

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092820\
 Data File : VN063717.D
 Acq On : 28 Sep 2020 19:56
 Operator : JC/MD
 Sample : L4178-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-5

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.772	299	321	347	rBV	1632974	3402795	9.06%	0.999%
2	3.107	406	425	456	rBV	1171369	2613340	6.96%	0.767%
3	3.766	608	630	656	rBV5	156530	509337	1.36%	0.150%
4	4.547	832	873	926	rVV2	2233119	11424256	30.42%	3.354%
5	5.004	987	1015	1048	rBV	1752497	6339269	16.88%	1.861%
6	5.332	1093	1117	1145	rVB3	254584	865424	2.30%	0.254%
7	5.521	1150	1176	1205	rBV	1546297	4844592	12.90%	1.422%
8	5.695	1210	1230	1252	rVB2	131597	422990	1.13%	0.124%
9	5.881	1264	1288	1311	rBV2	726209	2239467	5.96%	0.657%
10	6.023	1311	1332	1350	rBV	362449	1140647	3.04%	0.335%
11	6.325	1405	1426	1450	rVB4	376745	1197415	3.19%	0.352%
12	6.476	1454	1473	1486	rBV2	334745	1056582	2.81%	0.310%
13	6.582	1487	1506	1525	rBV	3675015	11353804	30.23%	3.333%
14	7.348	1731	1744	1763	rVB2	778356	2067379	5.50%	0.607%
15	7.611	1789	1826	1843	rBV6	3242383	10801157	28.76%	3.171%
16	7.708	1844	1856	1878	rVB	1099481	3072172	8.18%	0.902%
17	7.840	1878	1897	1910	rBV	1852000	4536376	12.08%	1.332%
18	7.991	1930	1944	1962	rBV3	269005	741568	1.97%	0.218%
19	8.126	1969	1986	1991	rBV4	977126	2291673	6.10%	0.673%
20	8.267	2020	2030	2052	rVB	955634	2385955	6.35%	0.700%
21	8.402	2052	2072	2093	rVB2	1020261	2318187	6.17%	0.681%
22	8.547	2099	2117	2141	rBV4	1639037	4666504	12.42%	1.370%
23	8.830	2197	2205	2235	rVB	538991	1246084	3.32%	0.366%
24	9.049	2242	2273	2286	rBV2	4449442	11175865	29.75%	3.281%
25	9.110	2286	2292	2307	rVB	495536	946117	2.52%	0.278%
26	9.238	2322	2332	2345	rVB	631775	1213938	3.23%	0.356%
27	9.335	2345	2362	2372	rBV2	220952	522059	1.39%	0.153%
28	9.402	2372	2383	2387	rBV2	290668	508730	1.35%	0.149%
29	9.486	2399	2409	2426	rVB2	1174957	2404606	6.40%	0.706%
30	9.585	2427	2440	2446	rBV	1461377	2782371	7.41%	0.817%
31	9.630	2448	2454	2466	rVB5	1187498	2342477	6.24%	0.688%
32	9.698	2466	2475	2482	rBV	501262	851353	2.27%	0.250%
33	9.836	2500	2518	2538	rBV3	711667	2040863	5.43%	0.599%
34	9.939	2538	2550	2560	rVV	1304523	2430495	6.47%	0.714%

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35	9.991	2560	2566	2575	rVV3	262435	466116	1.24%	0.137%
36	10.058	2575	2587	2617	rVV4	1605340	3769329	10.04%	1.107%
37	10.254	2632	2648	2661	rVB7	593514	1691570	4.50%	0.497%
38	10.402	2686	2694	2710	rVB2	505454	909641	2.42%	0.267%
39	10.492	2711	2722	2739	rBV5	774163	1653686	4.40%	0.485%
40	10.692	2765	2784	2792	rBV3	199599	466842	1.24%	0.137%
41	10.765	2793	2807	2815	rBV	779294	1529762	4.07%	0.449%
42	10.891	2839	2846	2850	rVV	323814	501117	1.33%	0.147%
43	10.923	2851	2856	2865	rVV	497755	752952	2.00%	0.221%
44	10.978	2865	2873	2883	rVB	395097	668487	1.78%	0.196%
45	11.090	2900	2908	2917	rVB6	205593	375816	1.00%	0.110%
46	11.177	2918	2935	2956	rVB8	453285	1224356	3.26%	0.359%
47	11.277	2956	2966	2984	rBV2	341909	775516	2.06%	0.228%
48	11.383	2987	2999	3010	rBV2	1039407	1906159	5.07%	0.560%
49	11.486	3011	3031	3048	rVB	10687205	19325787	51.45%	5.674%
50	12.093	3211	3220	3227	rBV4	288765	503728	1.34%	0.148%
51	12.225	3249	3261	3272	rBV	5807318	9420782	25.08%	2.766%
52	12.376	3299	3308	3317	rBV	746077	1142783	3.04%	0.336%
53	12.566	3352	3367	3381	rBV	13076525	22547893	60.03%	6.620%
54	12.865	3447	3460	3475	rBV	2493478	4150092	11.05%	1.218%
55	13.016	3498	3507	3516	rVB	386183	622636	1.66%	0.183%
56	13.142	3532	3546	3557	rBV3	944247	2244180	5.97%	0.659%
57	13.318	3591	3601	3604	rBV	982840	1405588	3.74%	0.413%
58	13.347	3605	3610	3619	rVB	1683482	2516306	6.70%	0.739%
59	13.518	3642	3663	3673	rBV2	17947821	37559787	100.00%	11.027%
60	13.566	3673	3678	3694	rVV2	2113613	4656867	12.40%	1.367%
61	13.646	3694	3703	3713	rVV	1038050	1680331	4.47%	0.493%
62	13.714	3715	3724	3734	rVV	903271	1381900	3.68%	0.406%
63	13.775	3735	3743	3758	rVB	591743	966778	2.57%	0.284%
64	13.862	3758	3770	3781	rBV	9541583	15272342	40.66%	4.484%
65	13.920	3781	3788	3794	rVV	1295690	1996807	5.32%	0.586%
66	13.968	3794	3803	3816	rVB2	5774165	9534053	25.38%	2.799%
67	14.103	3835	3845	3853	rBV2	488279	807773	2.15%	0.237%
68	14.183	3859	3870	3878	rBV	5402056	8467021	22.54%	2.486%
69	14.222	3878	3882	3890	rVB	1334159	1713765	4.56%	0.503%
70	14.267	3890	3896	3903	rBV2	521766	669229	1.78%	0.196%
71	14.450	3943	3953	3963	rBV	4687203	7409016	19.73%	2.175%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
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Stop Thrs : 0 Peak Location: TOP

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72	14.498	3963	3968	3975	rVB	402277	530650	1.41%	0.156%
73	14.566	3975	3989	4001	rBV3	11076729	22118242	58.89%	6.494%
74	14.627	4001	4008	4019	rVB3	1019244	1720571	4.58%	0.505%
75	14.714	4025	4035	4046	rBV	1391458	2196200	5.85%	0.645%
76	14.823	4053	4069	4076	rBV5	1118684	2284791	6.08%	0.671%
77	14.929	4091	4102	4115	rVB2	1113331	2278595	6.07%	0.669%
78	15.103	4136	4156	4169	rBV	3602981	6426692	17.11%	1.887%
79	15.212	4181	4190	4198	rBV2	352738	633502	1.69%	0.186%
80	15.260	4199	4205	4230	rVB2	344887	681761	1.82%	0.200%
81	15.386	4233	4244	4257	rBV	552466	1021310	2.72%	0.300%
82	15.547	4283	4294	4321	rVB2	436583	1309058	3.49%	0.384%
83	15.736	4343	4353	4364	rVB2	248991	492659	1.31%	0.145%
84	16.122	4459	4473	4497	rVB	5103639	10399350	27.69%	3.053%
85	16.312	4516	4532	4558	rVB	3330330	7083401	18.86%	2.080%

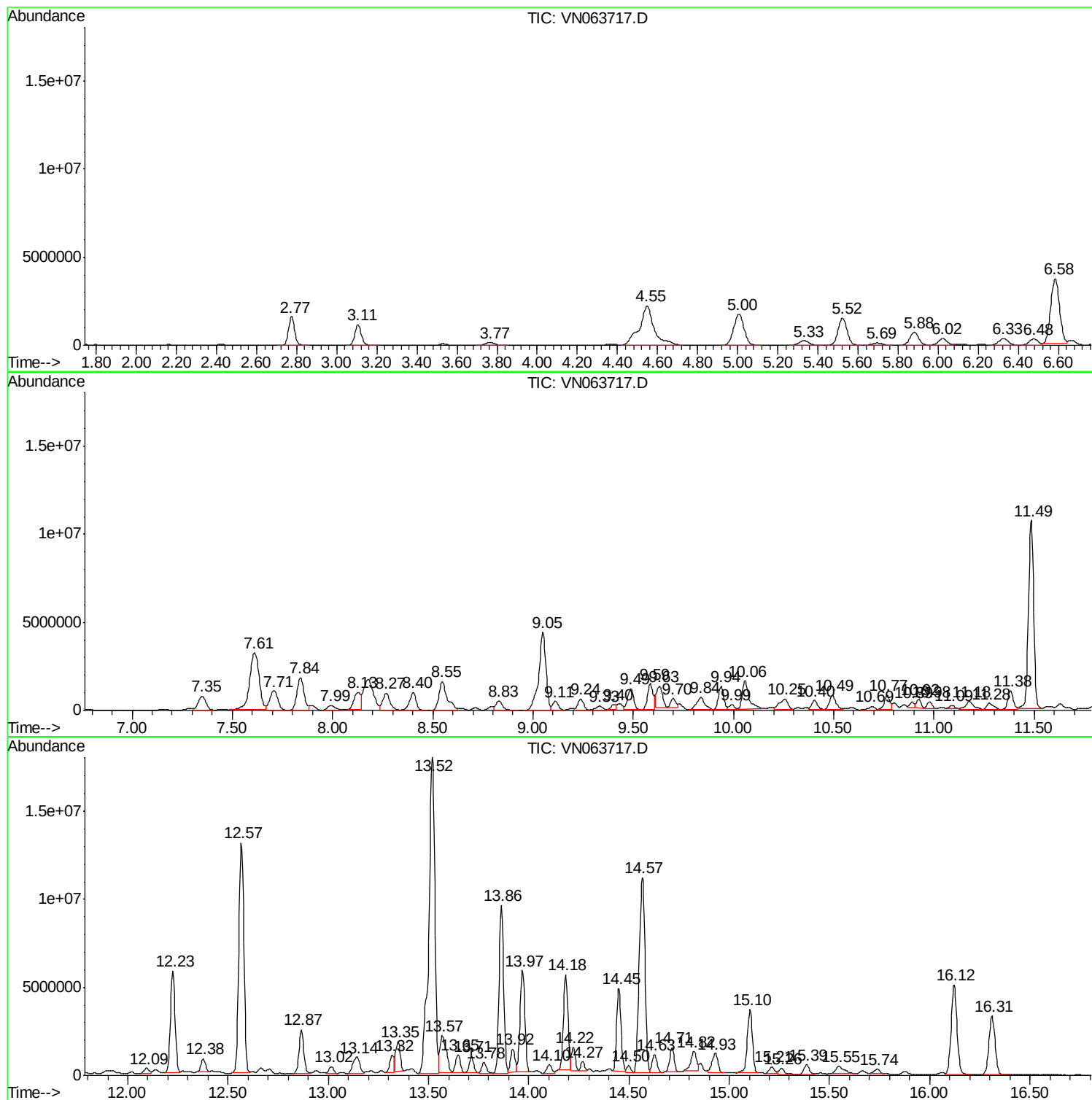
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampled :
TP-5

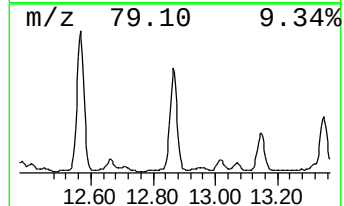
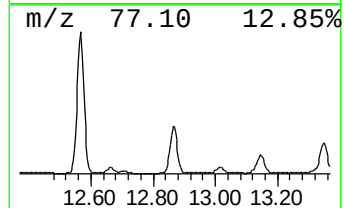
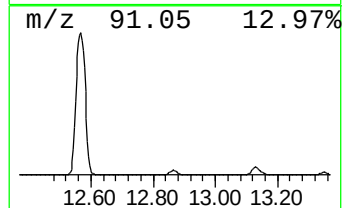
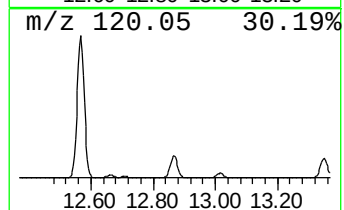
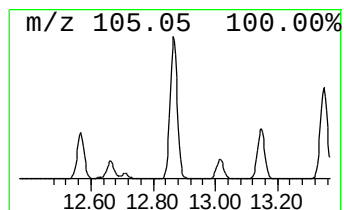
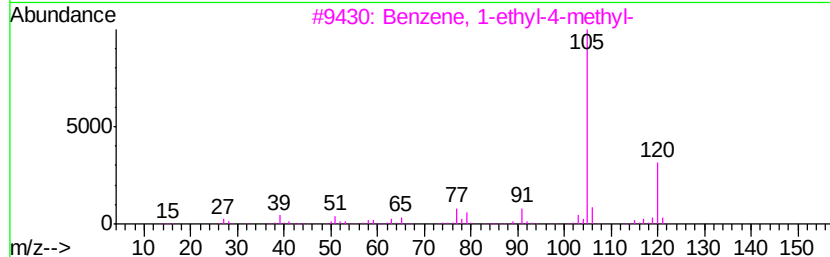
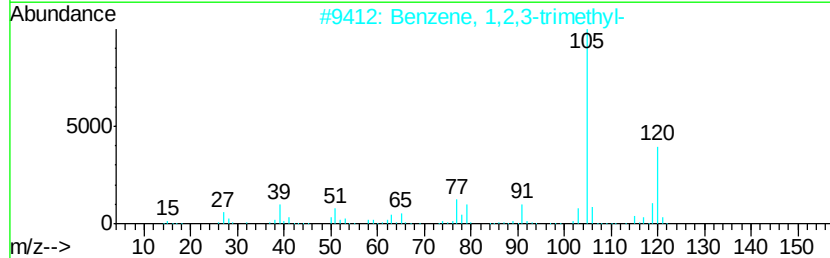
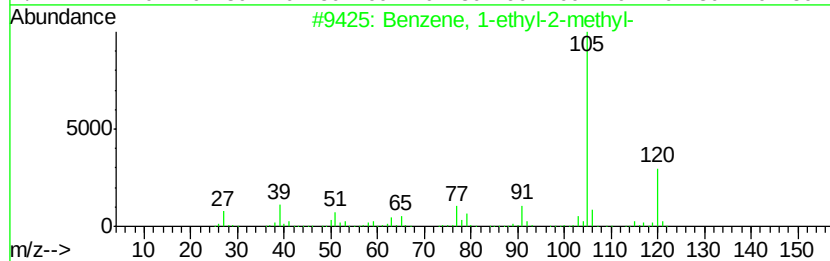
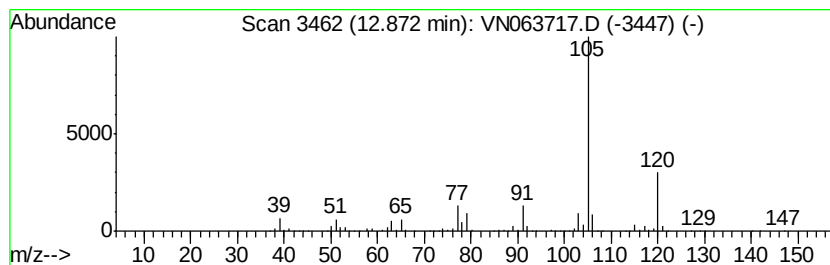
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	147.63 ug/l	4150090	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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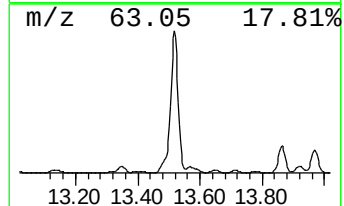
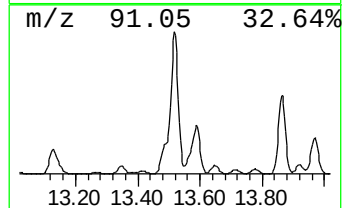
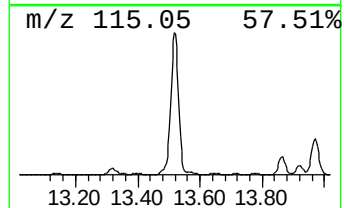
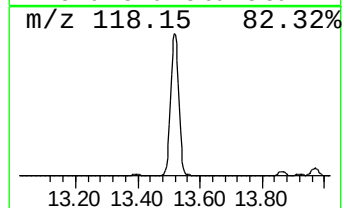
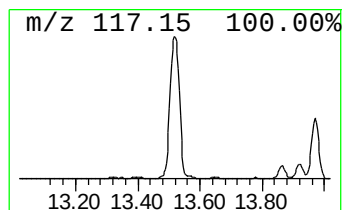
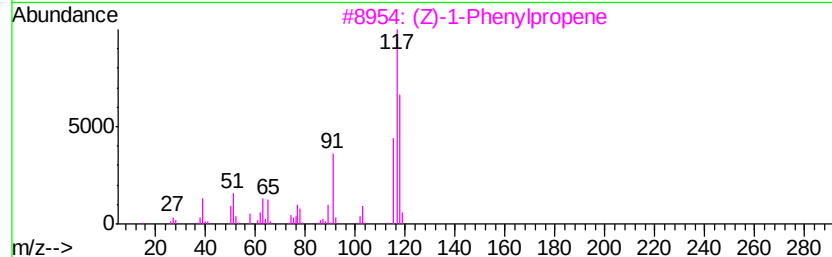
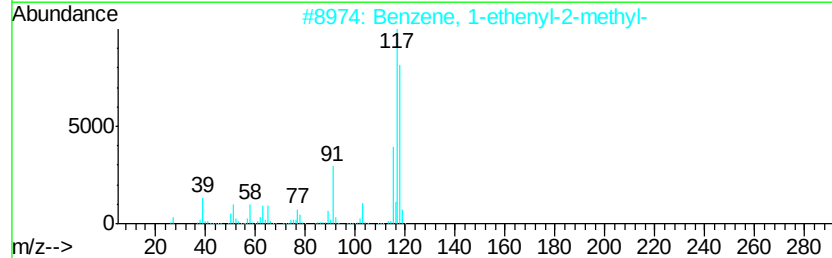
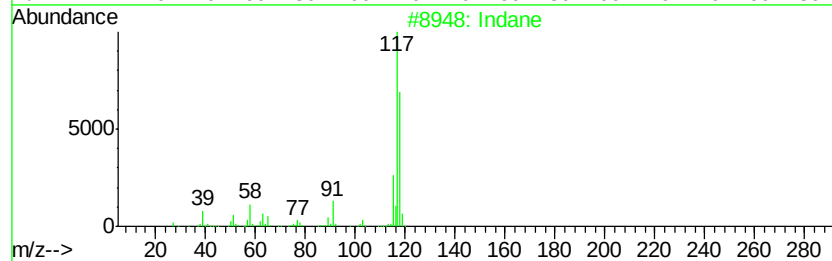
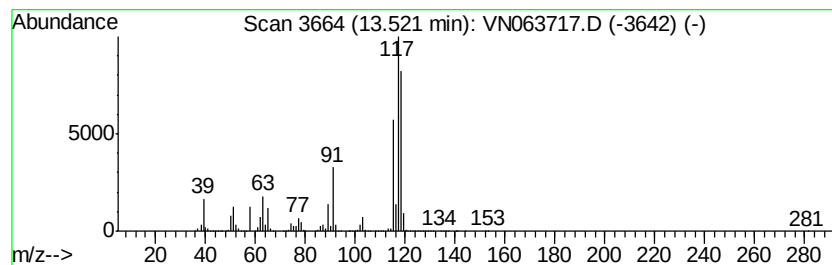
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	1336.09 ug/l	37559800	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	93
3	(Z)-1-Phenylpropene		118	C9H10	000766-90-5	90
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	81
5	Benzene, 1-propenyl-		118	C9H10	000637-50-3	81



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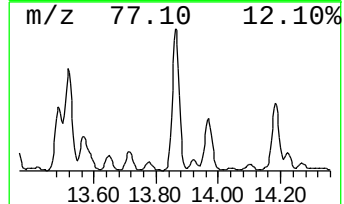
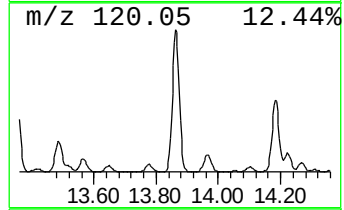
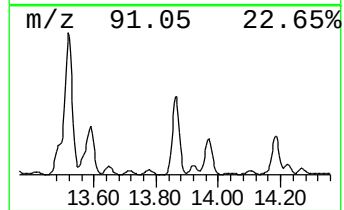
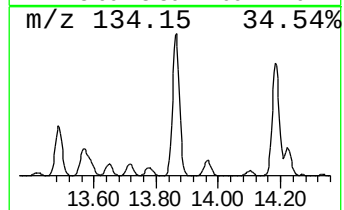
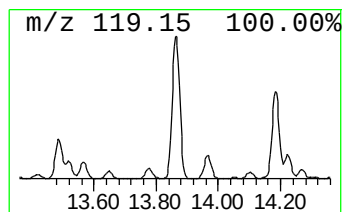
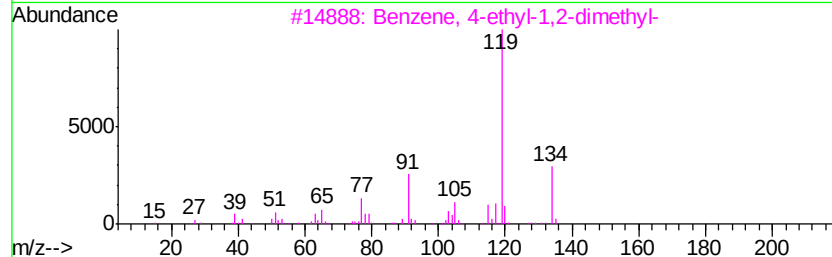
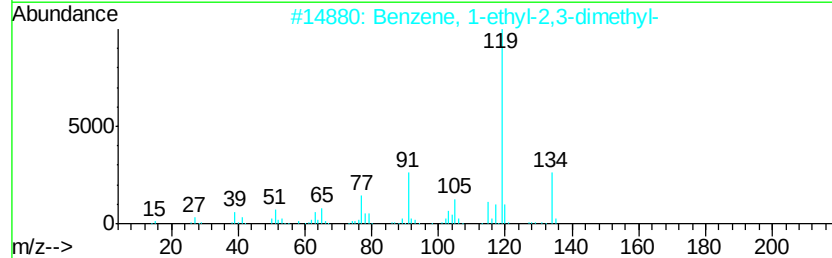
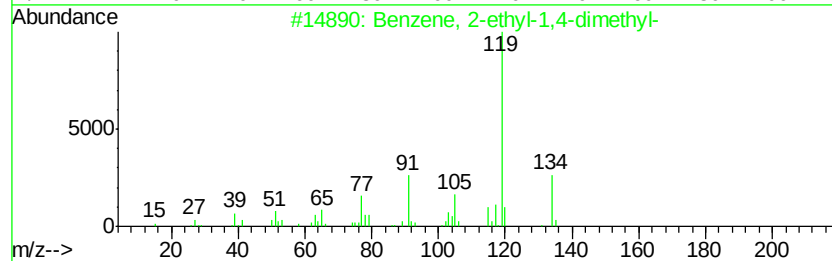
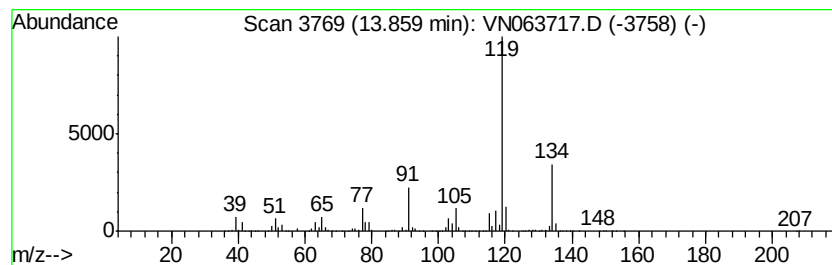
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	543.27 ug/l	15272300	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96	
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96	
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96	
4	o-Cymene	134	C10H14	000527-84-4	95	
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95	



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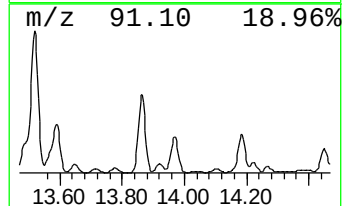
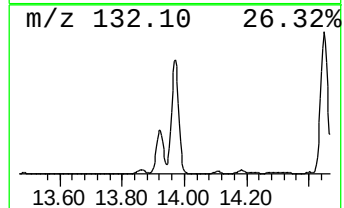
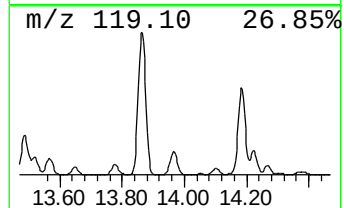
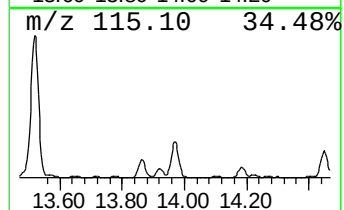
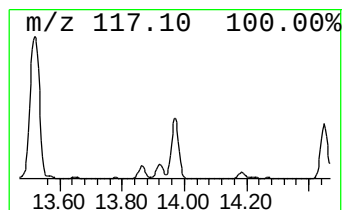
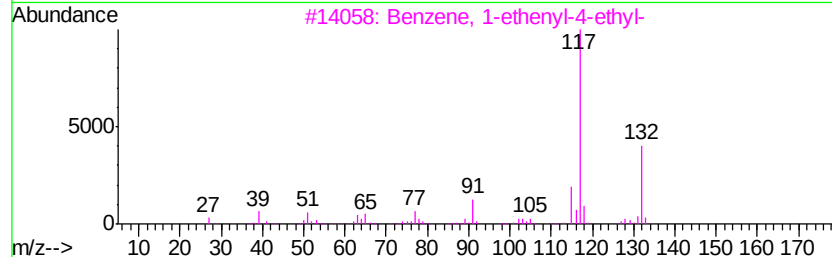
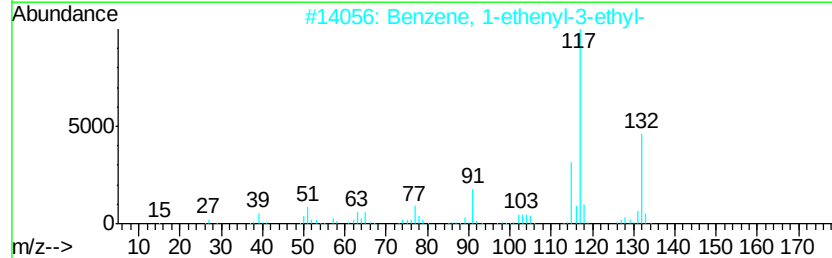
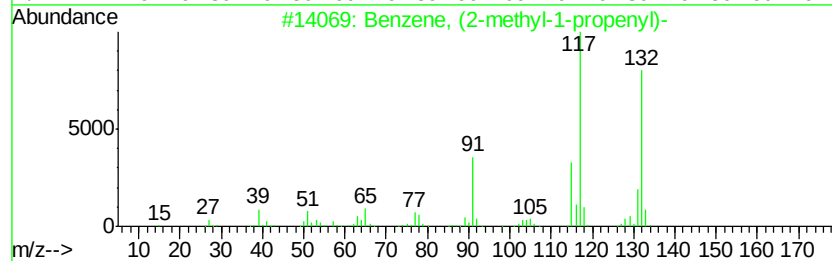
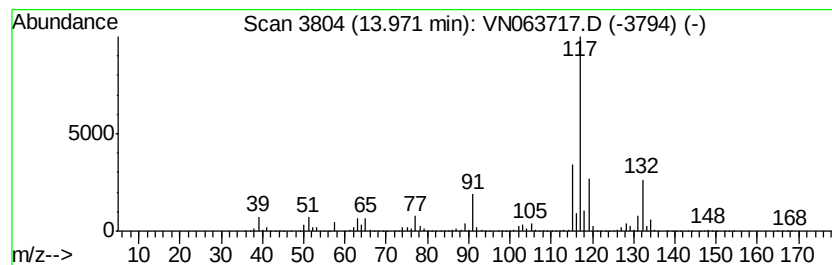
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, (2-methyl-1-propen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	339.15 ug/l	9534050	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
2		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
4		Indan, 1-methyl-	132	C10H12	000767-58-8	76
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TP-5

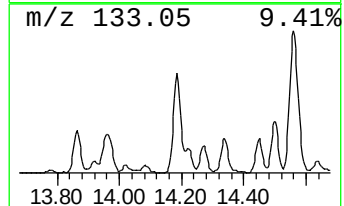
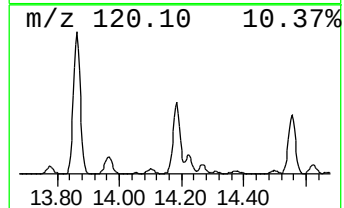
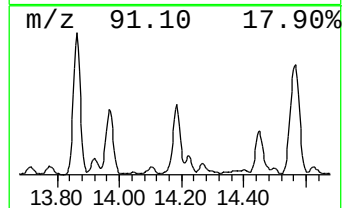
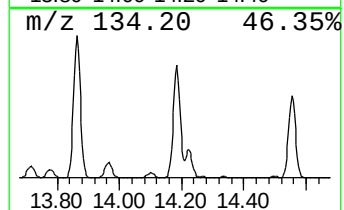
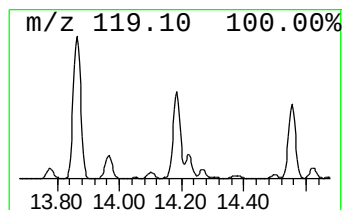
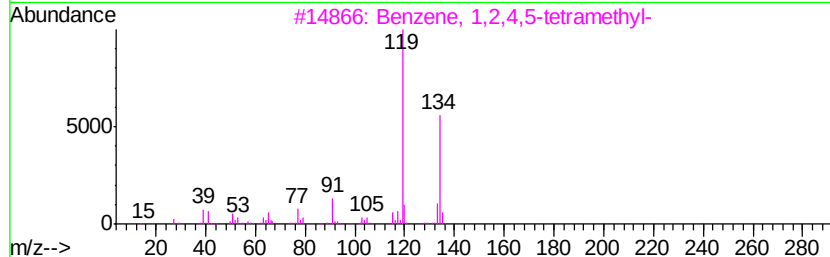
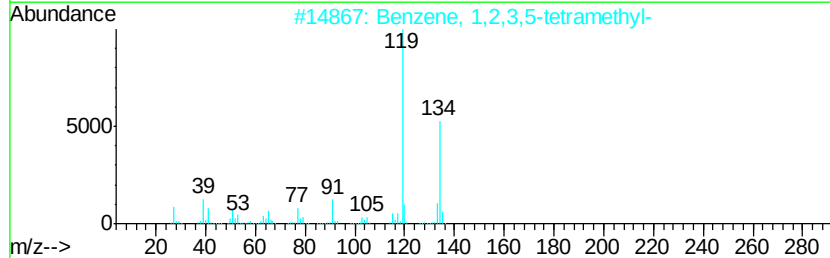
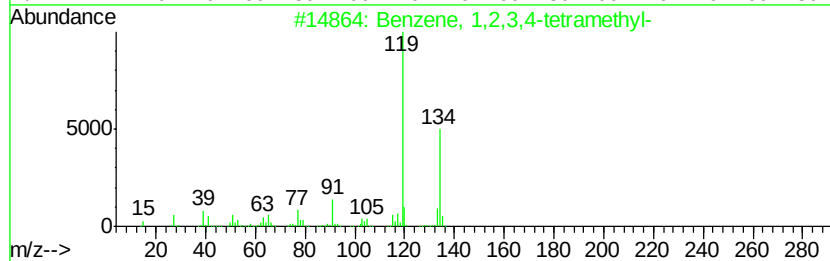
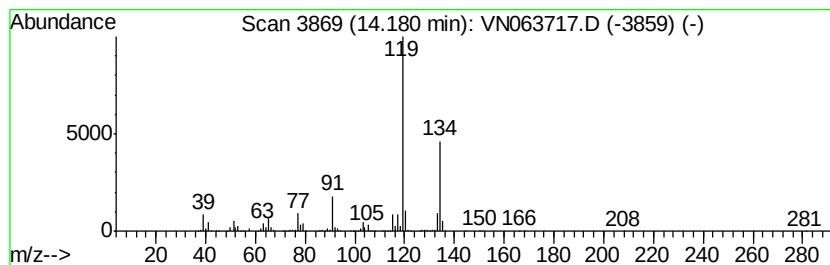
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	301.19 ug/l	8467020	1,4-Dichlorobenzene-d4	13.32

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-5

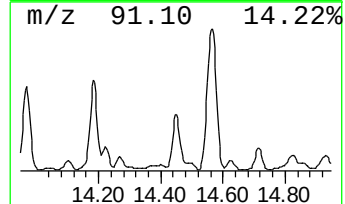
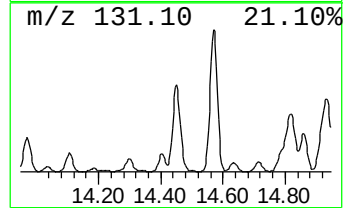
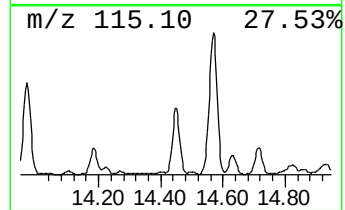
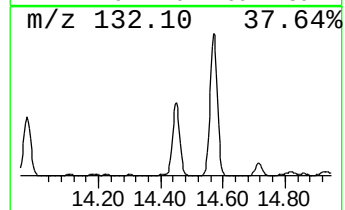
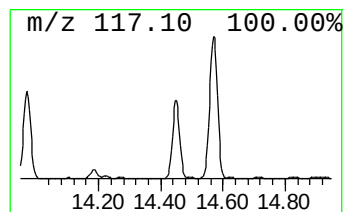
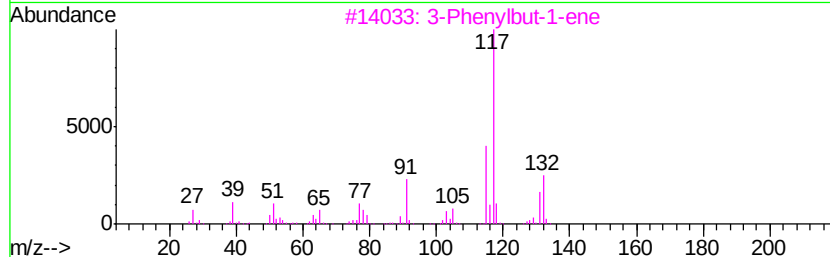
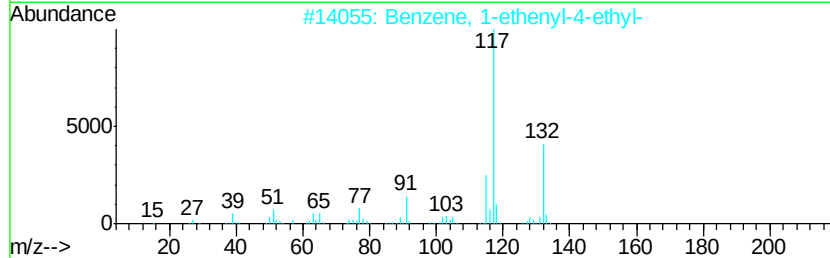
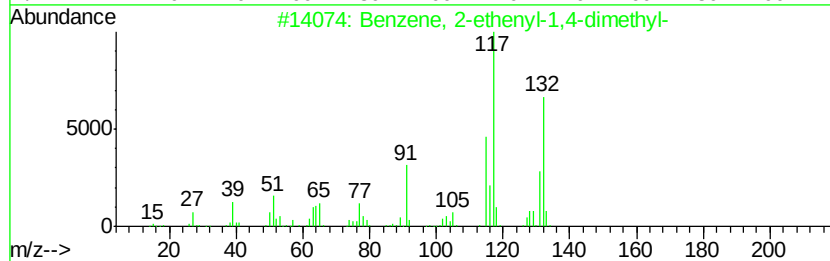
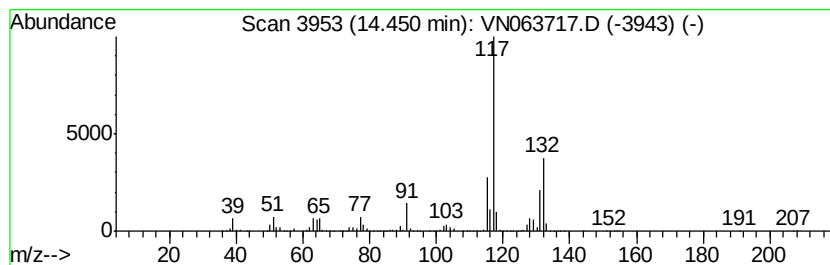
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	263.56 ug/l	7409020	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TP-5

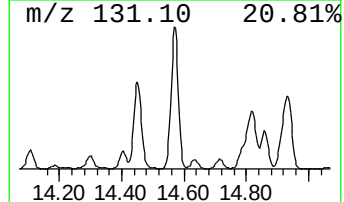
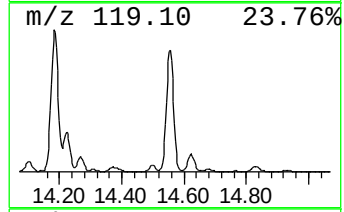
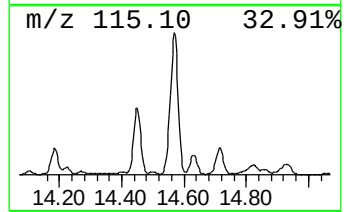
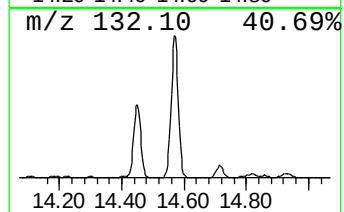
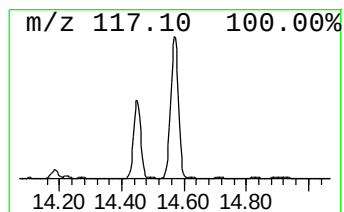
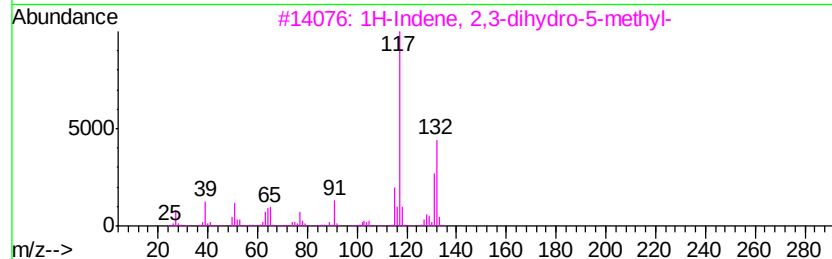
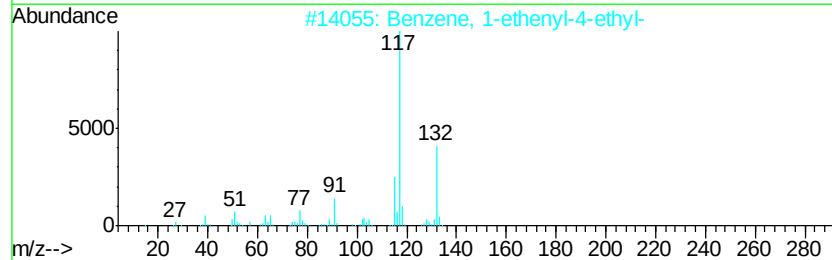
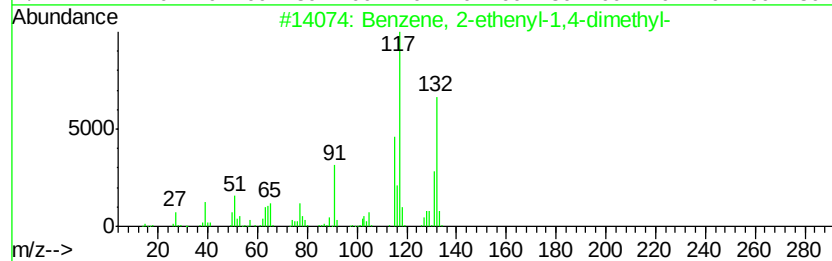
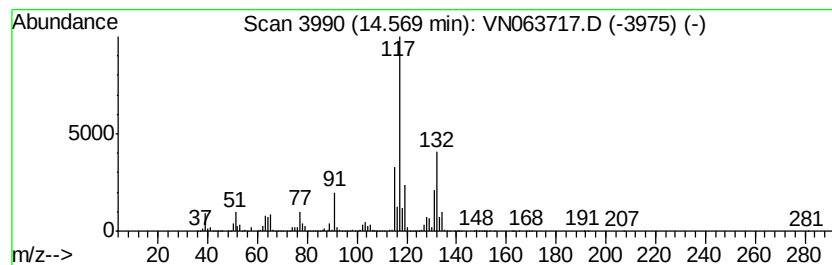
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	786.80 ug/l	22118200	1,4-Dichlorobenzene-d4	13.32

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	93
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	90
5		2,4-Dimethylstyrene	132	C10H12	002234-20-0	90



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

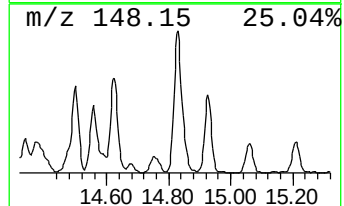
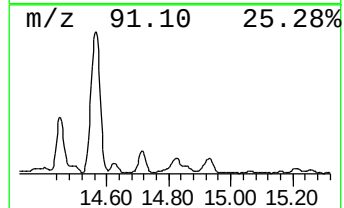
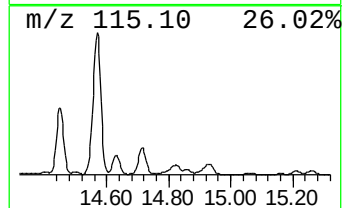
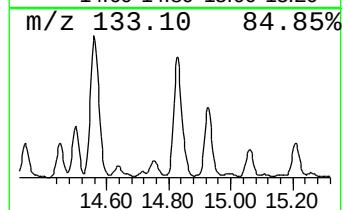
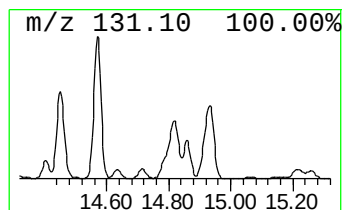
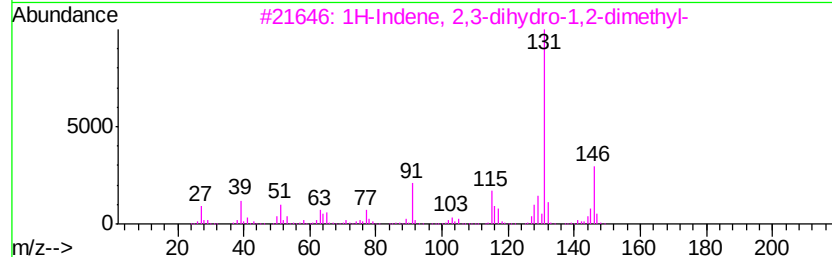
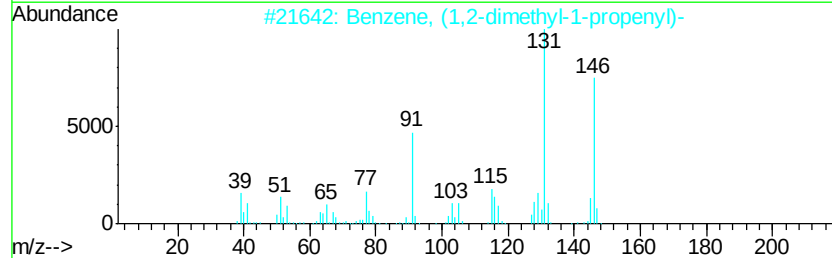
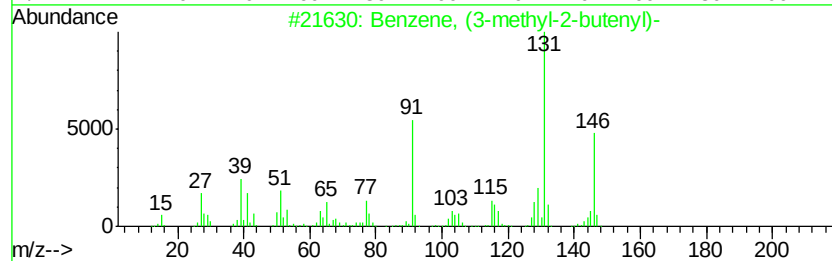
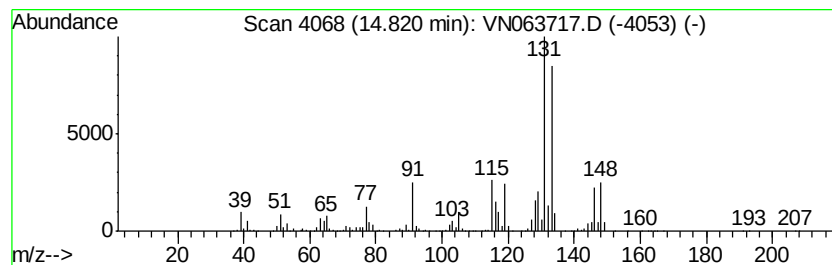
Instrument :
MSVOA_N
ClientSampleId :
TP-5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, (3-methyl-2-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	81.28 ug/l	2284790	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 64
2		Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3 64
3		1H-Indene, 2,3-dihydro-1,2-dimethyl-	146 C11H14	017057-82-8 64
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 50
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3 50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-5

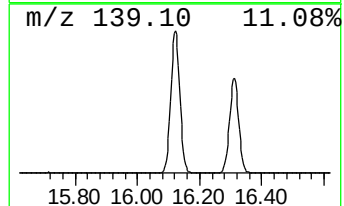
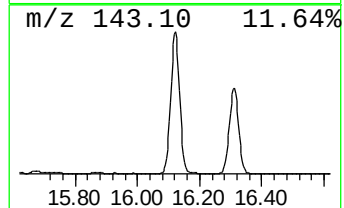
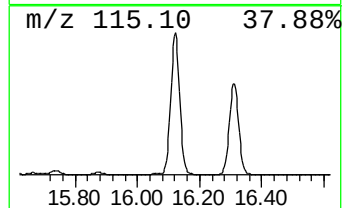
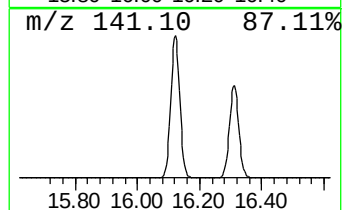
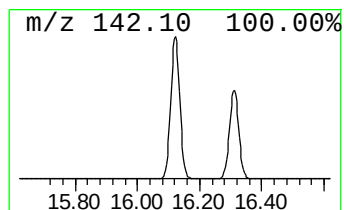
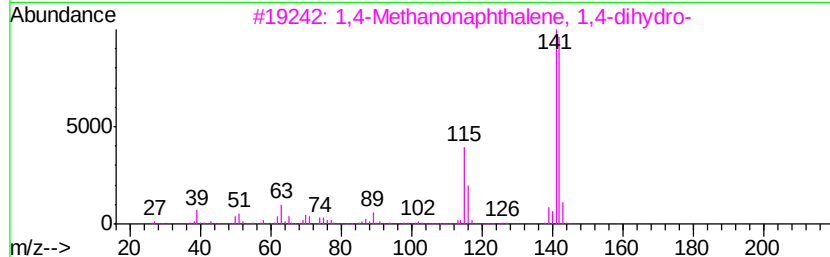
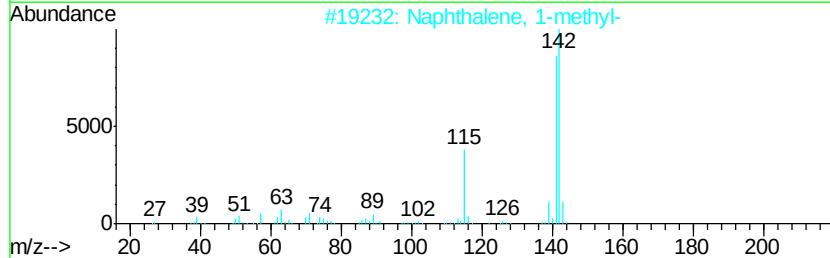
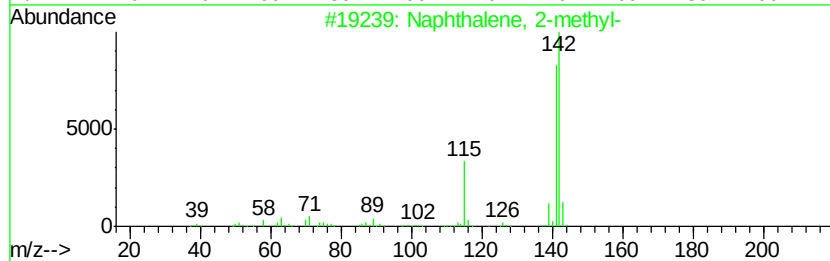
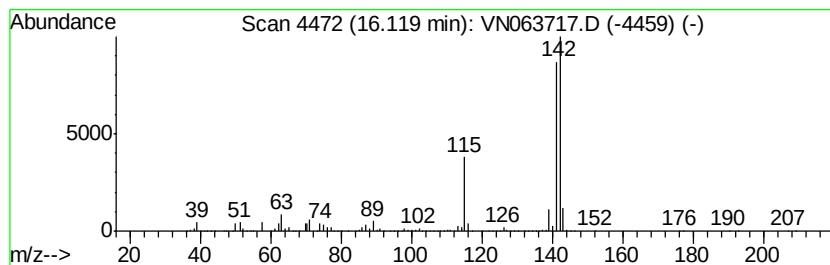
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,4-Methanonaphthalene, 1,4... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.12	369.93 ug/l	10399400	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-dihv...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 21 Sample Multiplier: 1

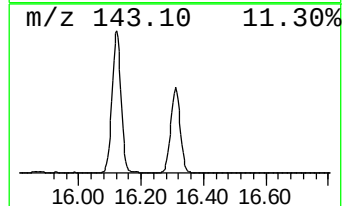
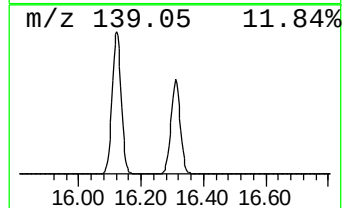
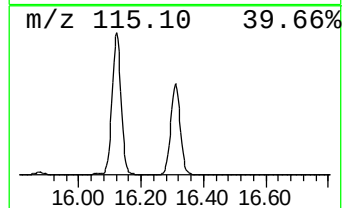
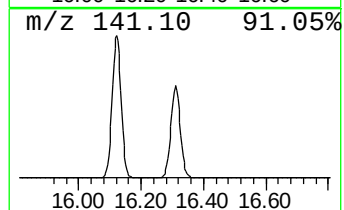
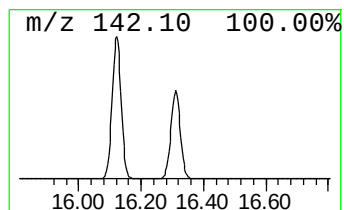
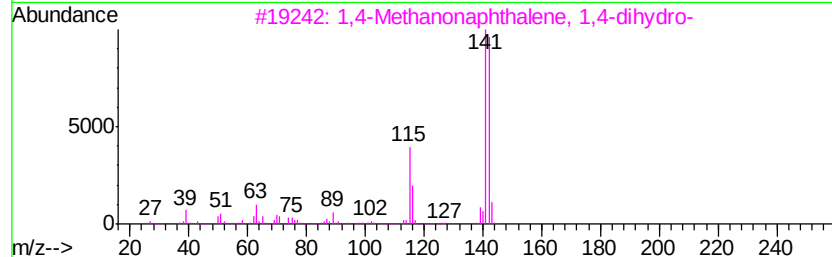
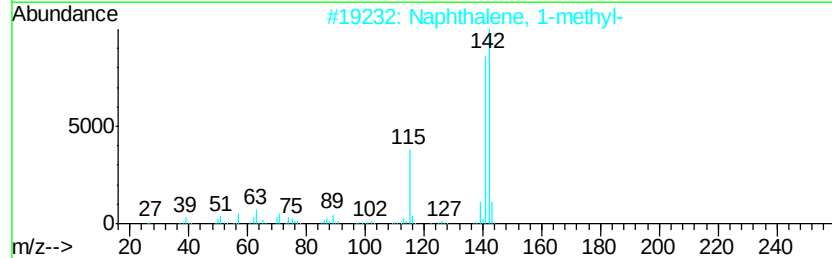
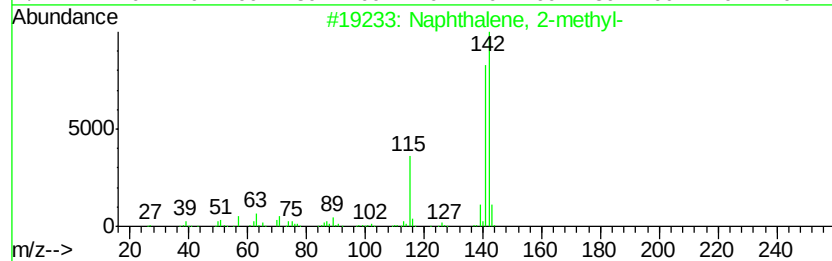
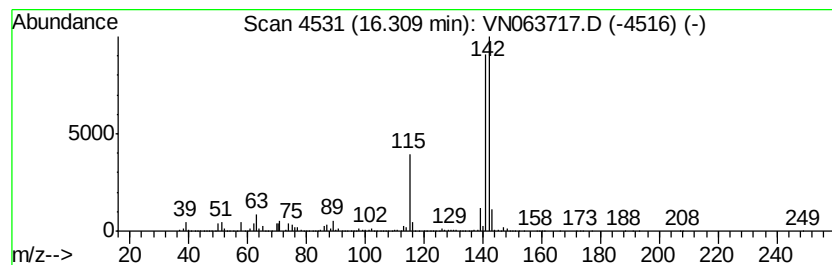
Instrument :
MSVOA_N
ClientSampleId :
TP-5

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
16.31	251.97 ug/l	7083400	1,4-Dichlorobenzene-d4	13.32		
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-dihv...	142	C11H10	004453-90-1	94
4		Benzocycloheptatriene	142	C11H10	000264-09-5	91
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN092820\
Data File : VN063717.D
Acq On : 28 Sep 2020 19:56
Operator : JC/MD
Sample : L4178-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	12.87	147.6	ug/l	4150090	4	13.32	1405590	50.0
Indane	13.52	1336.1	ug/l	37559800	4	13.32	1405590	50.0
Benzene, 2-ethyl-...	13.86	543.3	ug/l	15272300	4	13.32	1405590	50.0
Benzene, (2-methy...	13.97	339.1	ug/l	9534050	4	13.32	1405590	50.0
Benzene, 1,2,3,4-...	14.18	301.2	ug/l	8467020	4	13.32	1405590	50.0
Benzene, 2-etheny...	14.45	263.6	ug/l	7409020	4	13.32	1405590	50.0
Benzene, 1-etheny...	14.57	786.8	ug/l	22118200	4	13.32	1405590	50.0
Benzene, (3-methy...	14.82	81.3	ug/l	2284790	4	13.32	1405590	50.0
1,4-Methanonaphth...	16.12	369.9	ug/l	10399400	4	13.32	1405590	50.0
Naphthalene, 1-me...	16.31	252.0	ug/l	7083400	4	13.32	1405590	50.0