

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111620\
 Data File : VN064666.D
 Acq On : 16 Nov 2020 17:36
 Operator : JC/MD
 Sample : L4772-05
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-3R

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\82N110420W.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	7.627	1819	1831	1849	rVB	340667	823213	1.42%	0.419%
2	7.987	1926	1943	1965	rBV	252821	625518	1.08%	0.318%
3	8.550	2103	2118	2143	rVB	568995	1196368	2.06%	0.609%
4	10.058	2571	2587	2614	rBV	1035279	1992271	3.43%	1.014%
5	11.380	2985	2998	3015	rBV	972489	1726157	2.98%	0.878%
6	11.482	3015	3030	3047	rBV	1775110	3244190	5.59%	1.651%
7	11.598	3047	3066	3084	rBV	3448072	7217453	12.44%	3.672%
8	11.920	3144	3166	3182	rBV4	584815	1919039	3.31%	0.976%
9	12.093	3212	3220	3227	rBV	460364	674463	1.16%	0.343%
10	12.142	3227	3235	3249	rVV5	412097	777501	1.34%	0.396%
11	12.222	3249	3260	3272	rVB	5085535	8055007	13.88%	4.099%
12	12.376	3296	3308	3319	rBV	876259	1519823	2.62%	0.773%
13	12.563	3350	3366	3378	rBV	7182714	12219006	21.06%	6.217%
14	12.630	3378	3387	3391	rVV	1156844	1801213	3.10%	0.916%
15	12.662	3391	3397	3403	rVV	1951234	3006038	5.18%	1.530%
16	12.707	3404	3411	3421	rVB	2534078	3904573	6.73%	1.987%
17	12.865	3446	3460	3471	rBV	370109	766474	1.32%	0.390%
18	12.939	3473	3483	3486	rBV4	468282	738419	1.27%	0.376%
19	13.016	3495	3507	3531	rVB	34064498	58018901	100.00%	29.521%
20	13.145	3532	3547	3557	rVV2	2509426	5379475	9.27%	2.737%
21	13.209	3557	3567	3575	rVV	1436335	2363880	4.07%	1.203%
22	13.264	3575	3584	3592	rVV	1434952	2180244	3.76%	1.109%
23	13.347	3592	3610	3623	rVV	9651598	17311135	29.84%	8.808%
24	13.418	3623	3632	3642	rVV	952876	1503113	2.59%	0.765%
25	13.486	3643	3653	3656	rVV	2752294	3763877	6.49%	1.915%
26	13.518	3656	3663	3671	rVV	6851568	11415091	19.67%	5.808%
27	13.566	3671	3678	3693	rVV3	2123642	5330441	9.19%	2.712%
28	13.646	3693	3703	3714	rVB2	1180859	2077081	3.58%	1.057%
29	13.775	3732	3743	3748	rBV	2030645	3247210	5.60%	1.652%
30	13.804	3748	3752	3761	rVV	2141641	3153055	5.43%	1.604%
31	13.862	3761	3770	3780	rVV	4165504	6327310	10.91%	3.219%
32	13.920	3780	3788	3794	rVV	1296104	2008064	3.46%	1.022%
33	13.968	3794	3803	3814	rVB2	4175957	6922580	11.93%	3.522%
34	14.103	3835	3845	3854	rBV2	1050915	1659200	2.86%	0.844%

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Stop Thrs : 0 Peak Location: TOP

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35	14.183	3854	3870	3876	rVV	1268998	2252597	3.88%	1.146%
36	14.222	3876	3882	3891	rVB	1772917	2567069	4.42%	1.306%
37	14.450	3945	3953	3962	rBV3	462869	732699	1.26%	0.373%
38	14.556	3975	3986	4000	rBV4	2187753	4352191	7.50%	2.214%
39	14.714	4025	4035	4045	rBV	1129520	1761003	3.04%	0.896%

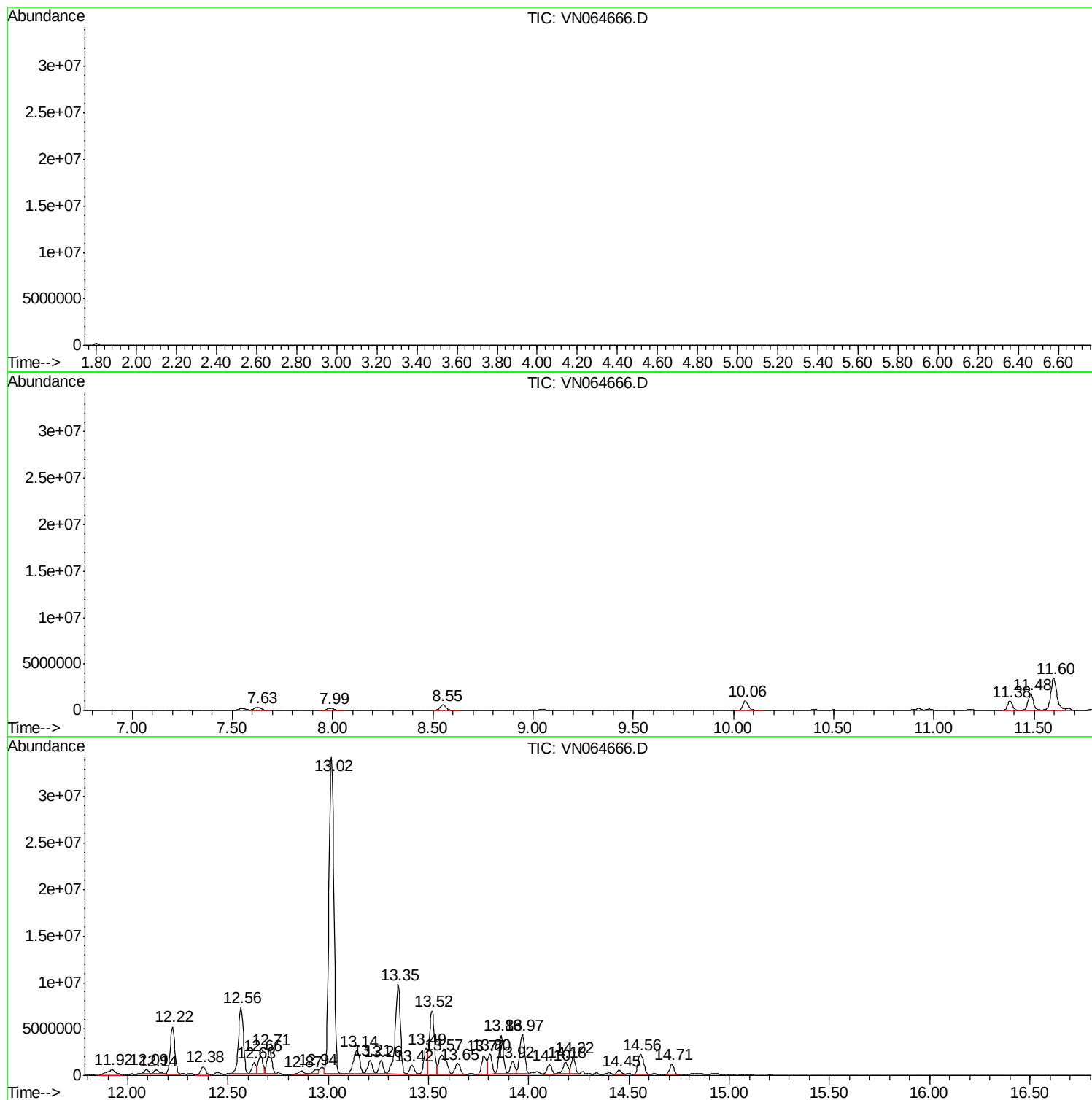
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Quant Title : SW846 8260

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



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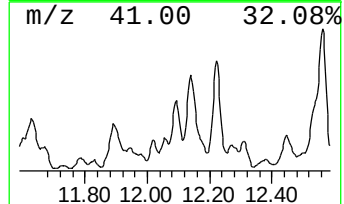
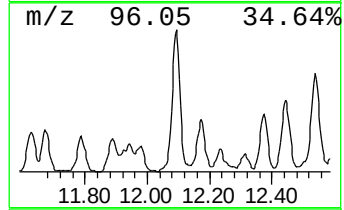
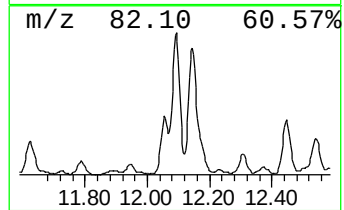
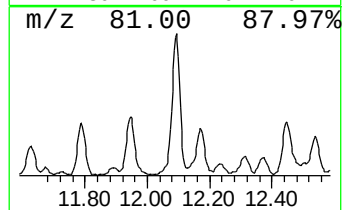
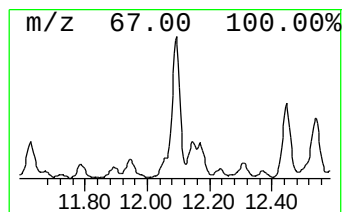
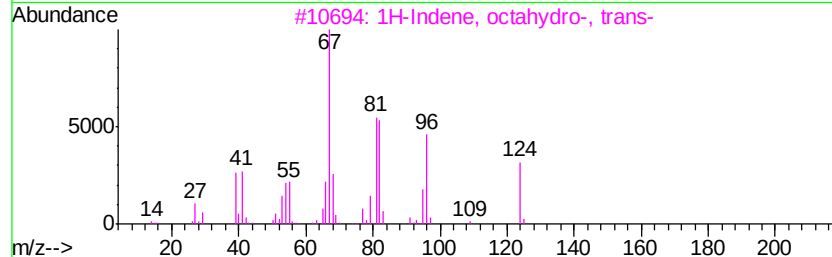
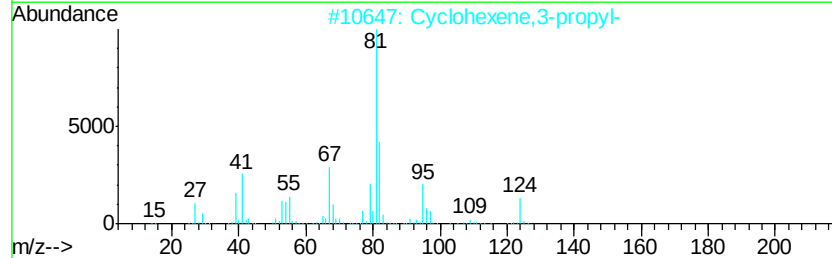
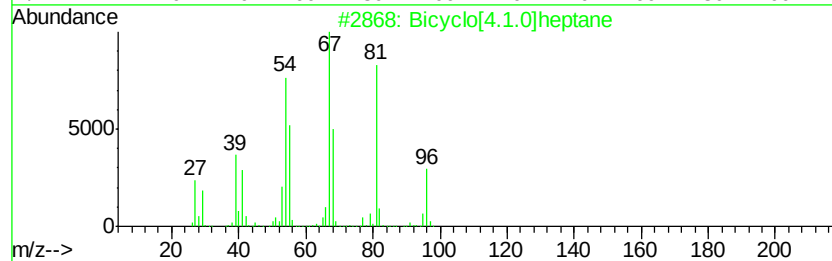
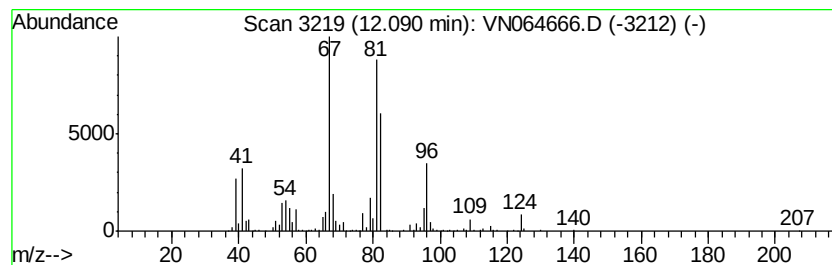
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Quant Title : SW846 8260

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

Peak Number 1 Bicyclo[4.1.0]heptane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	19.54 ug/l	674463	Chlorobenzene-d5	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[4.1.0]heptane	96	C7H12	000286-08-8	64
2		Cyclohexene,3-propyl-	124	C9H16	003983-06-0	49
3		1H-Indene, octahydro-, trans-	124	C9H16	003296-50-2	49
4		Isopropenylcyclopropane	82	C6H10	004663-22-3	43
5		2,4-Hexadiene, (E,E)-	82	C6H10	005194-51-4	43



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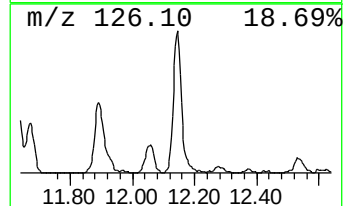
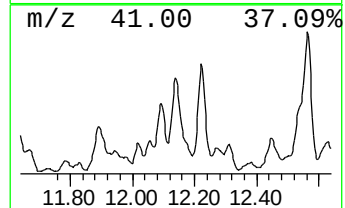
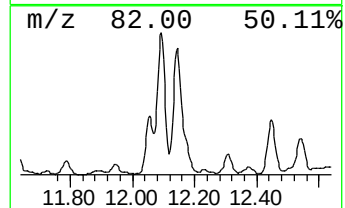
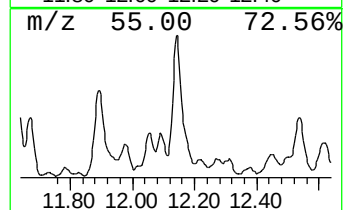
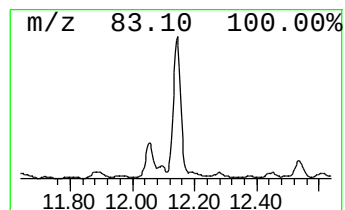
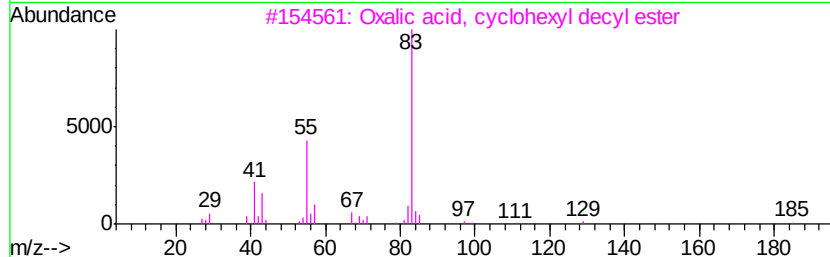
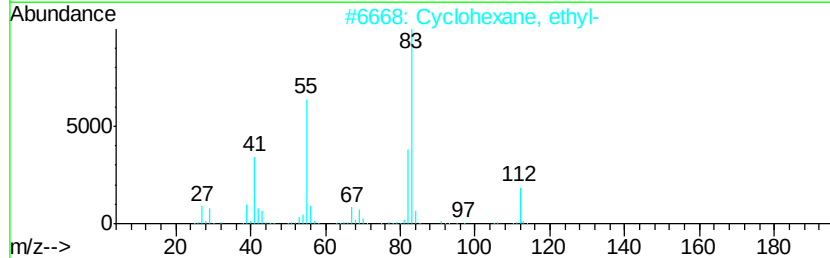
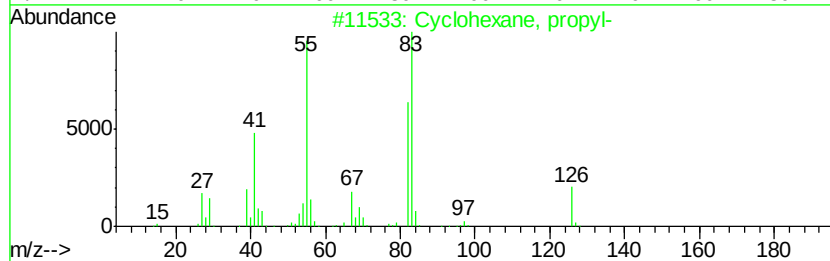
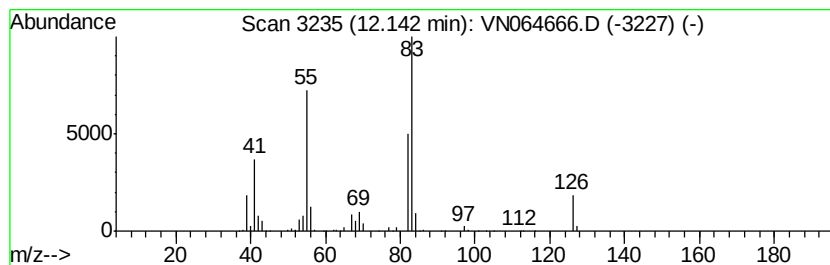
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	22.52 ug/l	777501	Chlorobenzene-d5	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, propyl-	126	C9H18	001678-92-8	94
2		Cyclohexane, ethyl-	112	C8H16	001678-91-7	78
3		Oxalic acid, cyclohexyl decyl ester	312	C18H32O4	1000309-31-2	59
4		Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	59
5		Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	53



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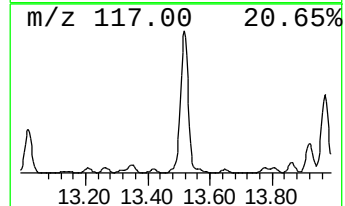
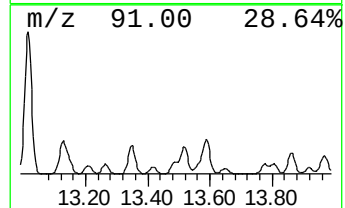
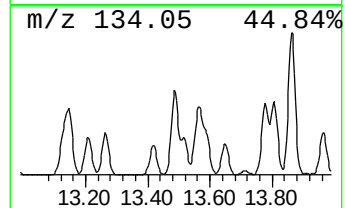
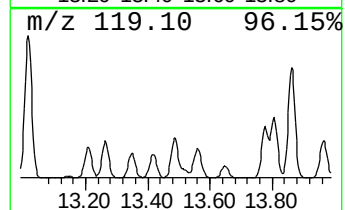
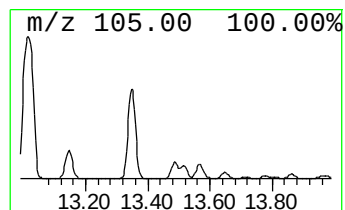
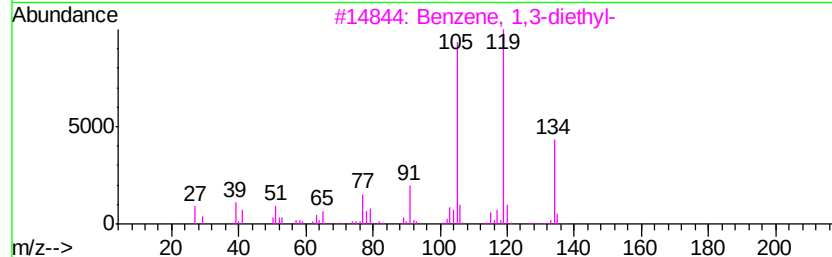
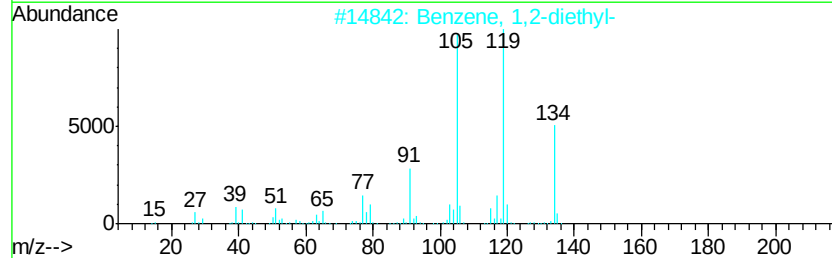
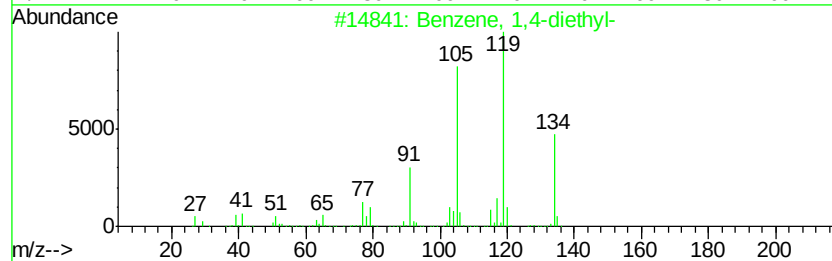
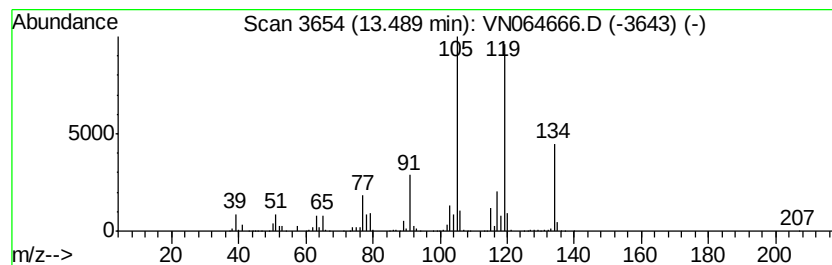
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Benzene, 1,4-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	10.87 ug/l	3763880	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
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-3,5-dimethyl-	134	C10H14	000934-74-7	90
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	83



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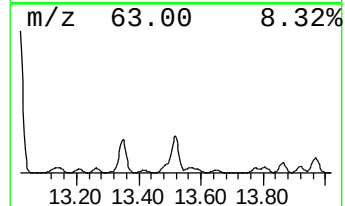
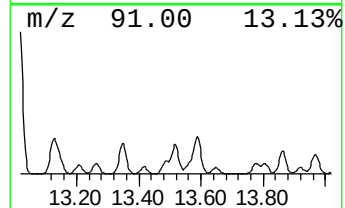
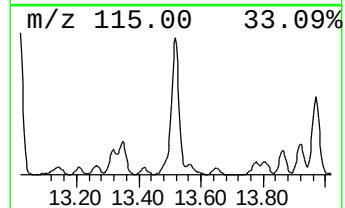
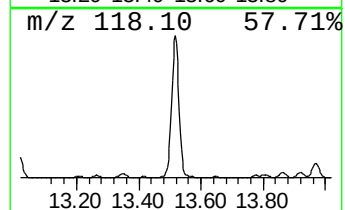
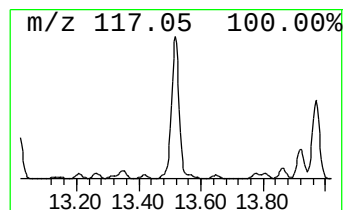
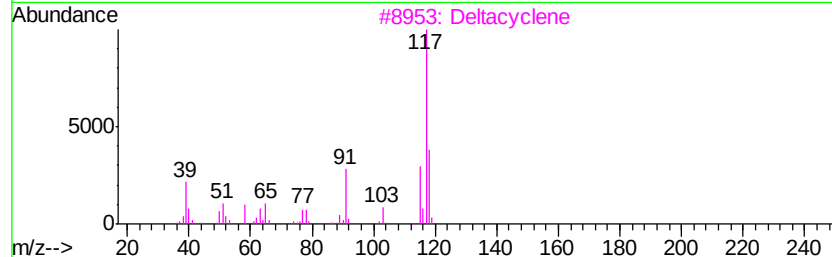
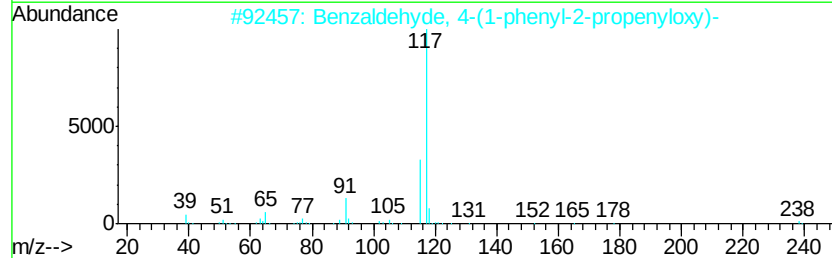
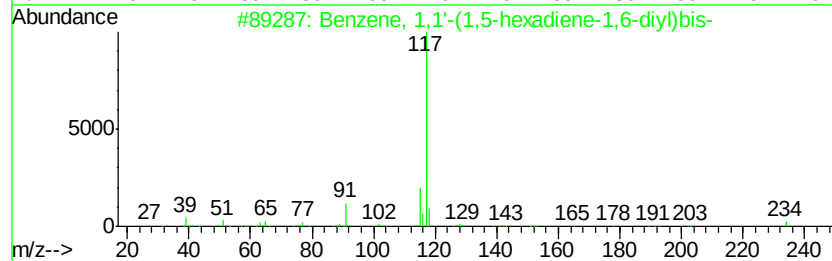
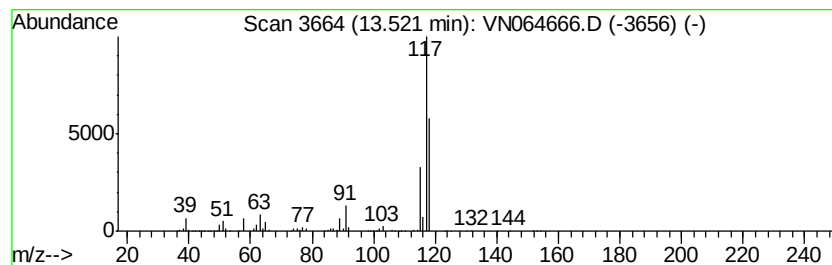
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Peak Number 4 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	32.97 ug/l	11415100	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2		Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	53
3		Deltacyclene	118	C9H10	007785-10-6	50
4		Indole	117	C8H7N	000120-72-9	49
5		5H-1-Pyridine	117	C8H7N	000270-91-7	43



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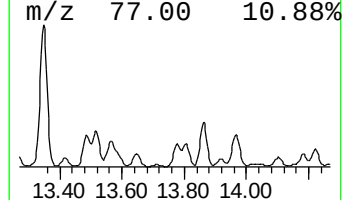
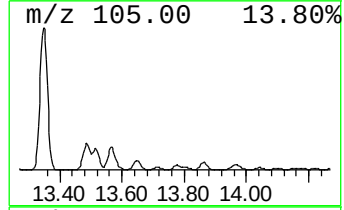
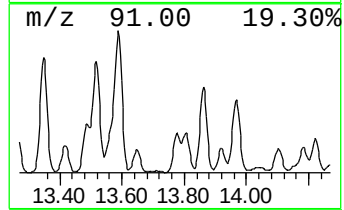
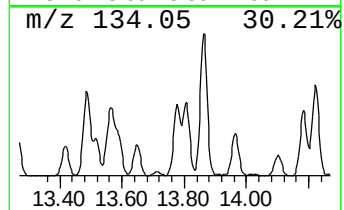
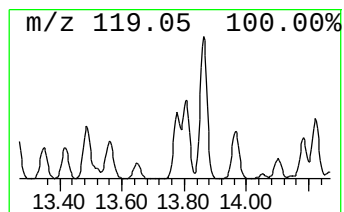
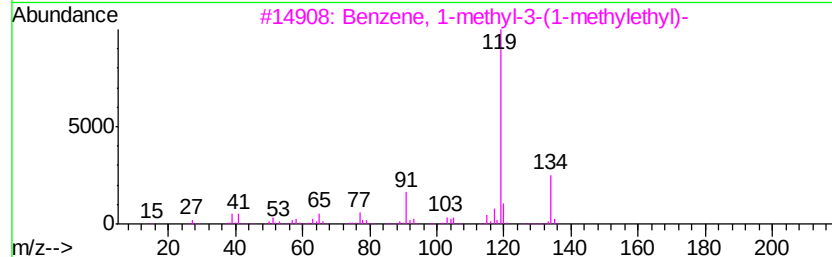
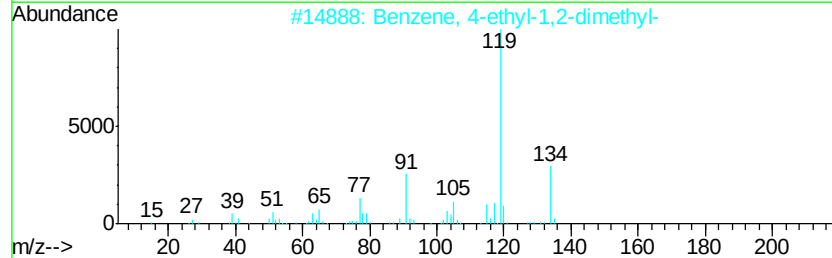
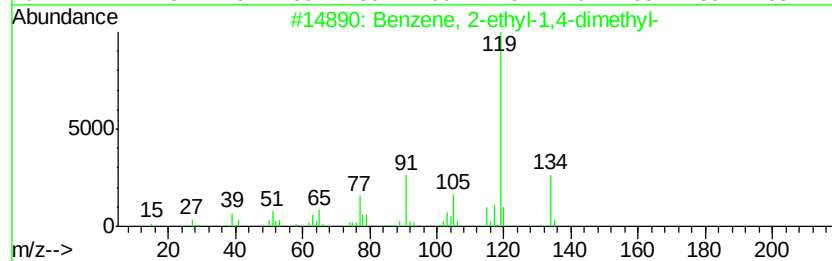
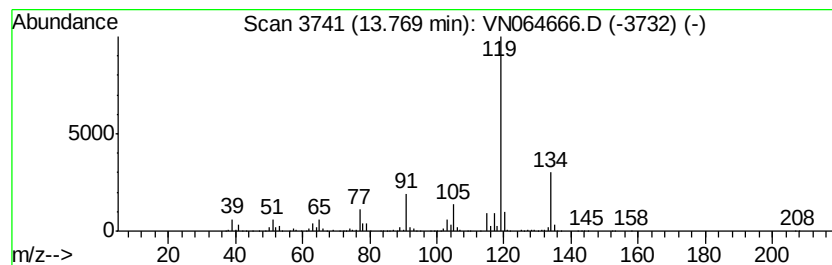
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Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	9.38 ug/l	3247210	1,4-Dichlorobenzene-d4	13.32

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



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064666.D
Acq On : 16 Nov 2020 17:36
Operator : JC/MD
Sample : L4772-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

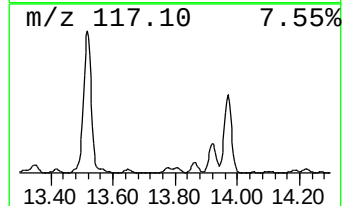
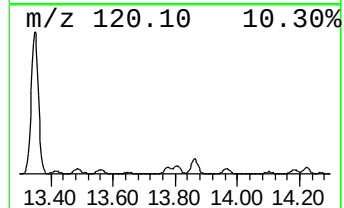
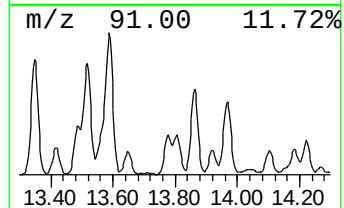
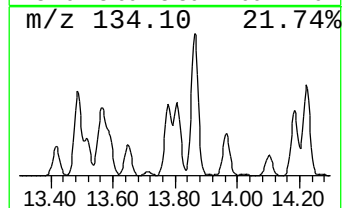
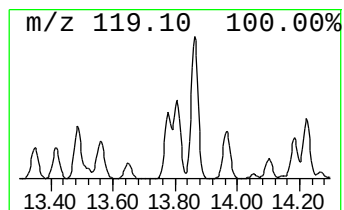
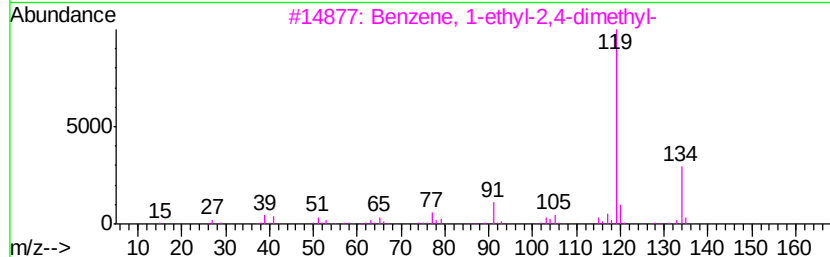
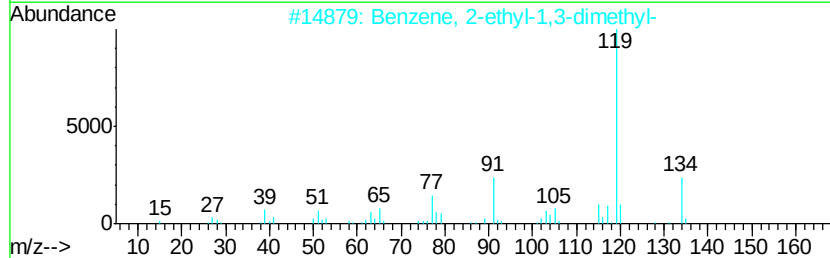
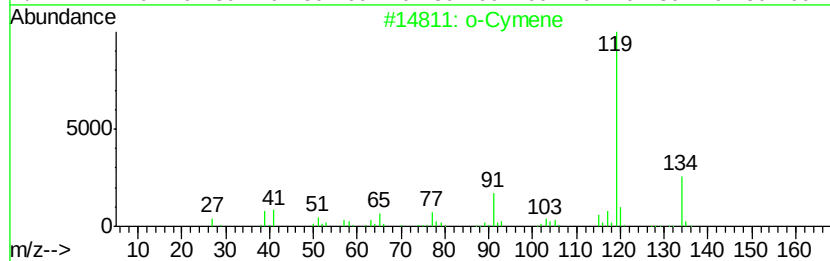
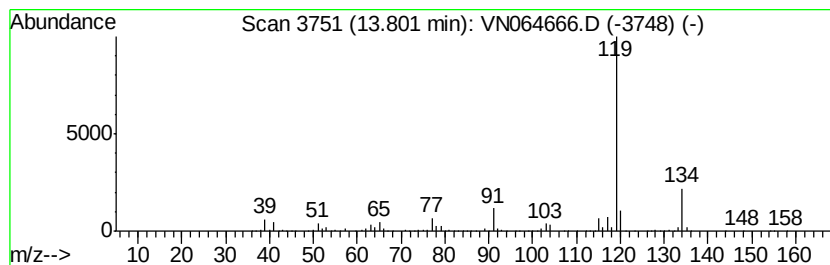
Instrument :
MSVOA_N
ClientSampleId :
MW-3R

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260

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

Peak Number 6 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.80	9.11 ug/l	3153060	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		o-Cymene	134 C10H14	000527-84-4 94
2		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 91
3		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 91
4		Benzene, 1-ethyl-3,5-dimethyl-	134 C10H14	000934-74-7 91
5		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064666.D
Acq On : 16 Nov 2020 17:36
Operator : JC/MD
Sample : L4772-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3R

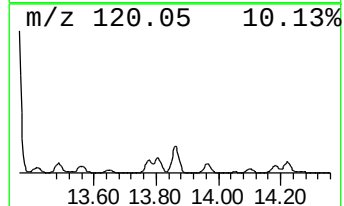
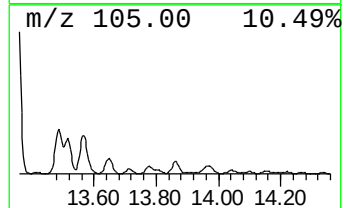
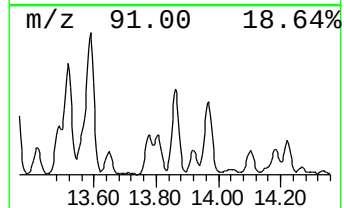
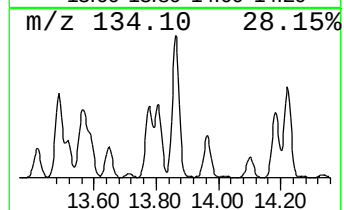
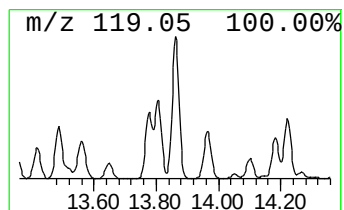
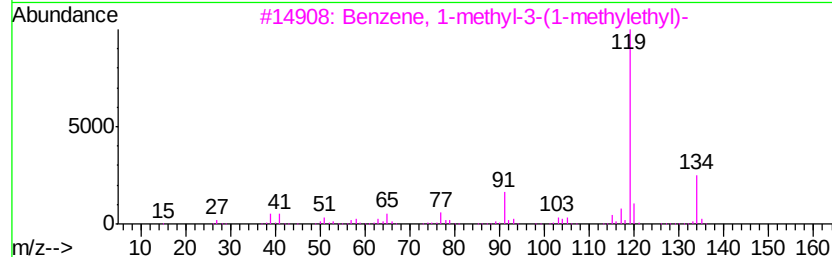
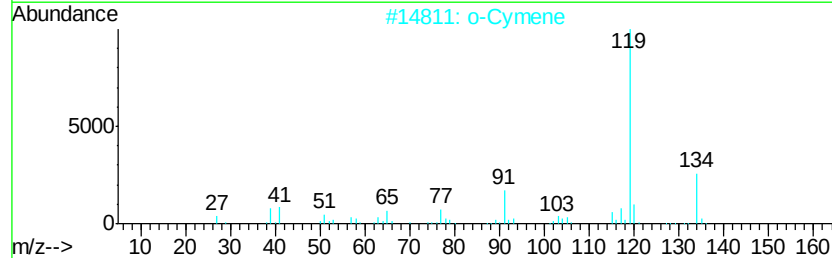
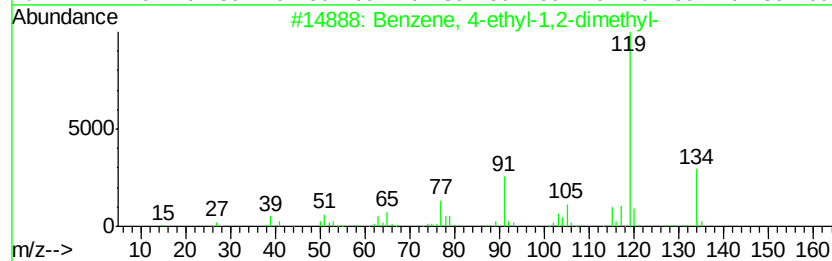
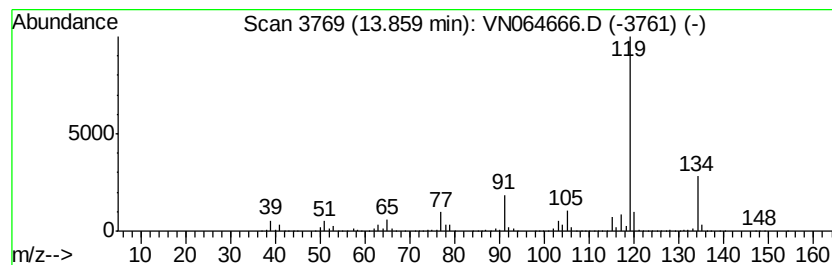
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	18.28 ug/l	6327310	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 1-methyl-3-(1-methylethyl-)	134	C10H14	000535-77-3	95
4		p-Cymene	134	C10H14	000099-87-6	95
5		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064666.D
Acq On : 16 Nov 2020 17:36
Operator : JC/MD
Sample : L4772-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3R

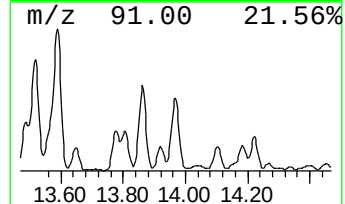
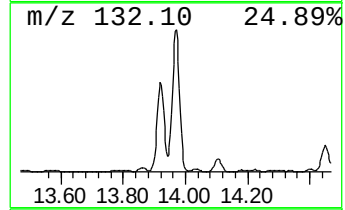
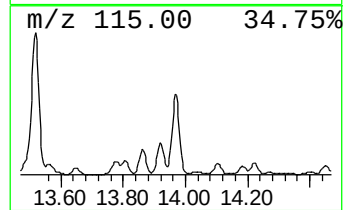
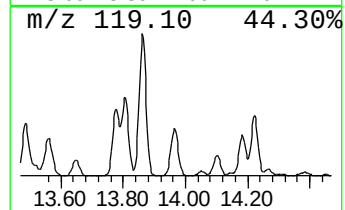
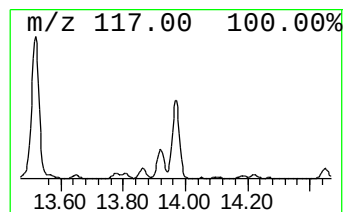
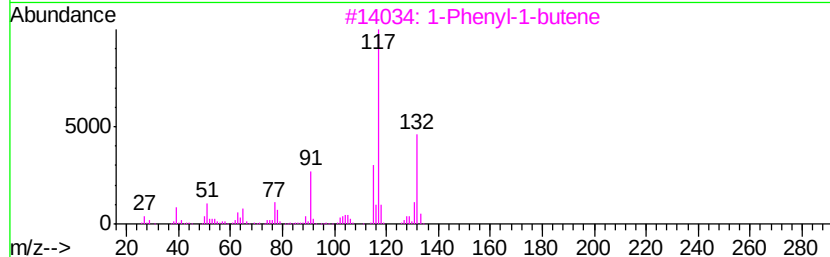
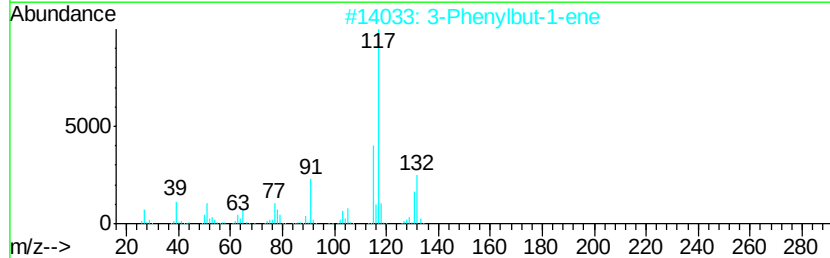
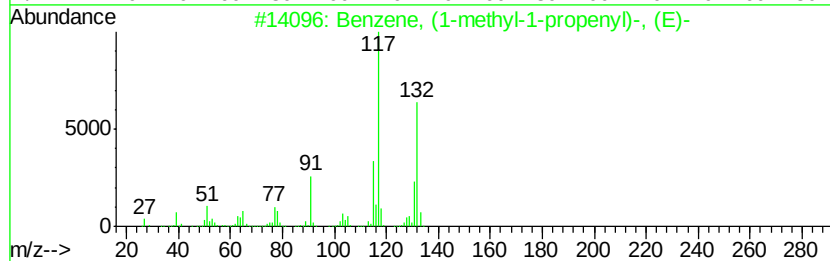
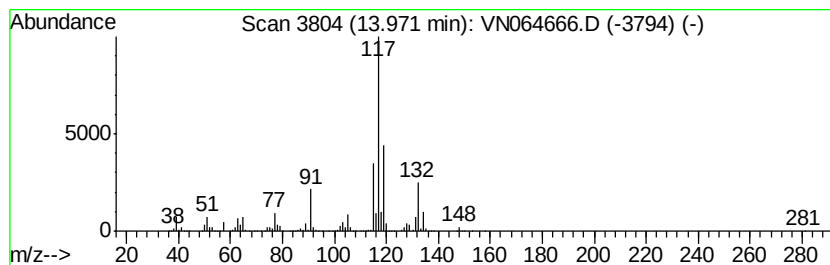
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (1-methyl-1-propen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	19.99 ug/l	6922580	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4		Indan, 1-methyl-	132	C10H12	000767-58-8	64
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064666.D
Acq On : 16 Nov 2020 17:36
Operator : JC/MD
Sample : L4772-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3R

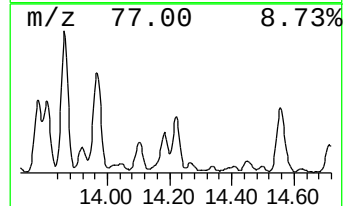
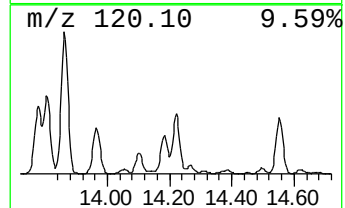
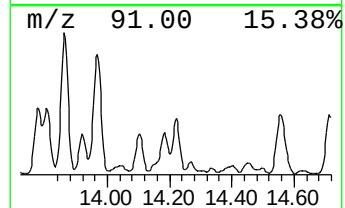
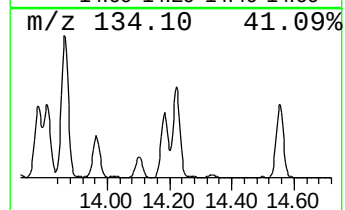
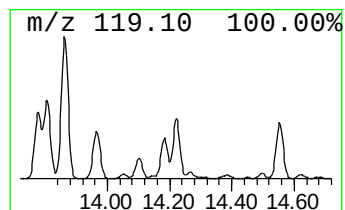
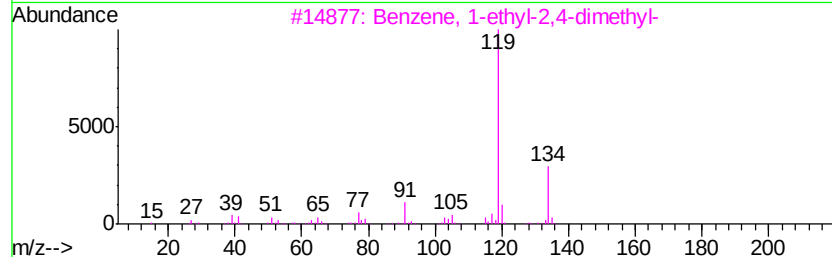
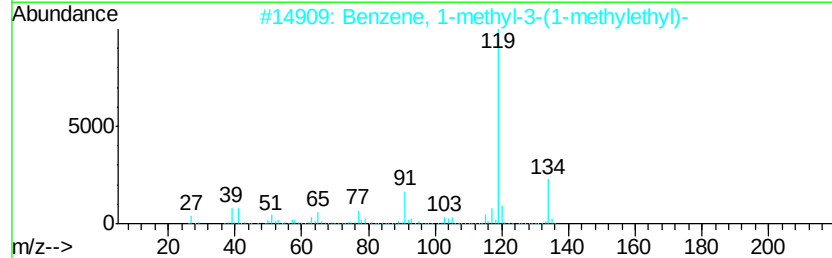
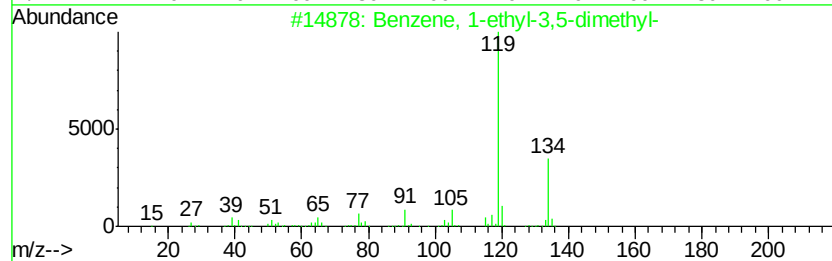
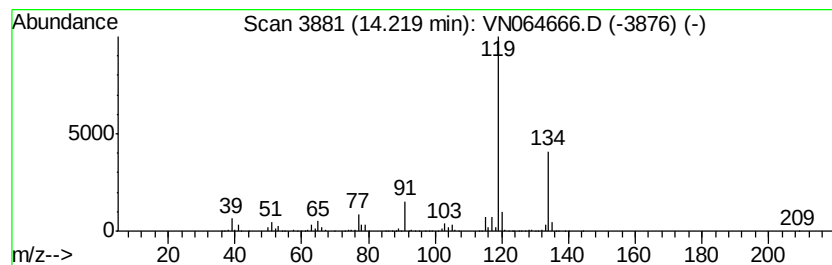
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	7.41 ug/l	2567070	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5		o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064666.D
Acq On : 16 Nov 2020 17:36
Operator : JC/MD
Sample : L4772-05
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3R

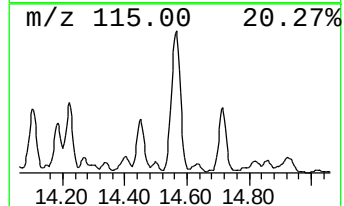
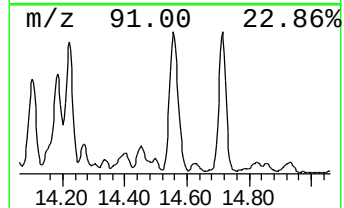
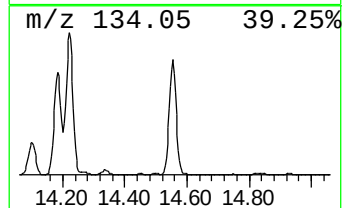
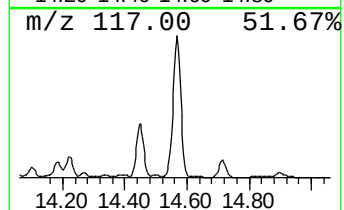
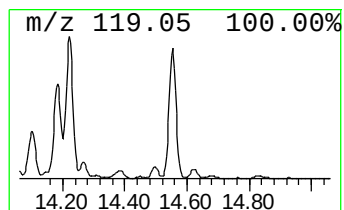
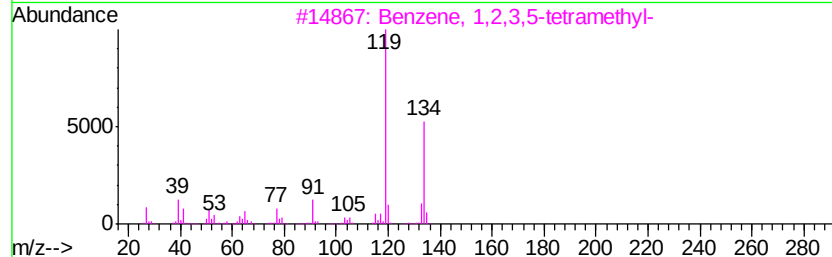
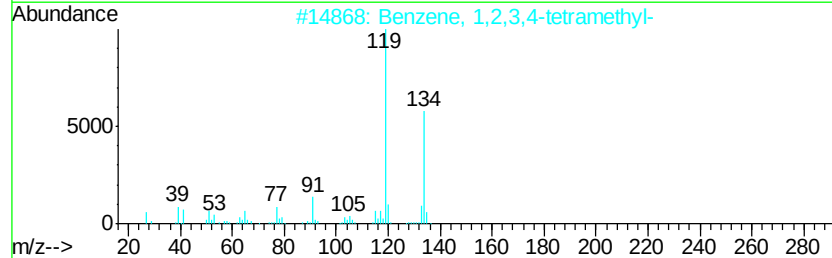
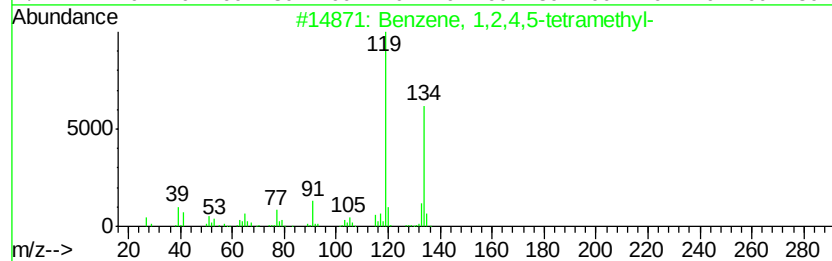
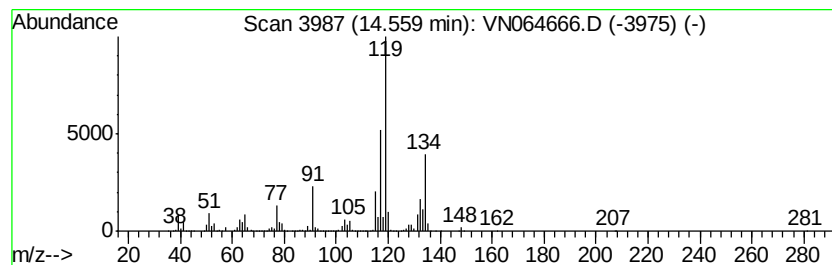
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	12.57 ug/l	4352190	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	87
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
4		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	76



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN111620\
Data File : VN064666.D
Acq On : 16 Nov 2020 17:36
Operator : JC/MD
Sample : L4772-05
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3R

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.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
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Cyclohexane, propyl-	12.14	22.5	ug/l	777501	3	11.38	1726160	50.0
Benzene, 1,4-diet...	13.49	10.9	ug/l	3763880	4	13.32	17311100	50.0
Benzene, 1,1'-(1,...	13.52	33.0	ug/l	11415100	4	13.32	17311100	50.0
Benzene, 2-ethyl-...	13.77	9.4	ug/l	3247210	4	13.32	17311100	50.0
o-Cymene	13.80	9.1	ug/l	3153060	4	13.32	17311100	50.0
Benzene, 4-ethyl-...	13.86	18.3	ug/l	6327310	4	13.32	17311100	50.0
Benzene, (1-methy...	13.97	20.0	ug/l	6922580	4	13.32	17311100	50.0
Benzene, 1-ethyl-...	14.22	7.4	ug/l	2567070	4	13.32	17311100	50.0
Benzene, 1,2,4,5-...	14.56	12.6	ug/l	4352190	4	13.32	17311100	50.0