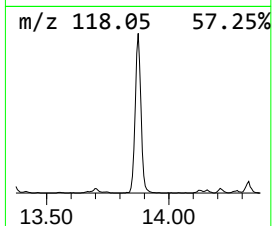
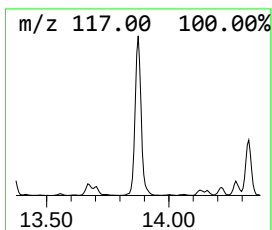
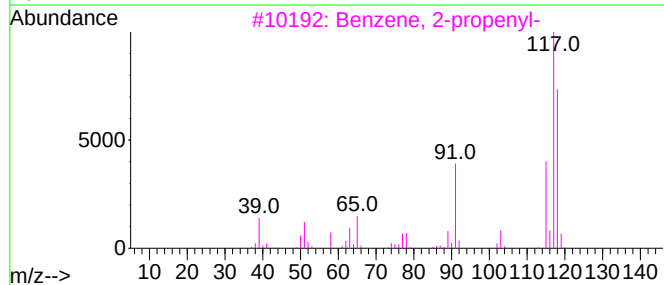
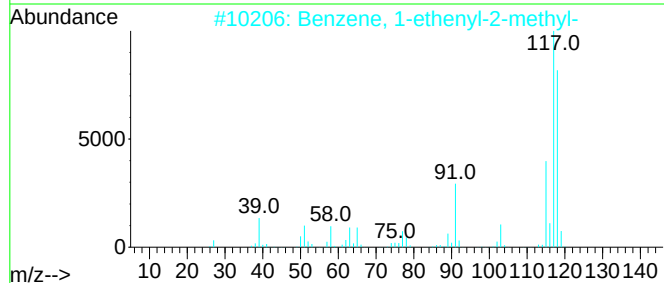
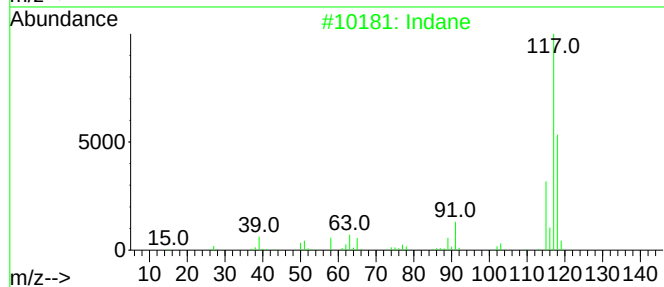
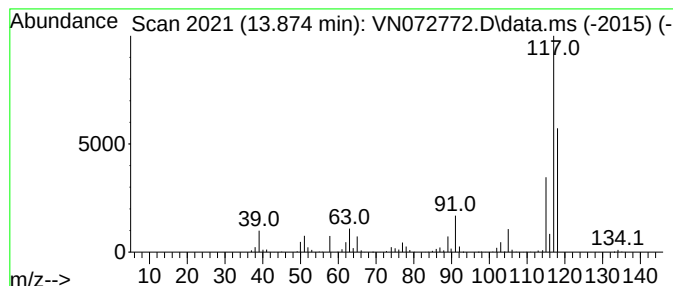
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

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

Peak Number 3 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	119.49 ug/l	2108180	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	59



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Instrument :
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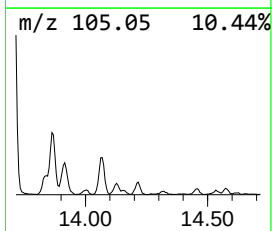
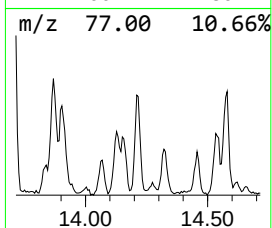
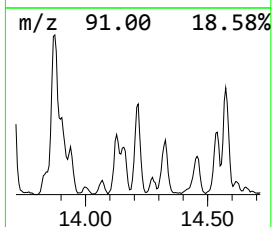
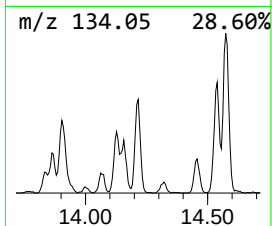
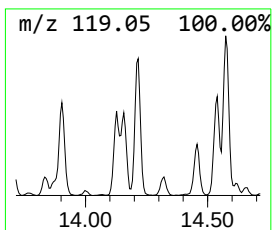
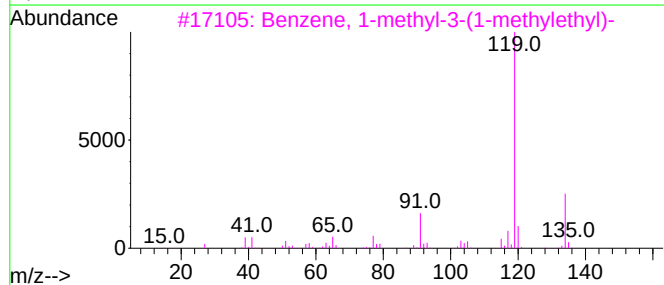
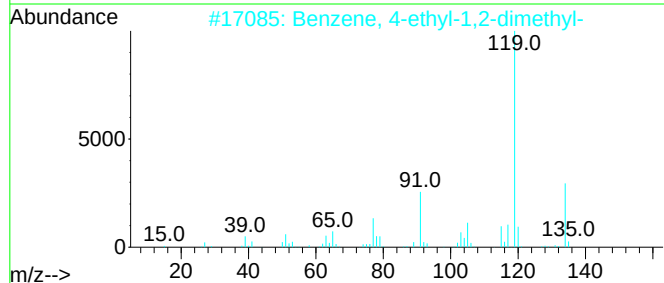
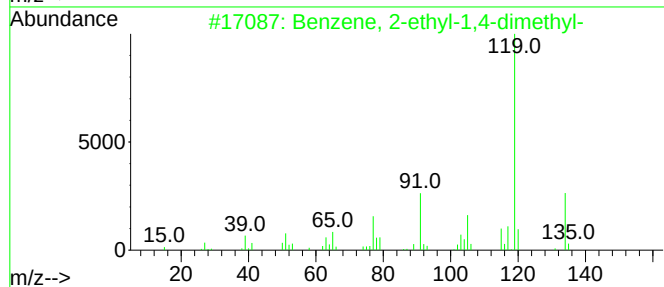
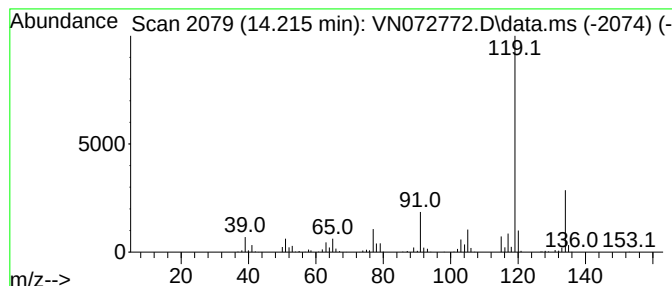
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.215	34.01 ug/l	600054	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			p-Cymene	134	C10H14	000099-87-6	95



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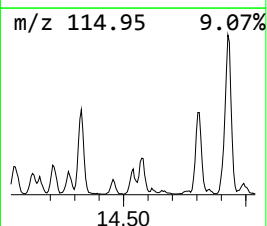
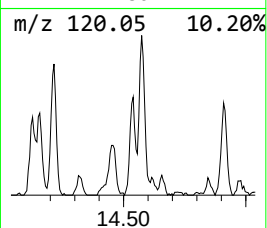
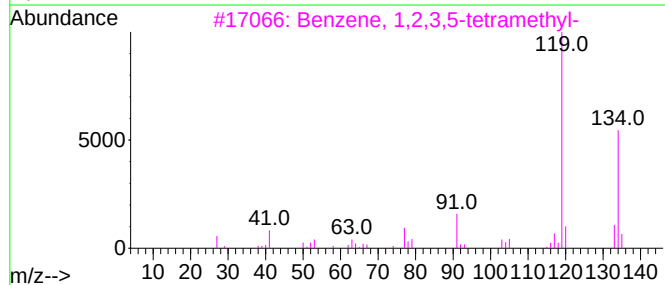
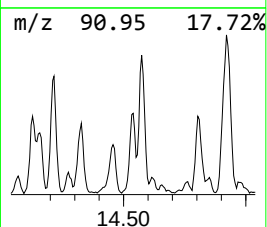
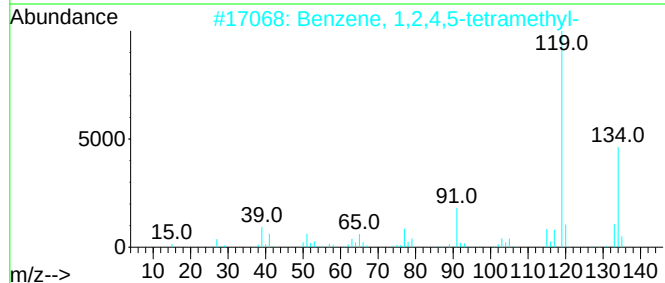
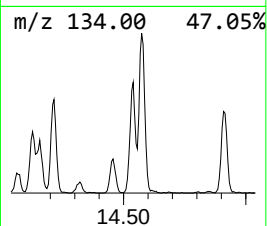
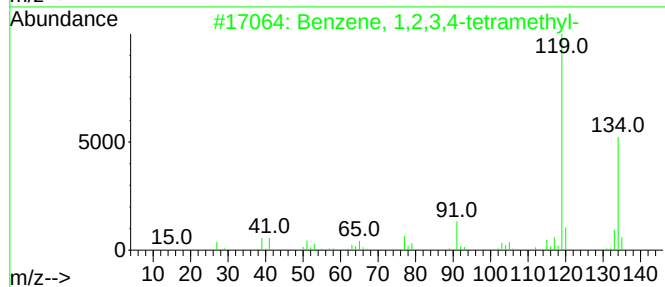
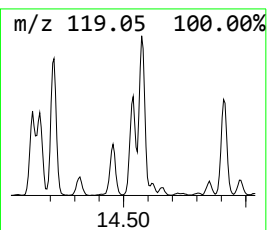
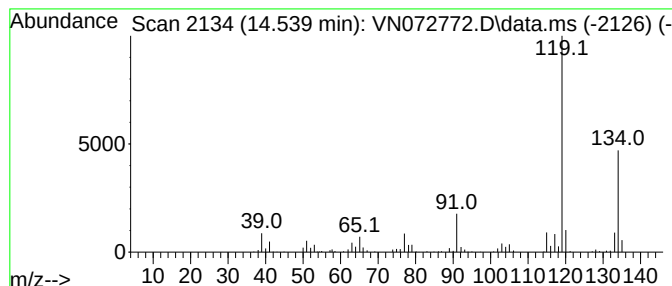
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.539	29.28 ug/l	516590	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5			o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072772.D
Acq On : 10 Jun 2022 18:12
Operator : JC\MD
Sample : N3293-13
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
40805

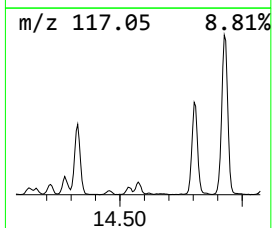
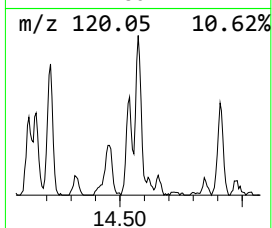
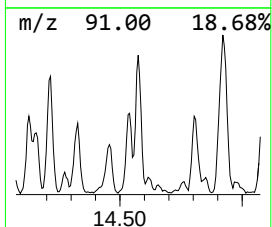
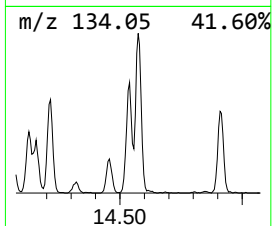
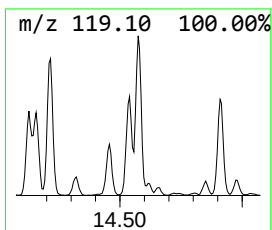
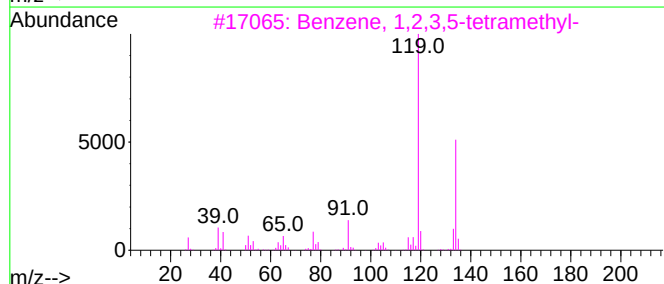
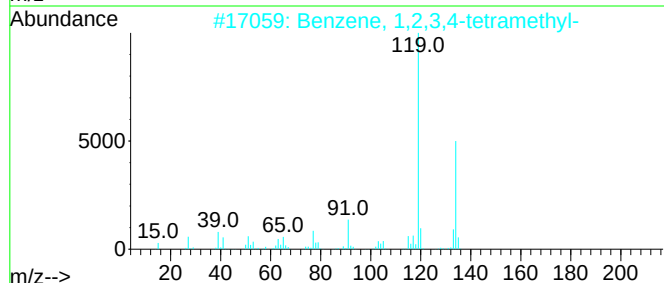
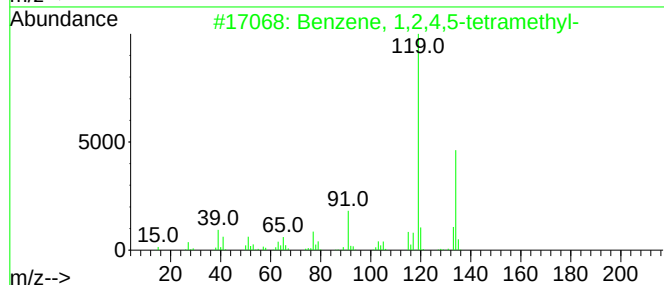
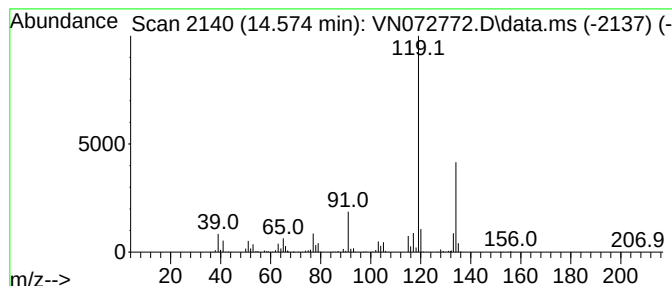
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.574	36.66 ug/l	646711	1,4-Dichlorobenzene-d4	13.668

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	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4			o-Cymene	134	C10H14	000527-84-4	93
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072772.D
Acq On : 10 Jun 2022 18:12
Operator : JC\MD
Sample : N3293-13
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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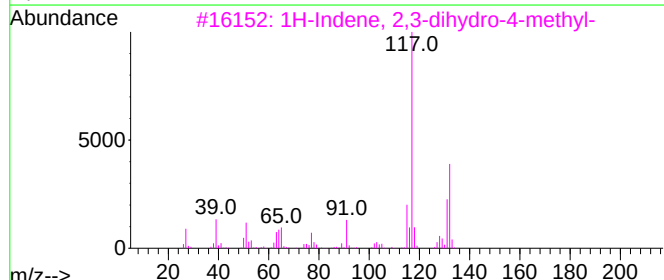
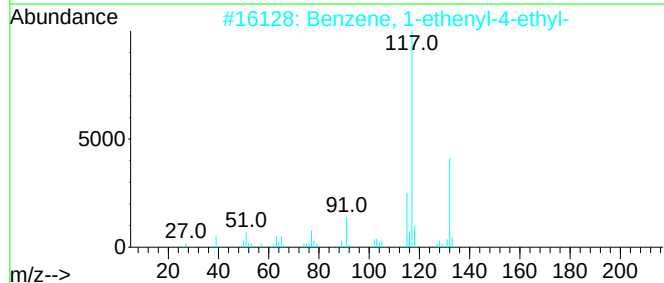
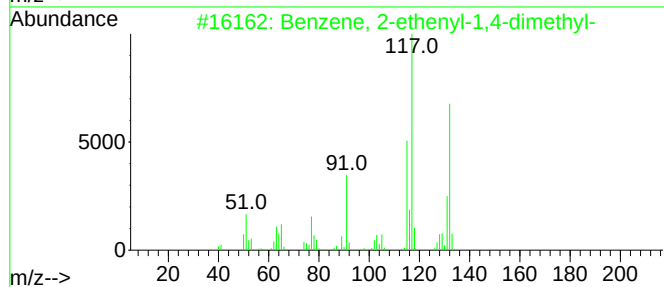
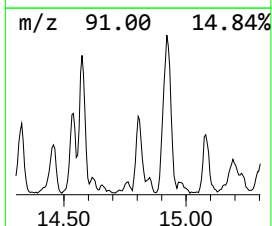
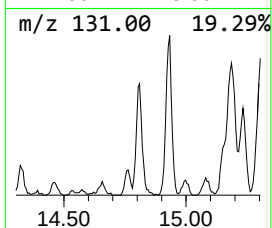
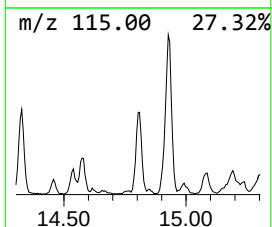
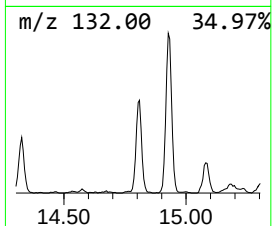
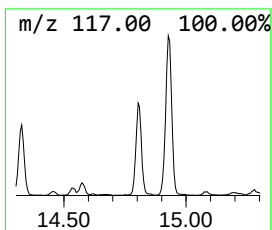
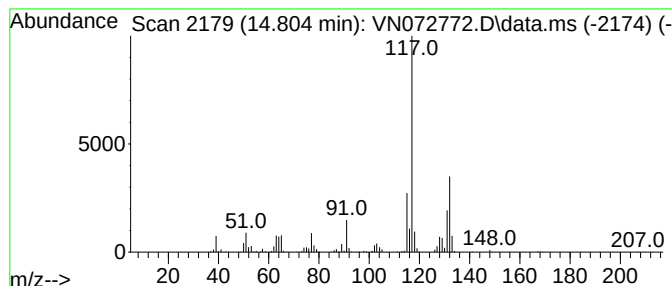
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	32.52 ug/l	573732	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072772.D
Acq On : 10 Jun 2022 18:12
Operator : JC\MD
Sample : N3293-13
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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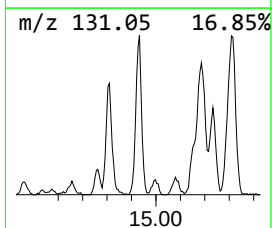
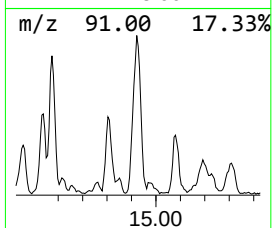
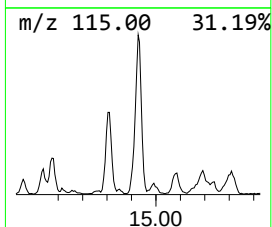
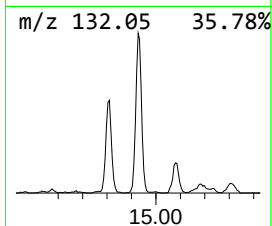
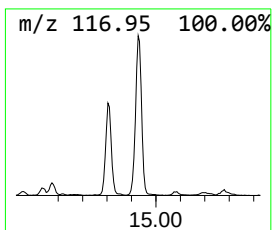
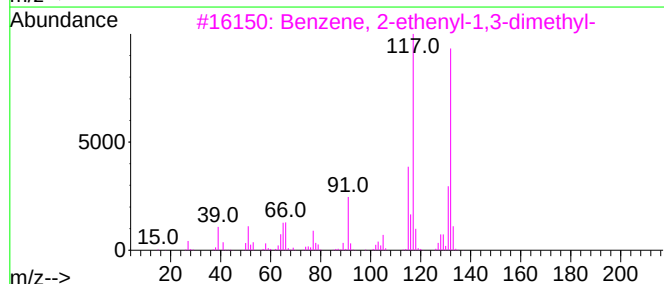
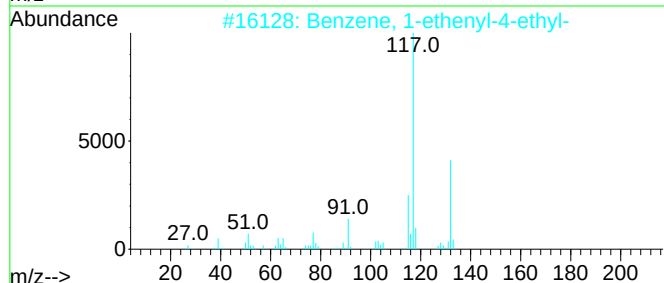
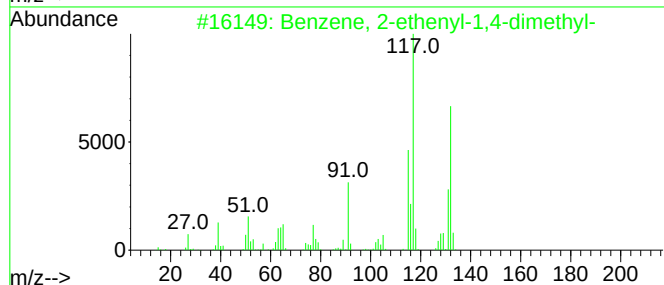
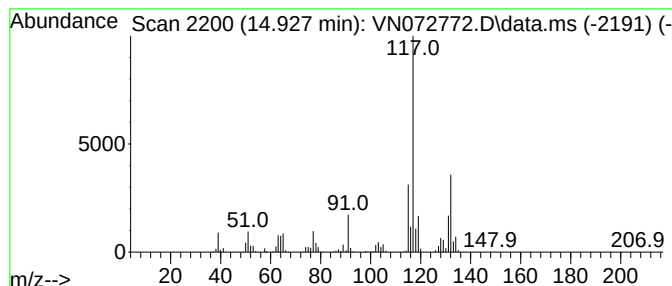
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	84.23 ug/l	1486120	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	94
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
5			2,4-Dimethylstyrene	132	C10H12	002234-20-0	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072772.D
Acq On : 10 Jun 2022 18:12
Operator : JC\MD
Sample : N3293-13
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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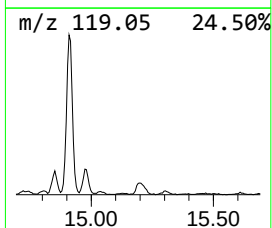
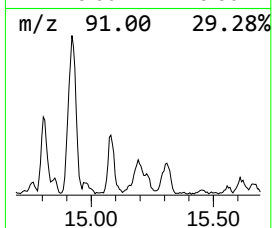
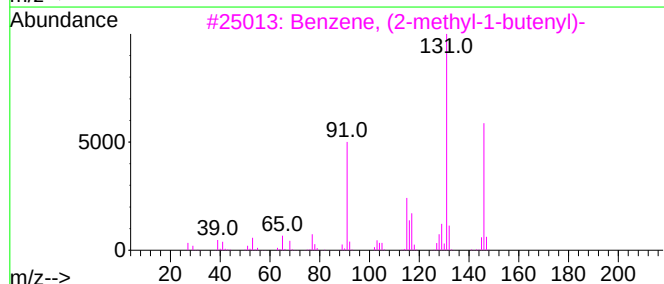
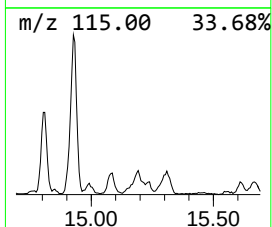
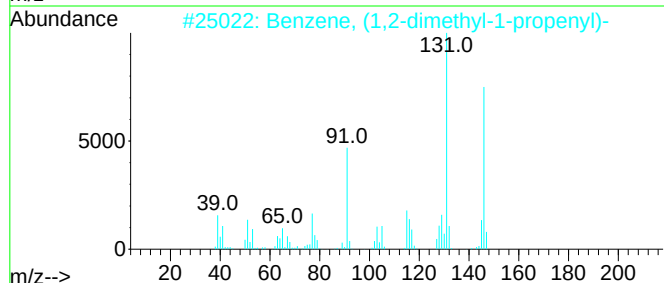
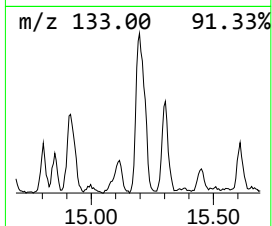
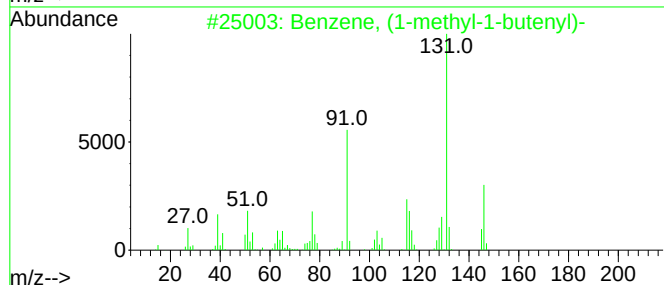
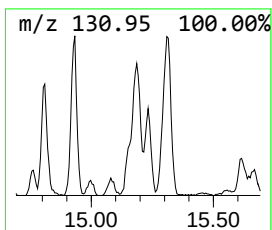
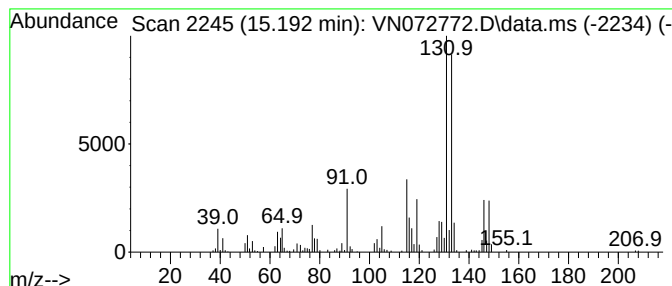
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1-methyl-1-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.192	30.98 ug/l	546622	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
3			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	62
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	50
5			Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072772.D
Acq On : 10 Jun 2022 18:12
Operator : JC\MD
Sample : N3293-13
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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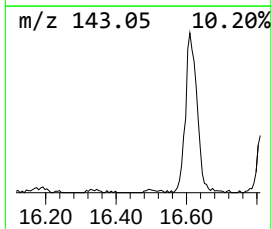
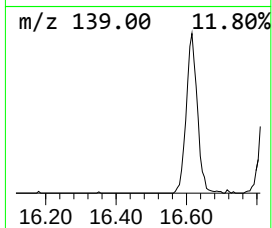
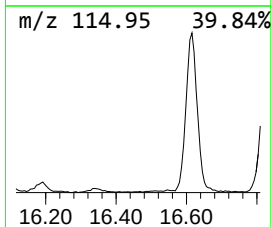
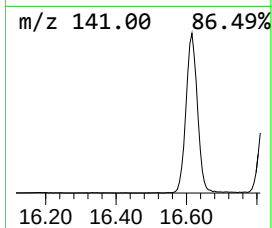
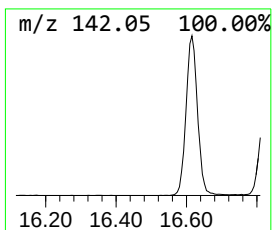
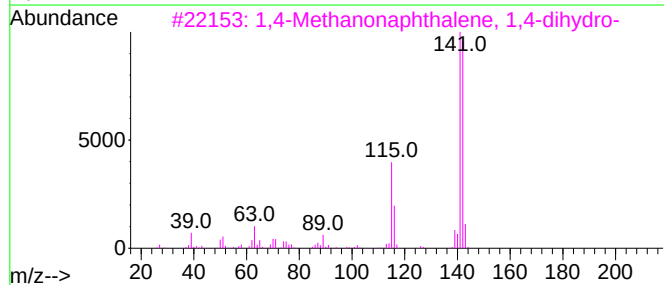
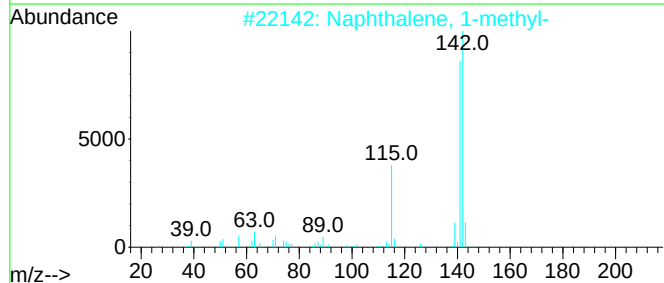
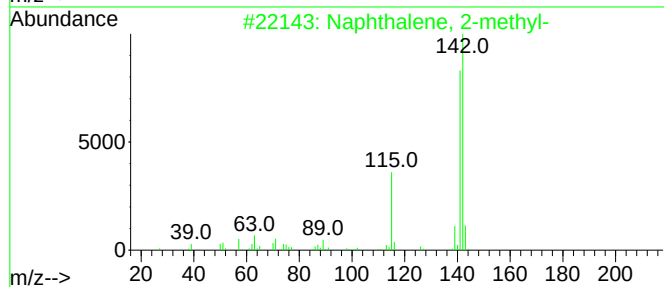
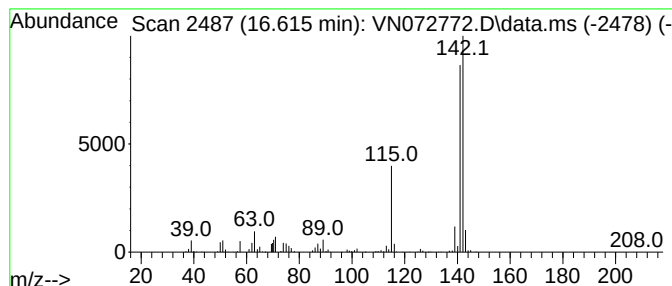
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.615	74.07 ug/l	1306770	1,4-Dichlorobenzene-d4	13.668

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			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061022\
Data File : VN072772.D
Acq On : 10 Jun 2022 18:12
Operator : JC\MD
Sample : N3293-13
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060822W.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
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Benzene, 1-ethy...	13.216	64.3	ug/l	1134060	4	13.668	882138	50.0
Indane	13.874	119.5	ug/l	2108180	4	13.668	882138	50.0
Benzene, 2-ethy...	14.215	34.0	ug/l	600054	4	13.668	882138	50.0
Benzene, 1,2,3,...	14.539	29.3	ug/l	516590	4	13.668	882138	50.0
Benzene, 1,2,4,...	14.574	36.7	ug/l	646711	4	13.668	882138	50.0
Benzene, 2-ethe...	14.804	32.5	ug/l	573732	4	13.668	882138	50.0
Benzene, 1-ethe...	14.927	84.2	ug/l	1486120	4	13.668	882138	50.0
Benzene, (1-met...	15.192	31.0	ug/l	546622	4	13.668	882138	50.0
Naphthalene, 2-...	16.615	74.1	ug/l	1306770	4	13.668	882138	50.0