

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
 Data File : VN071778.D
 Acq On : 02 Apr 2022 02:10
 Operator : JC\MD
 Sample : N2222-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-16

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\82N033022W.M
 Title : SW846 8260

Signal : TIC: VN071778.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.257	36	46	64	rBV2	397810	806233	14.88%	1.270%
2	2.440	68	77	87	rVB	592101	918975	16.97%	1.447%
3	3.134	186	195	207	rVB	319080	735969	13.59%	1.159%
4	3.516	249	260	273	rBV4	73452	209631	3.87%	0.330%
5	3.716	287	294	304	rVB2	26591	62524	1.15%	0.098%
6	3.998	332	342	355	rVB	50678	136839	2.53%	0.215%
7	5.081	514	526	532	rBV3	44376	157583	2.91%	0.248%
8	5.157	532	539	565	rVV6	63482	413479	7.63%	0.651%
9	5.363	565	574	585	rVB2	19255	66213	1.22%	0.104%
10	5.616	604	617	634	rVB3	102264	354551	6.55%	0.558%
11	6.122	692	703	717	rVB5	19825	65273	1.21%	0.103%
12	6.469	753	762	775	rVB3	39459	131080	2.42%	0.206%
13	6.610	775	786	795	rBV4	35169	110730	2.04%	0.174%
14	6.898	826	835	850	rVB5	21184	63423	1.17%	0.100%
15	7.145	865	877	886	rVV2	261110	777296	14.35%	1.224%
16	7.216	886	889	900	rVB3	31960	74929	1.38%	0.118%
17	7.845	989	996	1005	rVB	38094	93163	1.72%	0.147%
18	8.022	1015	1026	1031	rBV	399208	997733	18.42%	1.571%
19	8.086	1031	1037	1046	rVV2	885604	2111293	38.98%	3.325%
20	8.175	1046	1052	1061	rVB4	56682	157991	2.92%	0.249%
21	8.298	1064	1073	1078	rBV3	25267	56812	1.05%	0.089%
22	8.451	1088	1099	1109	rBV3	621565	1901742	35.11%	2.995%
23	8.581	1109	1121	1124	rBV10	62162	176656	3.26%	0.278%
24	8.704	1137	1142	1154	rVB2	59691	141908	2.62%	0.223%
25	8.963	1177	1186	1199	rBV	1187064	2578675	47.61%	4.061%
26	9.198	1219	1226	1229	rBV2	28275	58052	1.07%	0.091%
27	9.245	1229	1234	1244	rVB2	44212	98506	1.82%	0.155%
28	9.457	1256	1270	1278	rBV2	192235	571284	10.55%	0.900%
29	9.639	1295	1301	1308	rVB2	37429	72624	1.34%	0.114%
30	9.804	1322	1329	1330	rBV3	33721	59320	1.10%	0.093%
31	9.880	1337	1342	1350	rVB2	57324	116253	2.15%	0.183%
32	9.969	1350	1357	1361	rBV2	132179	266935	4.93%	0.420%
33	10.010	1361	1364	1372	rVB	111369	209854	3.87%	0.330%
34	10.198	1389	1396	1404	rBV9	23201	65354	1.21%	0.103%
35	10.322	1410	1417	1422	rBV	77909	150608	2.78%	0.237%
36	10.439	1429	1437	1444	rBV	1886050	3470132	64.06%	5.465%

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Integrator: RTE

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Max Peaks: 100

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Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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37	10.504	1444	1448	1460	rVV	85191	196050	3.62%	0.309%
38	10.639	1460	1471	1478	rVV4	131234	293566	5.42%	0.462%
39	10.710	1478	1483	1490	rVB	92294	163881	3.03%	0.258%
40	10.857	1501	1508	1518	rVB6	90428	218097	4.03%	0.343%
41	11.739	1651	1658	1669	rBV	1775982	3160097	58.34%	4.977%
42	11.839	1669	1675	1687	rVV	1102018	1885972	34.82%	2.970%
43	11.957	1687	1695	1707	rVB	191787	367477	6.78%	0.579%
44	12.274	1739	1749	1759	rBV	134582	240605	4.44%	0.379%
45	12.574	1791	1800	1809	rBV	613764	1007593	18.60%	1.587%
46	12.727	1818	1826	1836	rBV	1330927	2255642	41.64%	3.552%
47	12.916	1851	1858	1865	rBV	1065422	1739108	32.11%	2.739%
48	13.010	1865	1874	1879	rVV	264926	514699	9.50%	0.811%
49	13.057	1879	1882	1889	rVB	69801	99127	1.83%	0.156%
50	13.216	1901	1909	1919	rBV	1898739	3117681	57.56%	4.910%
51	13.363	1927	1934	1948	rVB	3411597	5416642	100.00%	8.530%
52	13.492	1948	1956	1963	rBV3	115238	258099	4.76%	0.406%
53	13.557	1963	1967	1972	rVB	36942	56648	1.05%	0.089%
54	13.674	1980	1987	1988	rBV	1508495	2293371	42.34%	3.612%
55	13.698	1988	1991	1999	rVV	2435475	4130198	76.25%	6.504%
56	13.768	1999	2003	2008	rVB2	42111	66932	1.24%	0.105%
57	13.833	2008	2014	2017	rBV	544553	920787	17.00%	1.450%
58	13.874	2017	2021	2026	rVV	1553609	2603315	48.06%	4.100%
59	13.915	2026	2028	2037	rVB2	205386	339575	6.27%	0.535%
60	13.998	2037	2042	2047	rBV	130842	201356	3.72%	0.317%
61	14.063	2047	2053	2059	rBV	264431	423360	7.82%	0.667%
62	14.127	2059	2064	2073	rVV	274504	546874	10.10%	0.861%
63	14.215	2073	2079	2084	rVV	821715	1278432	23.60%	2.013%
64	14.274	2084	2089	2092	rVV	207520	337429	6.23%	0.531%
65	14.321	2092	2097	2105	rVB2	1034213	1740558	32.13%	2.741%
66	14.457	2113	2120	2126	rBV2	285642	466930	8.62%	0.735%
67	14.533	2126	2133	2137	rBV	710419	1179127	21.77%	1.857%
68	14.574	2137	2140	2145	rVB	560331	805835	14.88%	1.269%
69	14.757	2164	2171	2174	rBV2	37558	69770	1.29%	0.110%
70	14.804	2174	2179	2185	rVV	552406	924862	17.07%	1.456%
71	14.927	2191	2200	2206	rVV2	1431423	2915978	53.83%	4.592%
72	14.992	2206	2211	2216	rVV4	84284	166446	3.07%	0.262%
73	15.080	2219	2226	2233	rVV2	172752	327667	6.05%	0.516%
74	15.192	2234	2245	2249	rVV3	193895	510600	9.43%	0.804%
75	15.227	2249	2251	2257	rVV	102390	177452	3.28%	0.279%
76	15.309	2257	2265	2274	rVB3	206809	518130	9.57%	0.816%

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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

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Peak separation: 5

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Title : SW846 8260

77	15.504	2291	2298	2305	rVB	318025	594013	10.97%	0.935%
78	15.609	2311	2316	2321	rBV3	42937	92395	1.71%	0.146%
79	15.668	2321	2326	2335	rVB2	47552	98576	1.82%	0.155%
80	15.798	2341	2348	2356	rBV2	63999	125203	2.31%	0.197%
81	15.974	2368	2378	2393	rBV2	45485	142999	2.64%	0.225%
82	16.186	2407	2414	2424	rVB3	28916	70367	1.30%	0.111%

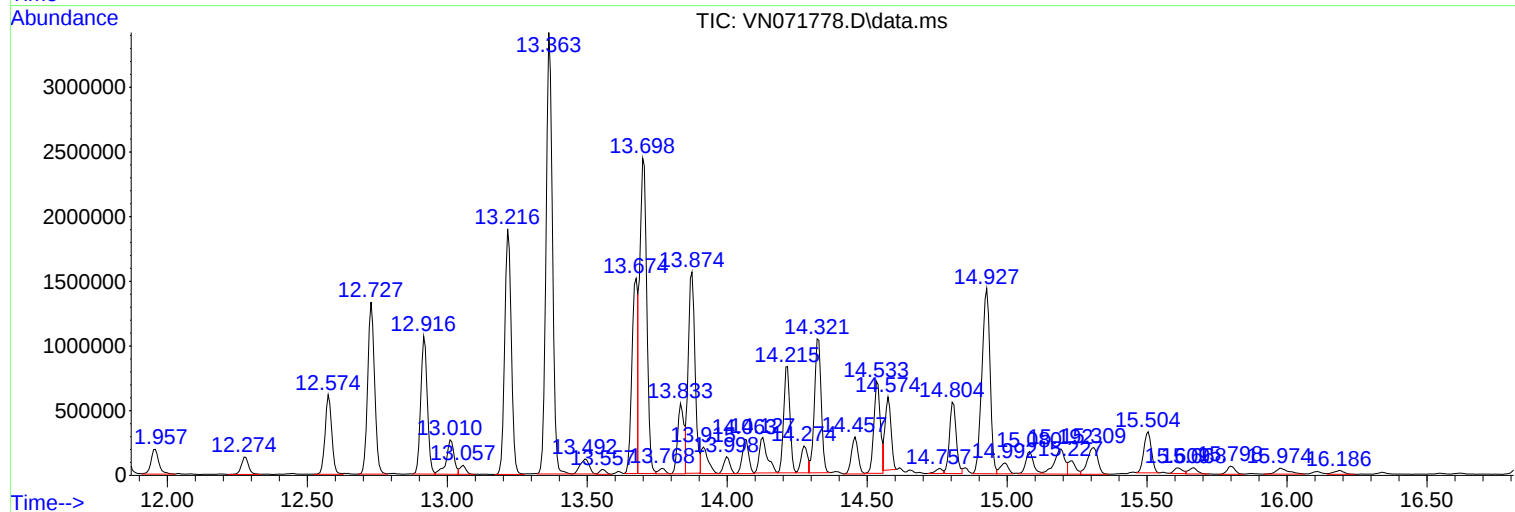
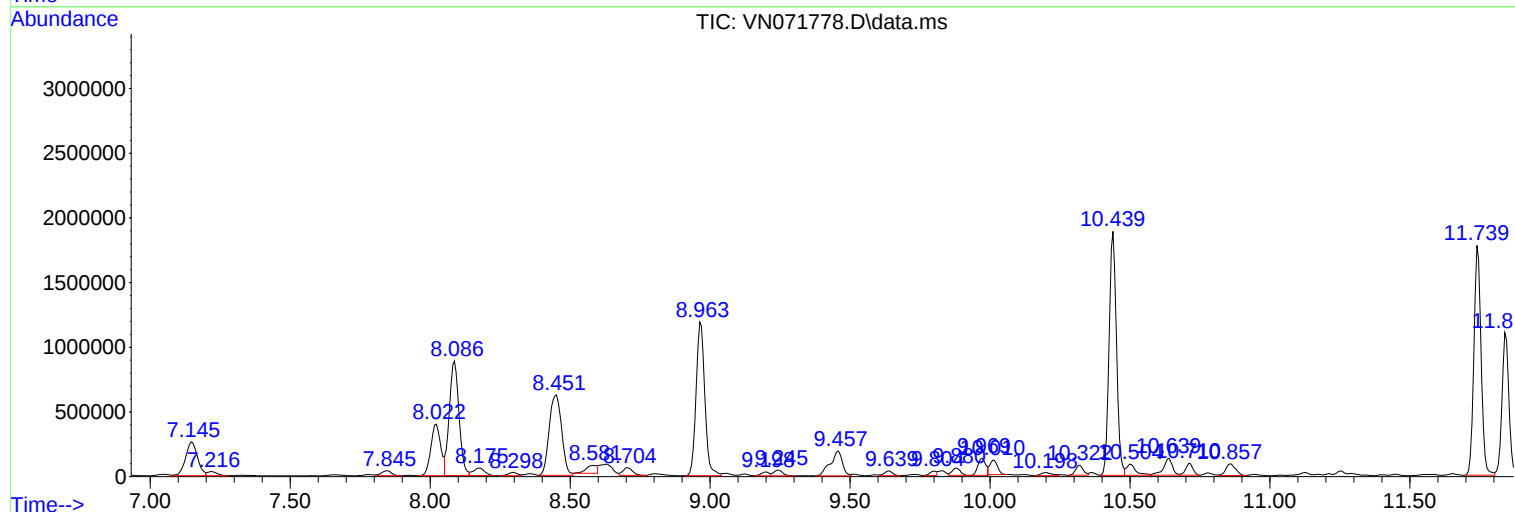
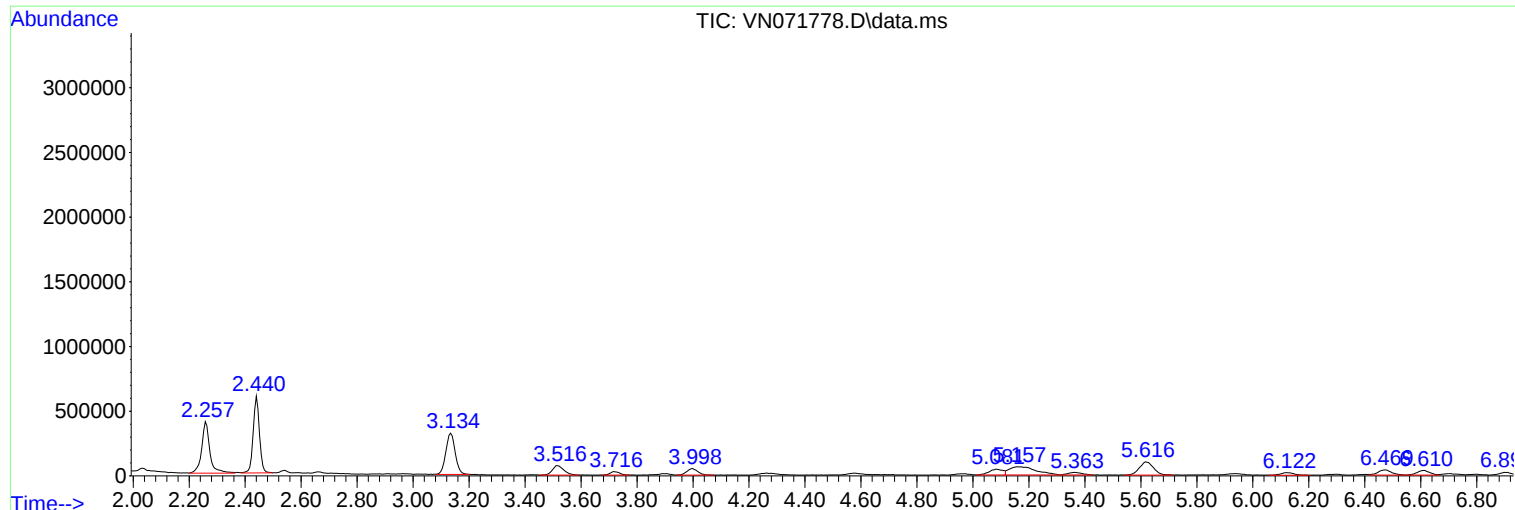
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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



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Instrument :
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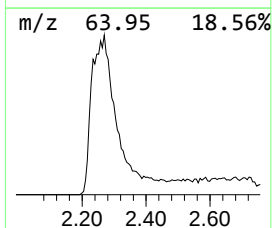
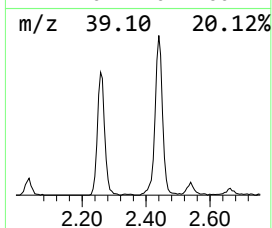
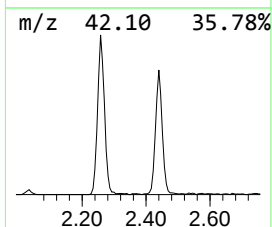
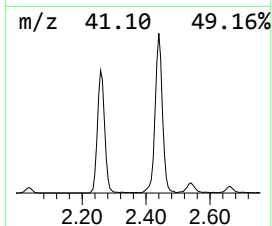
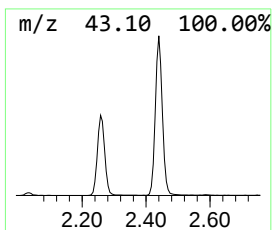
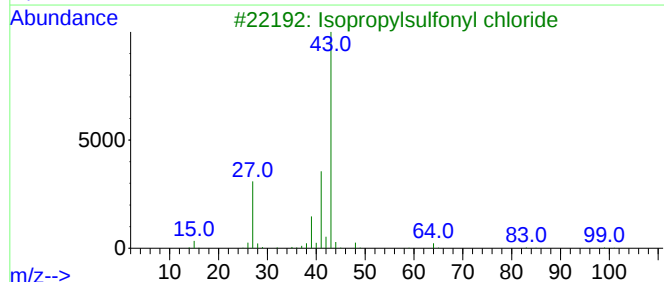
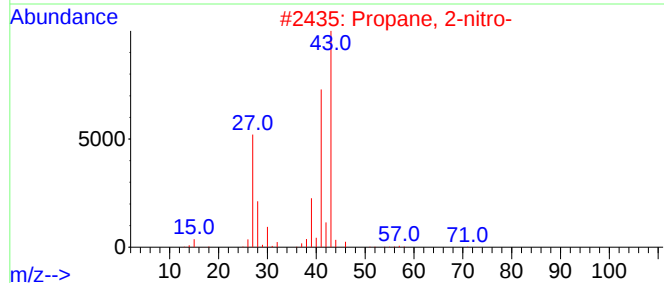
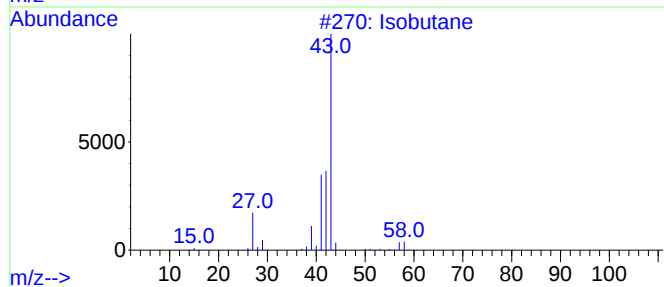
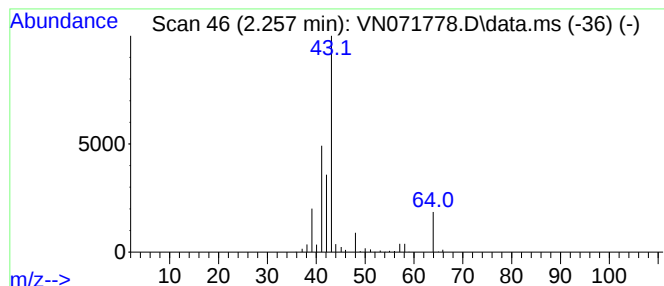
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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

Peak Number 1 Isobutane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.257	19.09 ug/l	806233	Pentafluorobenzene	8.086

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutane	58	C4H10	000075-28-5	58
2			Propane, 2-nitro-	89	C3H7NO2	000079-46-9	9
3			Isopropylsulfonyl chloride	142	C3H7ClO2S	010147-37-2	9
4			Propane, 1-chloro-2-methyl-	92	C4H9Cl	000513-36-0	4
5			5-Methyloxazolidine	87	C4H9NO	058328-22-6	4



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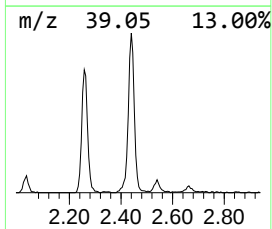
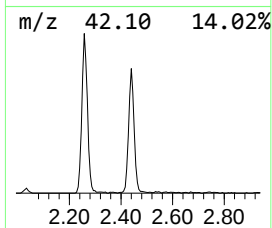
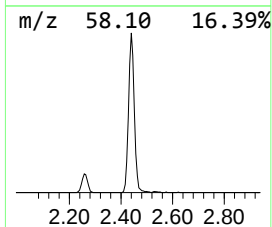
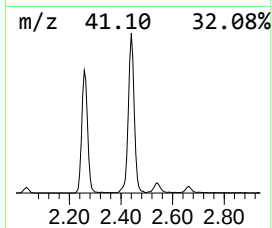
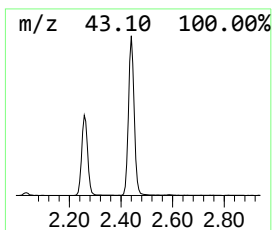
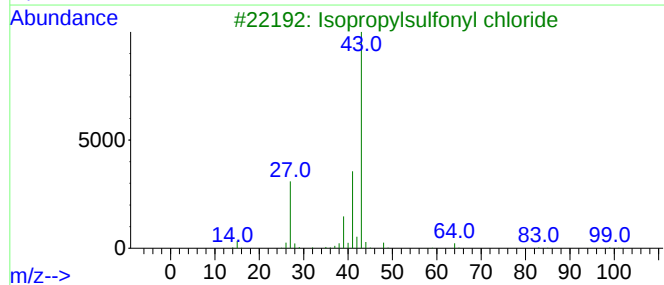
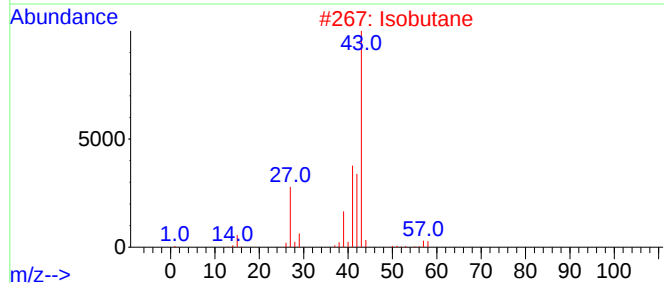
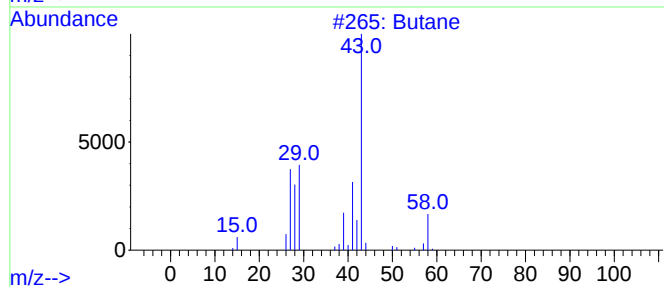
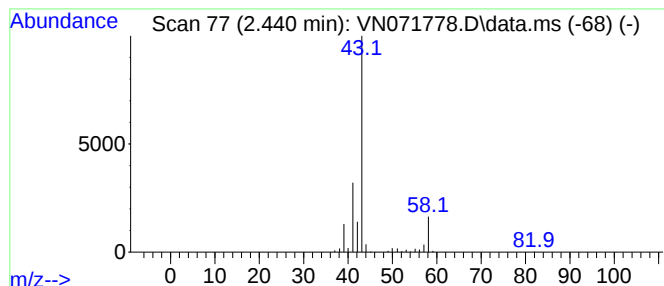
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.440	21.76 ug/l	918975	Pentafluorobenzene	8.086

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane		58	C4H10	000106-97-8	80
2	Isobutane		58	C4H10	000075-28-5	9
3	Isopropylsulfonyl chloride		142	C3H7ClO2S	010147-37-2	9
4	Propane, 1-nitro-		89	C3H7NO2	000108-03-2	9
5	Acetone		58	C3H6O	000067-64-1	4



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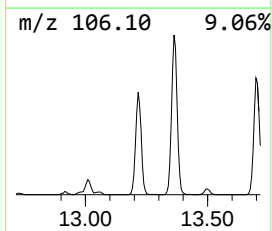
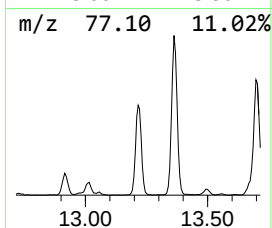
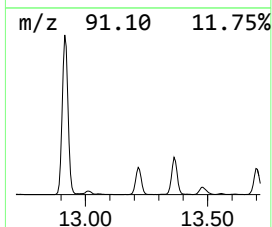
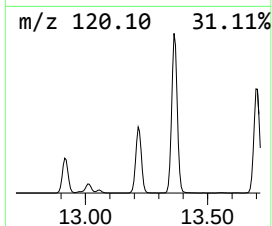
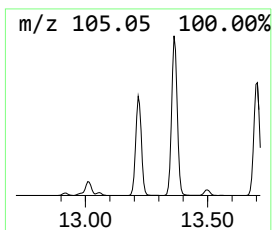
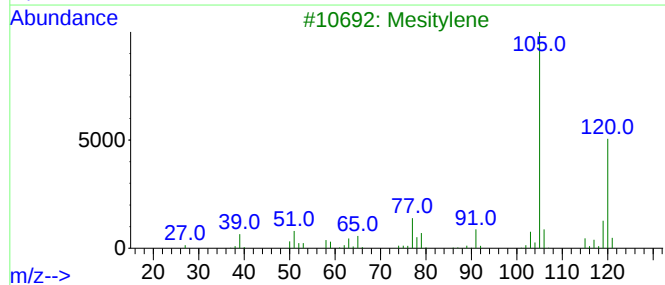
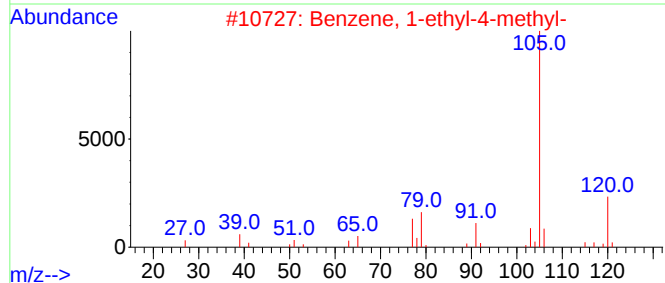
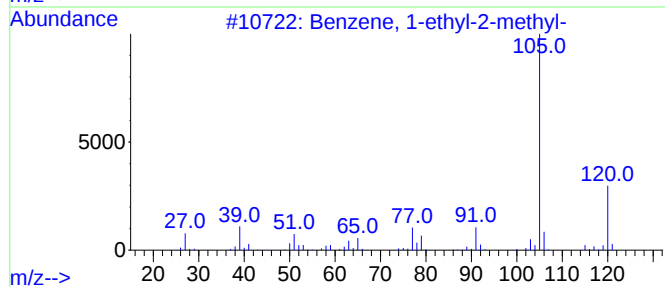
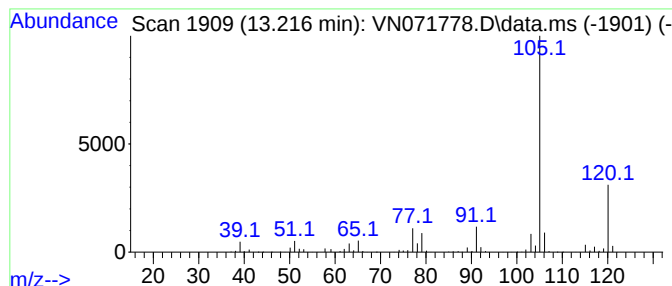
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	67.97 ug/l	3117680	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-4-methyl-	120	C9H12	000622-96-8	94
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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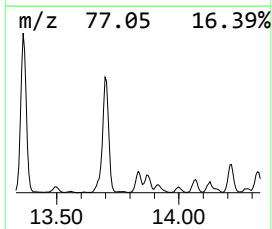
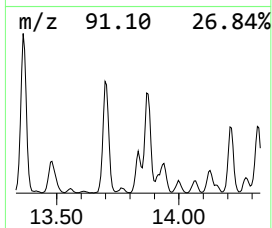
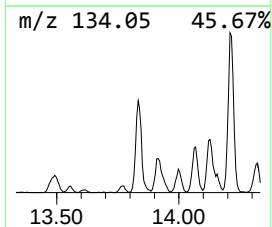
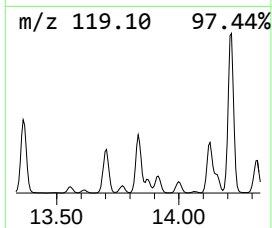
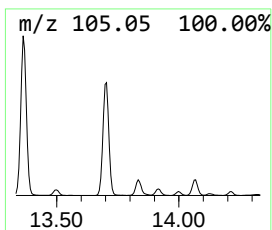
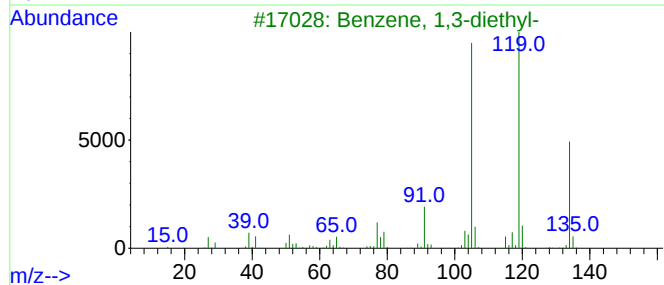
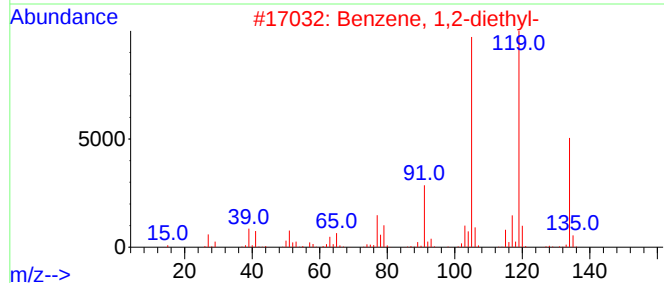
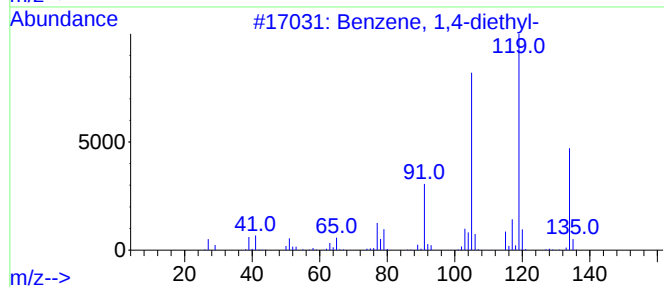
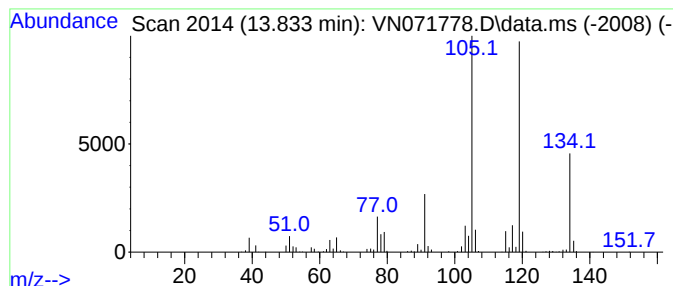
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,4-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.833	20.07 ug/l	920787	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-16

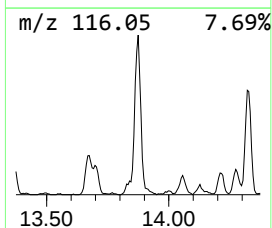
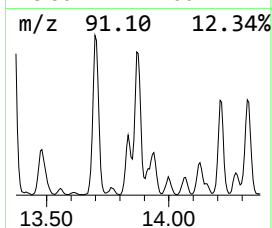
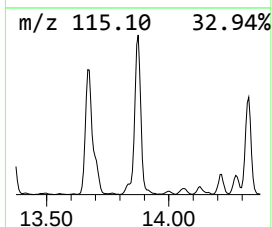
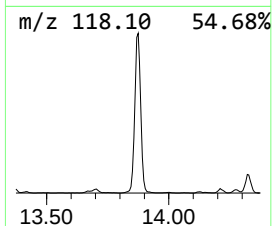
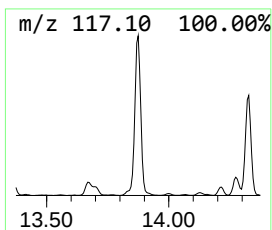
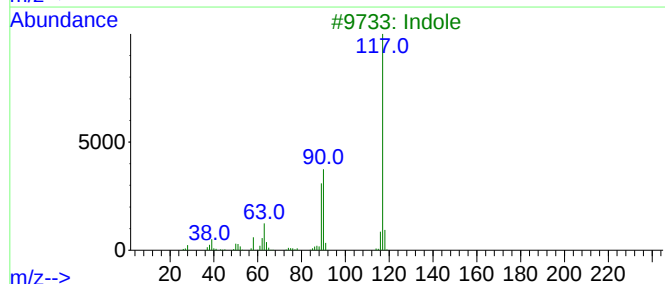
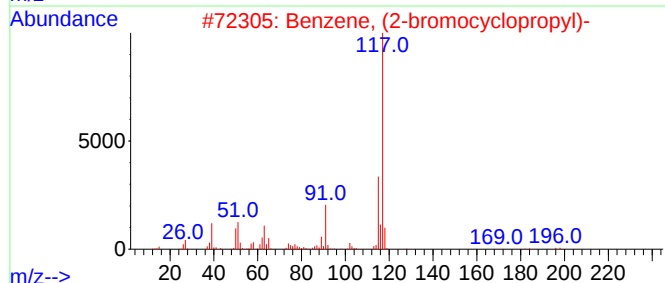
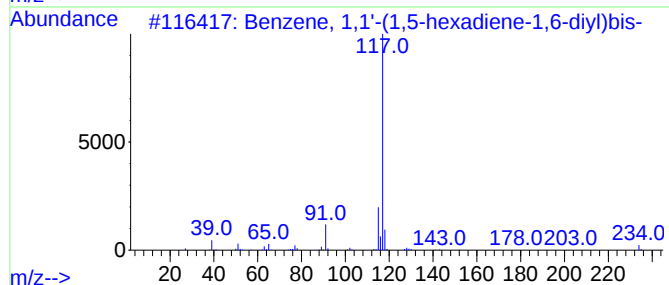
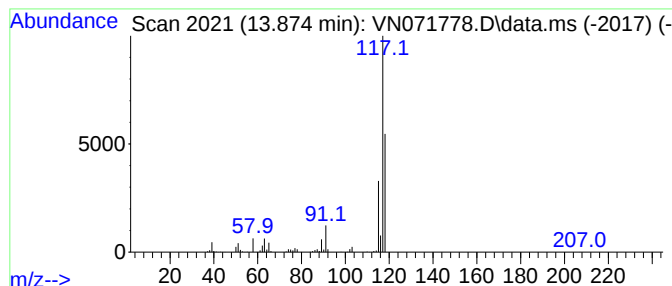
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.874	56.76 ug/l	2603320	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	50
2			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
3			Indole	117	C8H7N	000120-72-9	43
4			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42
5			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-16

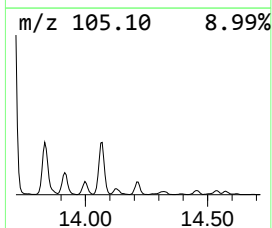
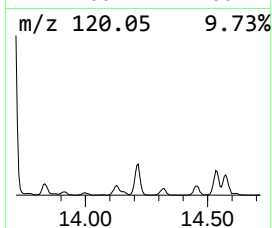
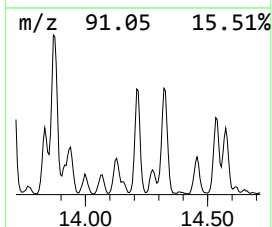
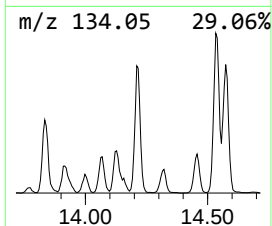
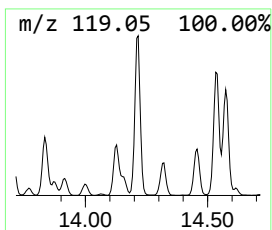
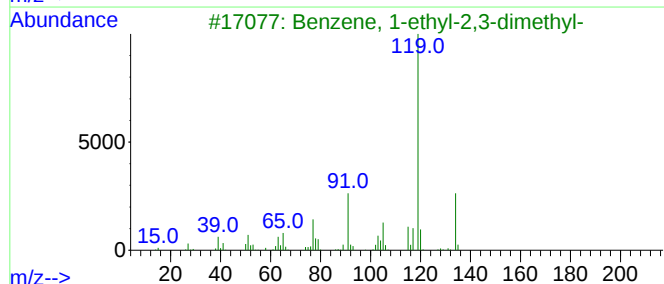
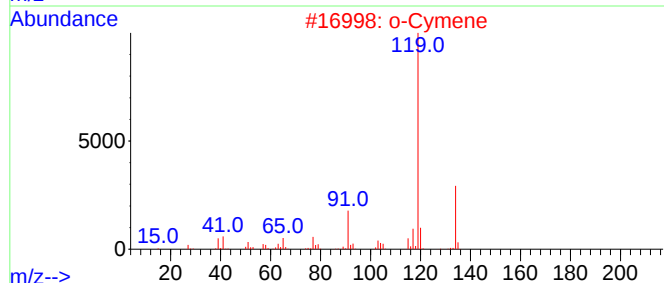
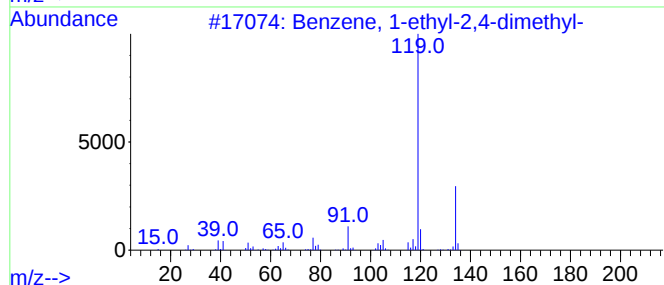
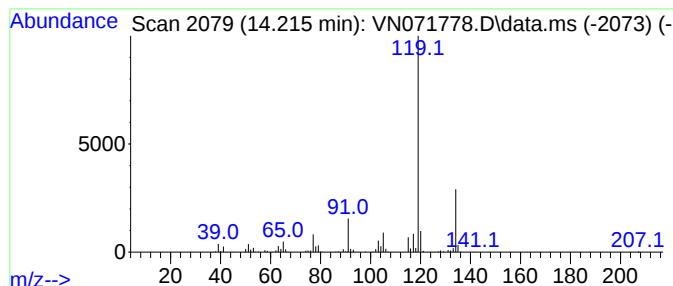
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.216	27.87 ug/l	1278430	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			o-Cymene	134	C10H14	000527-84-4	97
3			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-16

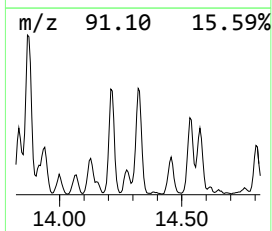
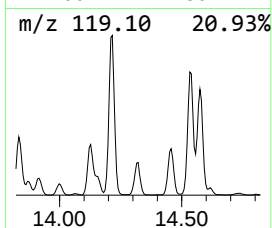
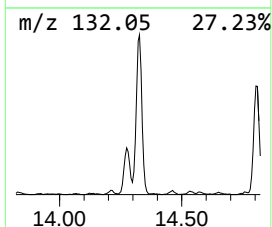
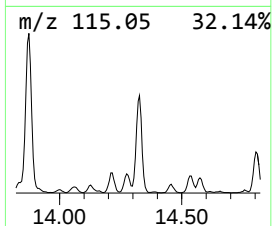
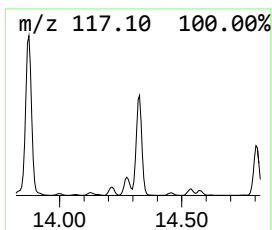
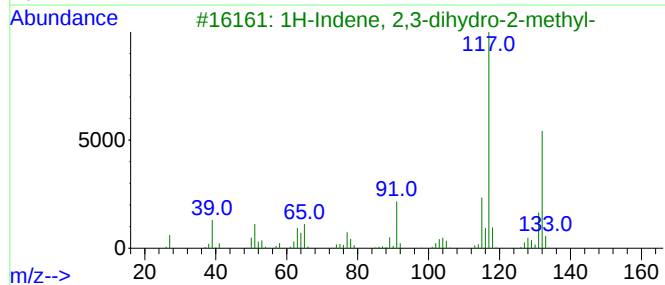
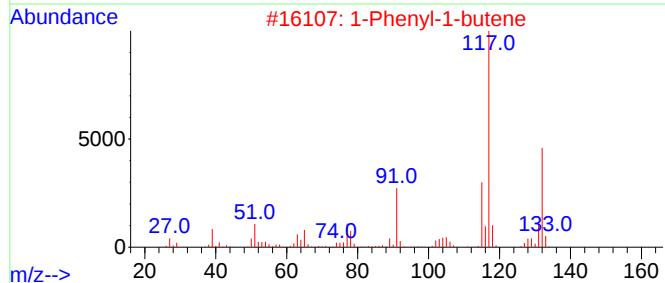
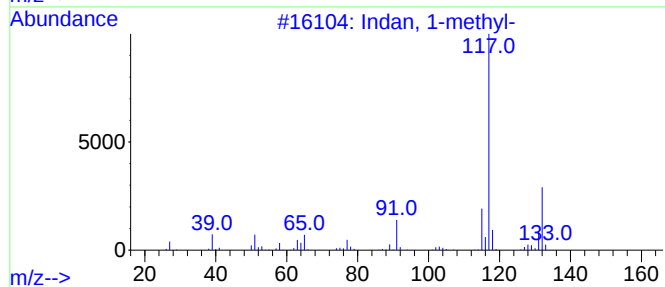
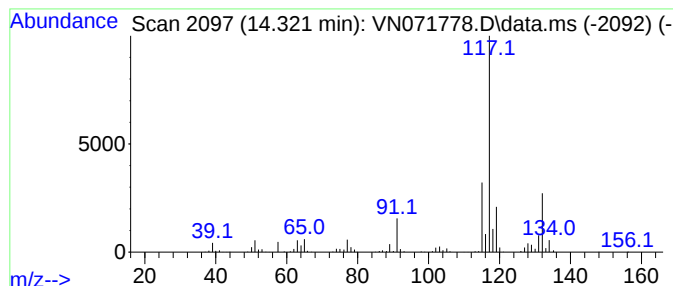
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.321	37.95 ug/l	1740560	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	76
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	74
3			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	72
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	72
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-16

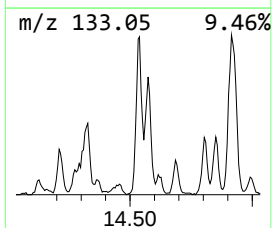
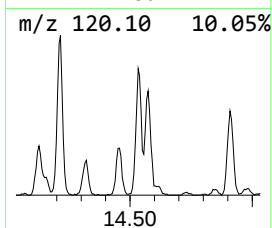
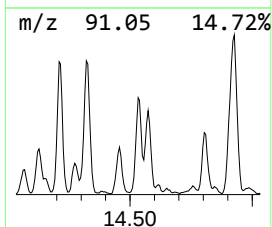
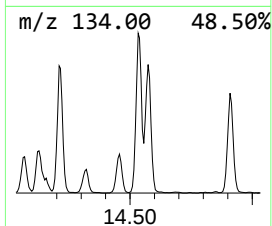
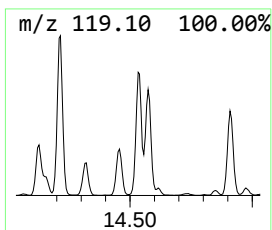
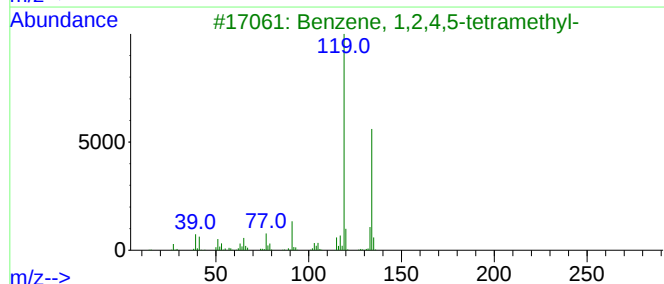
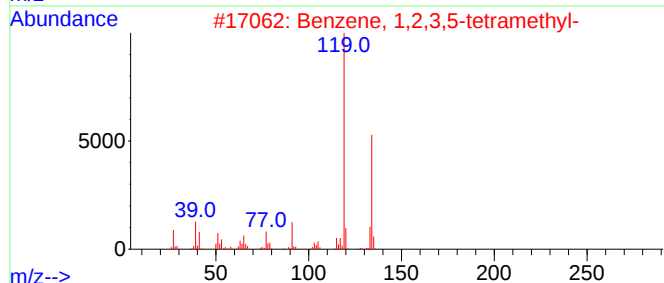
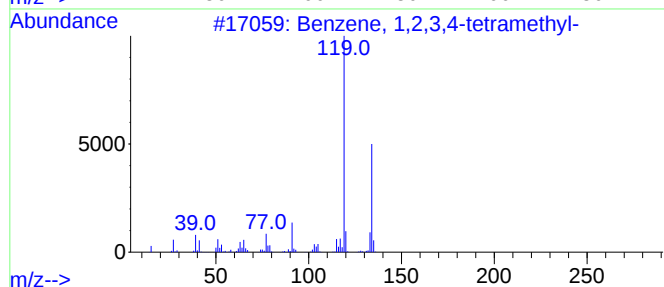
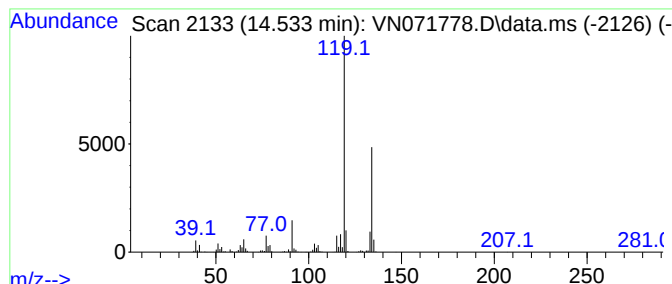
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.533	25.71 ug/l	1179130	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-16

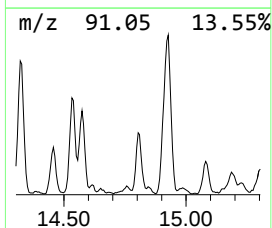
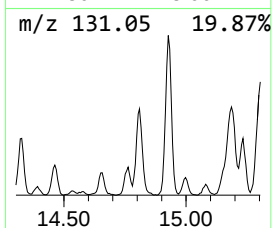
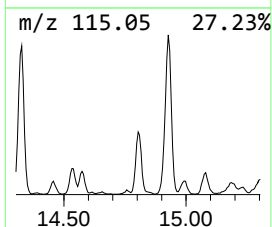
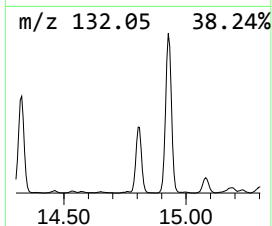
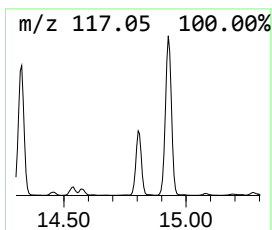
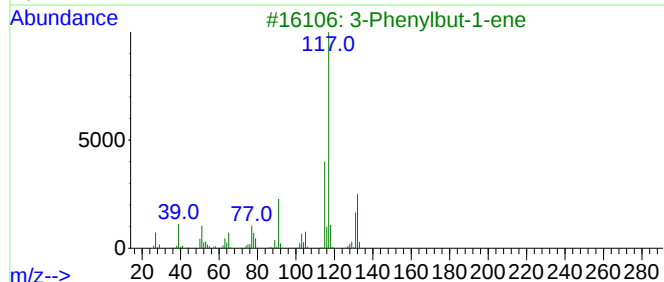
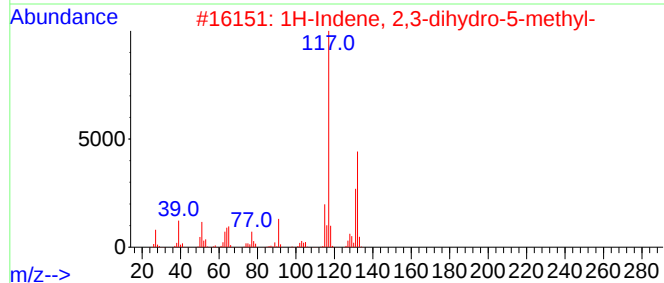
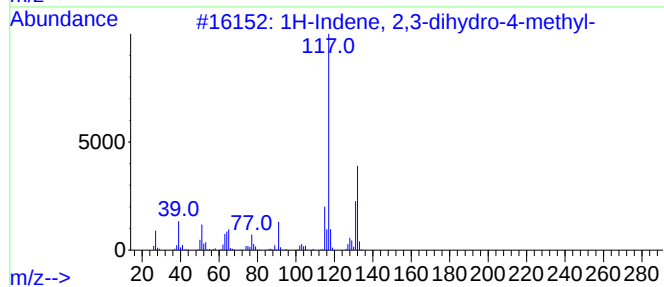
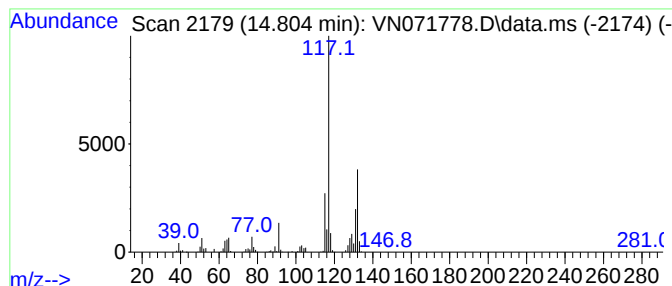
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.804	20.16 ug/l	924862	1,4-Dichlorobenzene-d4	13.668

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5			Indan, 1-methyl-	132	C10H12	000767-58-8	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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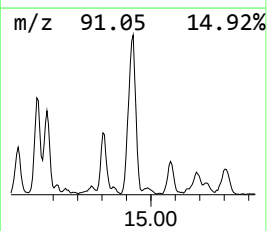
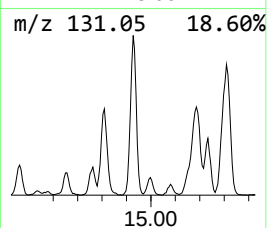
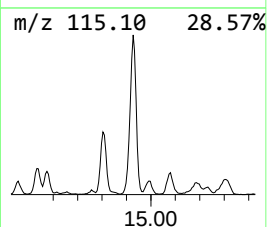
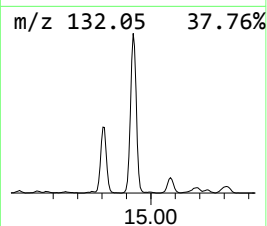
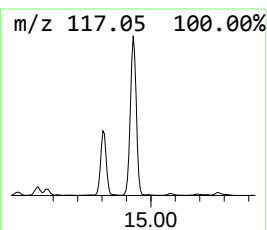
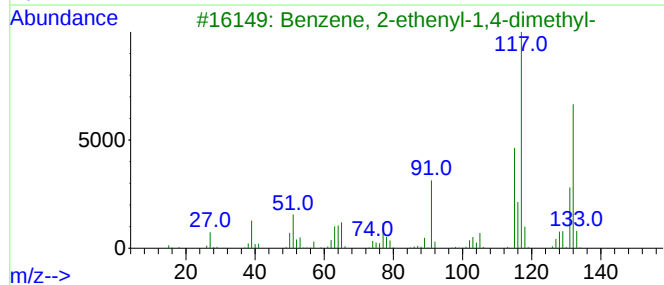
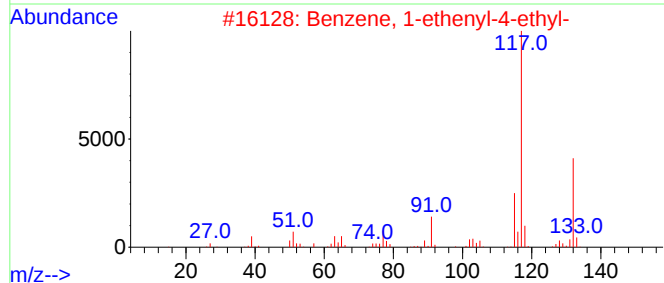
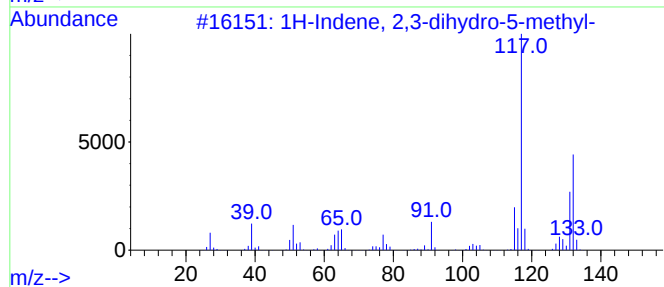
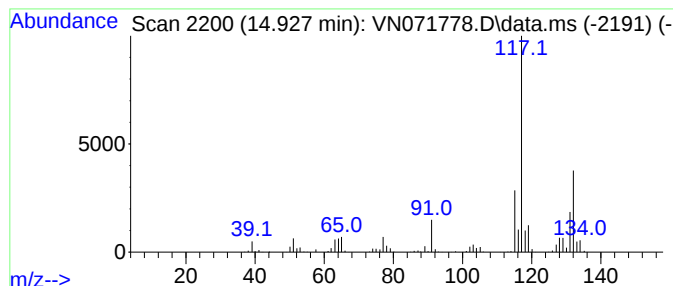
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	63.57 ug/l	2915980	1,4-Dichlorobenzene-d4	13.668

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040122\
Data File : VN071778.D
Acq On : 02 Apr 2022 02:10
Operator : JC\MD
Sample : N2222-05
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 39 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-16

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.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
Isobutane	2.257	19.1	ug/l	806233	1	8.086	2111290	50.0
Butane	2.440	21.8	ug/l	918975	1	8.086	2111290	50.0
Benzene, 1-ethy...	13.216	68.0	ug/l	3117680	4	13.668	2293370	50.0
Benzene, 1,4-di...	13.833	20.1	ug/l	920787	4	13.668	2293370	50.0
Benzene, 1,1'-(...	13.874	56.8	ug/l	2603320	4	13.668	2293370	50.0
Benzene, 1-ethy...	14.216	27.9	ug/l	1278430	4	13.668	2293370	50.0
Indan, 1-methyl-	14.321	38.0	ug/l	1740560	4	13.668	2293370	50.0
Benzene, 1,2,3,...	14.533	25.7	ug/l	1179130	4	13.668	2293370	50.0
1H-Indene, 2,3-...	14.804	20.2	ug/l	924862	4	13.668	2293370	50.0
1H-Indene, 2,3-...	14.927	63.6	ug/l	2915980	4	13.668	2293370	50.0