

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
 Data File : VN068087.D
 Acq On : 12 Aug 2021 14:12
 Operator : JC/MD
 Sample : M3374-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DEWATERING-DISCHARGE

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\82N072921W.M
 Title : SW846 8260

Signal : TIC: VN068087.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.437	162	167	180	rVB3	16472	22488	1.06%	0.093%
2	2.746	271	282	297	rVB4	19963	45738	2.16%	0.189%
3	3.132	412	426	445	rVB2	87000	192108	9.07%	0.793%
4	3.505	554	565	586	rVB3	21605	53307	2.52%	0.220%
5	3.982	722	743	761	rBV2	17840	46577	2.20%	0.192%
6	4.293	821	859	889	rBV	55183	178444	8.43%	0.737%
7	4.594	957	971	1001	rVB2	11824	36141	1.71%	0.149%
8	5.077	1126	1151	1153	rBV4	19214	43772	2.07%	0.181%
9	5.149	1162	1178	1180	rBV4	39972	69399	3.28%	0.287%
10	5.613	1325	1351	1381	rBV4	55924	188742	8.91%	0.779%
11	6.898	1810	1830	1849	rBV6	12094	36535	1.73%	0.151%
12	7.040	1867	1883	1900	rBV6	8727	25299	1.19%	0.104%
13	7.147	1902	1923	1943	rVV2	111654	321526	15.19%	1.328%
14	7.697	2108	2128	2148	rBV5	21691	58136	2.75%	0.240%
15	7.844	2165	2183	2201	rVB6	18633	47331	2.24%	0.195%
16	8.024	2225	2250	2260	rBV2	241096	587122	27.73%	2.425%
17	8.088	2261	2274	2294	rVV4	556156	1397566	66.01%	5.771%
18	8.174	2296	2306	2329	rVB4	42834	108652	5.13%	0.449%
19	8.295	2332	2351	2365	rBV3	64525	148860	7.03%	0.615%
20	8.445	2387	2407	2432	rBV5	347712	1008068	47.61%	4.163%
21	8.568	2436	2453	2467	rVV9	48963	145885	6.89%	0.602%
22	8.644	2471	2481	2493	rVV6	41620	94565	4.47%	0.391%
23	8.705	2493	2504	2526	rVB6	40606	99068	4.68%	0.409%
24	8.968	2585	2602	2630	rBV	661378	1432034	67.64%	5.914%
25	9.462	2755	2786	2801	rBV4	159035	448681	21.19%	1.853%
26	9.639	2839	2852	2867	rVB4	25304	50201	2.37%	0.207%
27	9.735	2867	2888	2902	rVB9	11921	26301	1.24%	0.109%
28	9.880	2935	2942	2955	rVB7	13866	24805	1.17%	0.102%
29	9.974	2962	2977	2988	rBV5	37205	75554	3.57%	0.312%
30	10.019	2989	2994	3005	rVB7	14730	24342	1.15%	0.101%
31	10.212	3047	3066	3094	rBV10	29994	97043	4.58%	0.401%
32	10.325	3094	3108	3135	rVB5	53929	125236	5.92%	0.517%
33	10.443	3135	3152	3182	rBV	1093371	2117182	100.00%	8.743%
34	10.620	3205	3218	3224	rBV	12650	29736	1.40%	0.123%
35	10.784	3267	3279	3292	rVB5	16341	31012	1.46%	0.128%
36	10.875	3292	3313	3327	rBV8	31381	76073	3.59%	0.314%

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Stop Thrs : 0

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37	11.296	3463	3470	3481	rVB4	17214	26784	1.27%	0.111%
38	11.529	3540	3557	3568	rBV9	10777	28876	1.36%	0.119%
39	11.747	3620	3638	3665	rBV	1009468	1866709	88.17%	7.709%
40	11.956	3700	3716	3727	rBV3	28094	66985	3.16%	0.277%
41	12.283	3830	3838	3855	rVB7	26426	59430	2.81%	0.245%
42	12.583	3940	3950	3962	rVB3	19825	32471	1.53%	0.134%
43	12.734	3991	4006	4027	rBV	743934	1292100	61.03%	5.336%
44	13.015	4102	4111	4118	rBV7	15362	22111	1.04%	0.091%
45	13.061	4120	4128	4139	rVB3	29515	46302	2.19%	0.191%
46	13.224	4175	4189	4205	rBV	215595	364427	17.21%	1.505%
47	13.369	4234	4243	4253	rVB2	19459	29147	1.38%	0.120%
48	13.501	4281	4292	4303	rBV4	19993	33489	1.58%	0.138%
49	13.616	4325	4335	4343	rBV4	23597	35353	1.67%	0.146%
50	13.678	4344	4358	4384	rVB2	881625	1651756	78.02%	6.821%
51	13.774	4387	4394	4405	rVB3	24500	34118	1.61%	0.141%
52	13.841	4406	4419	4423	rBV2	220213	317713	15.01%	1.312%
53	13.879	4424	4433	4445	rVB	886533	1398157	66.04%	5.774%
54	14.005	4466	4480	4491	rBV5	81983	145530	6.87%	0.601%
55	14.066	4491	4503	4517	rVB3	61002	114047	5.39%	0.471%
56	14.131	4517	4527	4542	rBV3	65892	106473	5.03%	0.440%
57	14.219	4546	4560	4573	rBV	575269	903759	42.69%	3.732%
58	14.281	4573	4583	4590	rBV	113661	175834	8.31%	0.726%
59	14.329	4590	4601	4615	rVB2	573135	979065	46.24%	4.043%
60	14.466	4639	4652	4661	rBV3	42036	71464	3.38%	0.295%
61	14.541	4663	4680	4688	rVV	356125	607992	28.72%	2.511%
62	14.581	4689	4695	4704	rVV	187836	274869	12.98%	1.135%
63	14.624	4704	4711	4720	rVV2	69580	98236	4.64%	0.406%
64	14.812	4770	4781	4792	rBV2	244640	388902	18.37%	1.606%
65	14.933	4807	4826	4839	rBV3	838816	1763587	83.30%	7.283%
66	14.989	4840	4847	4871	rVB6	58631	124958	5.90%	0.516%
67	15.086	4872	4883	4899	rBV3	80819	141896	6.70%	0.586%
68	15.198	4901	4925	4934	rBV5	182374	436906	20.64%	1.804%
69	15.314	4956	4968	4984	rVB4	175847	410087	19.37%	1.693%
70	15.458	5011	5022	5029	rBV4	18738	33717	1.59%	0.139%
71	15.507	5031	5040	5052	rVB	47119	82173	3.88%	0.339%
72	15.566	5053	5062	5070	rBV3	23623	38261	1.81%	0.158%
73	15.619	5073	5082	5089	rBV8	30691	56233	2.66%	0.232%
74	15.810	5138	5153	5169	rBV2	63935	125394	5.92%	0.518%
75	15.984	5204	5218	5225	rBV3	39843	82676	3.91%	0.341%
76	16.113	5255	5266	5278	rBV6	16163	31353	1.48%	0.129%

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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77	16.193	5281	5296	5312	rVB5	25830	60187	2.84%	0.249%
78	16.346	5334	5353	5368	rBV7	48553	109236	5.16%	0.451%
79	16.558	5411	5432	5443	rBV7	13811	33342	1.57%	0.138%

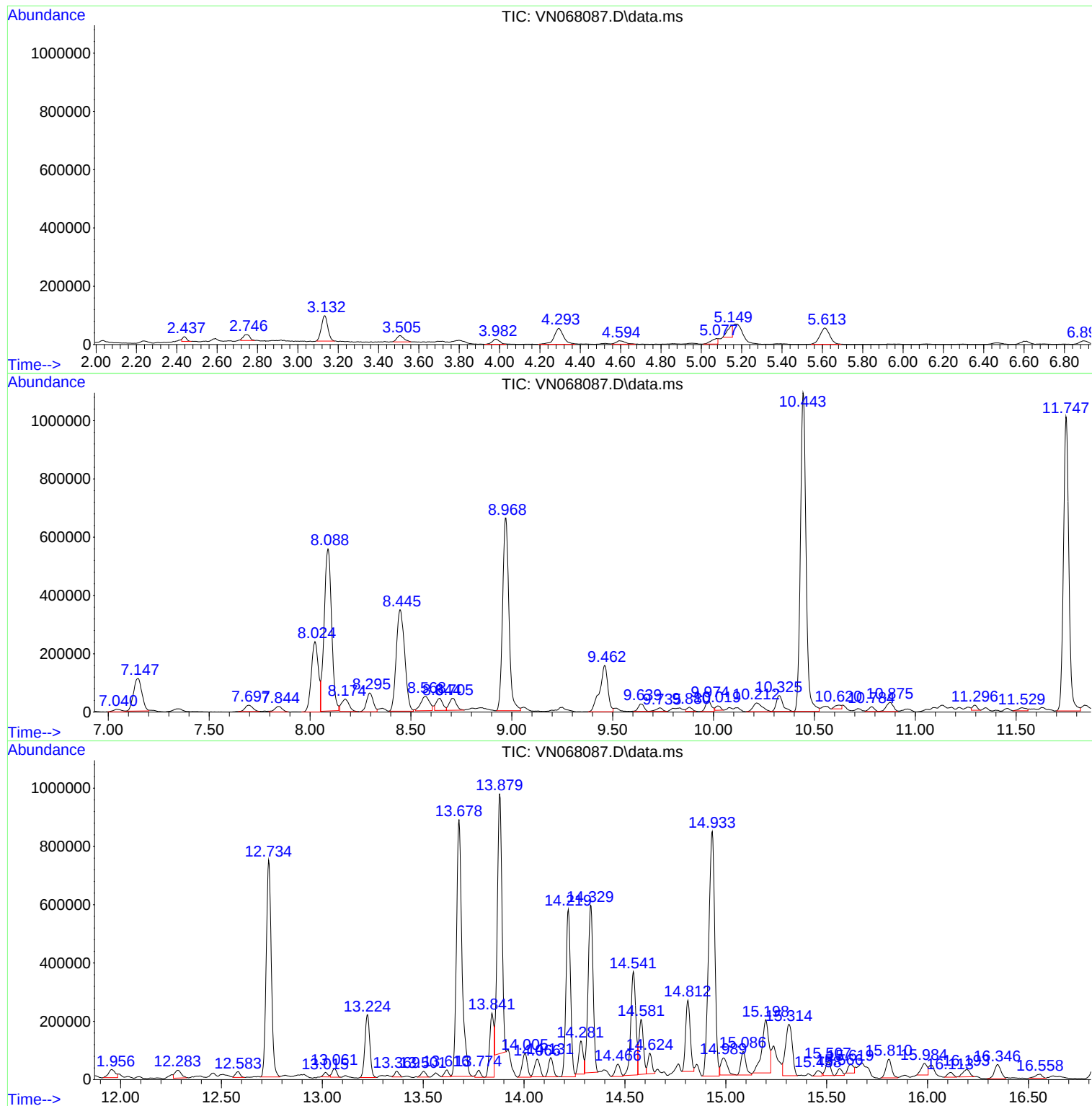
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
Quant Title : SW846 8260

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



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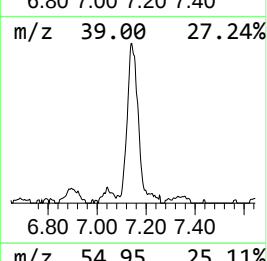
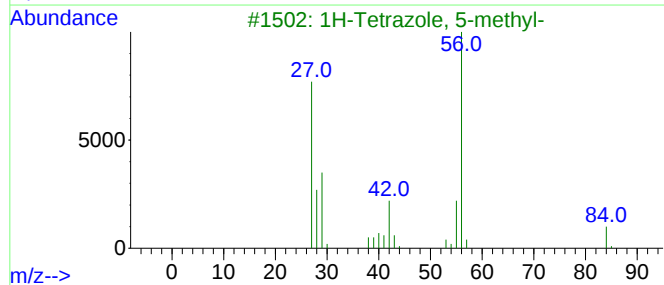
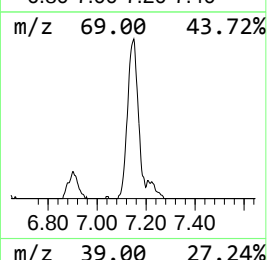
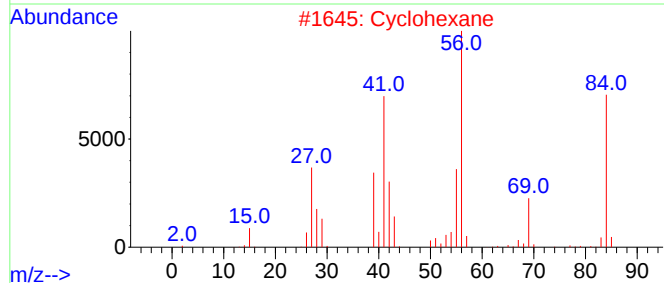
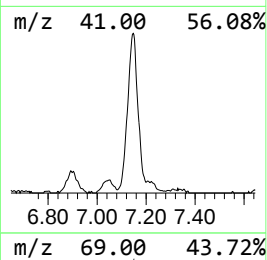
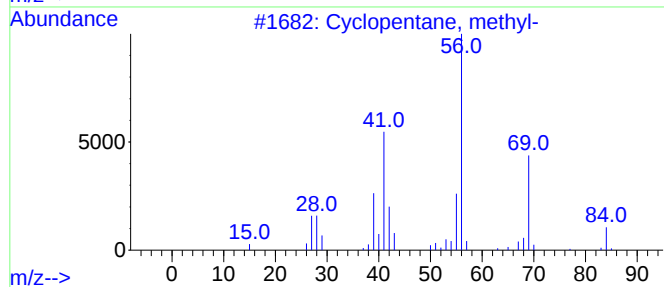
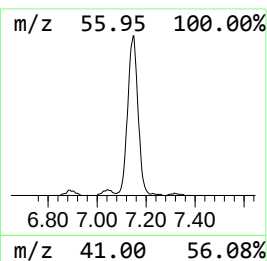
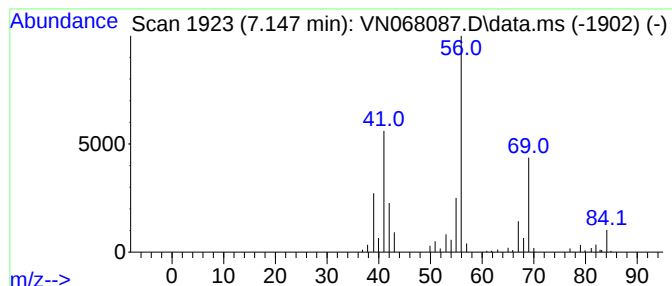
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.147	11.50 ug/l	321526	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2			Cyclohexane	84	C6H12	000110-82-7	80
3			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4			Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
5			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59



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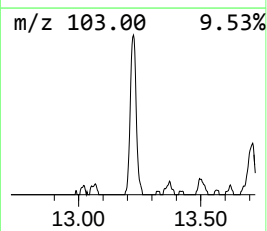
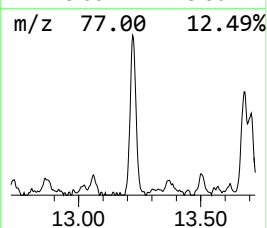
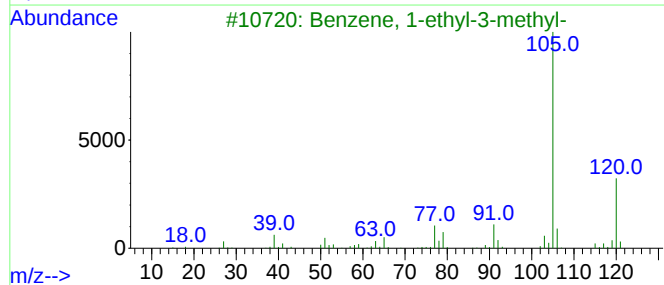
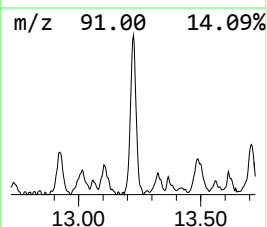
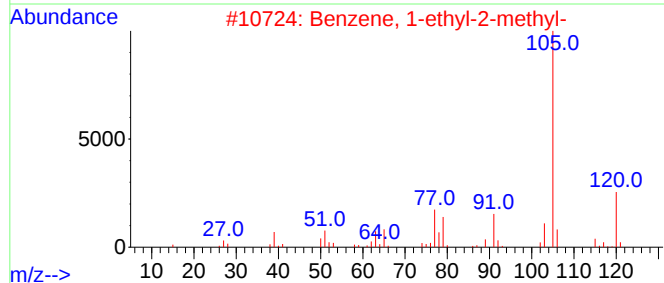
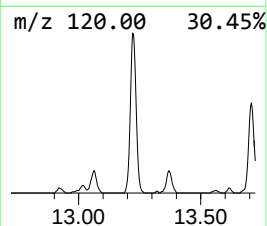
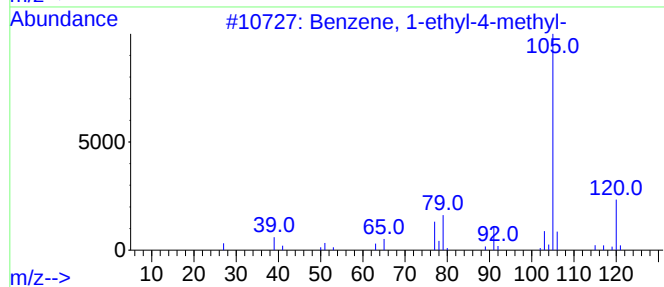
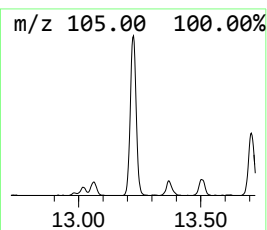
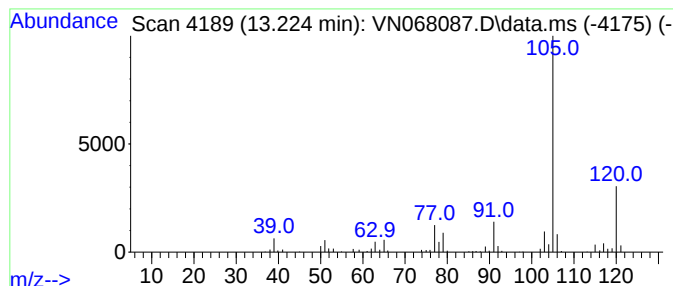
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.224	11.03 ug/l	364427	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	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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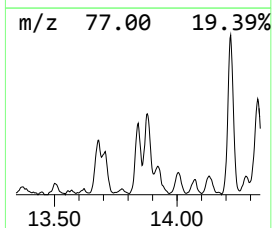
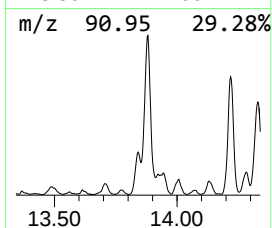
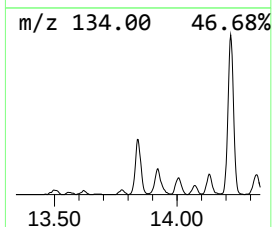
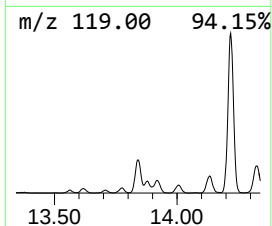
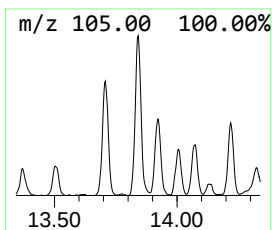
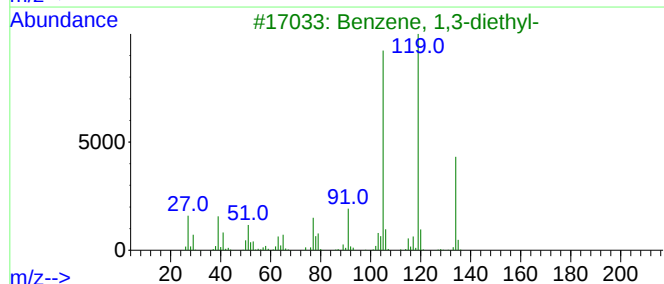
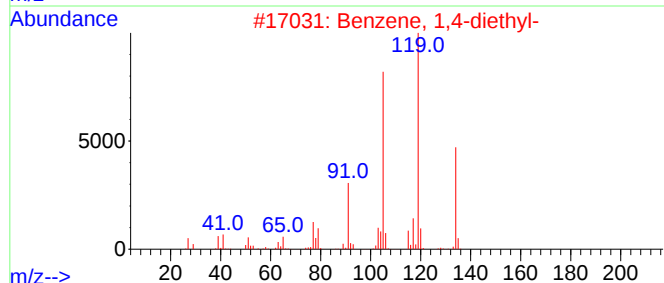
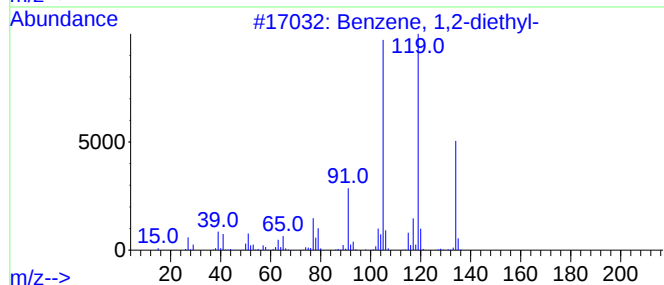
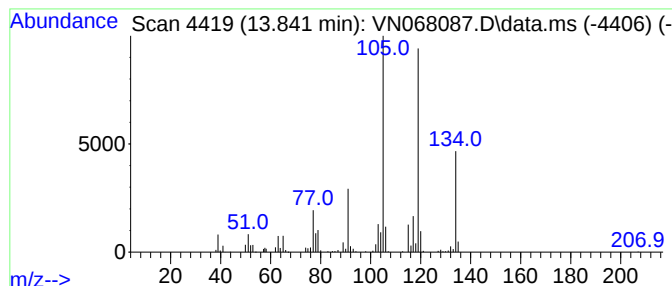
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.841	9.62 ug/l	317713	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	81
5			o-Cymene	134	C10H14	000527-84-4	76



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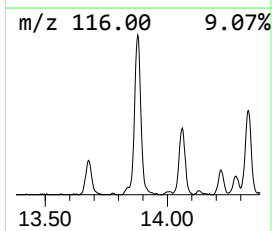
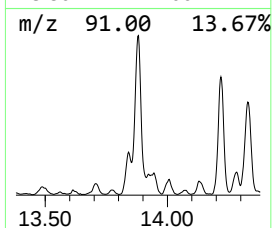
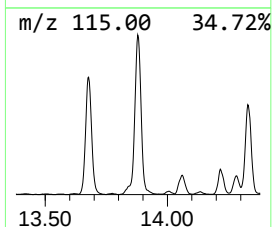
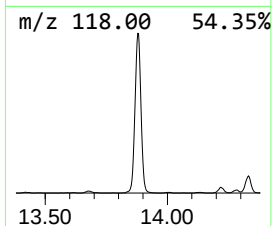
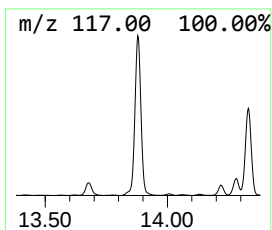
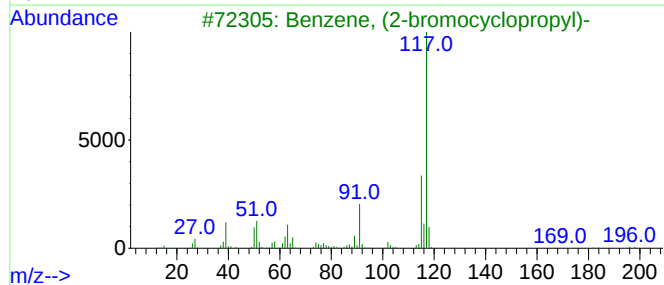
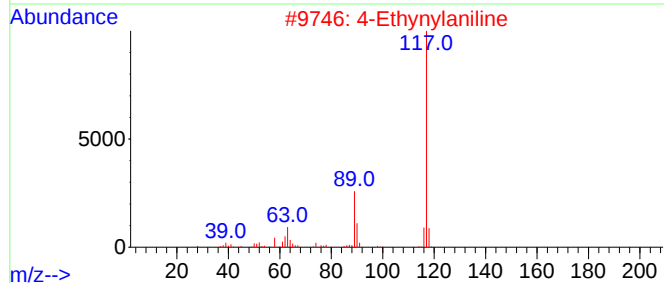
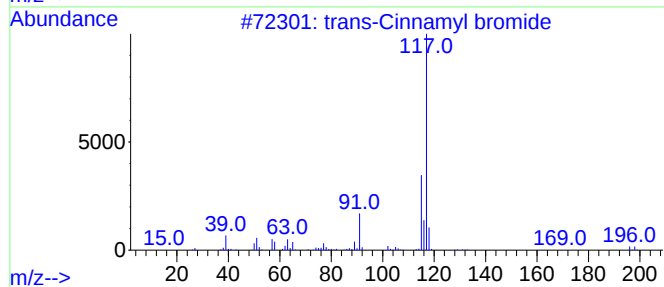
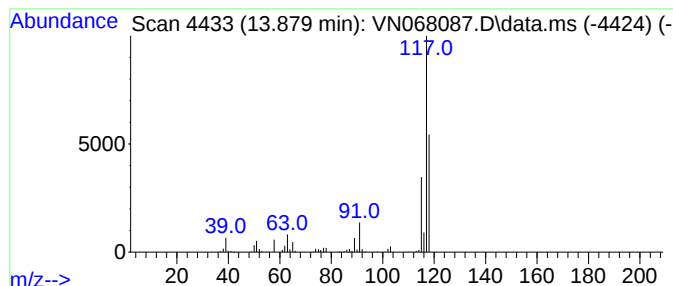
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Quant Title : SW846 8260

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

Peak Number 4 trans-Cinnamyl bromide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.879	42.32 ug/l	1398160	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	53
2			4-Ethynylaniline	117	C8H7N	014235-81-5	47
3			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
4			Indole	117	C8H7N	000120-72-9	43
5			Deltacyclene	118	C9H10	007785-10-6	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

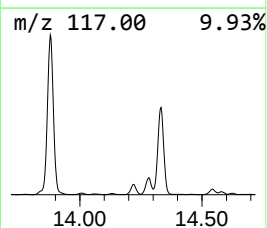
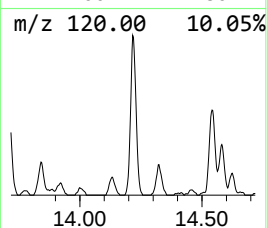
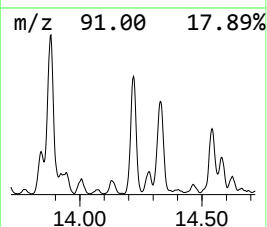
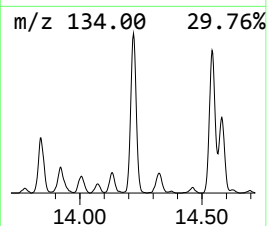
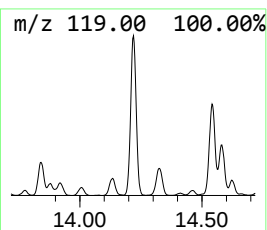
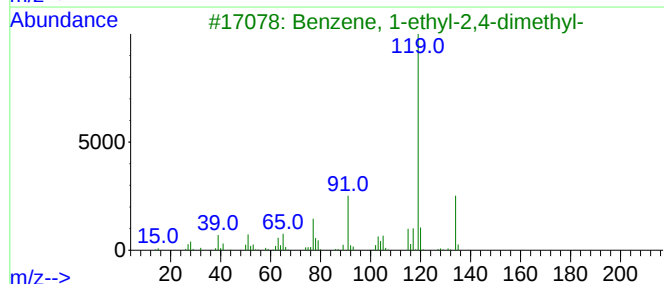
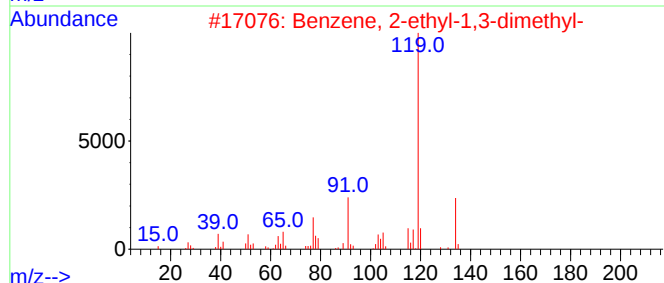
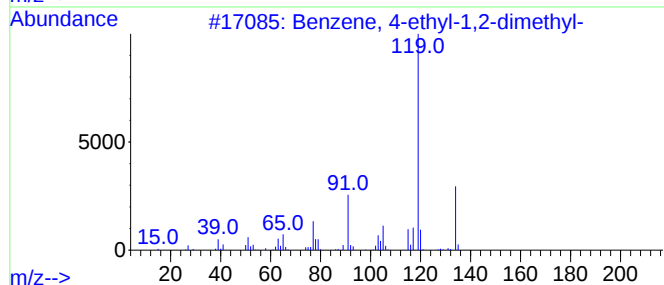
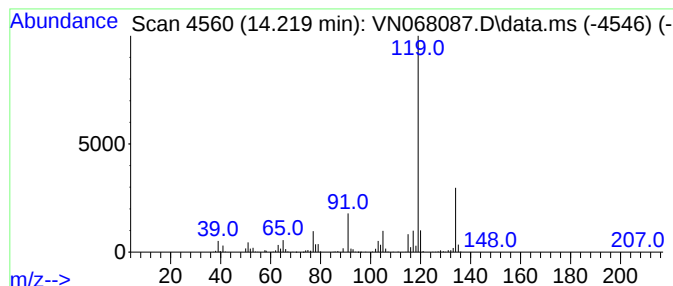
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.219	27.36 ug/l	903759	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

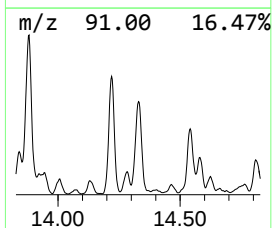
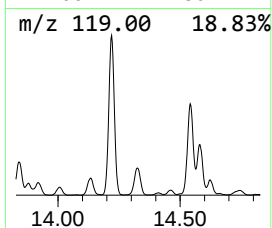
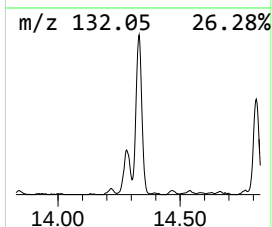
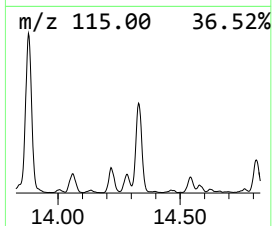
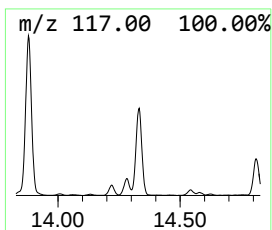
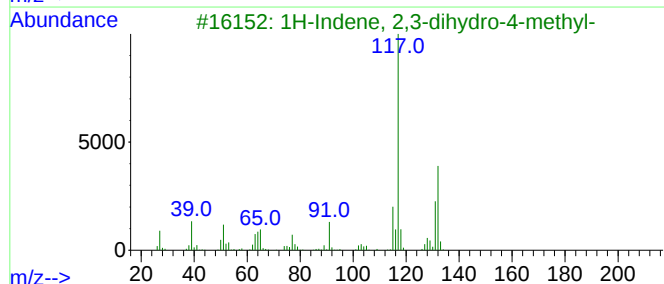
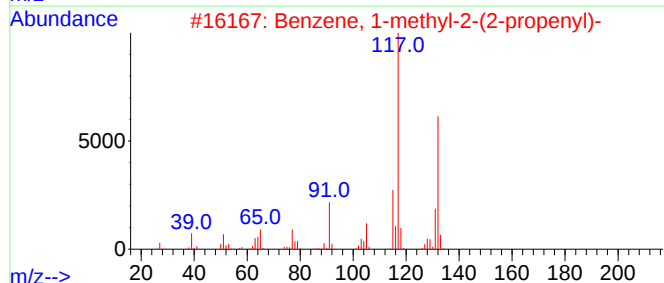
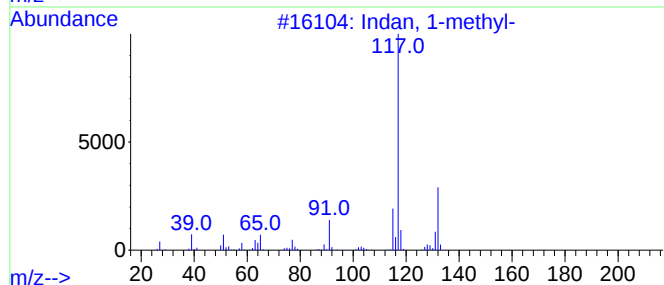
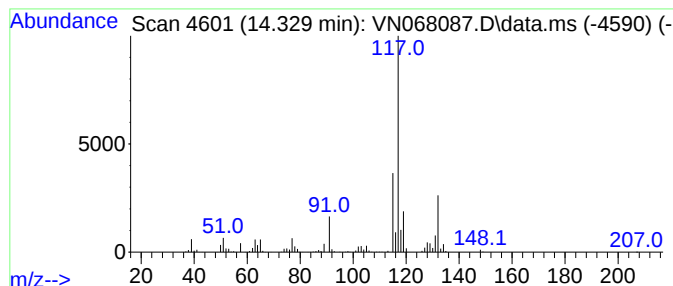
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Quant Title : SW846 8260

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

Peak Number 6 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.329	29.64 ug/l	979065	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	80
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	80
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

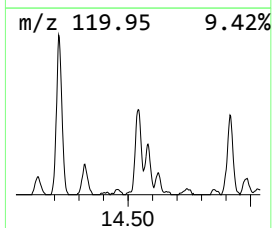
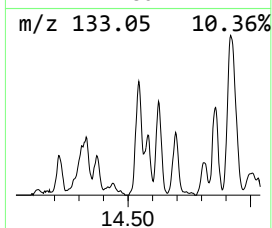
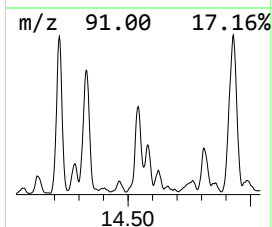
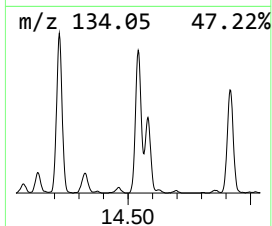
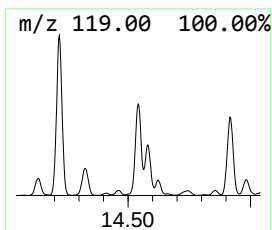
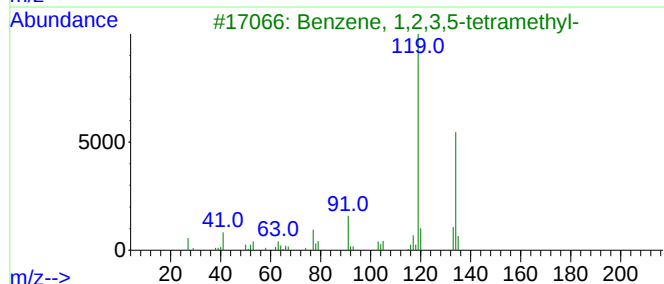
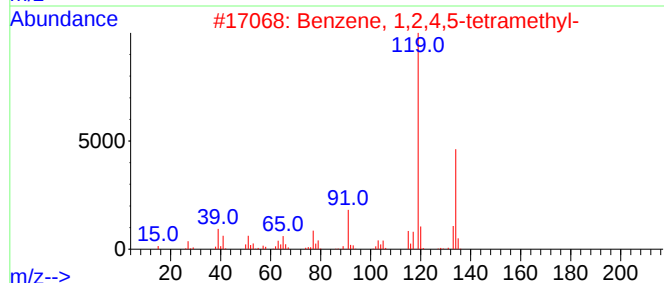
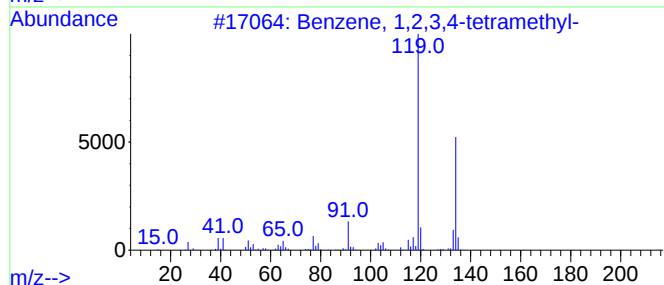
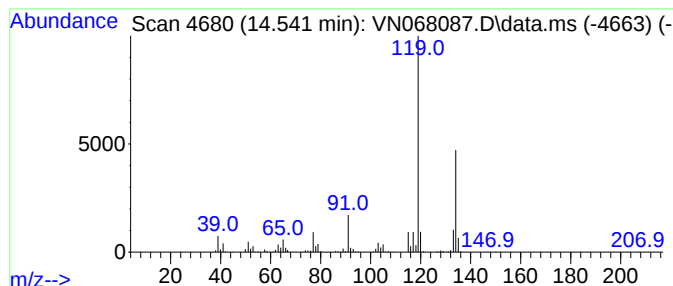
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.541	18.40 ug/l	607992	1,4-Dichlorobenzene-d4	13.678

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-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

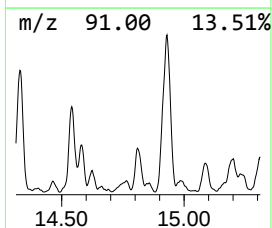
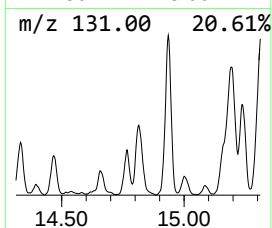
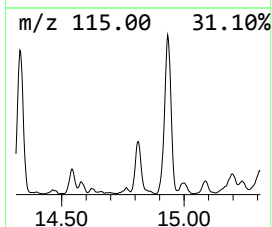
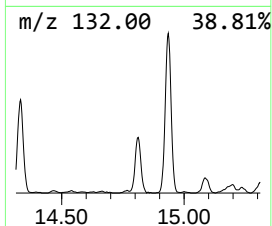
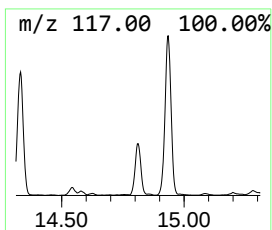
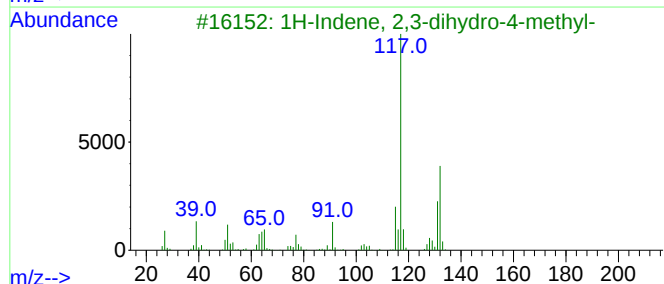
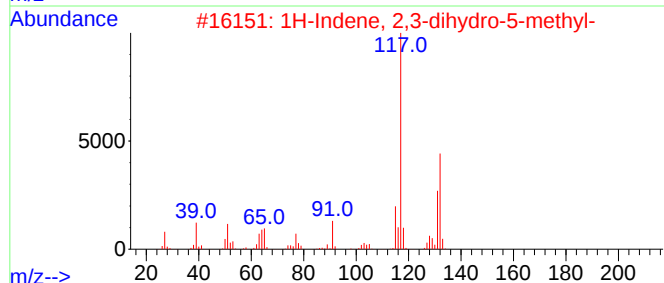
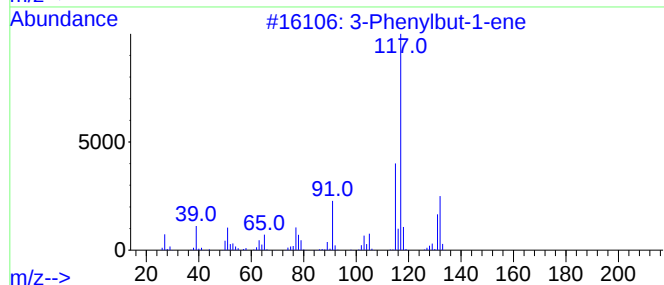
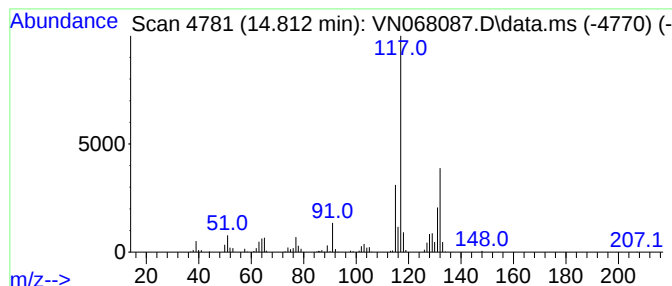
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Quant Title : SW846 8260

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

Peak Number 8 3-Phenylbut-1-ene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.812	11.77 ug/l	388902	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

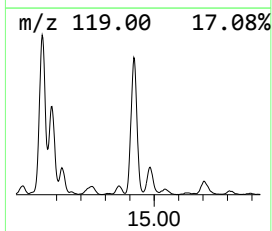
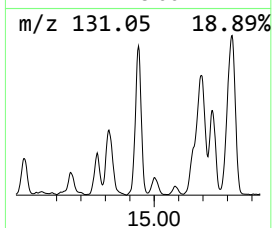
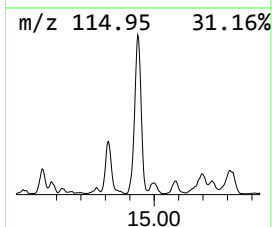
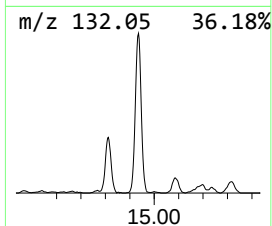
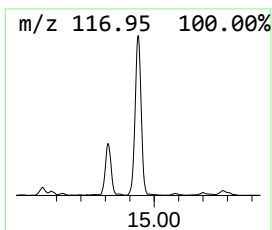
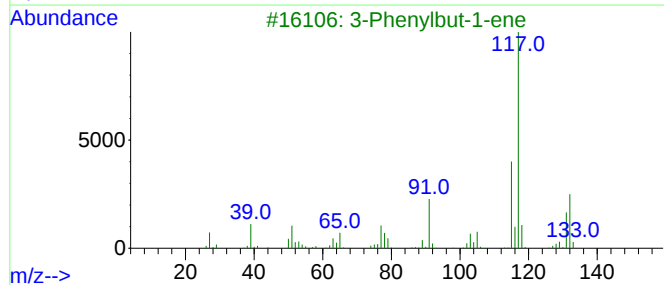
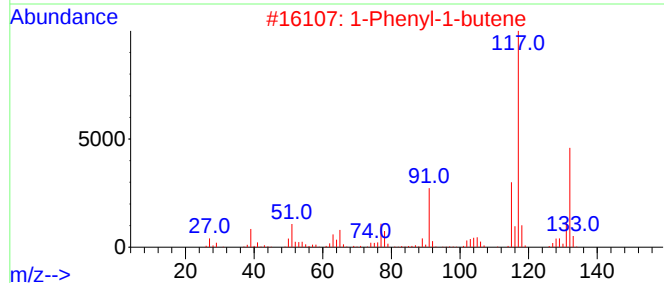
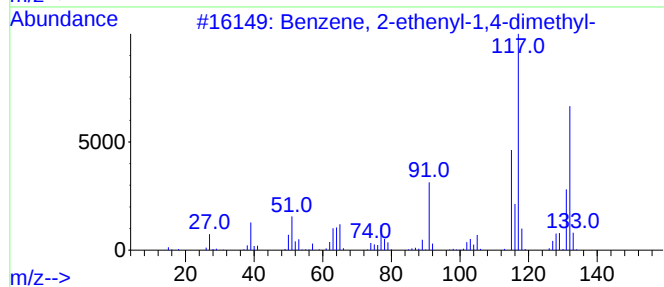
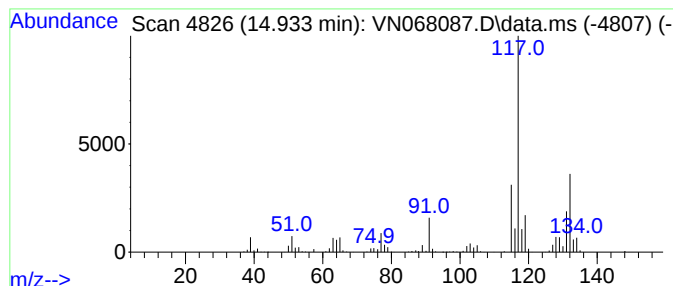
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Quant Title : SW846 8260

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.
14.933	53.39 ug/l	1763590	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	93
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

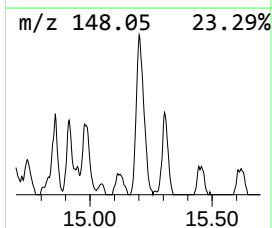
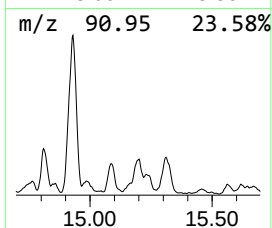
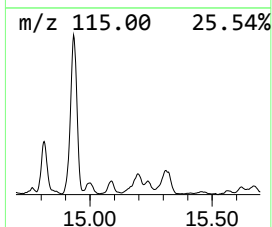
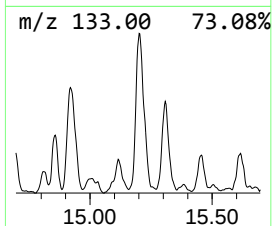
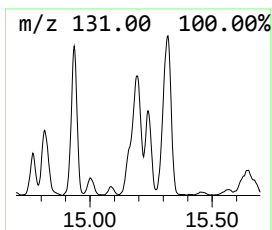
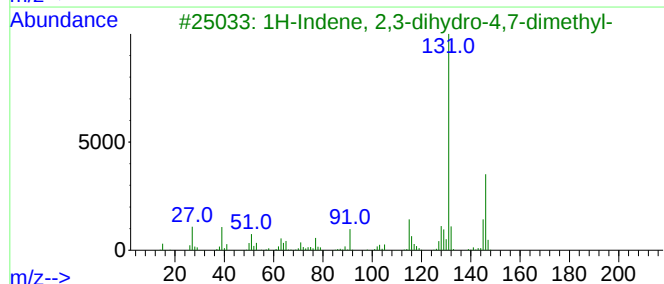
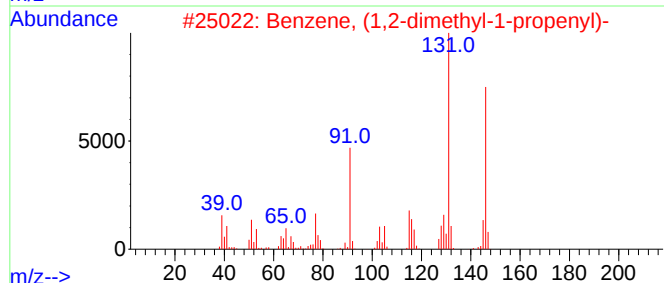
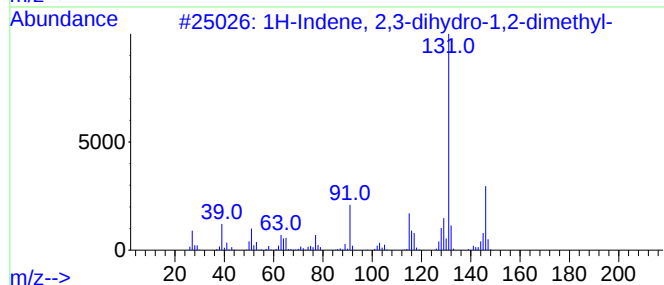
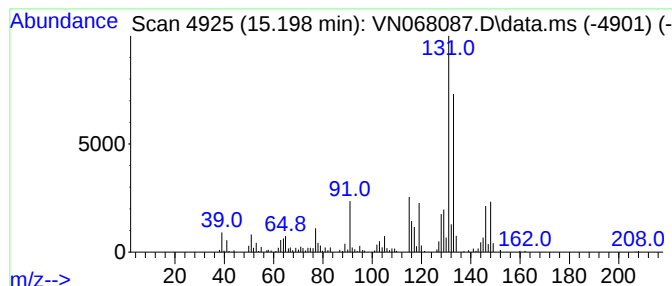
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.198	13.23 ug/l	436906	1,4-Dichlorobenzene-d4	13.678

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
2			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	91
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081221\
Data File : VN068087.D
Acq On : 12 Aug 2021 14:12
Operator : JC/MD
Sample : M3374-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DEWATERING-DISCHARGE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.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
Cyclopentane, m...	7.147	11.5	ug/l	321526	1	8.088	1397570	50.0
Benzene, 1-ethy...	13.224	11.0	ug/l	364427	4	13.678	1651760	50.0
Benzene, 1,2-di...	13.841	9.6	ug/l	317713	4	13.678	1651760	50.0
trans-Cinnamyl ...	13.879	42.3	ug/l	1398160	4	13.678	1651760	50.0
Benzene, 4-ethy...	14.219	27.4	ug/l	903759	4	13.678	1651760	50.0
Indan, 1-methyl-	14.329	29.6	ug/l	979065	4	13.678	1651760	50.0
Benzene, 1,2,3,...	14.541	18.4	ug/l	607992	4	13.678	1651760	50.0
3-Phenylbut-1-ene	14.812	11.8	ug/l	388902	4	13.678	1651760	50.0
Benzene, 2-ethe...	14.933	53.4	ug/l	1763590	4	13.678	1651760	50.0
1H-Indene, 2,3-...	15.198	13.2	ug/l	436906	4	13.678	1651760	50.0