

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070820\
 Data File : VN062433.D
 Acq On : 9 Jul 2020 2:47
 Operator : JC/MD
 Sample : L3215-10
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-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\82N070820W.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	2.773	299	311	325	rBB2	27756	55052	1.68%	0.274%
2	4.548	852	863	891	rVB5	34953	129440	3.95%	0.644%
3	4.995	977	1002	1033	rBV2	127403	470979	14.37%	2.345%
4	5.876	1257	1276	1297	rVB6	10695	35256	1.08%	0.176%
5	6.580	1476	1495	1515	rVV3	75528	233397	7.12%	1.162%
6	7.551	1778	1797	1806	rBV	133110	343826	10.49%	1.712%
7	7.622	1807	1819	1834	rVV2	308097	807284	24.63%	4.019%
8	7.709	1835	1846	1865	rVB3	59210	161864	4.94%	0.806%
9	7.837	1865	1886	1898	rBV3	54564	140393	4.28%	0.699%
10	7.992	1918	1934	1956	rBV5	230750	612579	18.69%	3.050%
11	8.130	1957	1977	1985	rVV9	94817	282250	8.61%	1.405%
12	8.198	1987	1998	2011	rVV2	250763	639868	19.53%	3.185%
13	8.265	2012	2019	2039	rVV4	46952	103204	3.15%	0.514%
14	8.371	2039	2052	2073	rVB6	11840	33110	1.01%	0.165%
15	8.551	2091	2108	2129	rBV	443302	1008009	30.76%	5.018%
16	8.789	2168	2182	2190	rBV3	22934	49929	1.52%	0.249%
17	9.046	2235	2262	2276	rBV5	171444	546549	16.68%	2.721%
18	9.239	2311	2322	2335	rVB3	39426	79007	2.41%	0.393%
19	9.333	2335	2351	2361	rBV4	16191	40381	1.23%	0.201%
20	9.400	2362	2372	2376	rBV4	36527	60948	1.86%	0.303%
21	9.487	2390	2399	2413	rVB3	47452	93547	2.85%	0.466%
22	9.583	2418	2429	2435	rBV2	73450	139822	4.27%	0.696%
23	9.821	2488	2503	2511	rBV2	26679	70035	2.14%	0.349%
24	9.937	2524	2539	2548	rVB5	49103	112947	3.45%	0.562%
25	10.059	2563	2577	2605	rBV	662460	1299079	39.64%	6.467%
26	10.230	2622	2630	2635	rBV4	30300	52482	1.60%	0.261%
27	10.403	2676	2684	2693	rVV5	22284	39050	1.19%	0.194%
28	10.490	2697	2711	2726	rVV4	118926	252657	7.71%	1.258%
29	10.596	2728	2744	2756	rVB10	22390	63757	1.95%	0.317%
30	10.763	2779	2796	2806	rBV3	52390	112969	3.45%	0.562%
31	10.889	2828	2835	2843	rVV3	28911	42929	1.31%	0.214%
32	11.291	2946	2960	2973	rVB4	62549	126692	3.87%	0.631%
33	11.381	2974	2988	3007	rBV	605220	1088201	33.21%	5.417%
34	11.603	3039	3057	3076	rBV2	74590	176140	5.37%	0.877%

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35	11.921	3148	3156	3170	rVB2	18701	34496	1.05%	0.172%
36	12.223	3241	3250	3260	rBV	55244	88700	2.71%	0.442%
37	12.377	3285	3298	3315	rBV	438586	750651	22.91%	3.737%
38	12.567	3344	3357	3371	rVB	15165	37040	1.13%	0.184%
39	12.660	3371	3386	3395	rVV2	24015	46969	1.43%	0.234%
40	12.866	3438	3450	3464	rBV	235248	380469	11.61%	1.894%
41	13.139	3519	3535	3548	rBV3	37669	87609	2.67%	0.436%
42	13.320	3579	3591	3614	rBV	503591	901022	27.49%	4.486%
43	13.519	3631	3653	3663	rBV2	1674856	3277048	100.00%	16.314%
44	13.567	3663	3668	3684	rVB	107687	171152	5.22%	0.852%
45	13.648	3684	3693	3702	rVB2	32111	48195	1.47%	0.240%
46	13.715	3702	3714	3725	rVB3	50413	90270	2.75%	0.449%
47	13.786	3725	3736	3748	rVB2	44911	79182	2.42%	0.394%
48	13.863	3748	3760	3770	rBV	402736	629992	19.22%	3.136%
49	13.921	3770	3778	3784	rVV	135126	212608	6.49%	1.058%
50	13.969	3784	3793	3807	rVV	805338	1303568	39.78%	6.490%
51	14.104	3825	3835	3844	rBV5	24457	38750	1.18%	0.193%
52	14.184	3845	3860	3873	rBV	267242	431761	13.18%	2.149%
53	14.265	3874	3885	3891	rVV2	19063	34652	1.06%	0.173%
54	14.451	3933	3943	3953	rVV2	286456	468318	14.29%	2.331%
55	14.567	3965	3979	3991	rBV2	471755	858466	26.20%	4.274%
56	14.631	3991	3999	4010	rVB4	46031	77677	2.37%	0.387%
57	14.715	4016	4025	4035	rVB3	32286	52403	1.60%	0.261%
58	14.821	4035	4058	4066	rBV4	83376	201583	6.15%	1.004%
59	14.927	4078	4091	4105	rVB2	88145	201405	6.15%	1.003%
60	15.384	4223	4233	4244	rBV2	25488	43688	1.33%	0.217%
61	15.545	4270	4283	4302	rBV4	12027	35865	1.09%	0.179%

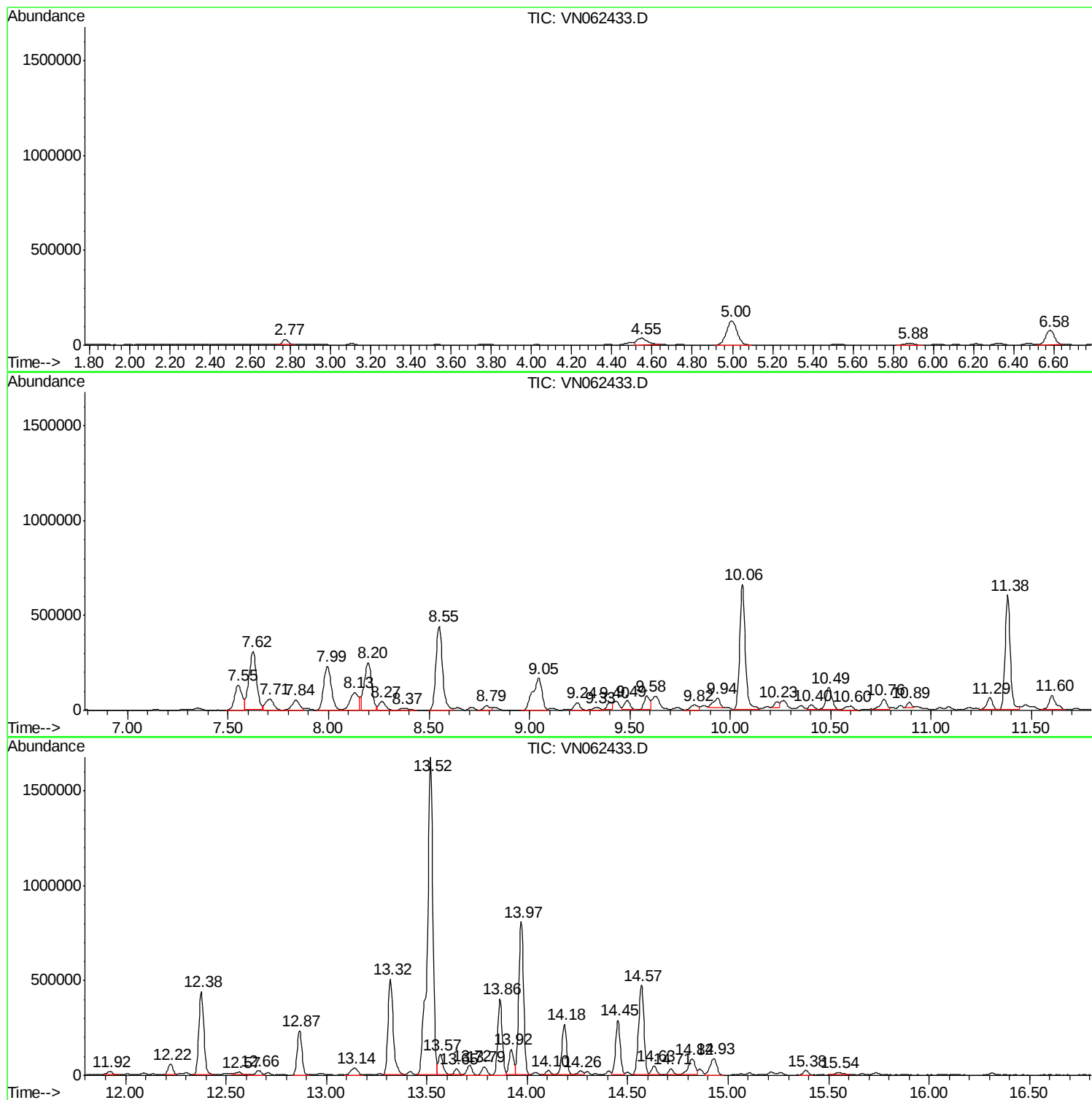
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TIC Integration Parameters: LSCINT.P



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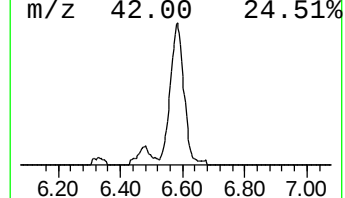
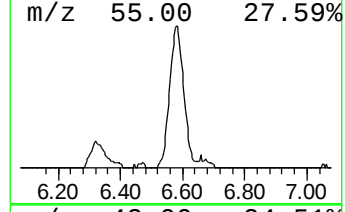
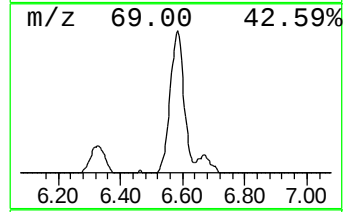
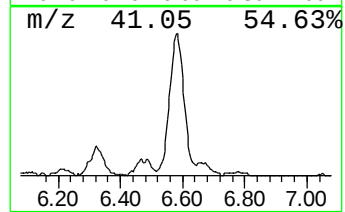
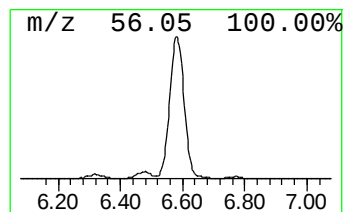
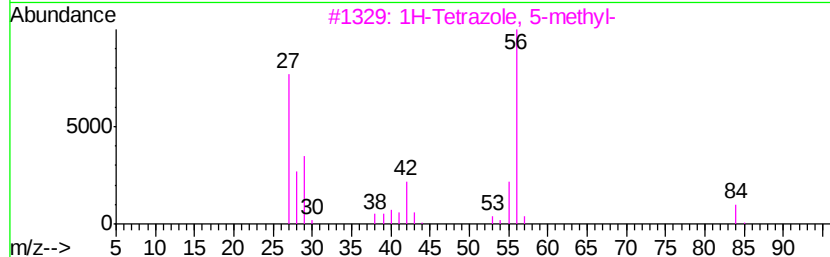
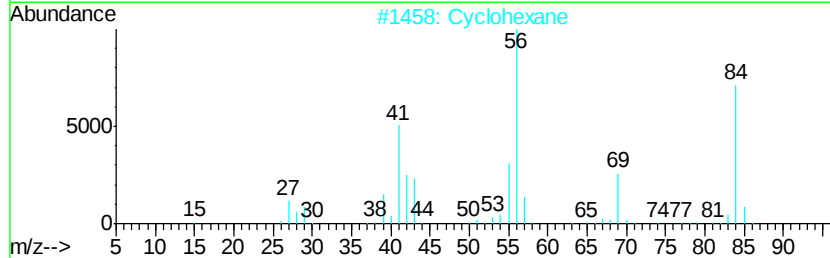
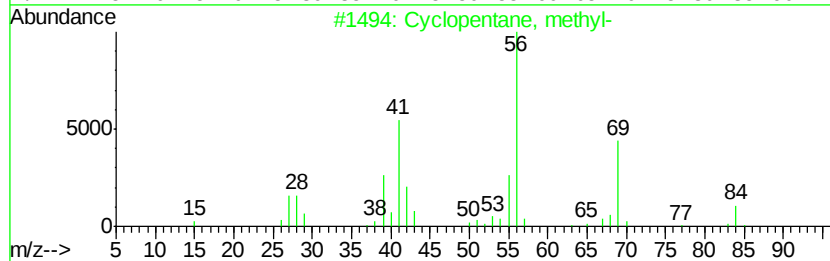
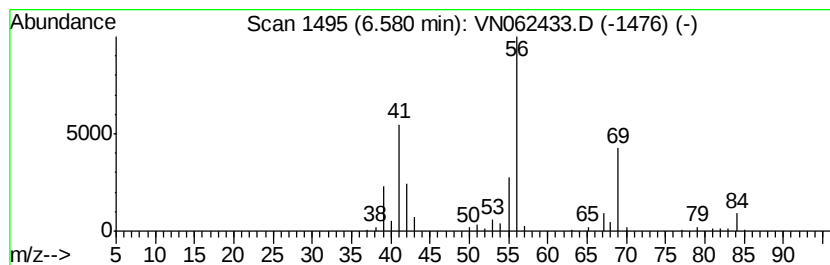
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	14.46 ug/l	233397	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		Cyclohexane	84	C6H12	000110-82-7	78
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50
4		Cyclopropane, propyl-	84	C6H12	002415-72-7	50
5		Cyclobutane	56	C4H8	000287-23-0	47



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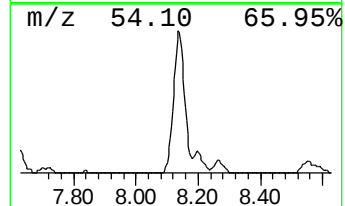
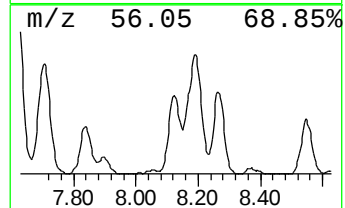
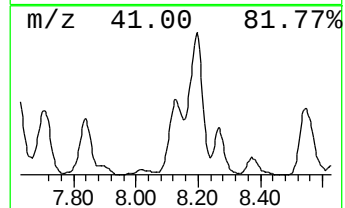
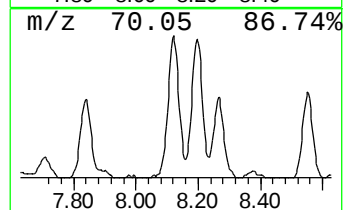
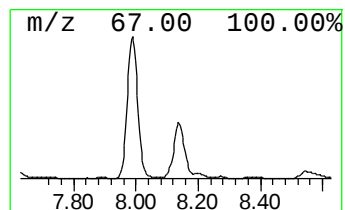
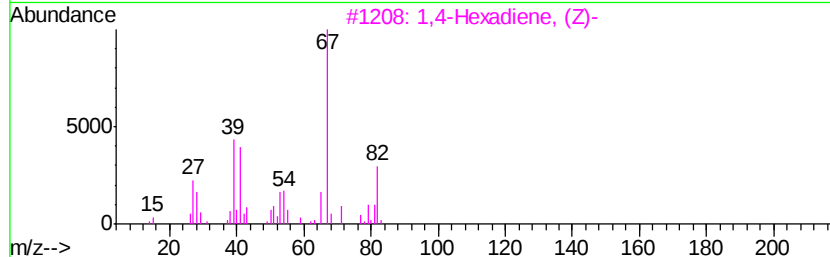
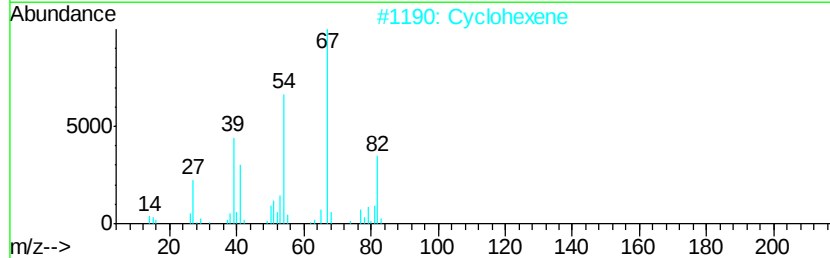
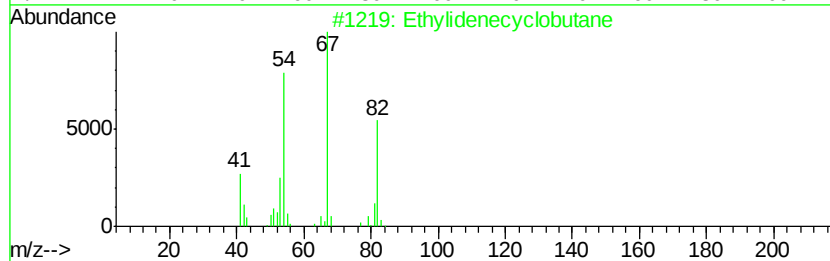
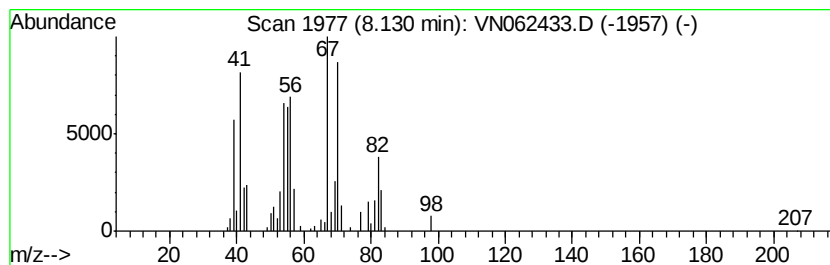
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethylidenecyclobutane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.13	14.00 ug/l	282250	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethylidenecvclobutane	82	C6H10	001528-21-8	55
2		Cyclohexene	82	C6H10	000110-83-8	55
3		1,4-Hexadiene, (Z)-	82	C6H10	007318-67-4	50
4		trans-1,4-Hexadiene	82	C6H10	007319-00-8	45
5		1,4-Hexadiene	82	C6H10	000592-45-0	45



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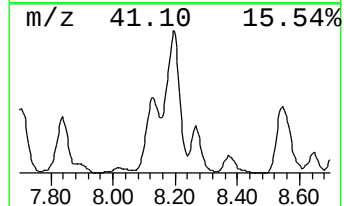
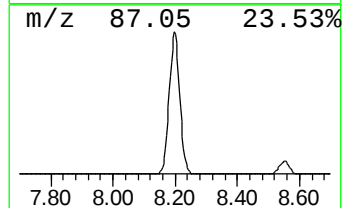
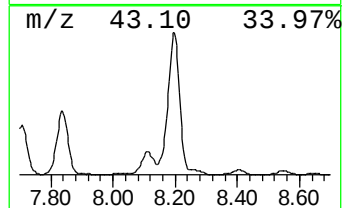
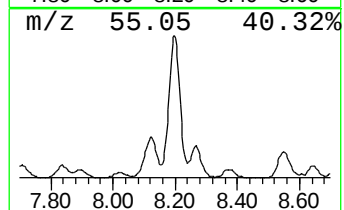
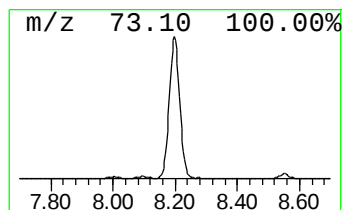
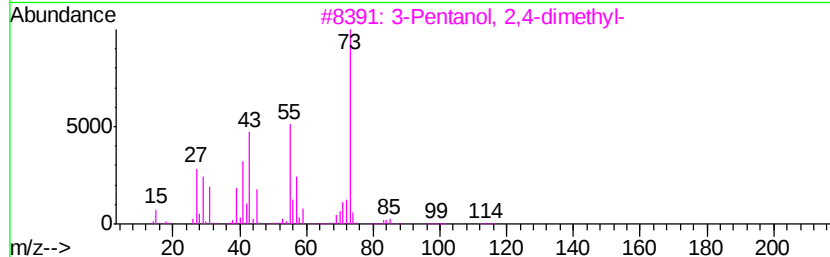
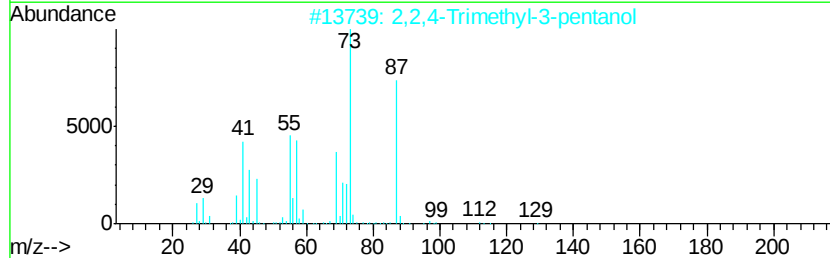
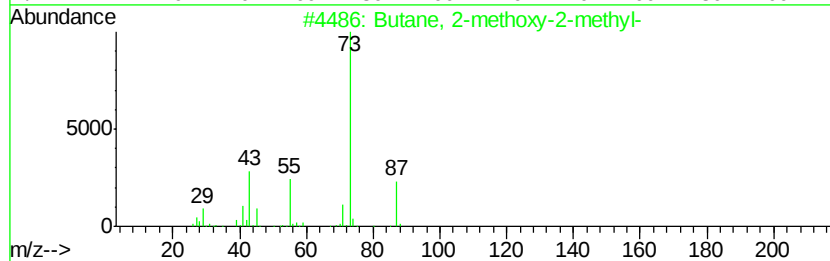
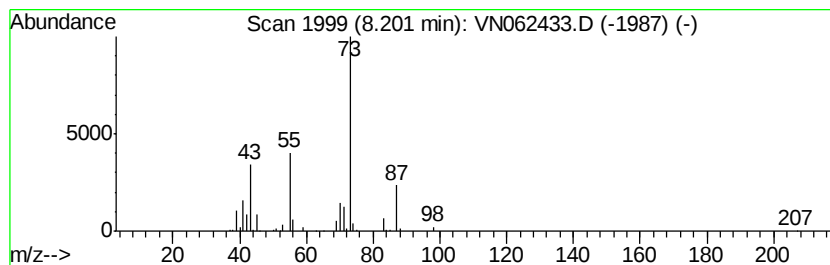
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	31.74 ug/l	639868	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	37
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
5		Silane, tetramethyl-	88	C4H12Si	000075-76-3	9



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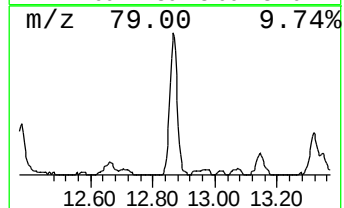
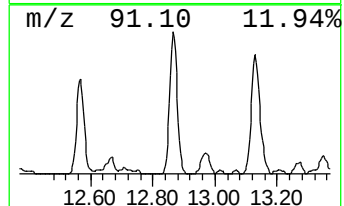
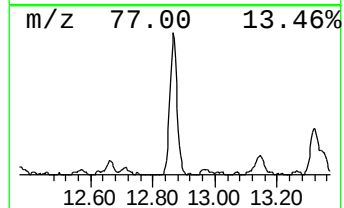
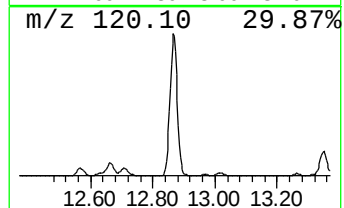
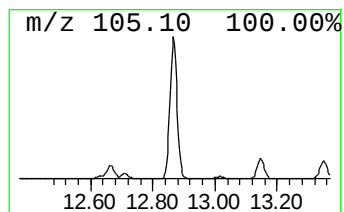
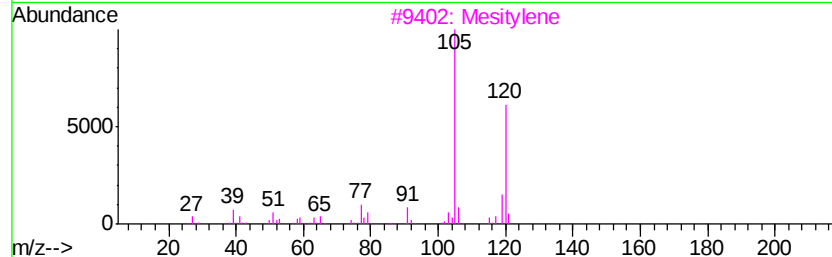
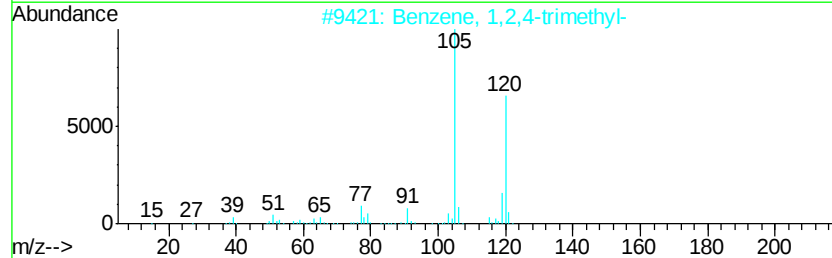
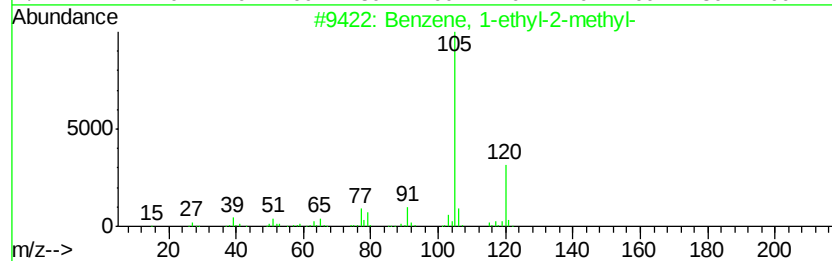
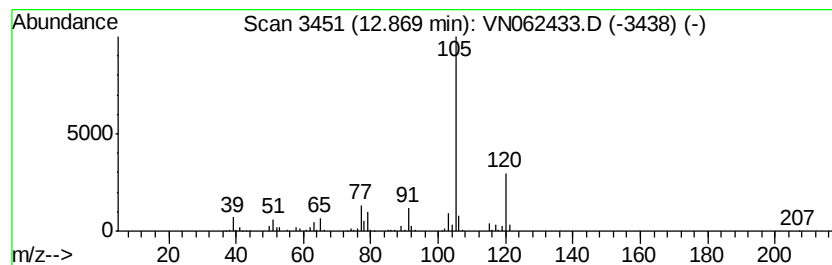
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Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	21.11 ug/l	380469	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	95
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
3		Mesitylene	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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MW-5

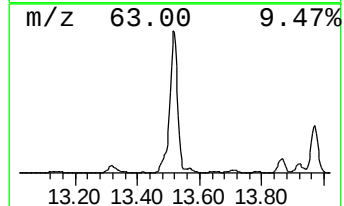
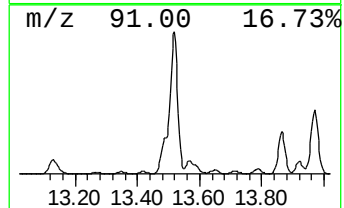
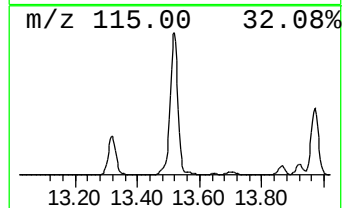
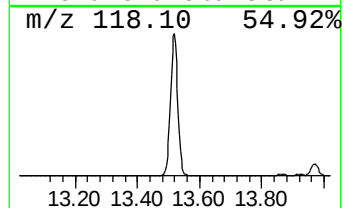
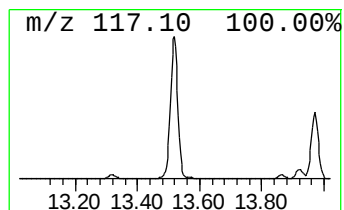
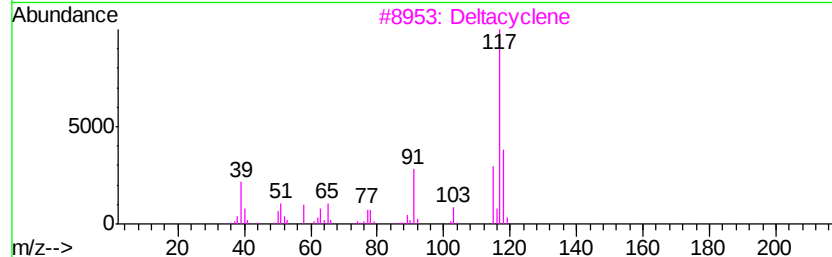
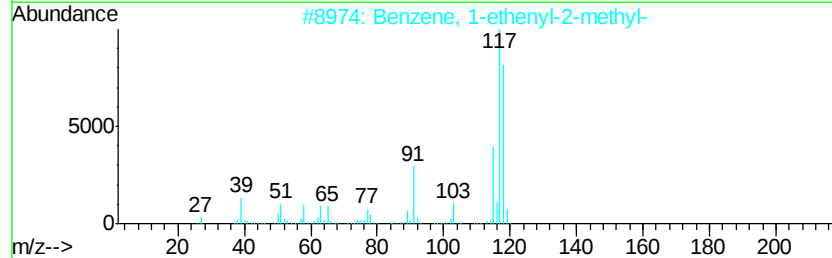
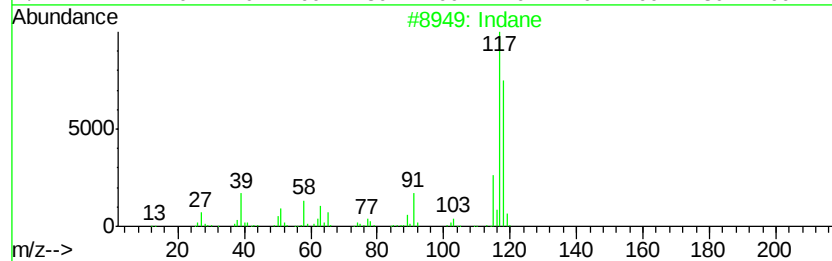
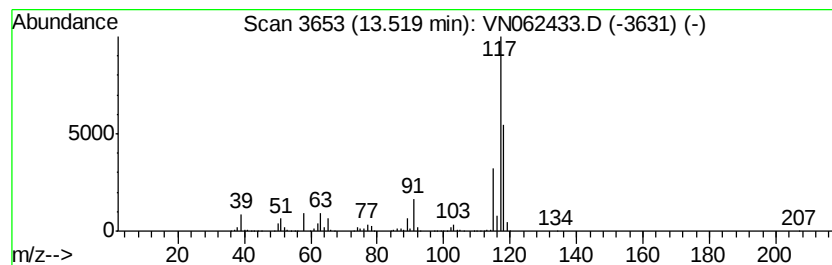
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Quant Title : SW846 8260

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

Peak Number 5 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	181.85 ug/l	3277050	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indane	118	C9H10	000496-11-7	96
2		Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	78
3		Deltacyclene	118	C9H10	007785-10-6	68
4		Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
5		Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN070820\
Data File : VN062433.D
Acq On : 9 Jul 2020 2:47
Operator : JC/MD
Sample : L3215-10
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

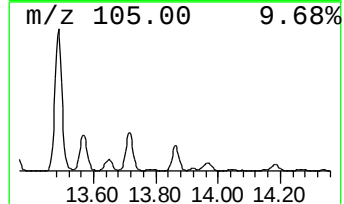
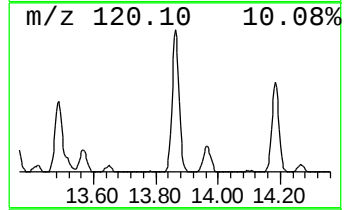
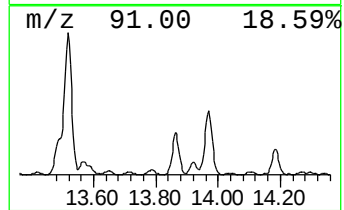
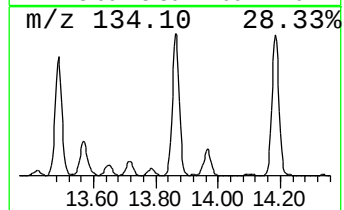
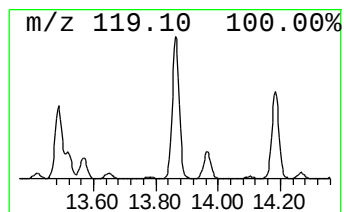
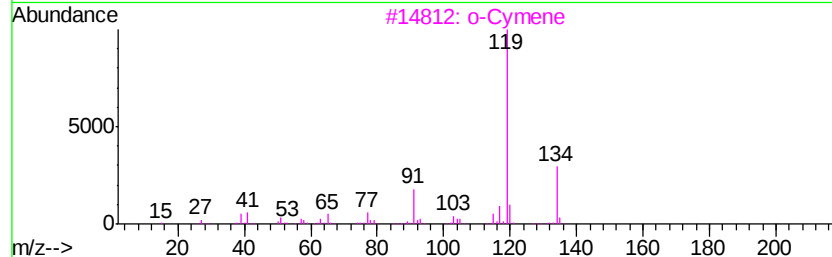
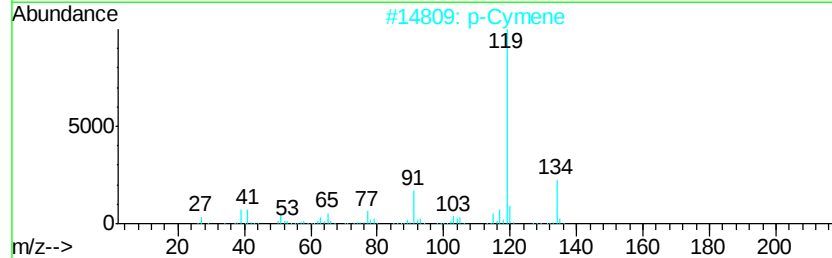
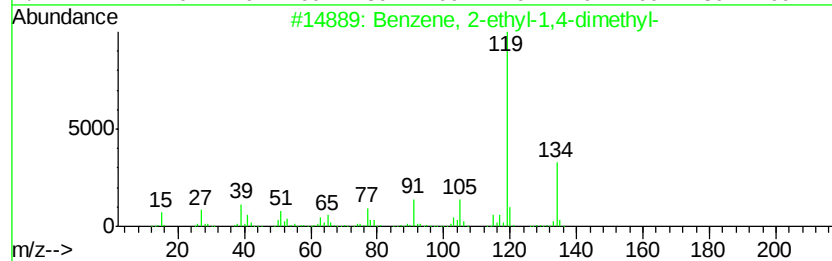
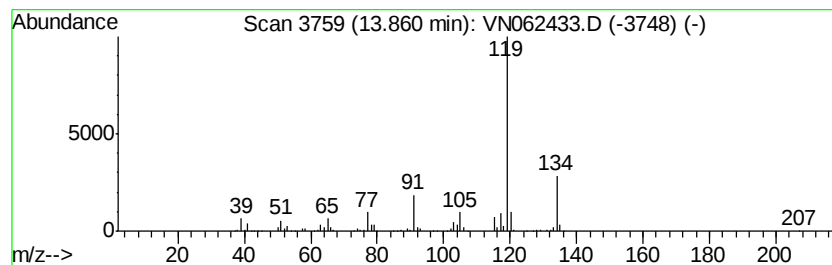
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	34.96 ug/l	629992	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2		p-Cymene	134	C10H14	000099-87-6	95
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN070820\
Data File : VN062433.D
Acq On : 9 Jul 2020 2:47
Operator : JC/MD
Sample : L3215-10
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

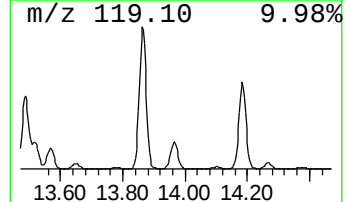
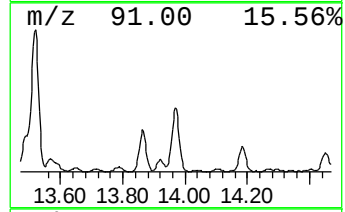
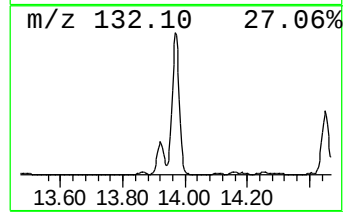
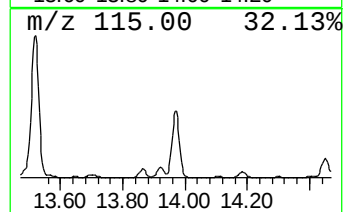
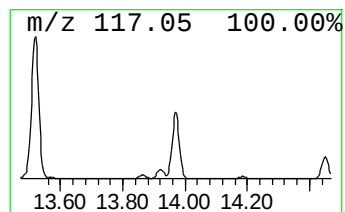
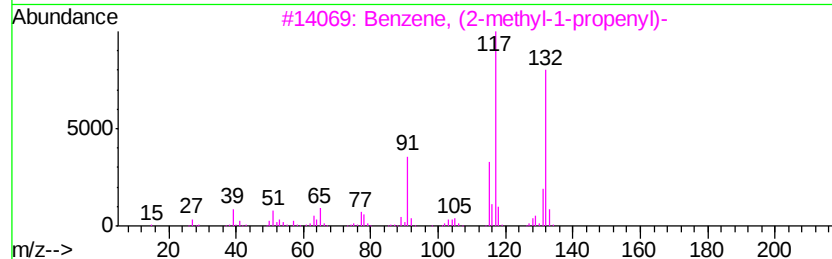
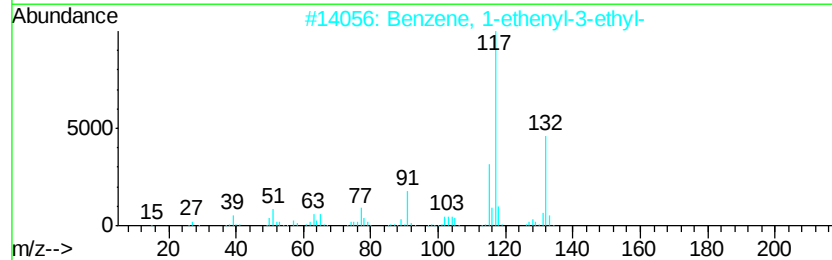
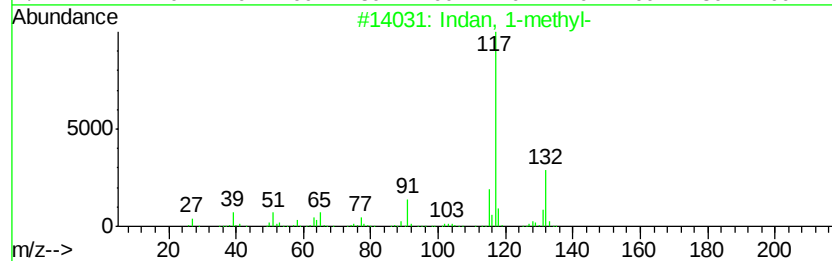
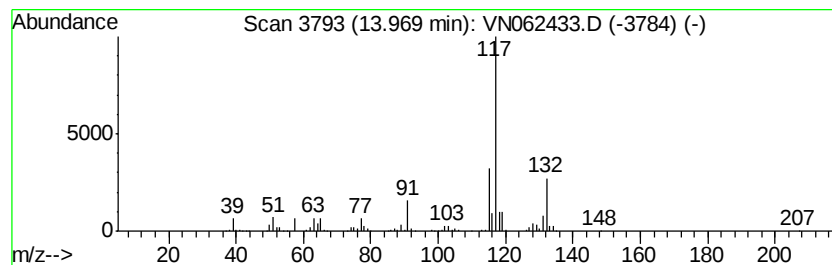
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Quant Title : SW846 8260

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

Peak Number 7 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	72.34 ug/l	1303570	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-		132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	90
3	Benzene, (2-methyl-1-propenyl)-		132	C10H12	000768-49-0	87
4	Benzene, 1-methyl-2-(2-propenyl)-		132	C10H12	001587-04-8	81
5	Benzene, (1-methyl-1-propenyl)-, ...		132	C10H12	000768-00-3	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN070820\
Data File : VN062433.D
Acq On : 9 Jul 2020 2:47
Operator : JC/MD
Sample : L3215-10
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
MW-5

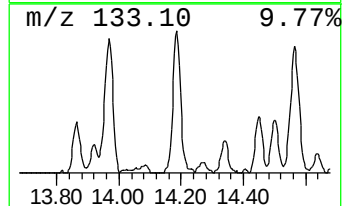
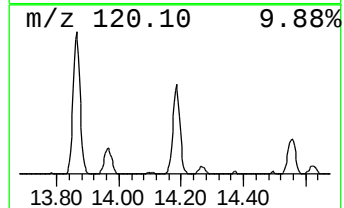
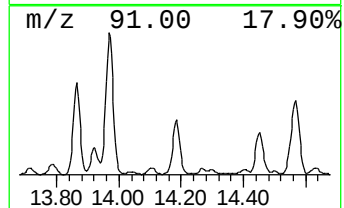
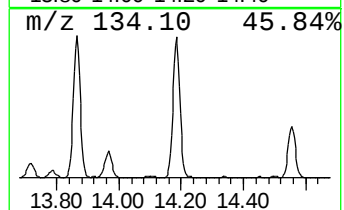
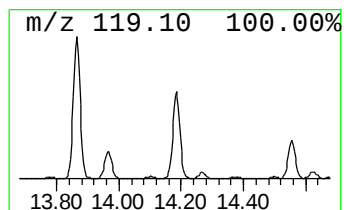
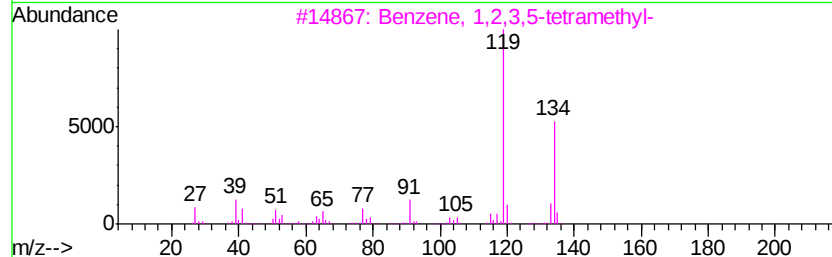
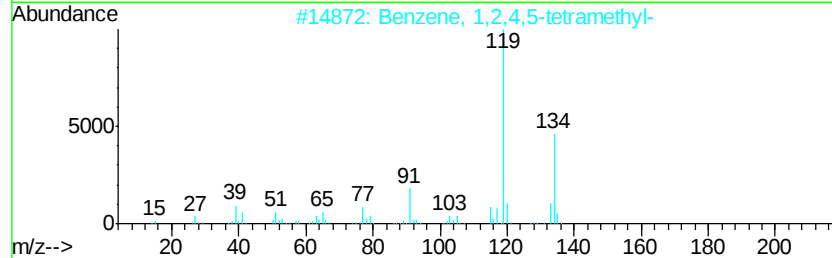
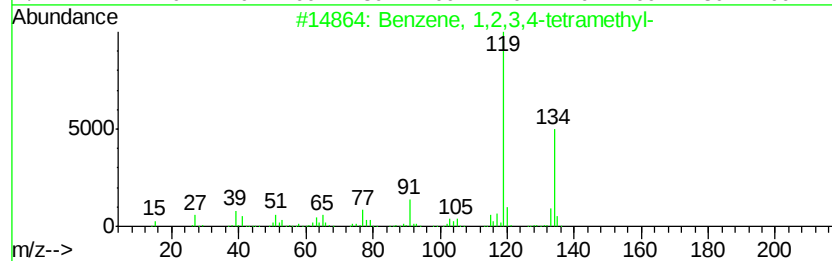
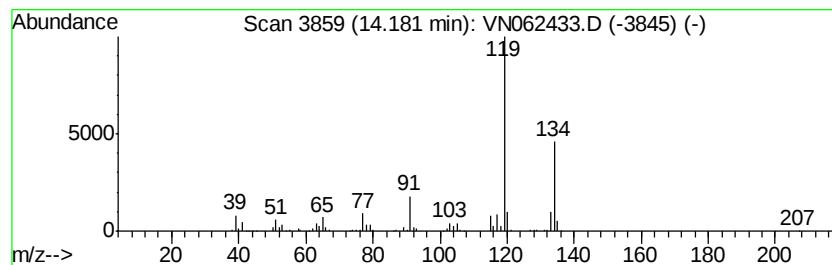
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	23.96 ug/l	431761	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	95
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:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN070820\
Data File : VN062433.D
Acq On : 9 Jul 2020 2:47
Operator : JC/MD
Sample : L3215-10
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

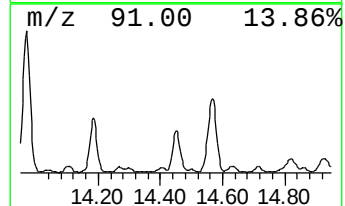
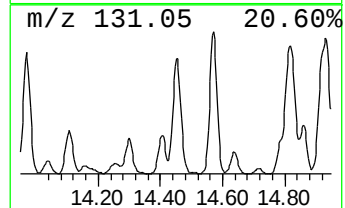
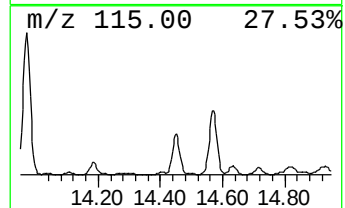
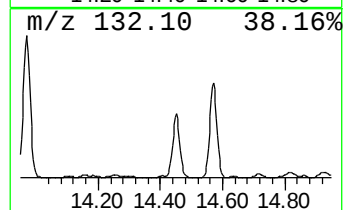
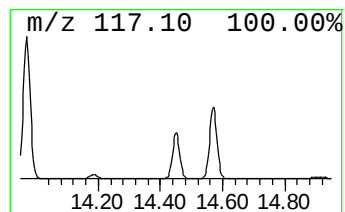
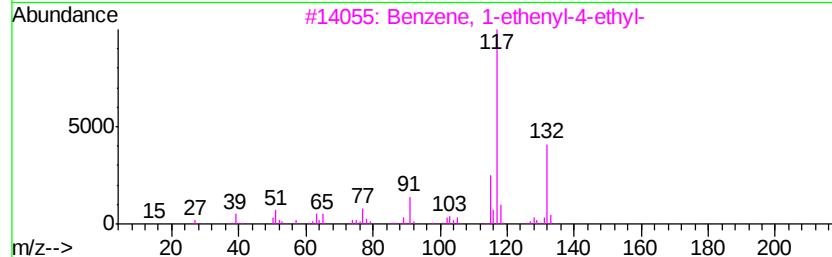
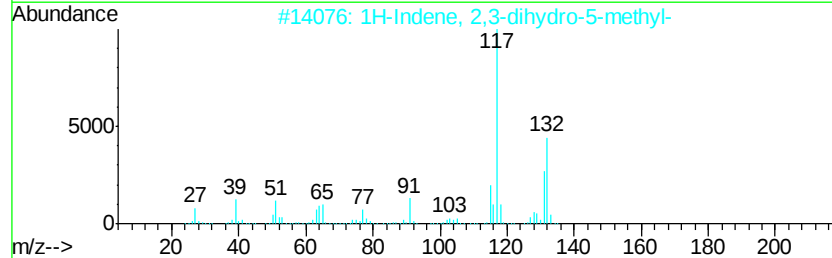
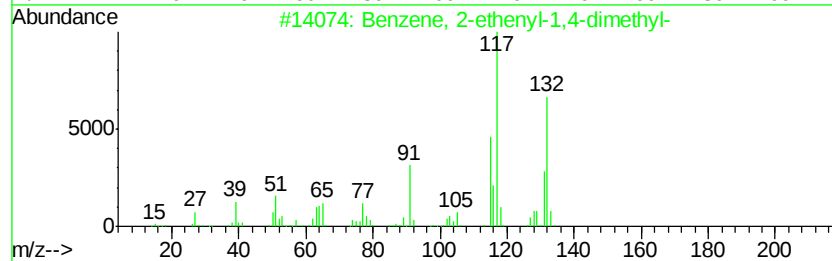
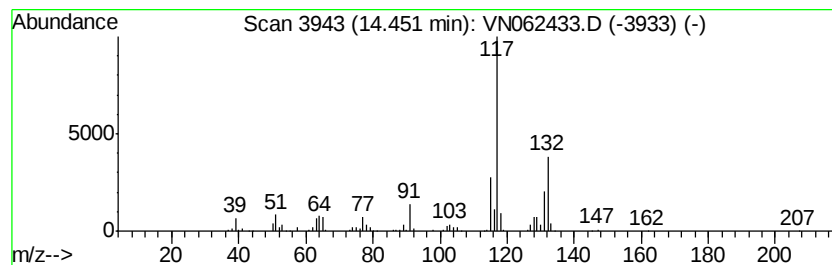
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	25.99 ug/l	468318	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		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062433.D
Acq On : 9 Jul 2020 2:47
Operator : JC/MD
Sample : L3215-10
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.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
Cyclopentane, met...	6.58	14.5	ug/l	233397	1	7.63	807284	50.0
Ethylidenecyclobu...	8.13	14.0	ug/l	282250	2	8.55	1008010	50.0
Butane, 2-methoxy...	8.20	31.7	ug/l	639868	2	8.55	1008010	50.0
Benzene, 1-ethyl-...	12.87	21.1	ug/l	380469	4	13.32	901022	50.0
Indane	13.52	181.8	ug/l	3277050	4	13.32	901022	50.0
Benzene, 2-ethyl-...	13.86	35.0	ug/l	629992	4	13.32	901022	50.0
Indan, 1-methyl-	13.97	72.3	ug/l	1303570	4	13.32	901022	50.0
Benzene, 1,2,3,4-...	14.18	24.0	ug/l	431761	4	13.32	901022	50.0
Benzene, 2-etheny...	14.45	26.0	ug/l	468318	4	13.32	901022	50.0