

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031319\
 Data File : VN054220.D
 Acq On : 13 Mar 2019 18:45
 Operator : JC/SP
 Sample : K1842-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PR-MW-05_030819

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\82N031219W.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	2.812	310	323	338	rVB	537592	1080228	3.51%	0.809%
2	3.143	415	426	443	rVB2	139686	308719	1.00%	0.231%
3	4.580	866	873	902	rVB3	161914	534328	1.74%	0.400%
4	5.047	996	1018	1039	rBV3	154797	547351	1.78%	0.410%
5	6.625	1489	1509	1529	rBV2	365280	1121613	3.65%	0.840%
6	7.593	1786	1810	1820	rBV	900063	2301617	7.49%	1.724%
7	7.664	1821	1832	1847	rVB	1371150	3192213	10.38%	2.391%
8	8.027	1928	1945	1971	rBV	836880	1935780	6.30%	1.450%
9	8.204	1972	2000	2019	rBV6	174750	849186	2.76%	0.636%
10	8.587	2103	2119	2159	rVB	2295094	5247647	17.07%	3.930%
11	10.091	2571	2587	2600	rBV	3497538	6887711	22.40%	5.158%
12	10.156	2602	2607	2635	rVB	414280	836461	2.72%	0.626%
13	11.406	2981	2996	3015	rBV	3560408	6701988	21.80%	5.019%
14	11.509	3015	3028	3048	rVV	9728356	17493719	56.91%	13.100%
15	11.622	3049	3063	3102	rVB	15364779	30741898	100.00%	23.022%
16	12.249	3245	3258	3282	rVB	1686743	2786961	9.07%	2.087%
17	12.403	3290	3306	3337	rBV	3324539	5822711	18.94%	4.360%
18	12.590	3350	3364	3374	rBV	1276966	2098584	6.83%	1.572%
19	12.654	3374	3384	3390	rVV	1913476	3257227	10.60%	2.439%
20	12.686	3390	3394	3402	rVV	1704968	2596342	8.45%	1.944%
21	12.731	3402	3408	3429	rVB	1501863	2406507	7.83%	1.802%
22	12.892	3445	3458	3471	rBV	203793	349283	1.14%	0.262%
23	13.040	3482	3504	3530	rBV	6947868	10968457	35.68%	8.214%
24	13.342	3586	3598	3620	rVB2	4064850	8560056	27.84%	6.410%
25	13.545	3640	3661	3669	rBV2	2222972	4047694	13.17%	3.031%
26	13.583	3670	3673	3691	rVB4	426587	743987	2.42%	0.557%
27	13.737	3709	3721	3731	rVB2	195871	368436	1.20%	0.276%
28	13.802	3731	3741	3746	rBV	326818	516644	1.68%	0.387%
29	13.889	3759	3768	3778	rVB	551175	836376	2.72%	0.626%
30	13.995	3792	3801	3811	rVB2	471344	754847	2.46%	0.565%
31	14.210	3853	3868	3874	rBV	536923	885441	2.88%	0.663%
32	14.249	3874	3880	3889	rVB	620086	901347	2.93%	0.675%
33	14.477	3942	3951	3960	rBV	243397	363461	1.18%	0.272%
34	14.586	3972	3985	3999	rBV3	833107	1761204	5.73%	1.319%

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ClientSampleId :
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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

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

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Title : SW846 8260

35	14.741	4023	4033	4050	rVB	298912	472116	1.54%	0.354%
36	15.133	4134	4155	4167	rBV	1372013	2440474	7.94%	1.828%
37	16.159	4463	4474	4488	rBV	195336	394006	1.28%	0.295%
38	16.352	4522	4534	4549	rBV2	196398	422615	1.37%	0.316%

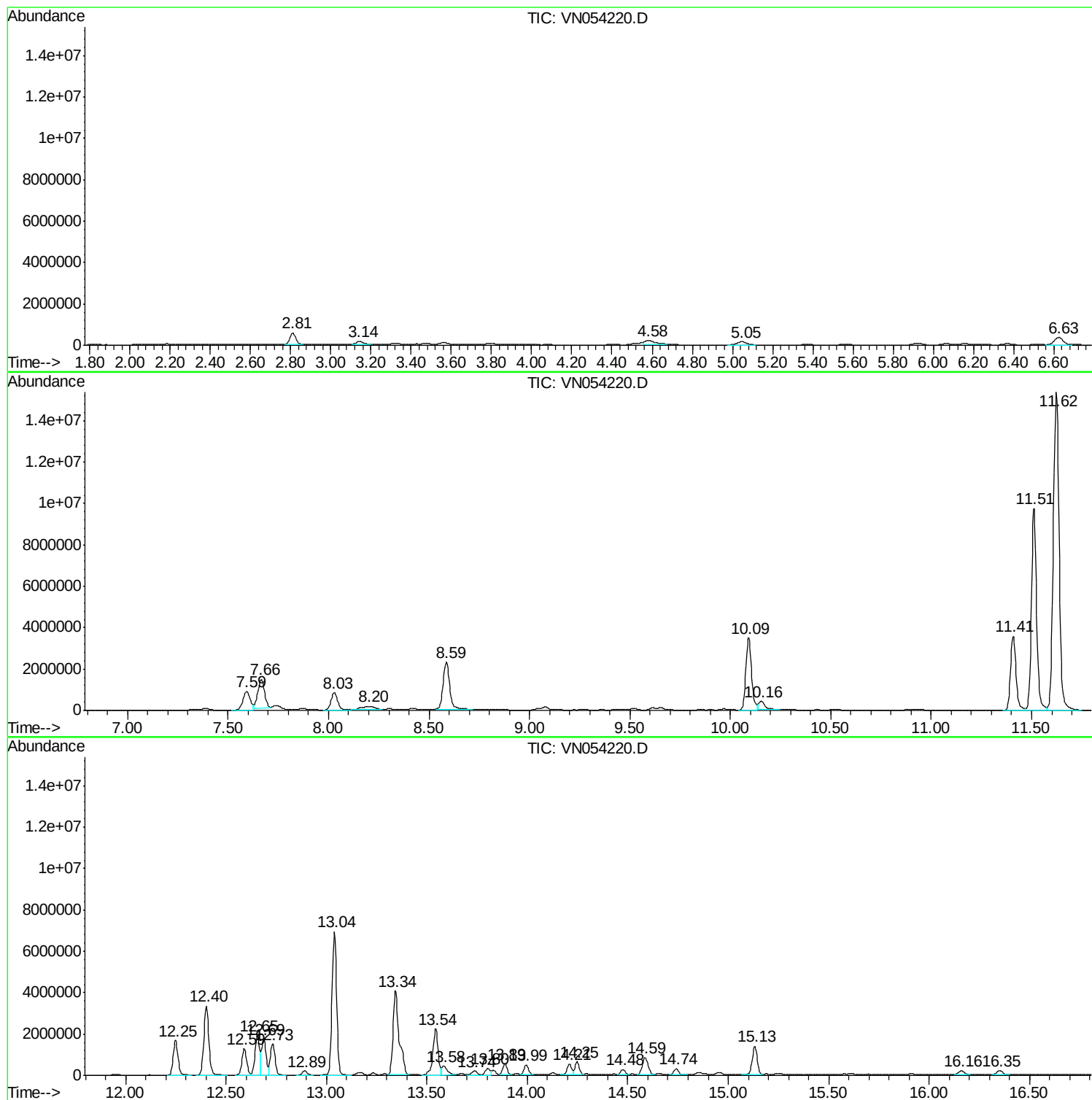
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Quant Title : SW846 8260

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



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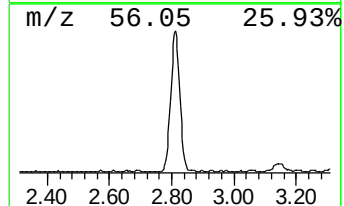
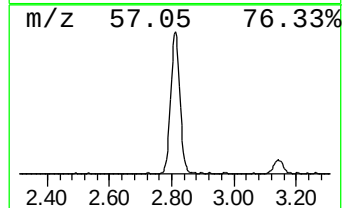
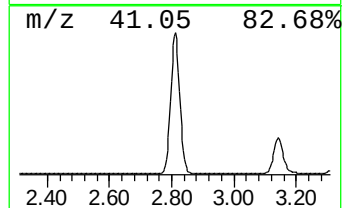
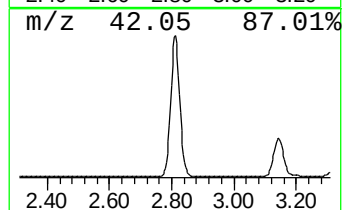
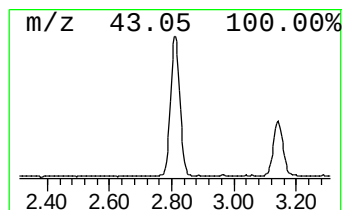
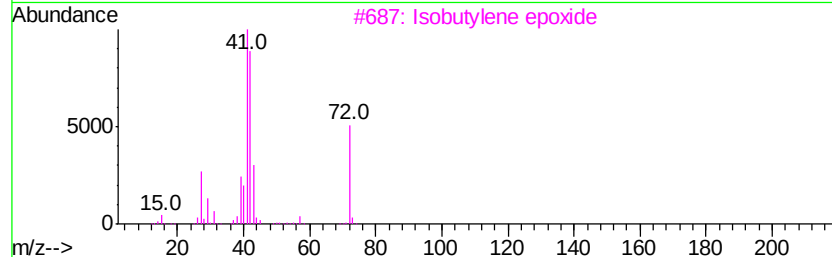
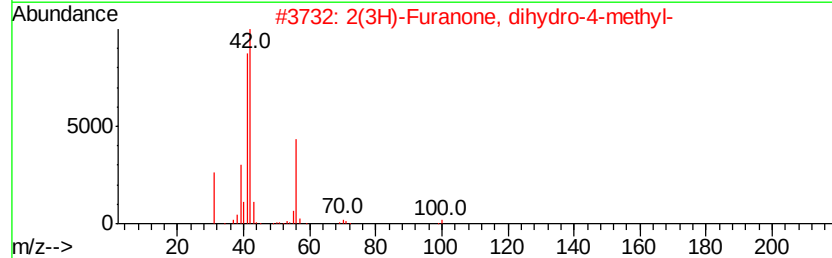
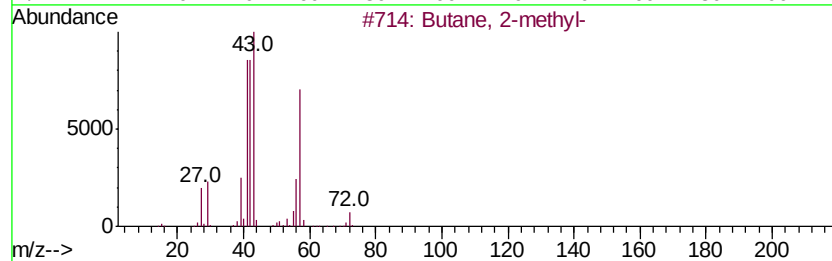
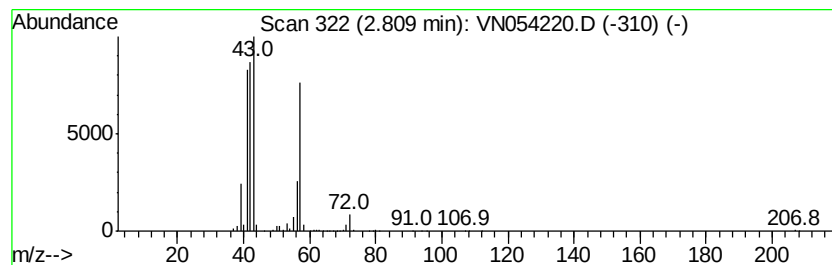
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.81	16.92 ug/l	1080230	Pentafluorobenzene	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2		2(3H)-Furanone, dihydro-4-methyl-	100	C5H8O2	001679-49-8	42
3		Isobutylene epoxide	72	C4H8O	000558-30-5	40
4		Pentane	72	C5H12	000109-66-0	37
5		3-Buten-1-ol	72	C4H8O	000627-27-0	25



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

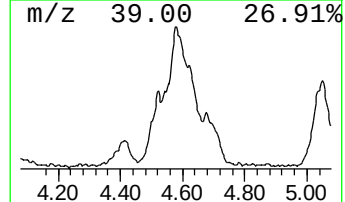
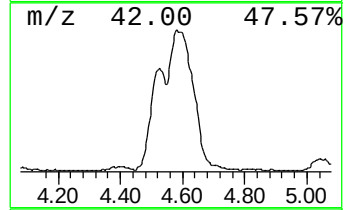
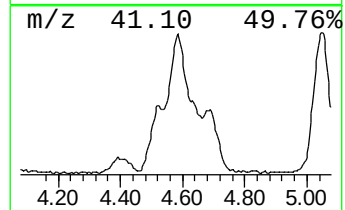
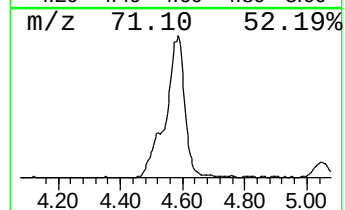
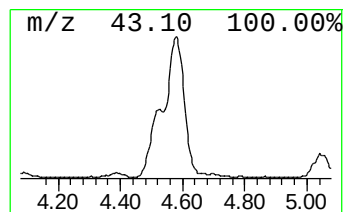
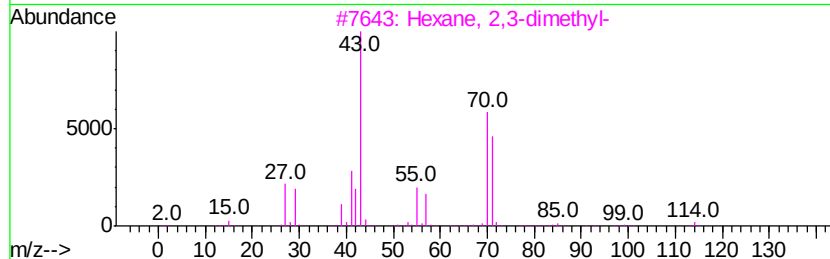
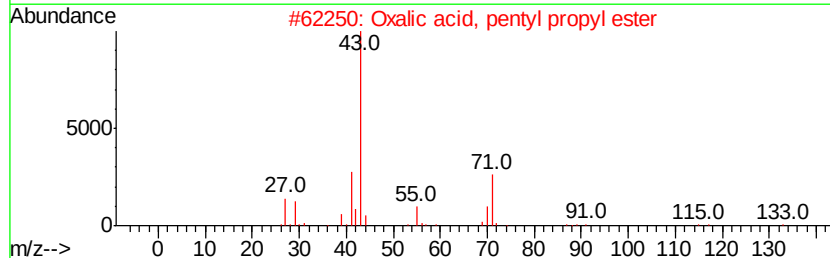
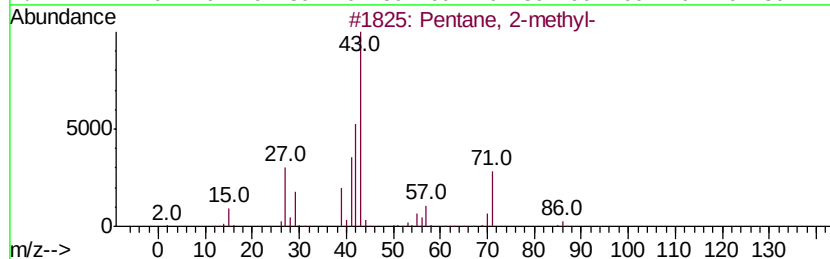
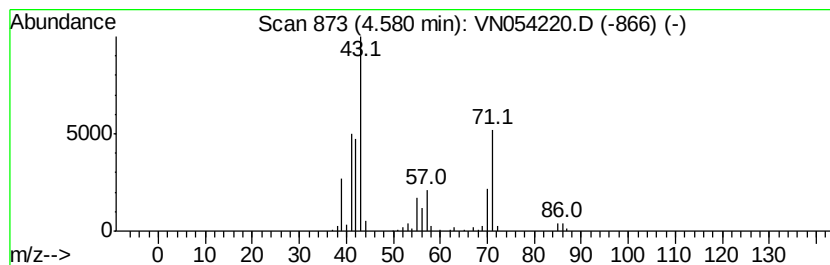
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.58	8.37 ug/l	534328	Pentafluorobenzene	7.66
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pentane, 2-methyl-	86 C6H14	000107-83-5	72
2	Oxalic acid, pentyl propyl ester	202 C10H18O4	1000309-25-7	38
3	Hexane, 2,3-dimethyl-	114 C8H18	000584-94-1	37
4	Furan, tetrahydro-2-methyl-	86 C5H10O	000096-47-9	28
5	1-Hexene, 4,5-dimethyl-	112 C8H16	016106-59-5	25



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ALS Vial : 22 Sample Multiplier: 1

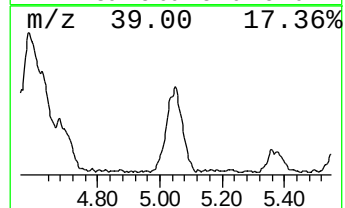
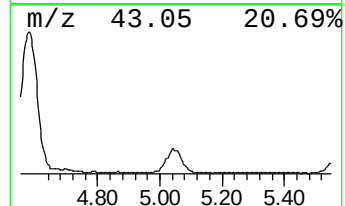
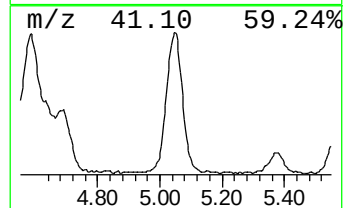
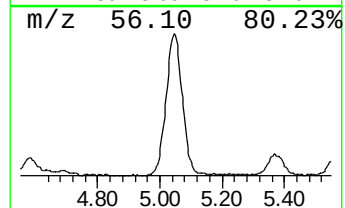
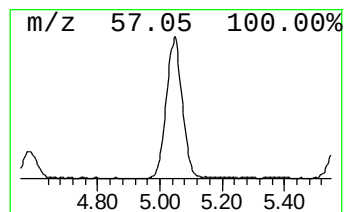
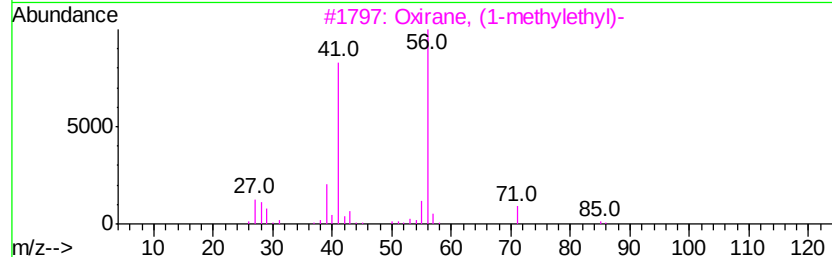
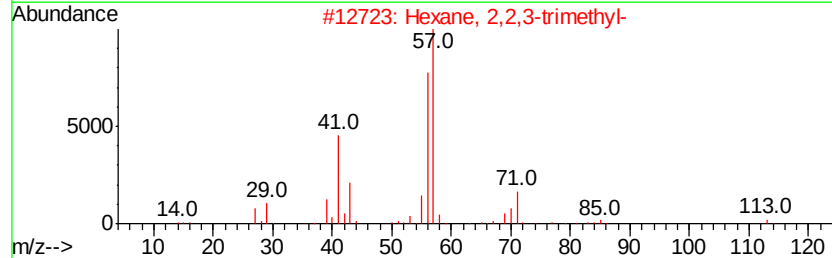
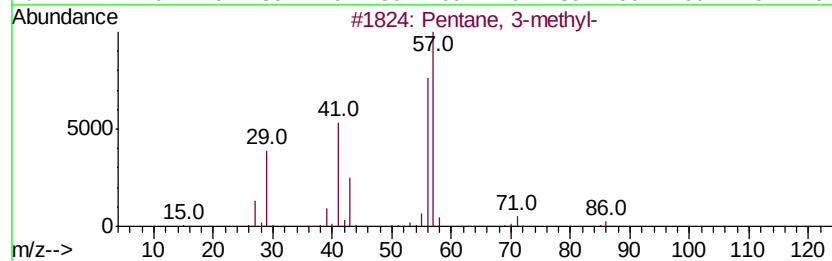
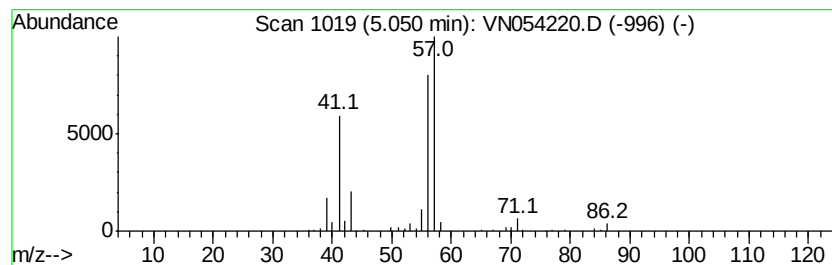
Instrument :
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Pentane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.05	8.57 ug/l	547351	Pentafluorobenzene	7.66	
Hit# of 5	Tentative ID		MW MolForm	CAS#	Qual
1	Pentane, 3-methyl-		86 C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-		128 C9H20	016747-25-4	64
3	Oxirane, (1-methylethyl)-		86 C5H10O	001438-14-8	50
4	Pentane, 2,2,4,4-tetramethyl-		128 C9H20	001070-87-7	45
5	Pentane, 3-ethyl-2,2-dimethyl-		128 C9H20	016747-32-3	43



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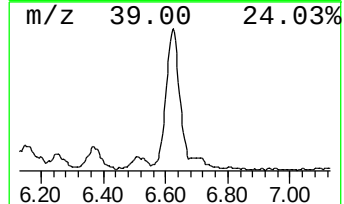
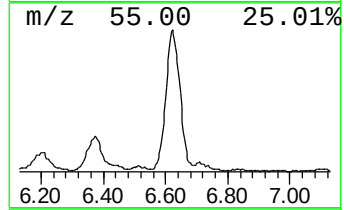
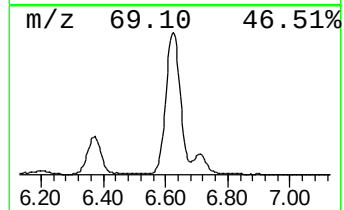
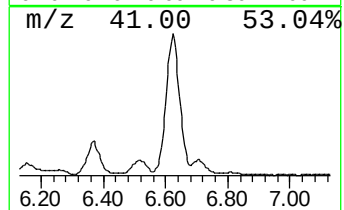
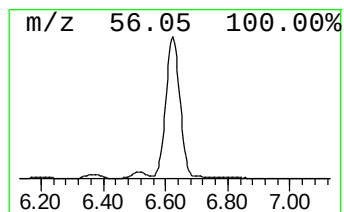
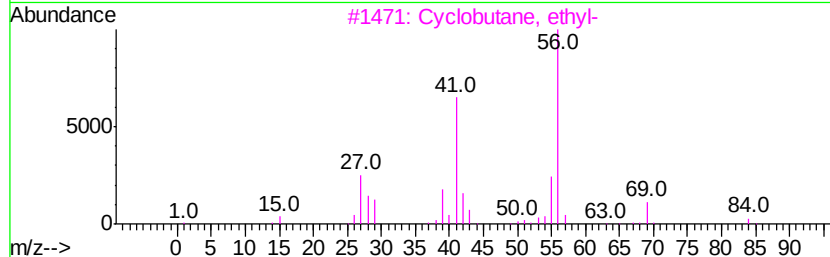
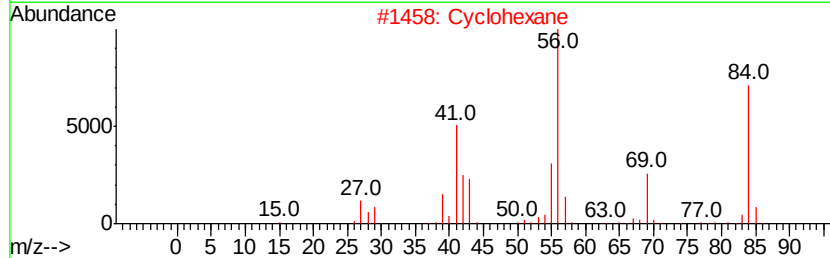
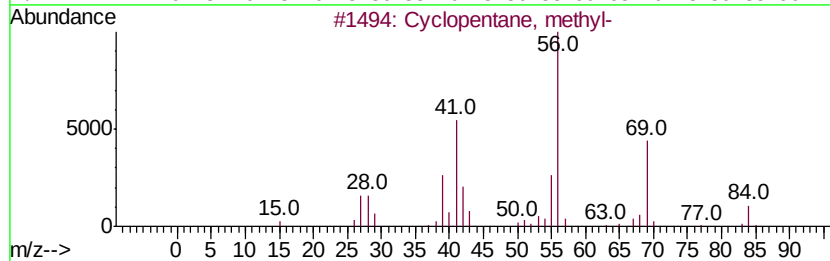
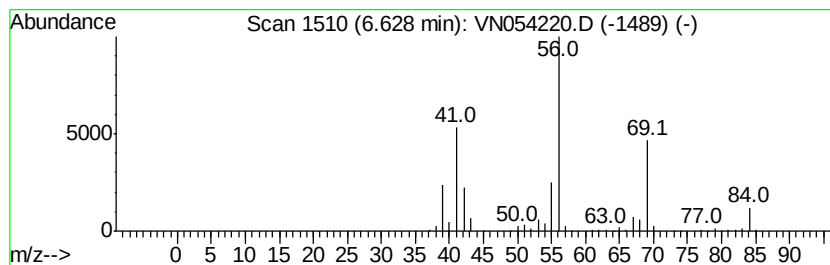
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	17.57 ug/l	1121610	Pentafluorobenzene	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2		Cyclohexane	84	C6H12	000110-82-7	78
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5		Cyclobutane	56	C4H8	000287-23-0	50



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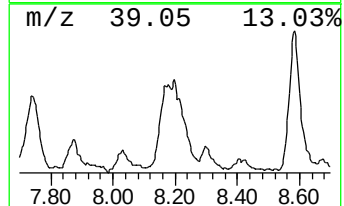
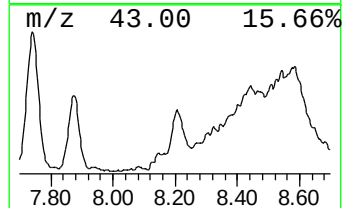
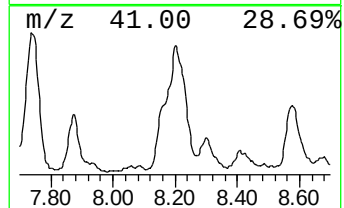
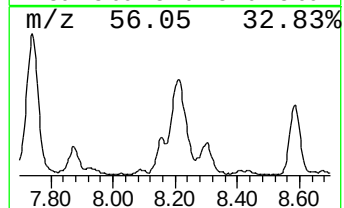
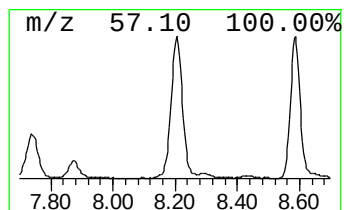
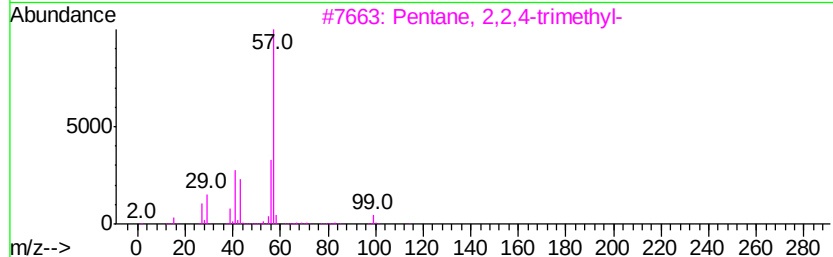
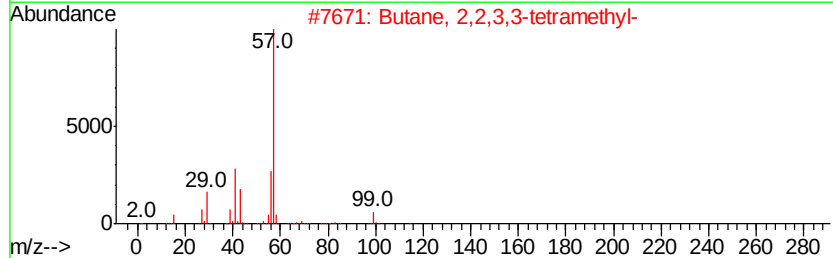
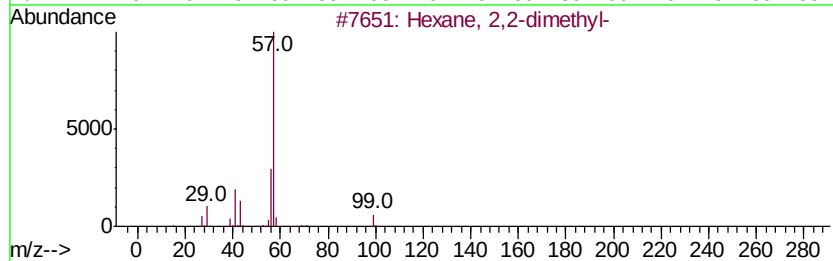
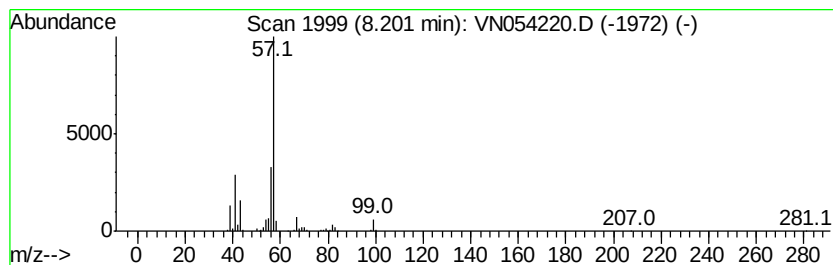
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 5 Hexane, 2,2-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.20	8.09 ug/l	849186	1,4-Difluorobenzene	8.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
2		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	64
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	59
4		Methane, isocyanato-	57	C2H3NO	000624-83-9	59
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN031319\
Data File : VN054220.D
Acq On : 13 Mar 2019 18:45
Operator : JC/SP
Sample : K1842-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PR-MW-05_030819

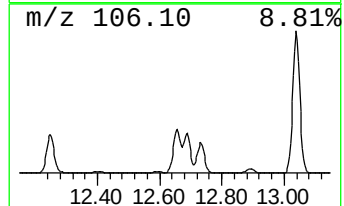
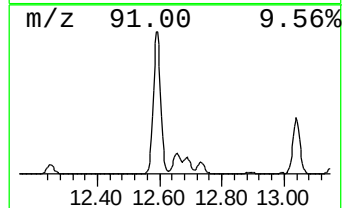
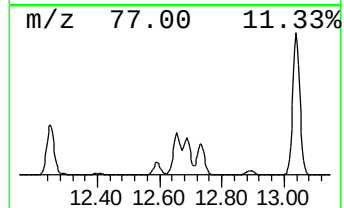
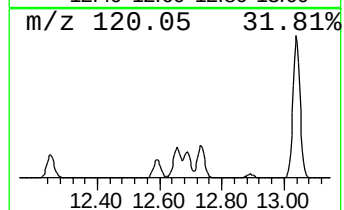
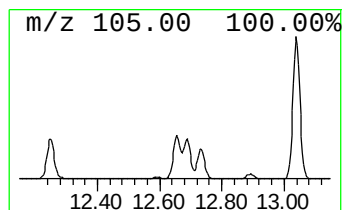
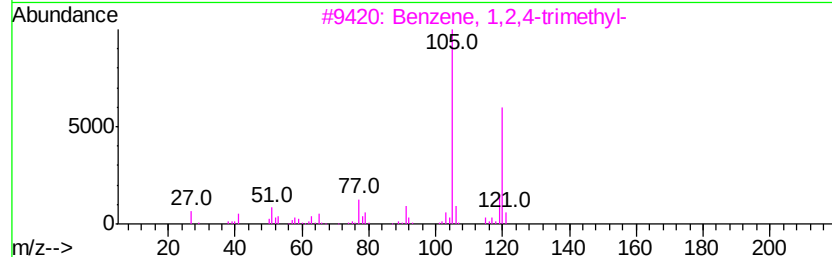
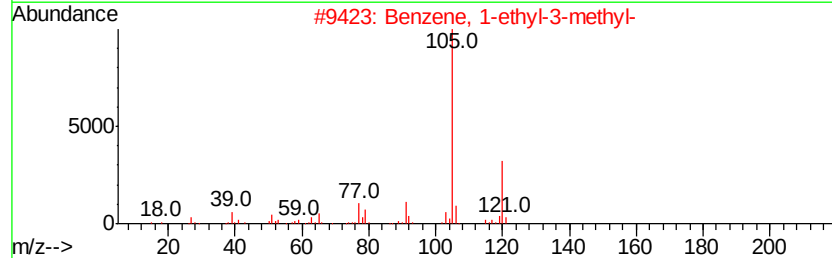
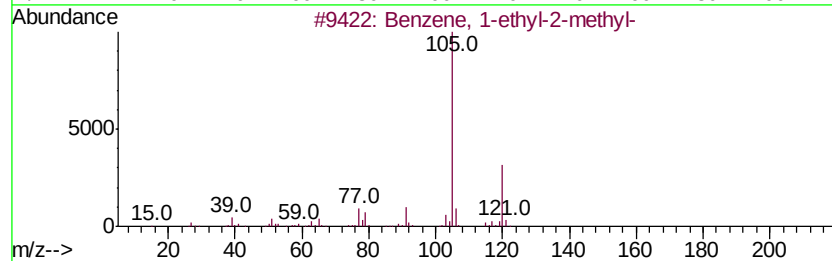
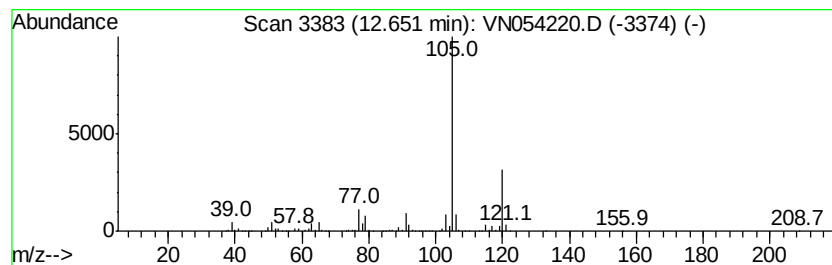
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	19.03 ug/l	3257230	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Mesitylene	120	C9H12	000108-67-8	91
5		Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031319\
Data File : VN054220.D
Acq On : 13 Mar 2019 18:45
Operator : JC/SP
Sample : K1842-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PR-MW-05_030819

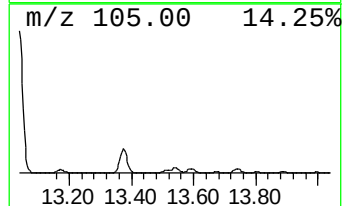
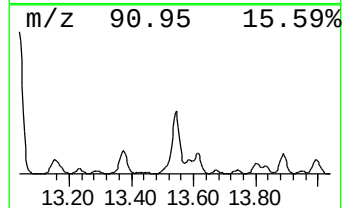
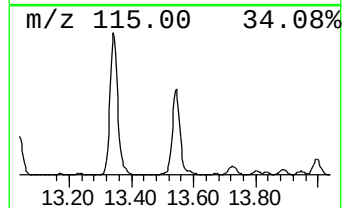
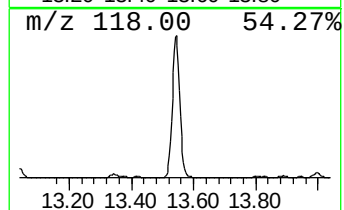
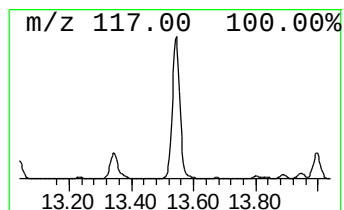
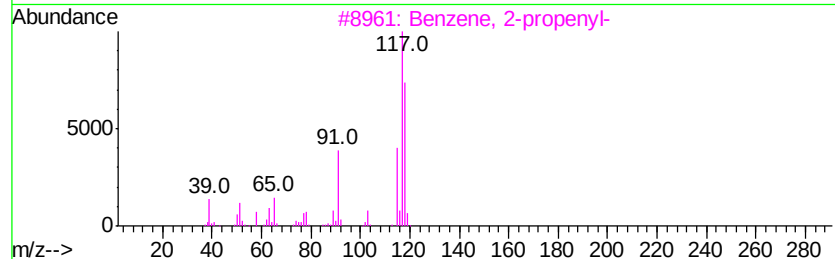
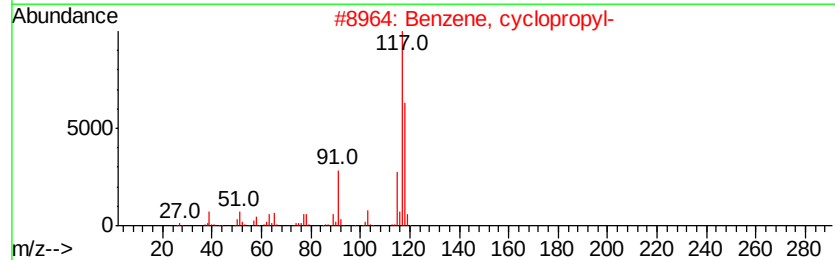
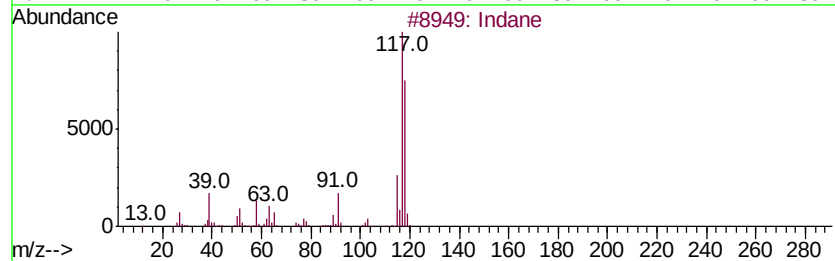
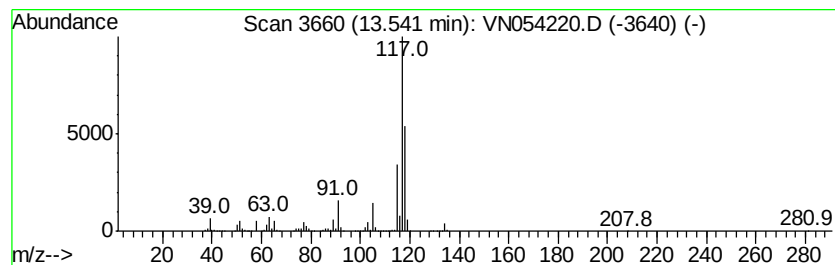
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Quant Title : SW846 8260

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

Peak Number 7 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	23.64 ug/l	4047690	1,4-Dichlorobenzene-d4	13.34

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	90
2	Benzene, cyclopropyl-		118	C9H10	000873-49-4	81
3	Benzene, 2-propenyl-		118	C9H10	000300-57-2	81
4	Benzene, 1-propenyl-		118	C9H10	000637-50-3	81
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...		118	C9H10	1000191-13-7	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN031319\
Data File : VN054220.D
Acq On : 13 Mar 2019 18:45
Operator : JC/SP
Sample : K1842-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PR-MW-05_030819

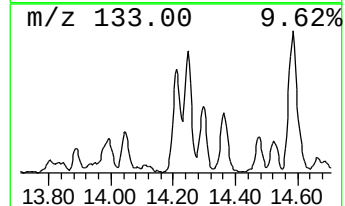
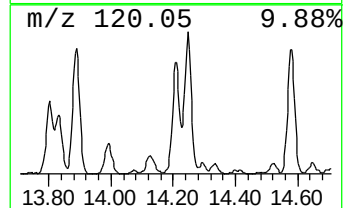
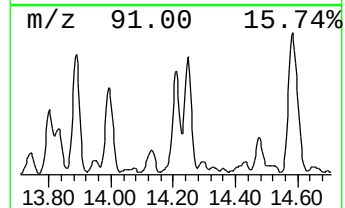
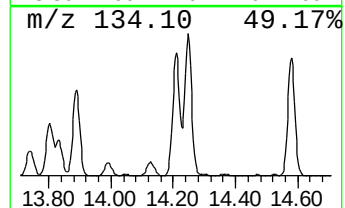
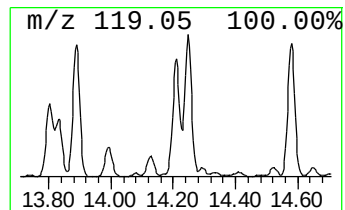
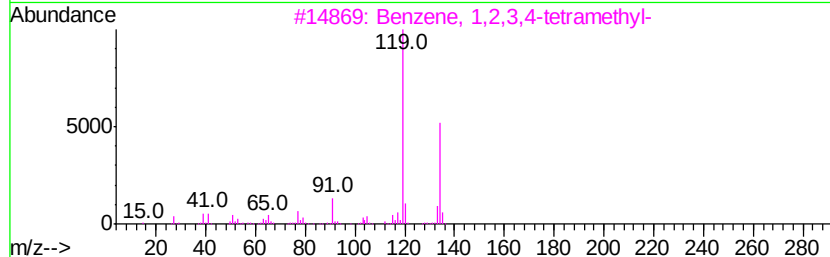
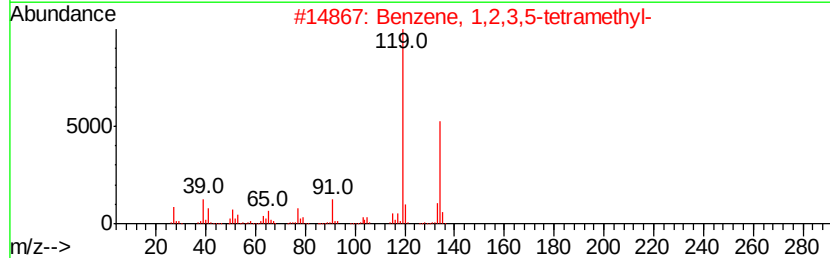
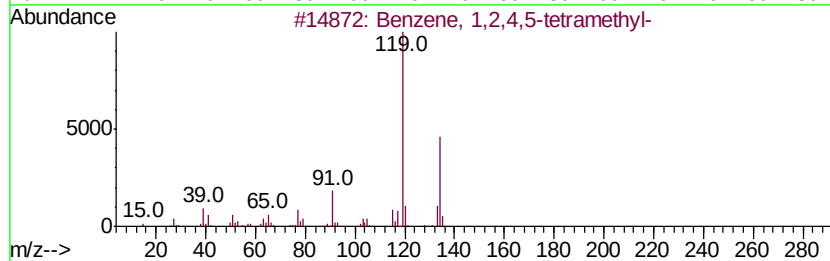
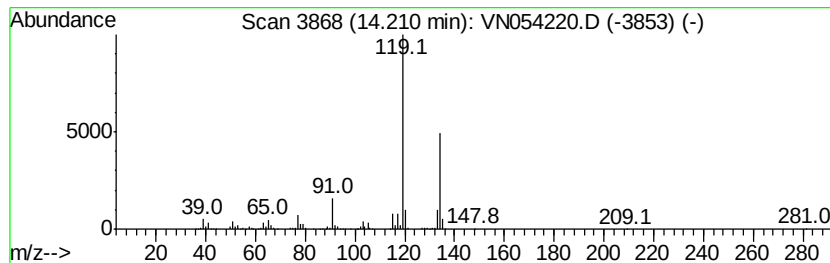
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	5.17 ug/l	885441	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		o-Cymene	134	C10H14	000527-84-4	91



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031319\
Data File : VN054220.D
Acq On : 13 Mar 2019 18:45
Operator : JC/SP
Sample : K1842-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

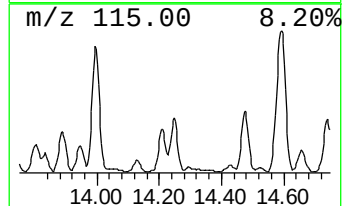
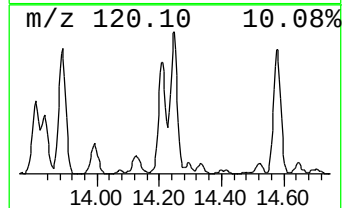
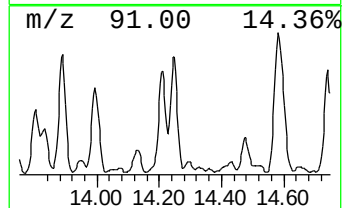
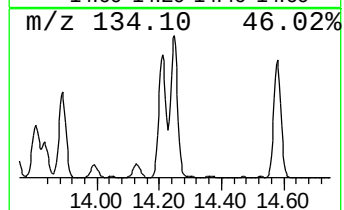
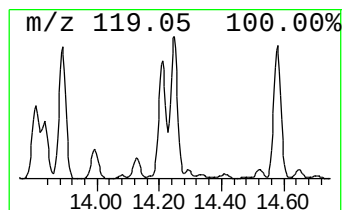
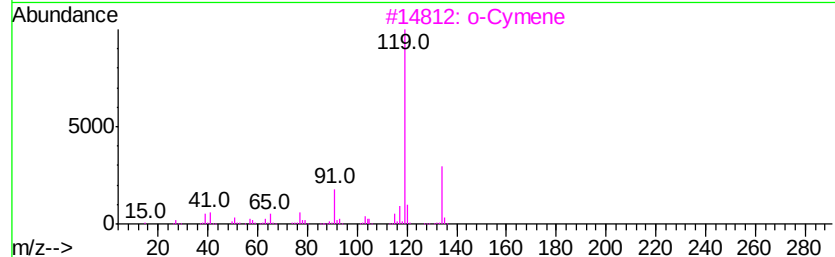
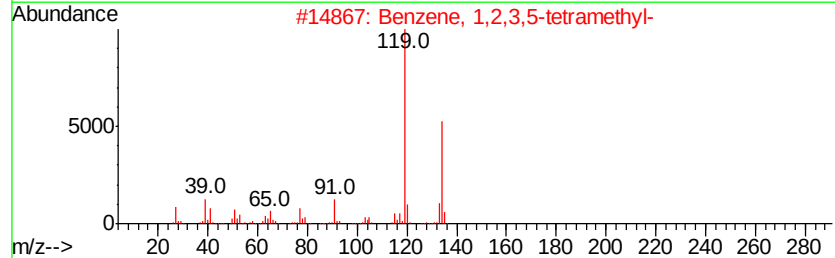
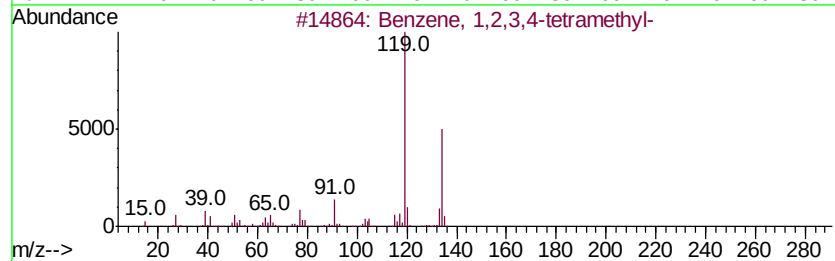
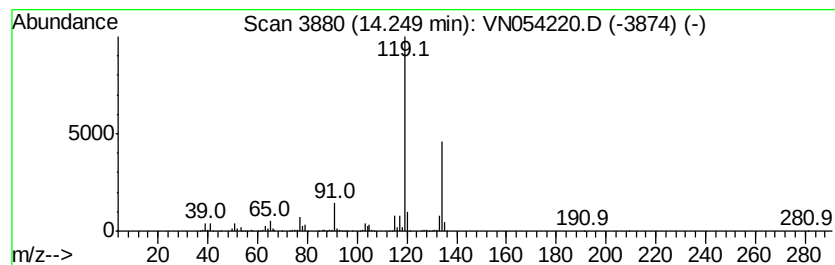
Instrument :
MSVOA_N
ClientSampleId :
PR-MW-05_030819

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	5.26 ug/l	901347	1,4-Dichlorobenzene-d4	13.34
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 97
2		Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7 97
3		o-Cymene	134 C10H14	000527-84-4 94
4		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 94
5		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN031319\
Data File : VN054220.D
Acq On : 13 Mar 2019 18:45
Operator : JC/SP
Sample : K1842-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
PR-MW-05_030819

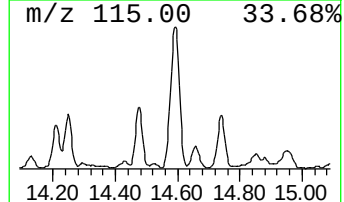
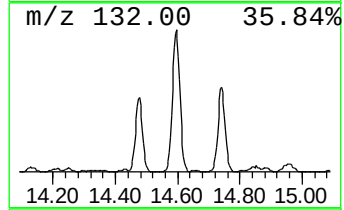
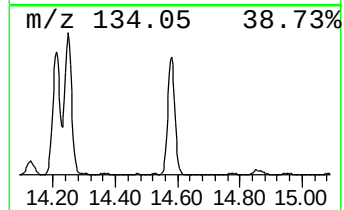
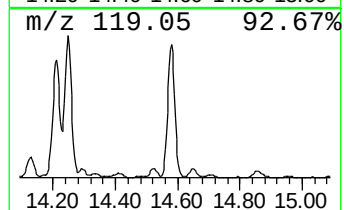
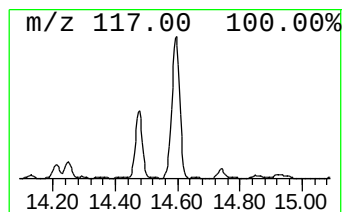
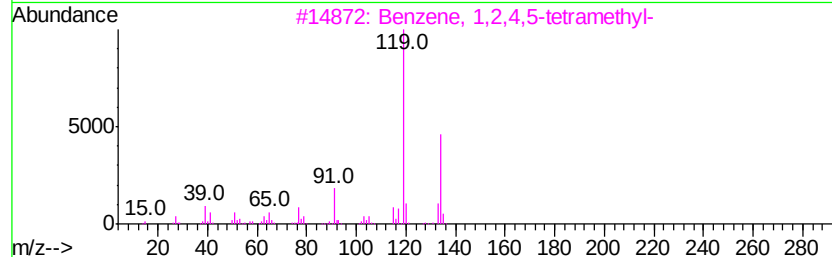
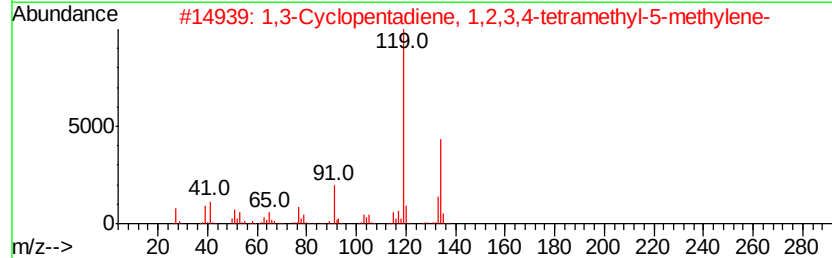
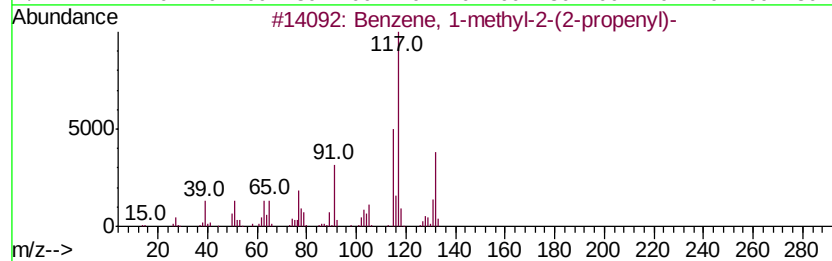
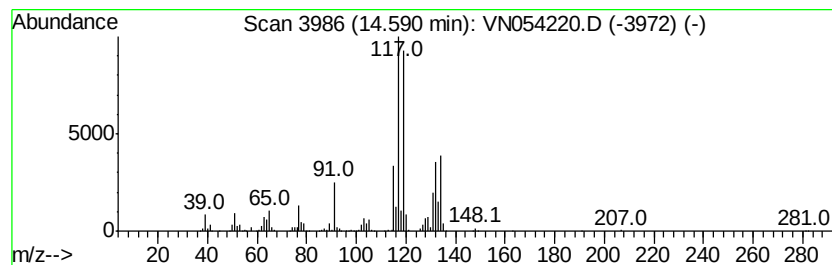
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-methyl-2-(2-prop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	10.29 ug/l	1761200	1,4-Dichlorobenzene-d4	13.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
2		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	64
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4		p-Cymene	134	C10H14	000099-87-6	60
5		o-Cymene	134	C10H14	000527-84-4	60



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031319\
 Data File : VN054220.D
 Acq On : 13 Mar 2019 18:45
 Operator : JC/SP
 Sample : K1842-12
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PR-MW-05_030819

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.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
Butane, 2-methyl-	2.81	16.9	ug/l	1080230	1	7.66	3192210	50.0
Pentane, 2-methyl-	4.58	8.4	ug/l	534328	1	7.66	3192210	50.0
Pentane, 3-methyl-	5.05	8.6	ug/l	547351	1	7.66	3192210	50.0
Cyclopentane, met...	6.63	17.6	ug/l	1121610	1	7.66	3192210	50.0
Hexane, 2,2-dimet...	8.20	8.1	ug/l	849186	2	8.59	5247650	50.0
Benzene, 1-ethyl-...	12.65	19.0	ug/l	3257230	4	13.34	8560060	50.0
Indane	13.54	23.6	ug/l	4047690	4	13.34	8560060	50.0
Benzene, 1,2,4,5-...	14.21	5.2	ug/l	885441	4	13.34	8560060	50.0
Benzene, 1,2,3,4-...	14.25	5.3	ug/l	901347	4	13.34	8560060	50.0
Benzene, 1-methyl...	14.59	10.3	ug/l	1761200	4	13.34	8560060	50.0