

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111620\
 Data File : VN064669.D
 Acq On : 16 Nov 2020 18:49
 Operator : JC/MD
 Sample : L4772-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 FD-111620

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	1853	rVB2	382791	919404	1.42%	0.405%
2	7.991	1924	1944	1965	rBV	281674	670313	1.04%	0.295%
3	8.550	2103	2118	2144	rVB	634681	1338349	2.07%	0.589%
4	10.058	2573	2587	2615	rBV	1138266	2202363	3.41%	0.969%
5	11.380	2985	2998	3015	rBV	1082559	1900415	2.94%	0.836%
6	11.482	3015	3030	3047	rBV	2065223	3782680	5.86%	1.665%
7	11.598	3047	3066	3084	rBV	3918866	8453356	13.09%	3.721%
8	11.920	3145	3166	3181	rBV5	725785	2570664	3.98%	1.131%
9	12.093	3212	3220	3227	rBV2	667468	998826	1.55%	0.440%
10	12.142	3227	3235	3249	rVB5	699841	1291414	2.00%	0.568%
11	12.222	3250	3260	3272	rBV	5721942	9127078	14.13%	4.017%
12	12.376	3297	3308	3319	rBV	1012587	1762556	2.73%	0.776%
13	12.447	3320	3330	3340	rBV4	366538	762508	1.18%	0.336%
14	12.566	3350	3367	3378	rBV	8090572	14108354	21.84%	6.209%
15	12.630	3378	3387	3391	rBV	1443604	2226209	3.45%	0.980%
16	12.662	3391	3397	3403	rVV	2306866	3608609	5.59%	1.588%
17	12.707	3404	3411	3421	rVB	3229807	4979743	7.71%	2.192%
18	12.865	3447	3460	3469	rBV2	434974	998174	1.55%	0.439%
19	12.936	3474	3482	3487	rBV4	824919	1301271	2.01%	0.573%
20	13.016	3495	3507	3520	rVB	37752929	64588447	100.00%	28.427%
21	13.145	3533	3547	3556	rBV2	3000868	6389858	9.89%	2.812%
22	13.209	3557	3567	3576	rVV2	1871870	3175144	4.92%	1.397%
23	13.264	3576	3584	3592	rVV	1770523	2671966	4.14%	1.176%
24	13.347	3592	3610	3623	rVB	10950948	19655551	30.43%	8.651%
25	13.418	3623	3632	3642	rVV	1063811	1614003	2.50%	0.710%
26	13.486	3643	3653	3656	rVV	3251166	4369932	6.77%	1.923%
27	13.518	3657	3663	3671	rVV	7828835	12350445	19.12%	5.436%
28	13.566	3671	3678	3693	rVV4	2535524	6268582	9.71%	2.759%
29	13.646	3693	3703	3712	rVB3	1445773	2509028	3.88%	1.104%
30	13.778	3733	3744	3748	rBV	2346802	3743325	5.80%	1.648%
31	13.807	3749	3753	3761	rVV	2590593	3439161	5.32%	1.514%
32	13.862	3761	3770	3780	rVV	4885480	7454763	11.54%	3.281%
33	13.920	3780	3788	3794	rVV	1479816	2307901	3.57%	1.016%
34	13.968	3794	3803	3814	rVB2	4878688	7983312	12.36%	3.514%

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Integration Parameters: RTEINT.P

Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	14.103	3835	3845	3854	rBV2	1238888	2002563	3.10%	0.881%
36	14.183	3855	3870	3876	rVV	1465760	2764113	4.28%	1.217%
37	14.222	3876	3882	3891	rVB	2027235	2946992	4.56%	1.297%
38	14.450	3945	3953	3962	rBV2	551285	869510	1.35%	0.383%
39	14.556	3975	3986	4000	rBV4	2540157	5038362	7.80%	2.218%
40	14.714	4025	4035	4045	rBV	1323231	2060814	3.19%	0.907%

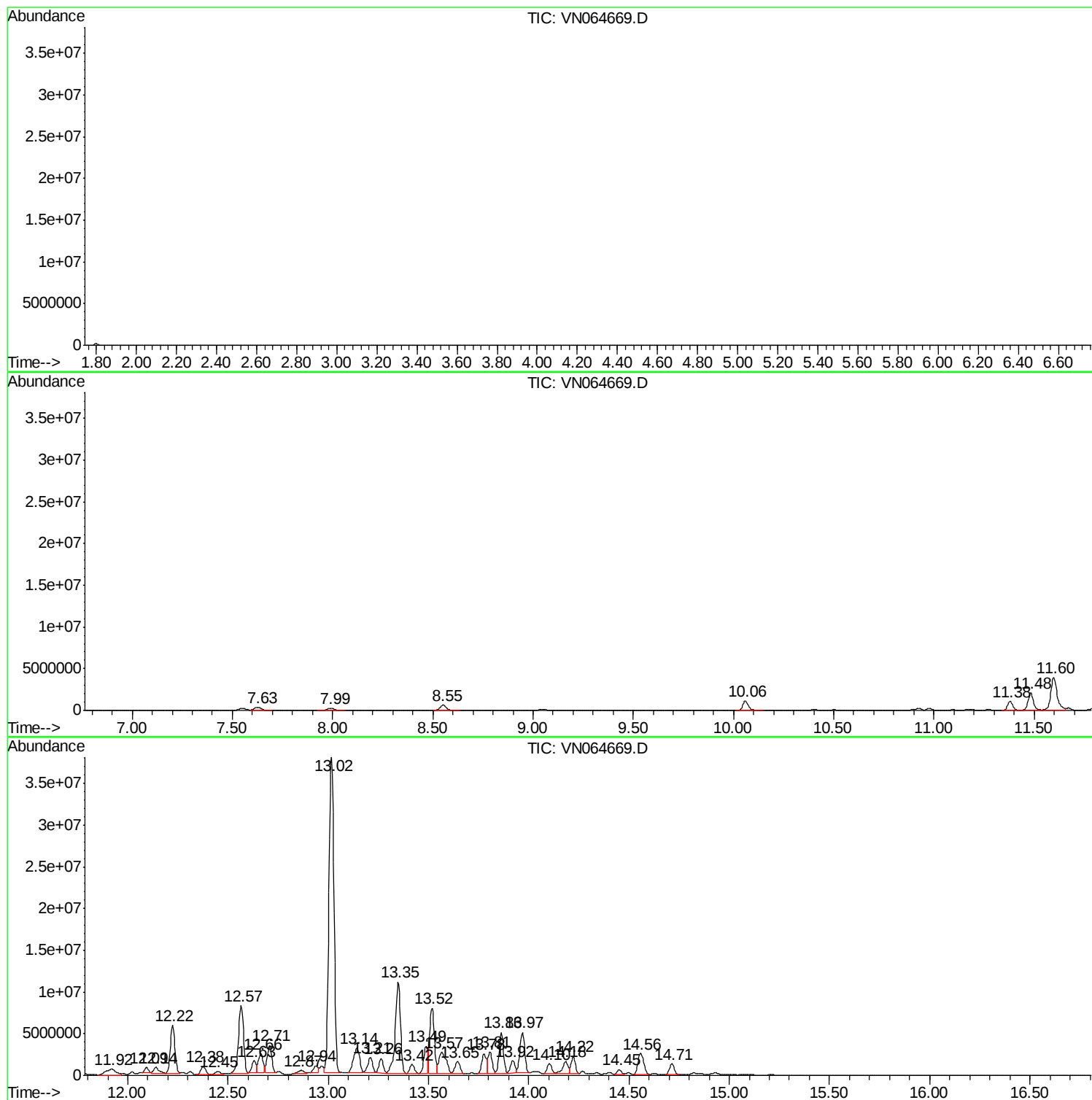
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Quant Title : SW846 8260

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



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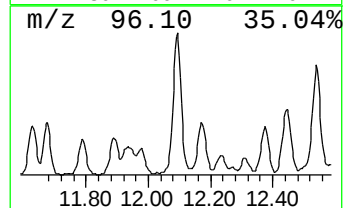
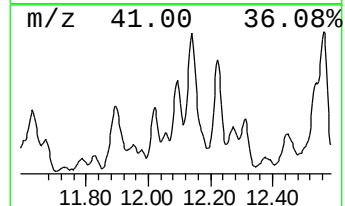
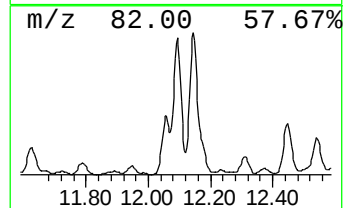
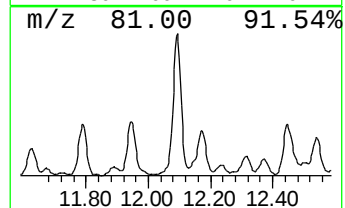
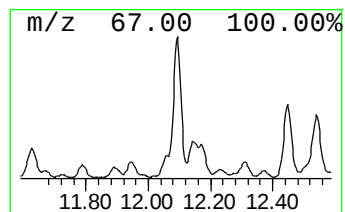
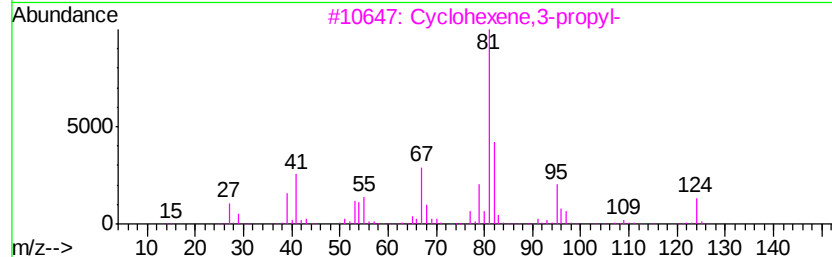
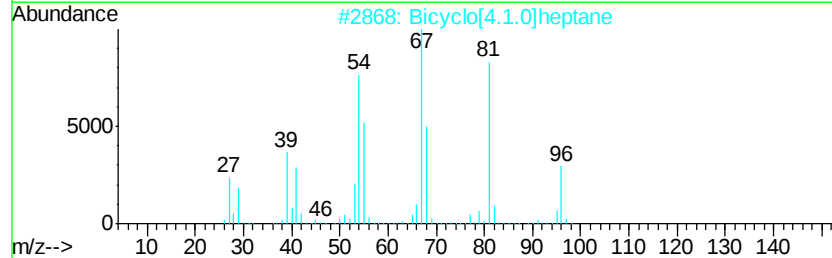
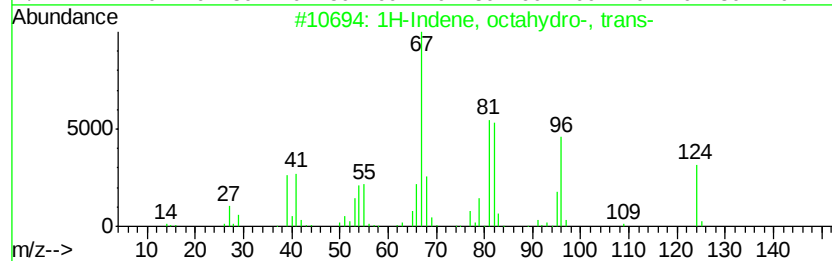
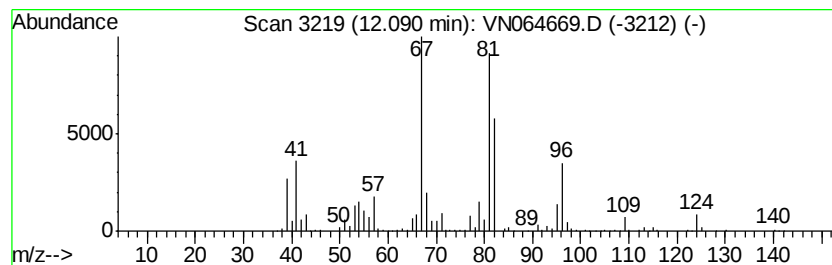
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N110420W.M
Quant Title : SW846 8260

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

Peak Number 1 1H-Indene, octahydro-, trans- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	26.28 ug/l	998826	Chlorobenzene-d5	11.38
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, octahydro-, trans-	124 C9H16	003296-50-2	70
2	Bicyclo[4.1.0]heptane	96 C7H12	000286-08-8	64
3	Cyclohexene, 3-propyl-	124 C9H16	003983-06-0	46
4	Cyclohexene, 3-methyl-	96 C7H12	000591-48-0	43
5	Cyclohexene, 1-methyl-	96 C7H12	000591-49-1	43



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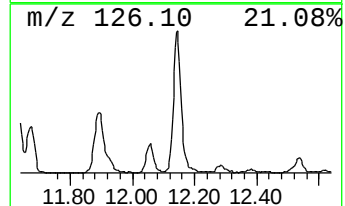
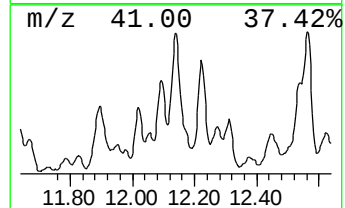
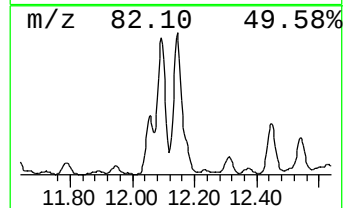
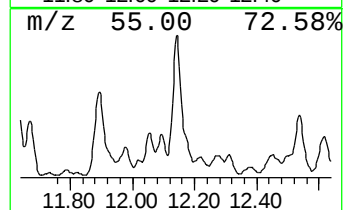
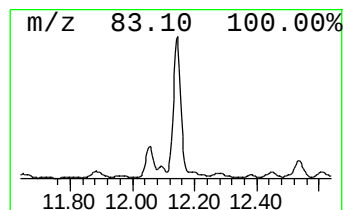
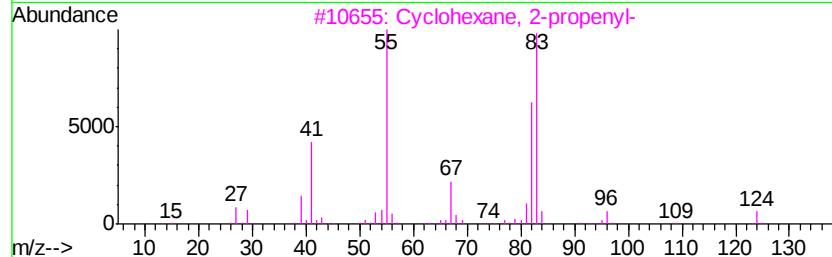
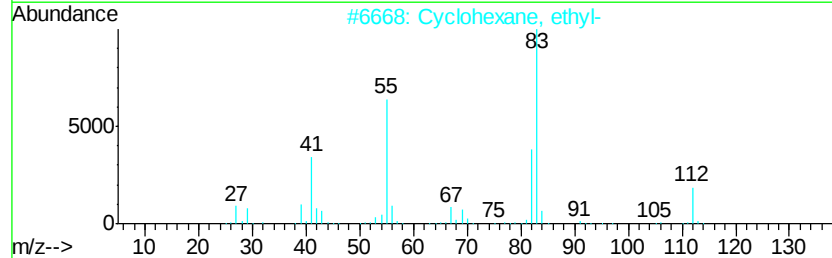
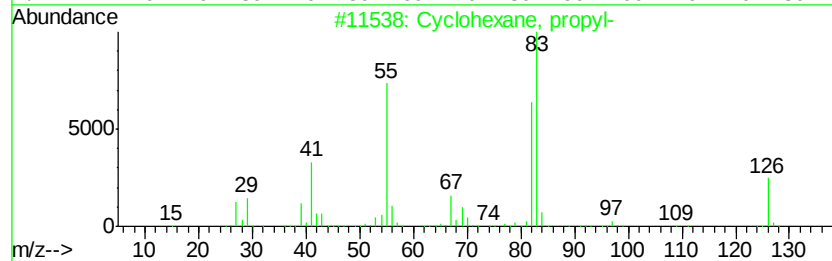
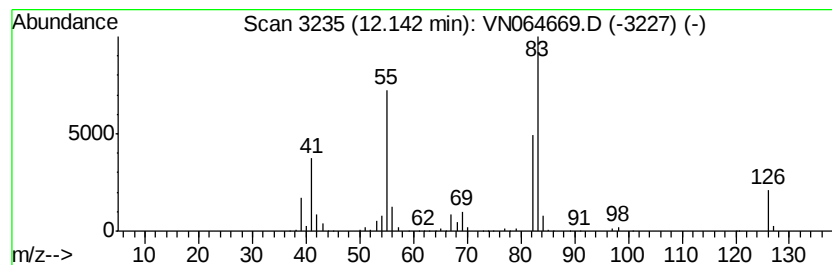
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclohexane, propyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	33.98 ug/l	1291410	Chlorobenzene-d5	11.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, propyl-	126	C9H18	001678-92-8	91
2		Cyclohexane, ethyl-	112	C8H16	001678-91-7	72
3		Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	59
4		Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	59
5		Lomustine	233	C9H16ClN3O2	013010-47-4	59



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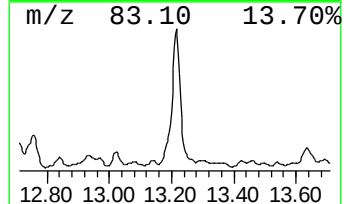
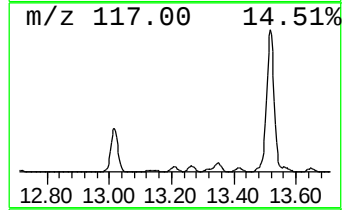
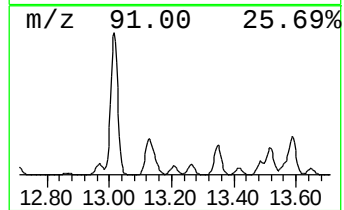
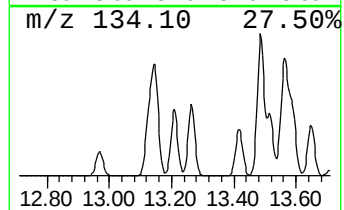
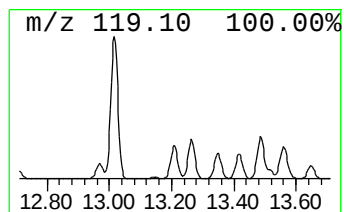
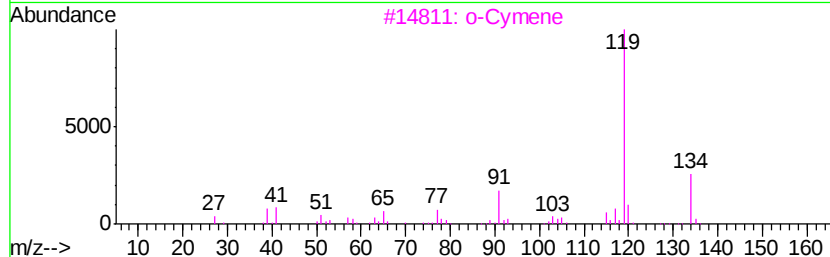
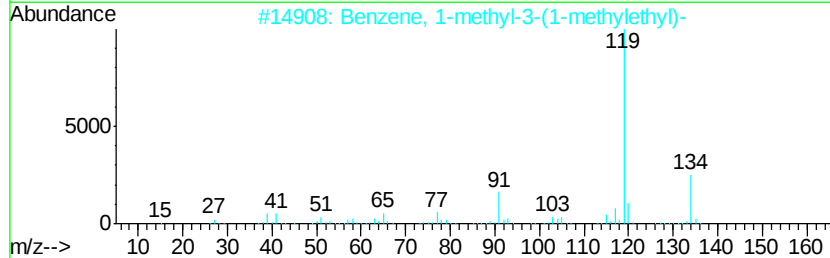
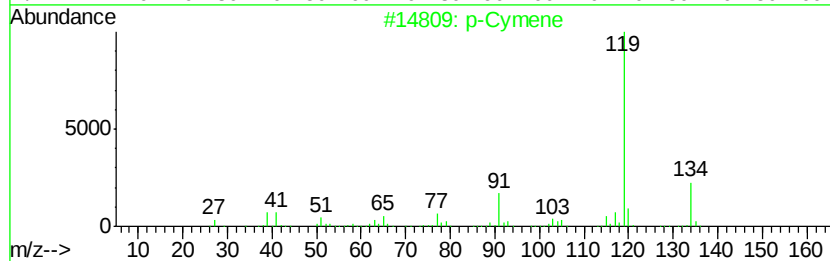
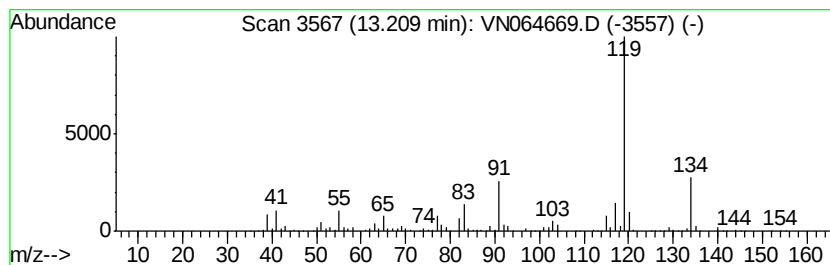
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	8.08 ug/l	3175140	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	93
2		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
3		o-Cymene	134	C10H14	000527-84-4	93
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	87
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87



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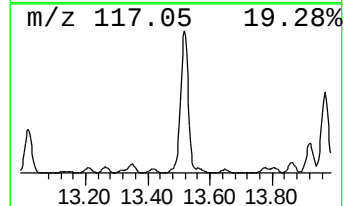
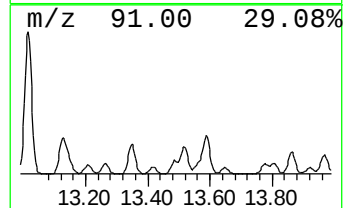
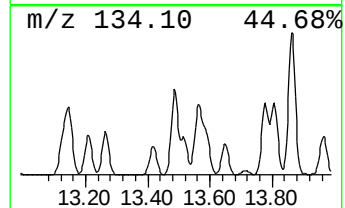
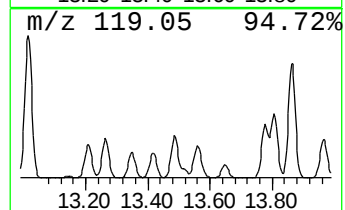
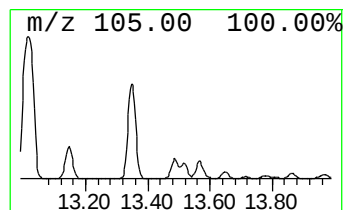
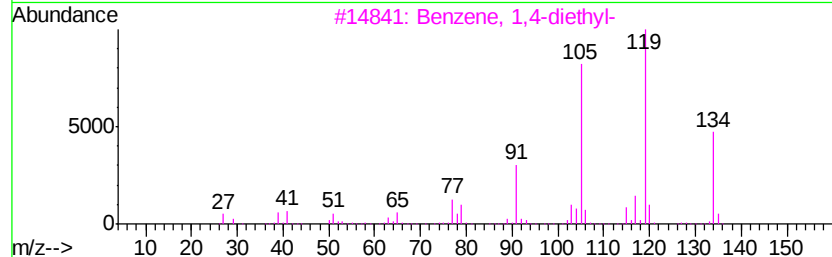
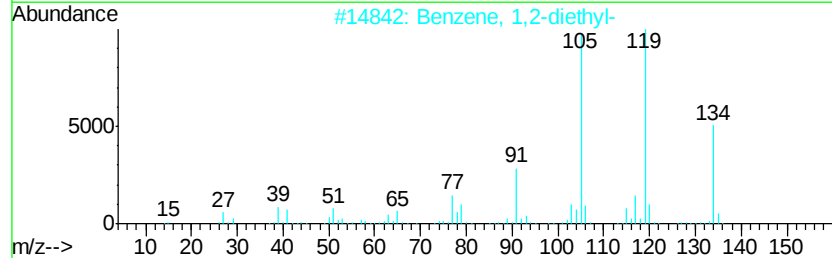
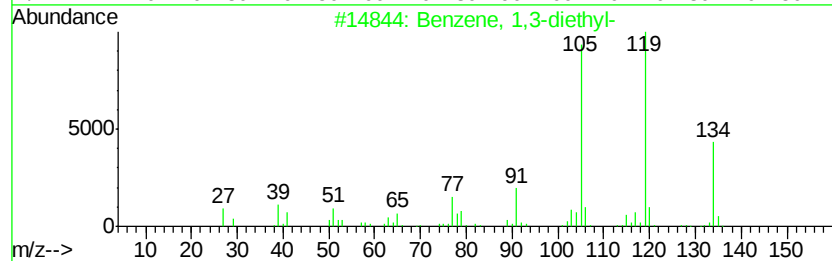
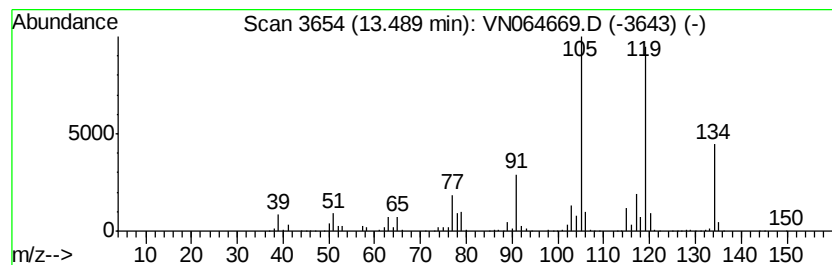
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	11.12 ug/l	4369930	1,4-Dichlorobenzene-d4	13.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



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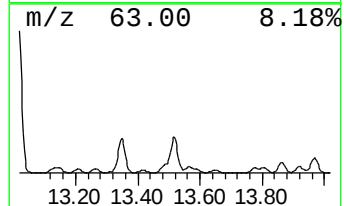
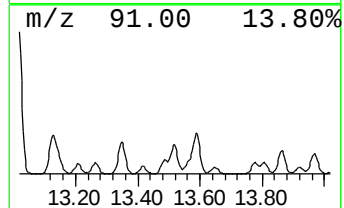
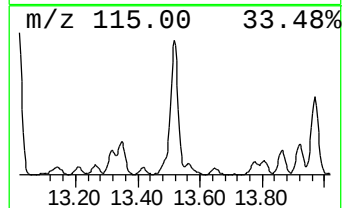
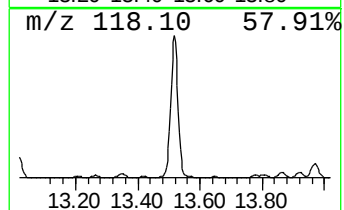
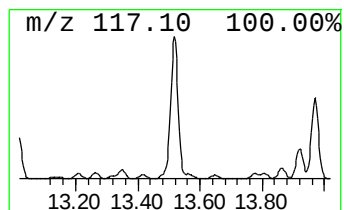
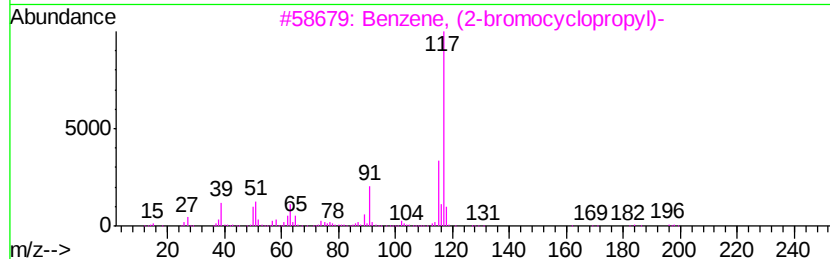
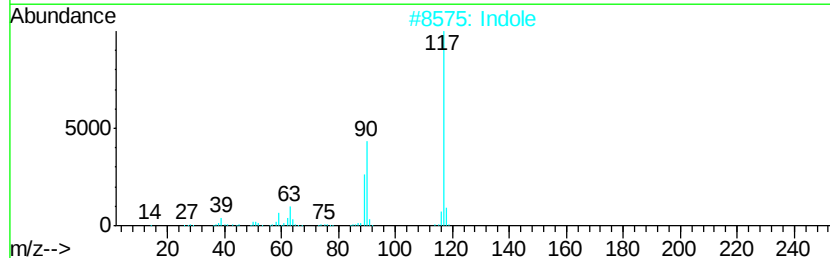
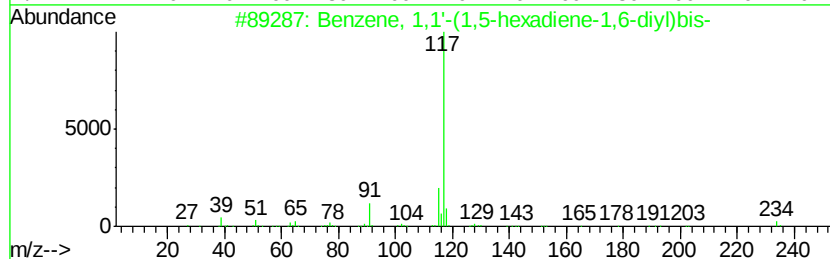
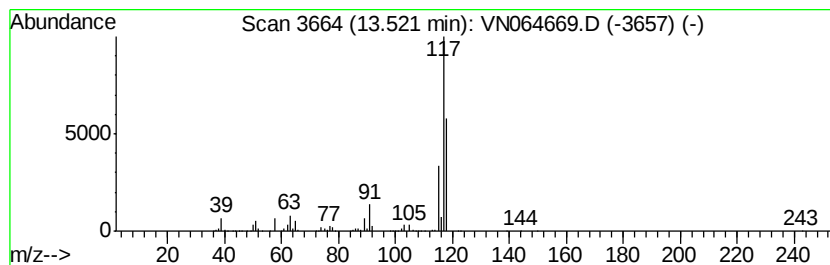
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	31.42 ug/l	12350400	1,4-Dichlorobenzene-d4	13.32

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2	Indole	117	C8H7N	000120-72-9	49
3	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
4	5H-1-Pyrindine	117	C8H7N	000270-91-7	43
5	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064669.D
Acq On : 16 Nov 2020 18:49
Operator : JC/MD
Sample : L4772-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FD-111620

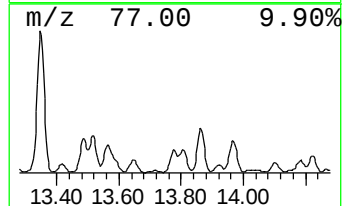
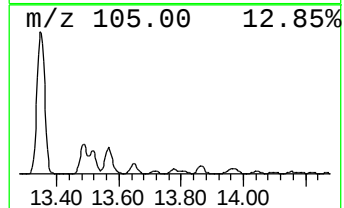
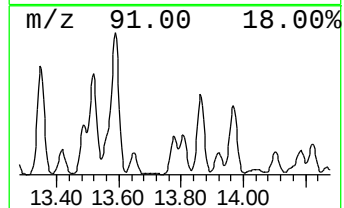
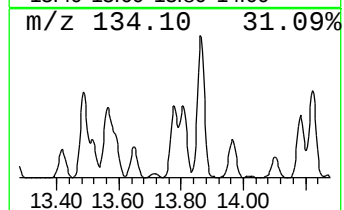
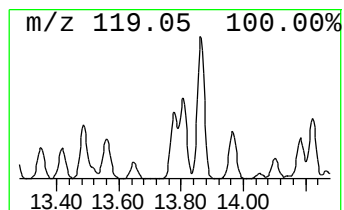
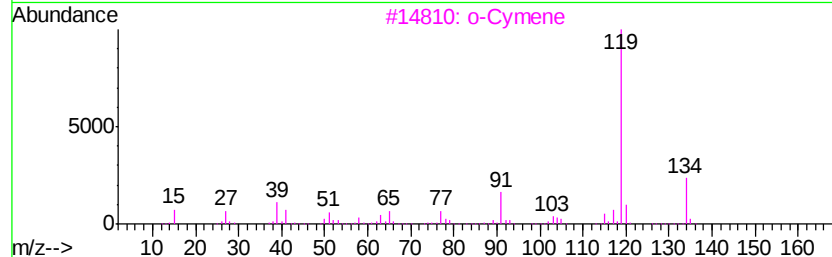
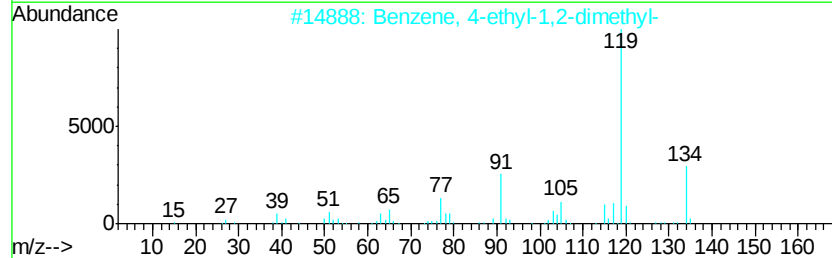
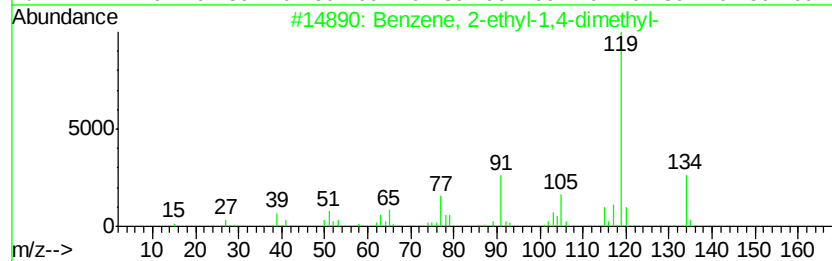
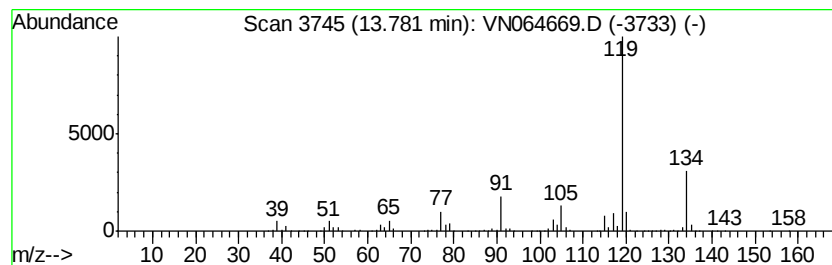
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	9.52 ug/l	3743330	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	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		o-Cymene	134	C10H14	000527-84-4	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064669.D
Acq On : 16 Nov 2020 18:49
Operator : JC/MD
Sample : L4772-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FD-111620

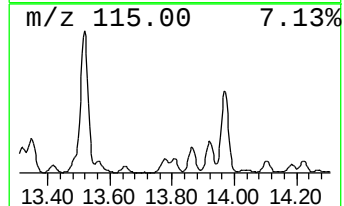
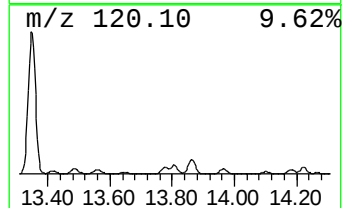
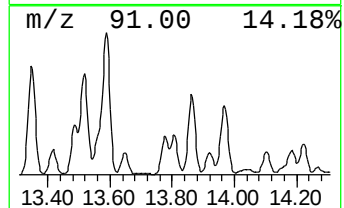
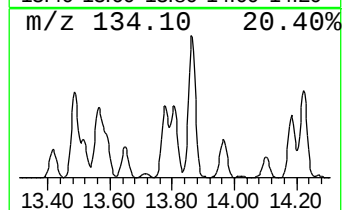
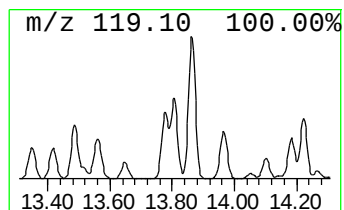
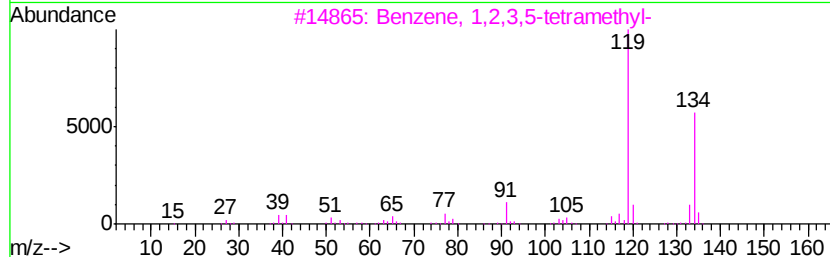
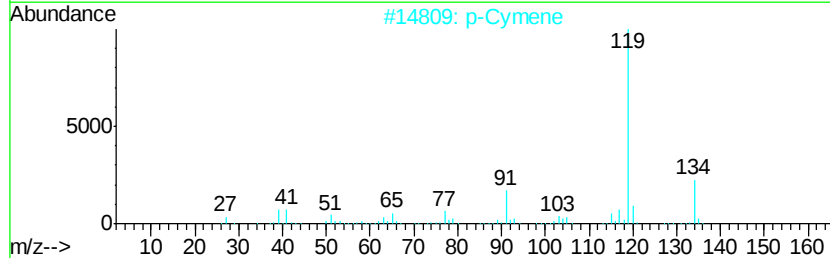
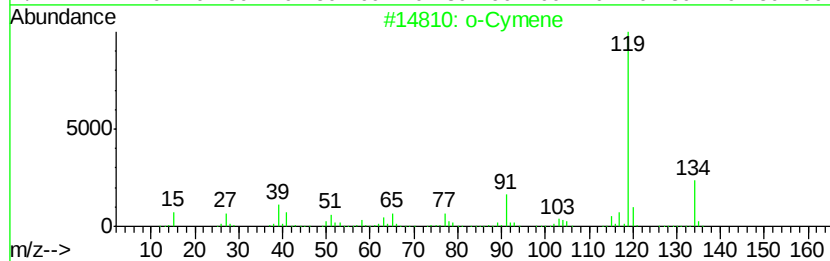
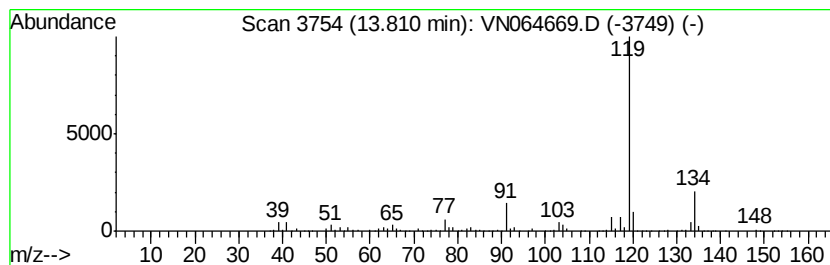
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Quant Title : SW846 8260

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

Peak Number 7 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	8.75 ug/l	3439160	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	94
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	91
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064669.D
Acq On : 16 Nov 2020 18:49
Operator : JC/MD
Sample : L4772-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FD-111620

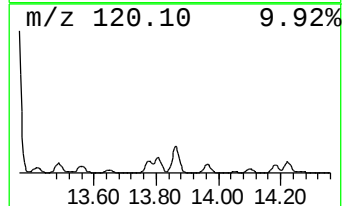
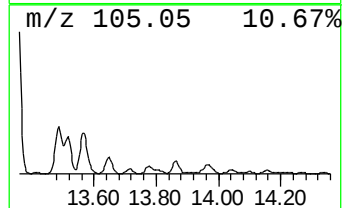
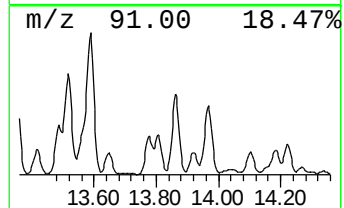
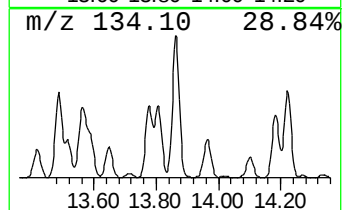
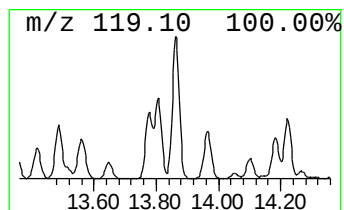
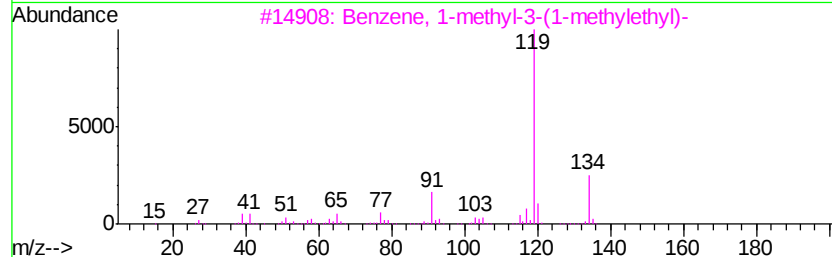
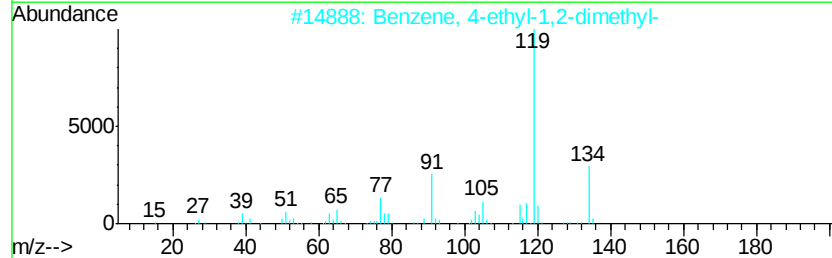
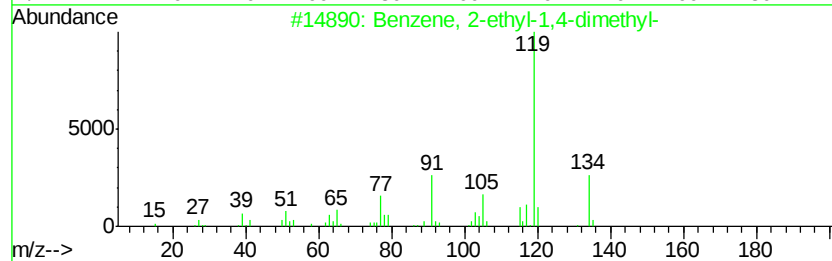
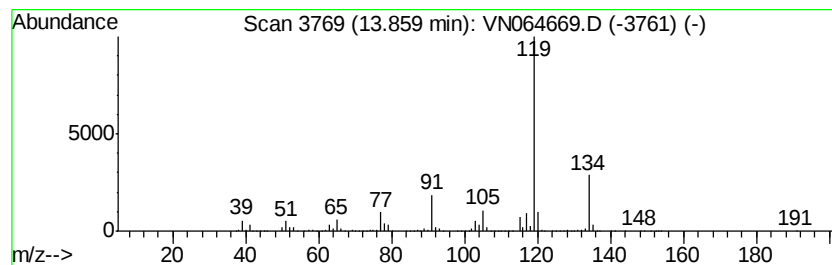
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	18.96 ug/l	7454760	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	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064669.D
Acq On : 16 Nov 2020 18:49
Operator : JC/MD
Sample : L4772-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FD-111620

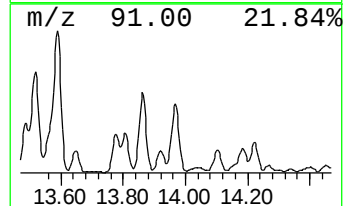
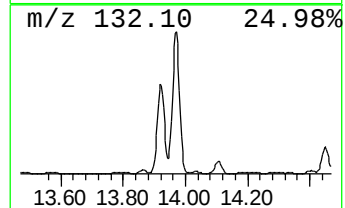
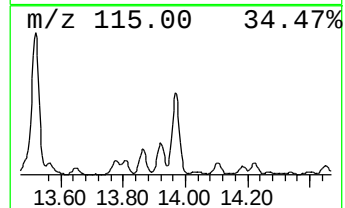
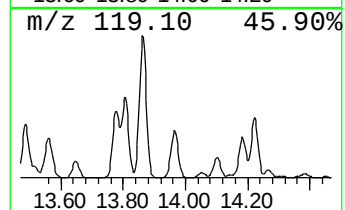
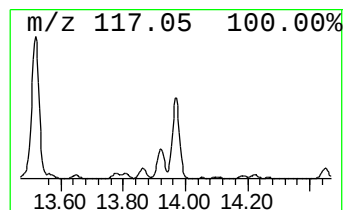
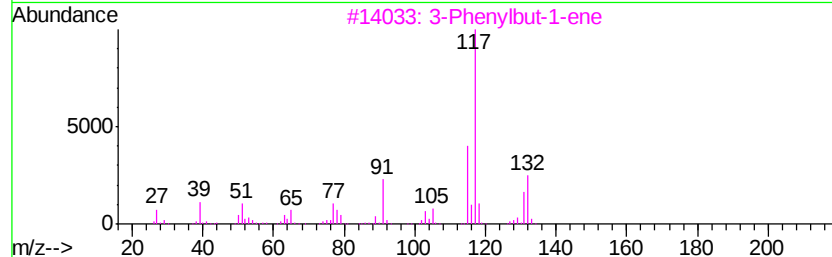
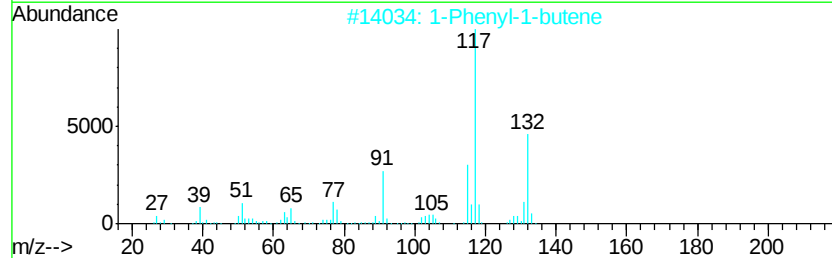
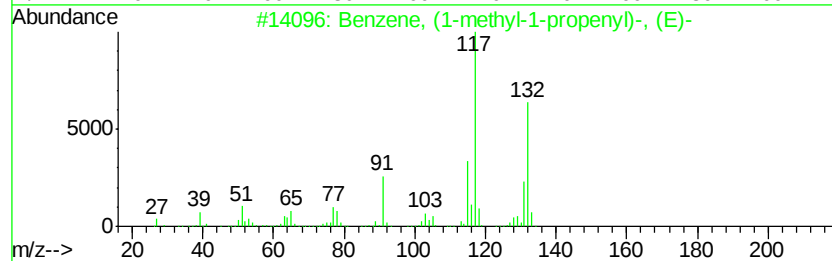
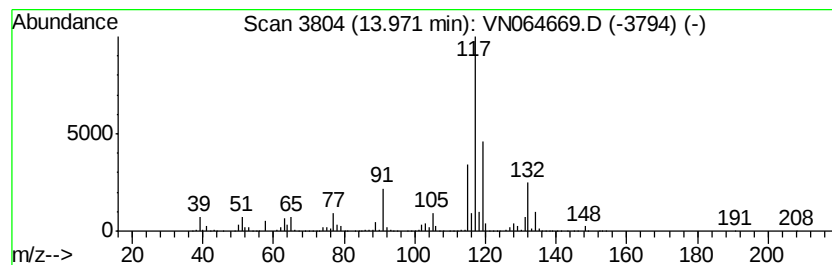
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1-methyl-1-propen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	20.31 ug/l	7983310	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	86
2		1-Phenyl-1-butene	132	C10H12	000824-90-8	86
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	83
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN111620\
Data File : VN064669.D
Acq On : 16 Nov 2020 18:49
Operator : JC/MD
Sample : L4772-08
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FD-111620

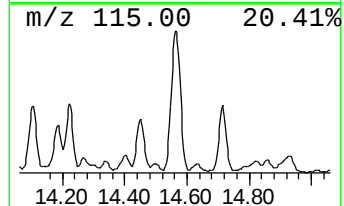
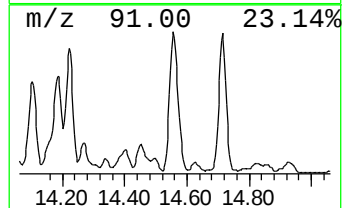
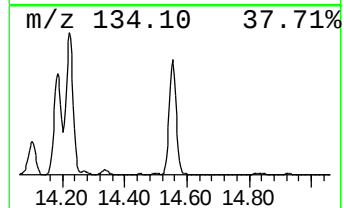
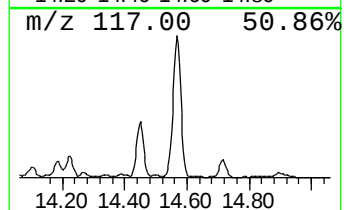
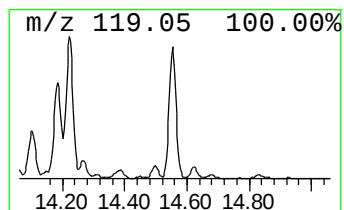
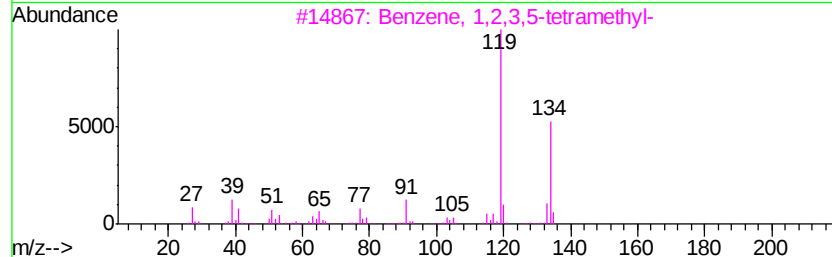
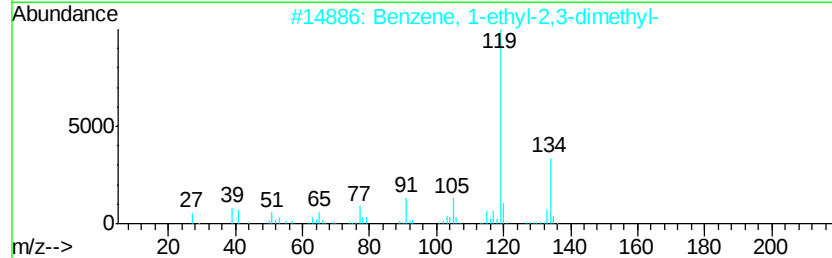
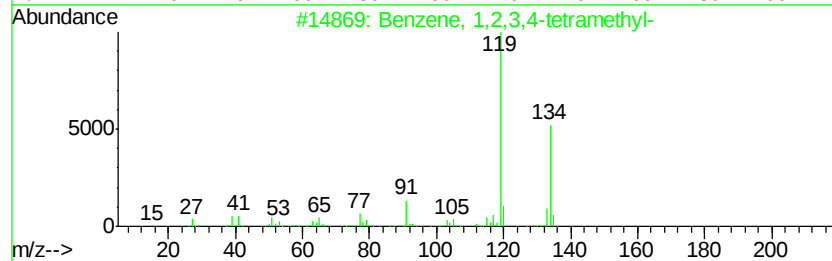
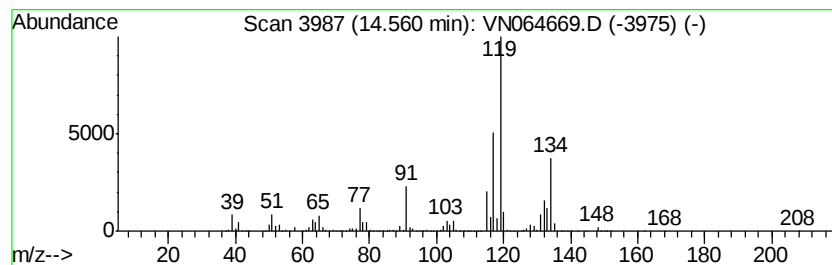
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	12.82 ug/l	5038360	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	81
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	81
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
4		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81
5		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	81



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN111620\
Data File : VN064669.D
Acq On : 16 Nov 2020 18:49
Operator : JC/MD
Sample : L4772-08
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
FD-111620

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	34.0	ug/l	1291410	3	11.38	1900420	50.0
p-Cymene	13.21	8.1	ug/l	3175140	4	13.32	19655600	50.0
Benzene, 1,3-diet...	13.49	11.1	ug/l	4369930	4	13.32	19655600	50.0
Benzene, 1,1'-(1,...	13.52	31.4	ug/l	12350400	4	13.32	19655600	50.0
Benzene, 2-ethyl-...	13.78	9.5	ug/l	3743330	4	13.32	19655600	50.0
o-Cymene	13.81	8.8	ug/l	3439160	4	13.32	19655600	50.0
Benzene, 4-ethyl-...	13.86	19.0	ug/l	7454760	4	13.32	19655600	50.0
Benzene, (1-methy...	13.97	20.3	ug/l	7983310	4	13.32	19655600	50.0
Benzene, 1,2,3,4-...	14.56	12.8	ug/l	5038360	4	13.32	19655600	50.0