

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062520\
 Data File : VN062115.D
 Acq On : 25 Jun 2020 16:53
 Operator : JC/MD
 Sample : L3020-17 20X
 Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-27(16.5-18.5)DL

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\82N062420W.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	2.423	184	202	203	rBV	17947	31781	1.34%	0.118%
2	7.548	1776	1796	1807	rBV2	172780	436781	18.46%	1.621%
3	7.625	1808	1820	1836	rVV	291824	712030	30.10%	2.643%
4	7.834	1873	1885	1900	rBV5	10025	25166	1.06%	0.093%
5	7.985	1911	1932	1959	rBV	186074	437610	18.50%	1.624%
6	8.168	1979	1989	2005	rVB2	24749	64453	2.72%	0.239%
7	8.397	2047	2060	2077	rVB5	15232	33121	1.40%	0.123%
8	8.548	2091	2107	2129	rBV	472500	999784	42.26%	3.711%
9	9.046	2239	2262	2274	rBV4	64707	166415	7.03%	0.618%
10	9.111	2274	2282	2292	rVB5	18266	35236	1.49%	0.131%
11	9.329	2333	2350	2365	rBV6	11344	26669	1.13%	0.099%
12	9.487	2386	2399	2419	rBV	64183	138037	5.83%	0.512%
13	9.622	2419	2441	2454	rBV4	87862	234012	9.89%	0.869%
14	9.699	2454	2465	2474	rBV	53204	106958	4.52%	0.397%
15	9.837	2489	2508	2528	rBV2	61859	156273	6.61%	0.580%
16	9.992	2545	2556	2565	rVV	97661	184143	7.78%	0.684%
17	10.059	2565	2577	2600	rVV	823773	1551482	65.58%	5.759%
18	10.181	2600	2615	2621	rVV6	17340	38316	1.62%	0.142%
19	10.255	2622	2638	2649	rVB5	89972	191539	8.10%	0.711%
20	10.403	2676	2684	2692	rBV5	21762	34434	1.46%	0.128%
21	10.493	2701	2712	2735	rBV8	55333	166024	7.02%	0.616%
22	10.596	2736	2744	2755	rVB5	26251	47937	2.03%	0.178%
23	10.693	2755	2774	2783	rBV3	41097	80544	3.40%	0.299%
24	10.792	2798	2805	2817	rVB2	48260	83139	3.51%	0.309%
25	10.889	2830	2835	2839	rVV3	22183	30058	1.27%	0.112%
26	10.924	2841	2846	2853	rVV	37075	50516	2.14%	0.188%
27	10.975	2854	2862	2872	rVB4	36046	61587	2.60%	0.229%
28	11.033	2873	2880	2889	rVB6	20902	35713	1.51%	0.133%
29	11.094	2890	2899	2907	rBV2	34750	55296	2.34%	0.205%
30	11.175	2907	2924	2943	rVB6	168832	409860	17.32%	1.521%
31	11.274	2944	2955	2973	rBV	133482	251470	10.63%	0.933%
32	11.381	2974	2988	3004	rBV	752527	1358278	57.41%	5.042%
33	11.483	3006	3020	3036	rVB	288939	555984	23.50%	2.064%
34	11.602	3037	3057	3075	rBV4	177055	483617	20.44%	1.795%

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35	11.786	3104	3114	3119	rBV8	15703	30848	1.30%	0.115%
36	11.831	3120	3128	3136	rVV4	42549	71374	3.02%	0.265%
37	11.901	3138	3150	3161	rVV7	61125	141678	5.99%	0.526%
38	12.020	3180	3187	3196	rVB	64174	91194	3.85%	0.339%
39	12.091	3201	3209	3216	rVV4	53454	95074	4.02%	0.353%
40	12.136	3217	3223	3232	rVB6	50091	80253	3.39%	0.298%
41	12.223	3243	3250	3256	rBV	51529	79584	3.36%	0.295%
42	12.374	3282	3297	3307	rVV2	570147	1130774	47.79%	4.198%
43	12.419	3307	3311	3319	rVB2	87417	110113	4.65%	0.409%
44	12.525	3336	3344	3346	rBV2	41123	56286	2.38%	0.209%
45	12.564	3348	3356	3368	rVB	276445	451423	19.08%	1.676%
46	12.663	3377	3387	3393	rBV	294029	461521	19.51%	1.713%
47	12.709	3393	3401	3424	rVB	516349	896964	37.91%	3.330%
48	12.866	3440	3450	3456	rBV	95490	169315	7.16%	0.629%
49	12.937	3467	3472	3476	rBV4	39518	49315	2.08%	0.183%
50	13.014	3486	3496	3507	rVB	1493300	2365898	100.00%	8.782%
51	13.069	3507	3513	3522	rVB3	69025	106972	4.52%	0.397%
52	13.130	3523	3532	3533	rBV4	61106	78351	3.31%	0.291%
53	13.178	3542	3547	3554	rVB2	66782	81978	3.46%	0.304%
54	13.262	3566	3573	3578	rBV4	80699	113127	4.78%	0.420%
55	13.319	3579	3591	3596	rBV	715859	1268996	53.64%	4.711%
56	13.487	3634	3643	3646	rBV	176661	245488	10.38%	0.911%
57	13.519	3647	3653	3659	rVV2	365525	621765	26.28%	2.308%
58	13.561	3659	3666	3684	rVB4	521492	1052550	44.49%	3.907%
59	13.647	3686	3693	3699	rBV4	45006	63741	2.69%	0.237%
60	13.715	3707	3714	3724	rVB	127067	197845	8.36%	0.734%
61	13.776	3724	3733	3738	rBV	240806	383057	16.19%	1.422%
62	13.863	3751	3760	3770	rBV	503510	769482	32.52%	2.856%
63	13.917	3772	3777	3784	rVB2	76940	106515	4.50%	0.395%
64	13.969	3784	3793	3803	rVB	328379	519763	21.97%	1.929%
65	14.036	3806	3814	3825	rVB6	46185	90065	3.81%	0.334%
66	14.101	3826	3834	3841	rBV2	106338	187753	7.94%	0.697%
67	14.184	3851	3860	3865	rBV	224768	363547	15.37%	1.350%
68	14.223	3866	3872	3880	rVV	291471	438112	18.52%	1.626%
69	14.271	3880	3887	3894	rVV3	197244	296655	12.54%	1.101%
70	14.307	3894	3898	3910	rVB2	110240	149863	6.33%	0.556%
71	14.409	3923	3930	3935	rBV2	77815	102420	4.33%	0.380%

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72	14.451	3935	3943	3952	rVV2	268515	452014	19.11%	1.678%
73	14.499	3953	3958	3966	rVB	117409	165327	6.99%	0.614%
74	14.567	3966	3979	3989	rBV5	453251	933256	39.45%	3.464%
75	14.622	3990	3996	4007	rVB4	125308	215667	9.12%	0.801%
76	14.824	4042	4059	4066	rBV3	204181	467160	19.75%	1.734%
77	14.930	4082	4092	4112	rVB3	173377	383161	16.20%	1.422%
78	15.104	4129	4146	4158	rBV	199792	437720	18.50%	1.625%
79	15.213	4171	4180	4188	rBV4	76457	135996	5.75%	0.505%
80	15.258	4189	4194	4202	rVB2	36093	52544	2.22%	0.195%
81	15.387	4223	4234	4249	rBV2	98030	202433	8.56%	0.751%
82	15.544	4271	4283	4302	rVB5	51938	160846	6.80%	0.597%
83	15.663	4314	4320	4330	rVB4	24164	41324	1.75%	0.153%
84	15.737	4332	4343	4352	rVB6	37168	74369	3.14%	0.276%
85	15.876	4376	4386	4395	rBV9	15845	29703	1.26%	0.110%
86	16.123	4452	4463	4490	rVB	119794	263539	11.14%	0.978%
87	16.313	4509	4522	4539	rVB3	58237	129801	5.49%	0.482%

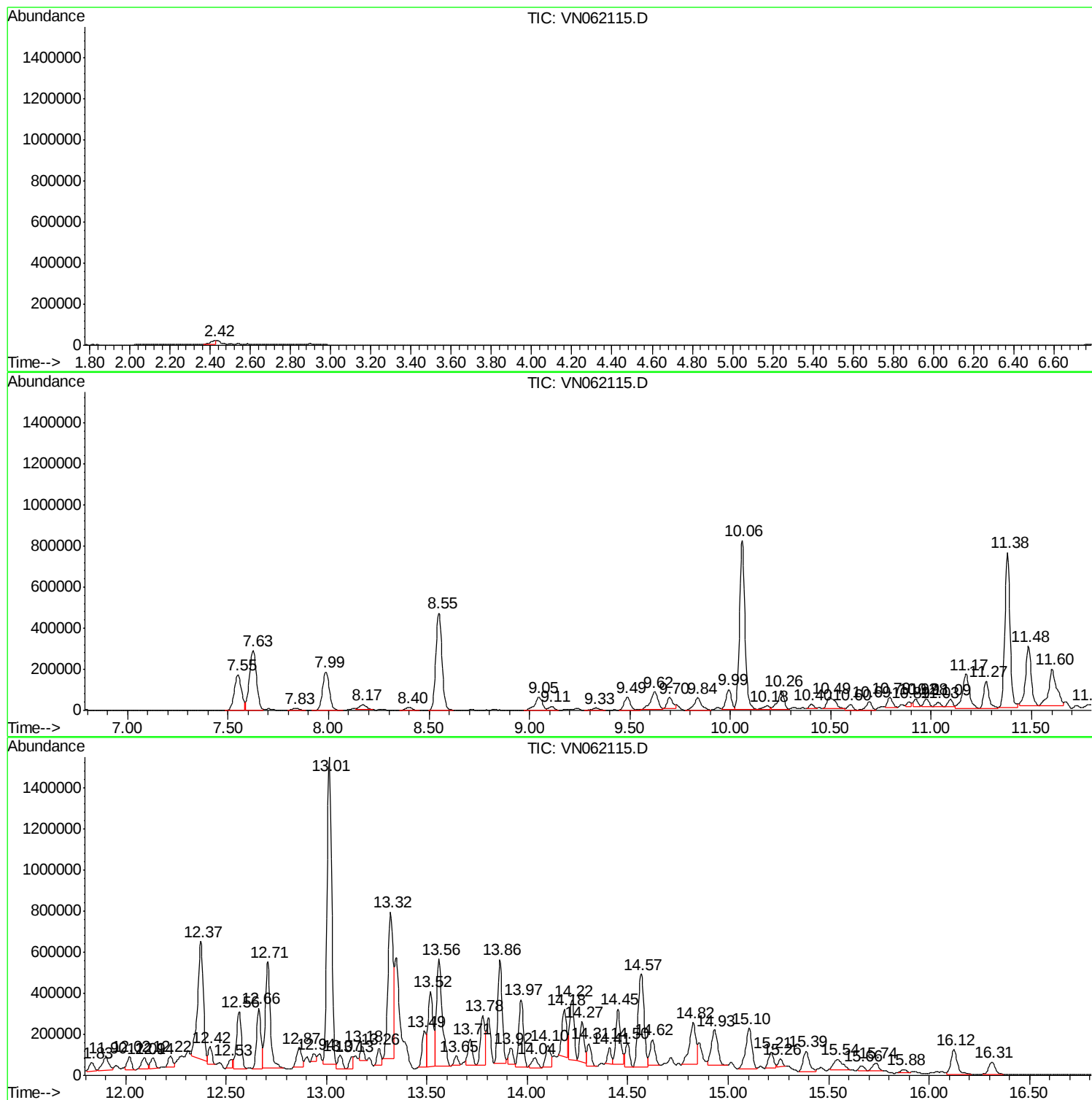
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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



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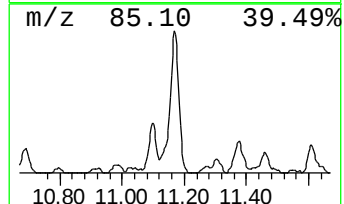
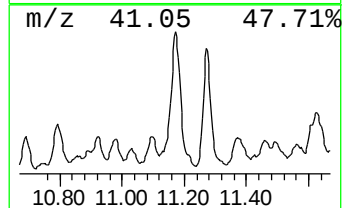
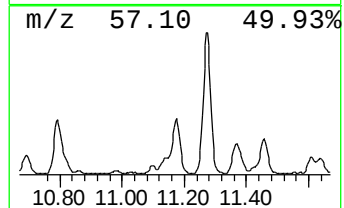
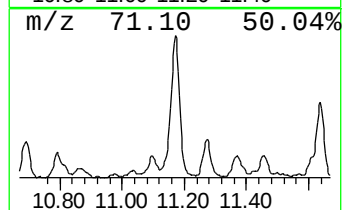
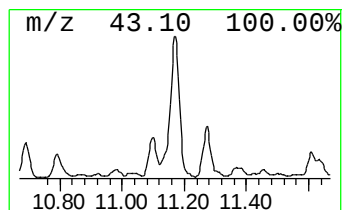
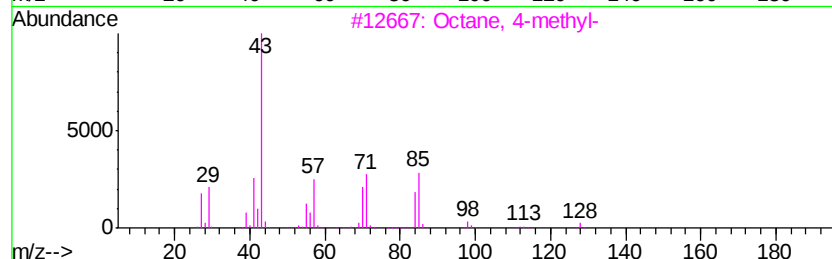
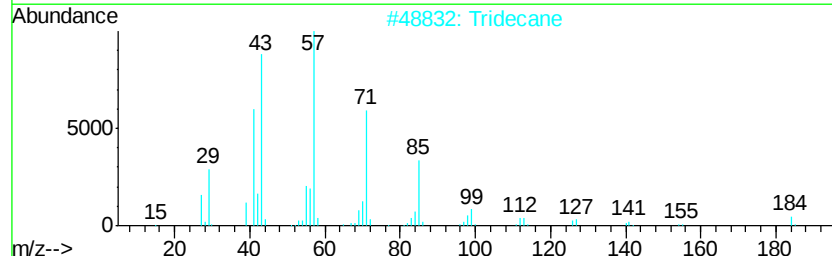
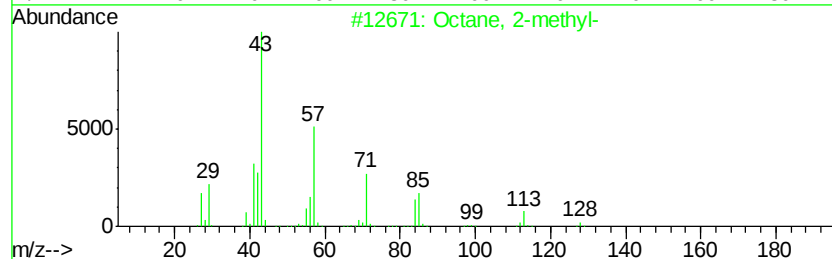
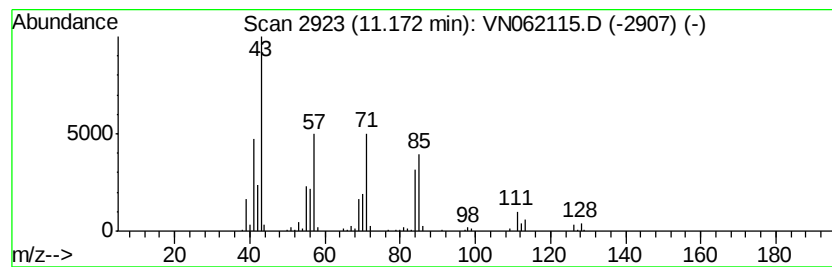
Instrument :
MSVOA_N
ClientSampleId :
MW-27(16.5-18.5)DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.17	15.09 ug/l	409860	Chlorobenzene-d5	11.38
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octane, 2-methyl-	128 C9H20	003221-61-2	83
2	Tridecane	184 C13H28	000629-50-5	59
3	Octane, 4-methyl-	128 C9H20	002216-34-4	53
4	Oxirane, 2-methyl-3-propyl-, cis-	100 C6H12O	006124-90-9	47
5	Sulfurous acid, 2-propyl tetradec...	320 C17H36O3S	1000309-12-5	47



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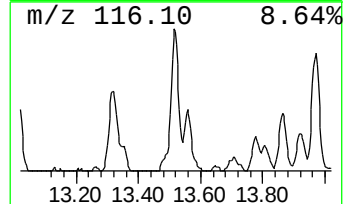
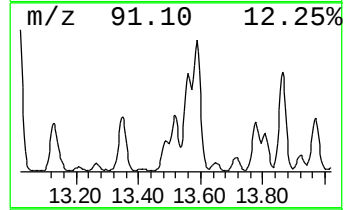
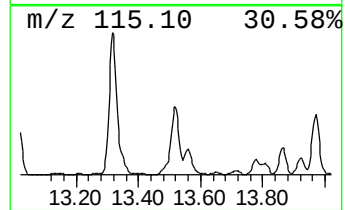
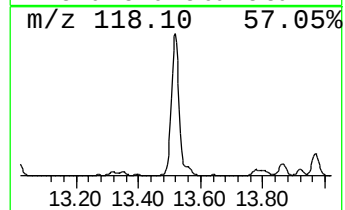
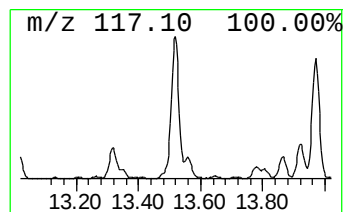
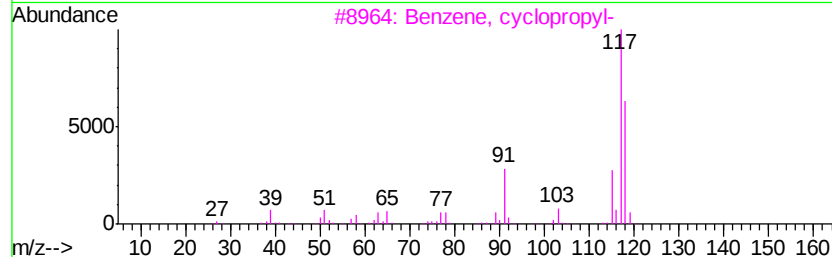
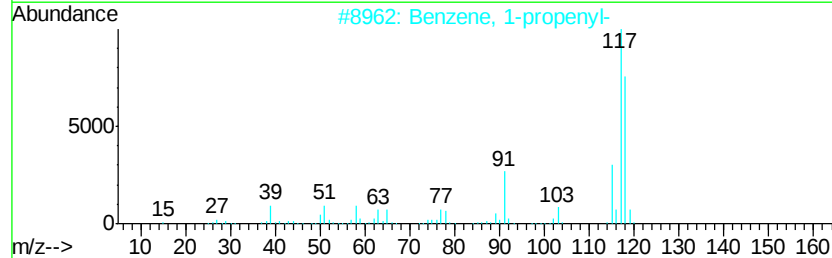
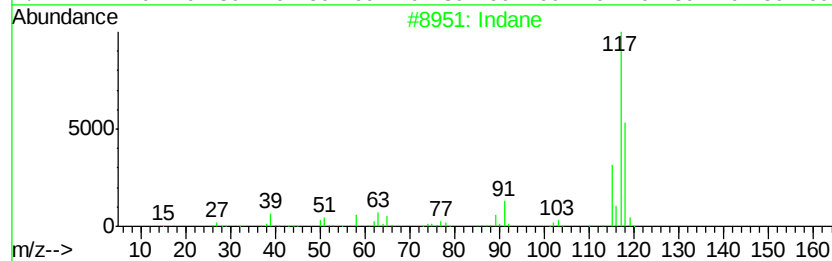
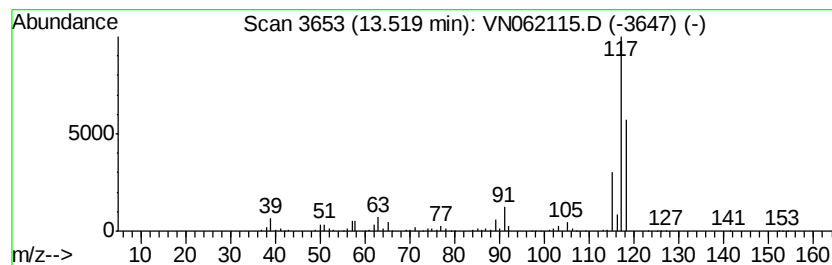
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	24.50 ug/l	621765	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 93
2	Benzene, 1-propenyl-		118 C9H10	000637-50-3 72
3	Benzene, cyclopropyl-		118 C9H10	000873-49-4 72
4	Benzene, 1-ethenyl-4-methyl-		118 C9H10	000622-97-9 72
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-		118 C9H10	1000191-13-7 64



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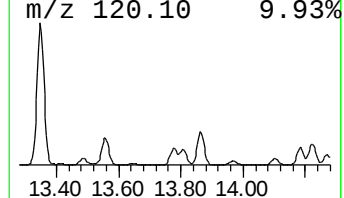
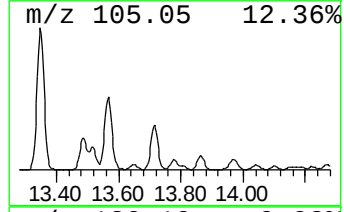
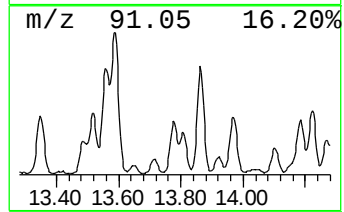
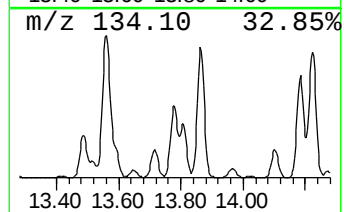
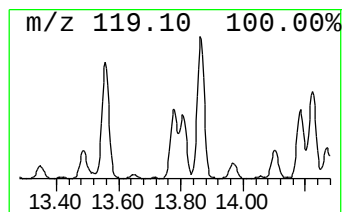
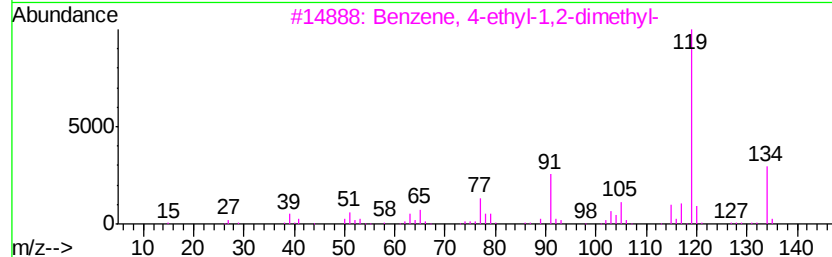
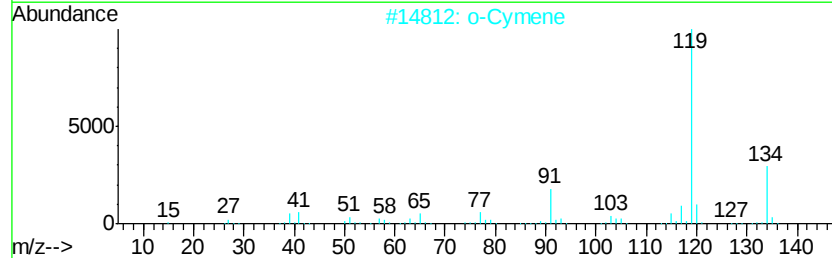
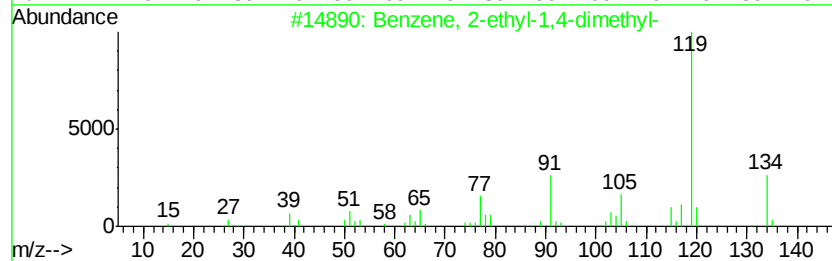
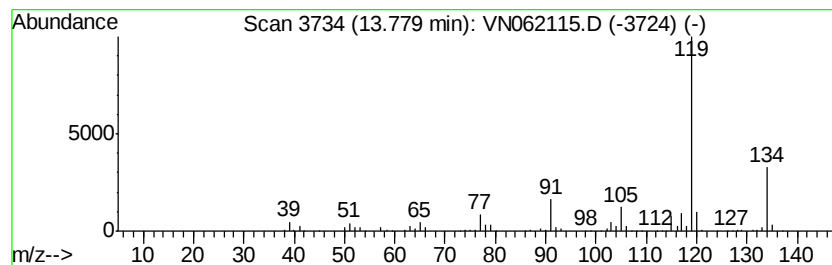
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.78	15.09 ug/l	383057	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		o-Cymene	134	C10H14	000527-84-4	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



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Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

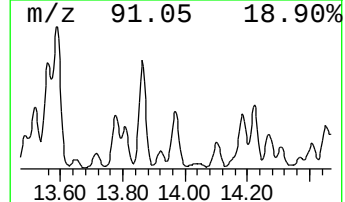
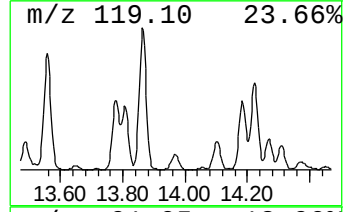
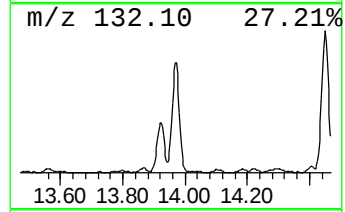
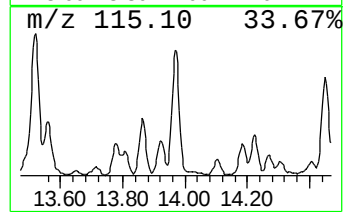
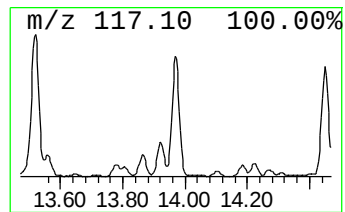
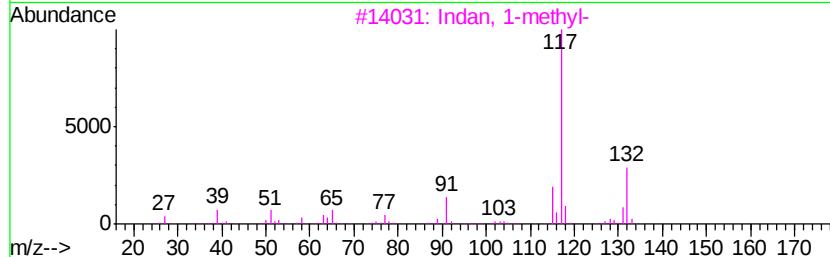
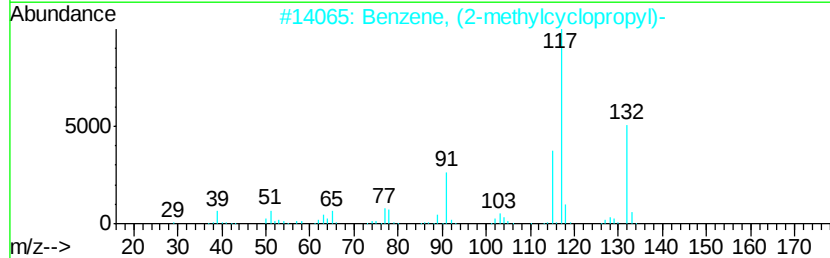
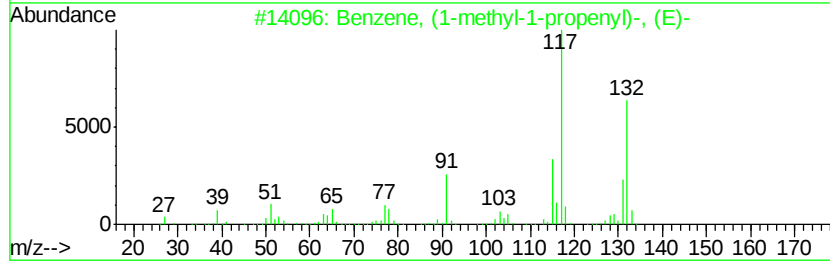
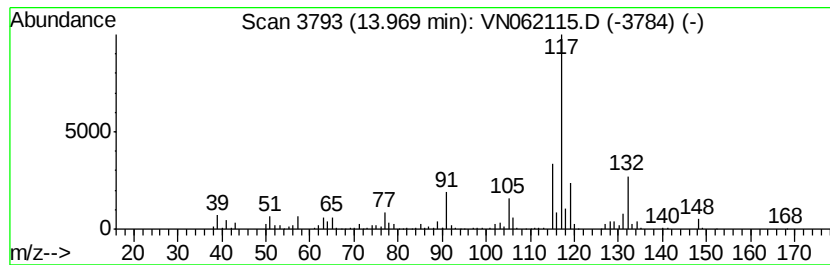
Instrument :
MSVOA_N
ClientSampleID :
MW-27(16.5-18.5)DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	20.48 ug/l	519763	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 81
2		Benzene, (2-methylcyclopropyl)-	132 C10H12	1000327-39-0 76
3		Indan, 1-methyl-	132 C10H12	000767-58-8 76
4		Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0 72
5		Benzene, 2-butenyl-	132 C10H12	001560-06-1 72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062115.D
Acq On : 25 Jun 2020 16:53
Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleID :
MW-27(16.5-18.5)DL

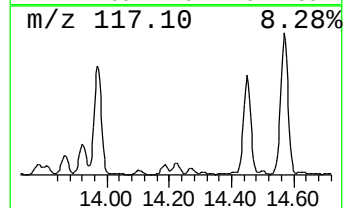
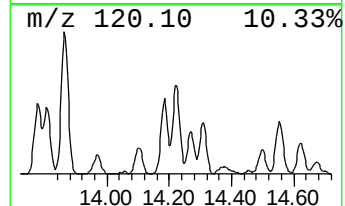
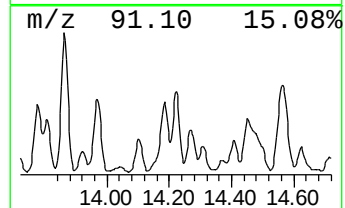
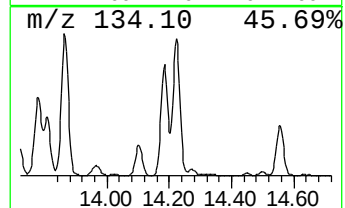
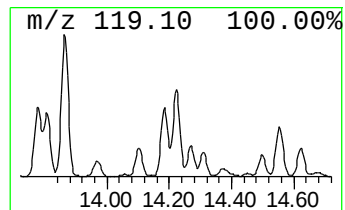
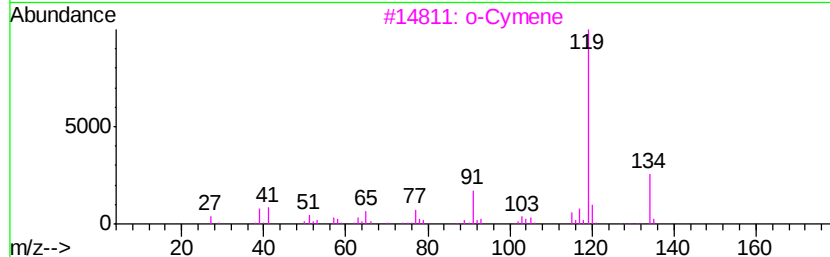
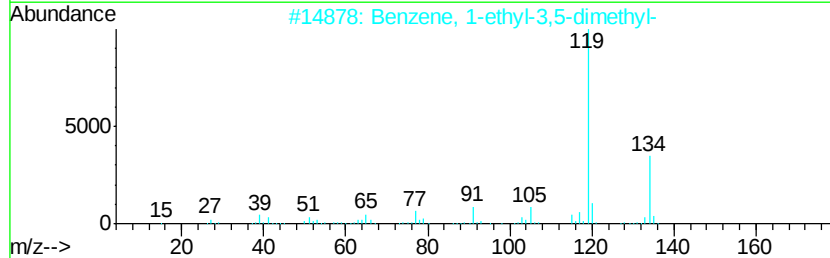
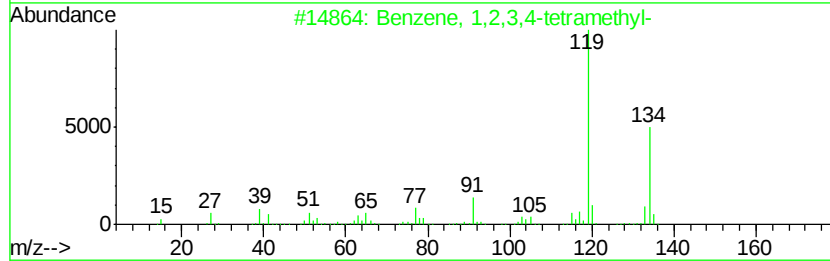
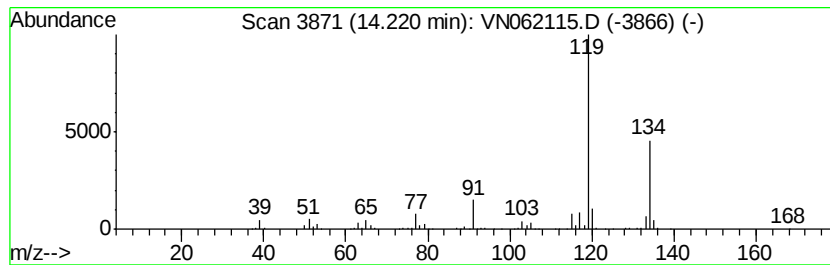
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.22	17.26 ug/l	438112	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	95
2		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
3		o-Cymene	134	C10H14	000527-84-4	94
4		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062115.D
Acq On : 25 Jun 2020 16:53
Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

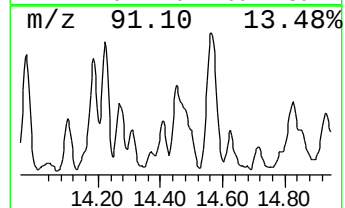
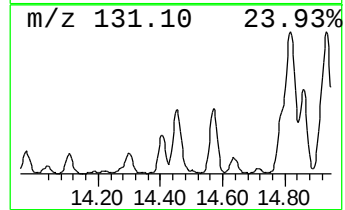
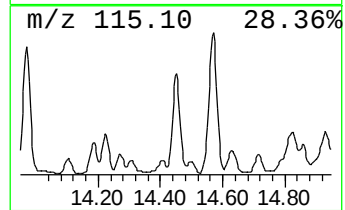
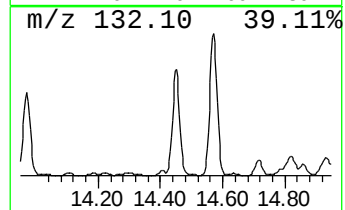
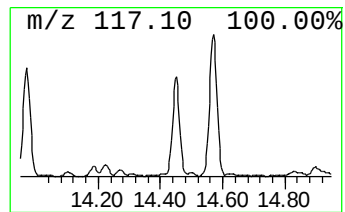
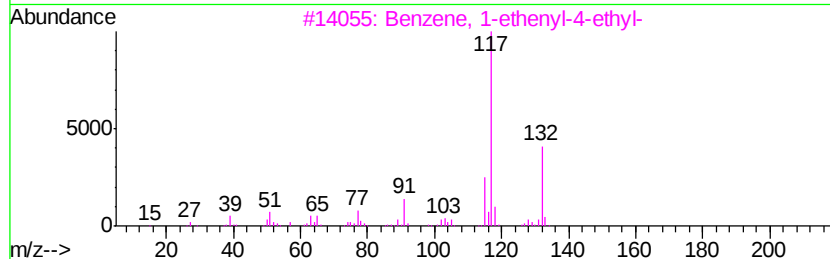
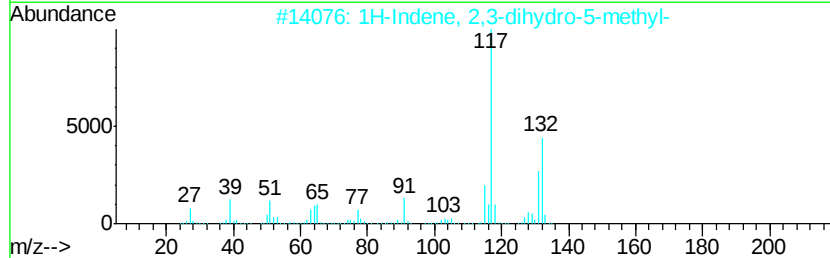
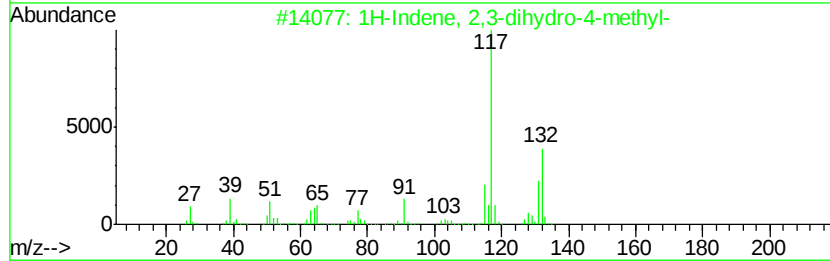
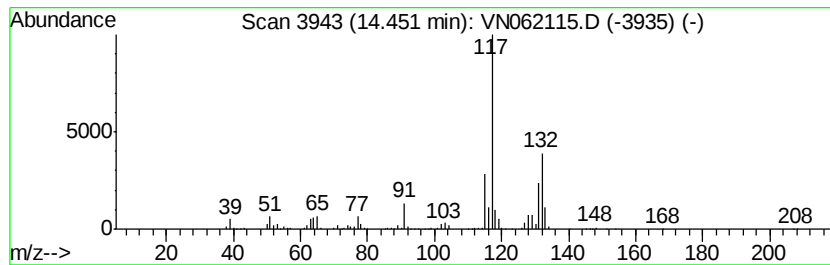
Instrument :
MSVOA_N
ClientSampleID :
MW-27(16.5-18.5)DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	17.81 ug/l	452014	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	93
2	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	87
3	Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7	76
4	Benzene, (2-methyl-1-propenyl)-	132 C10H12	000768-49-0	72
5	Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062115.D
Acq On : 25 Jun 2020 16:53
Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleID :
MW-27(16.5-18.5)DL

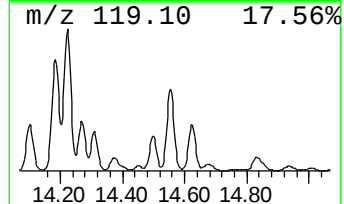
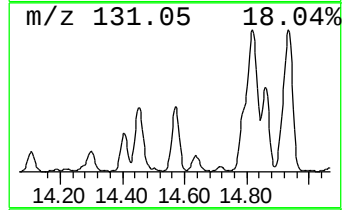
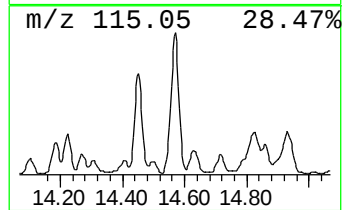
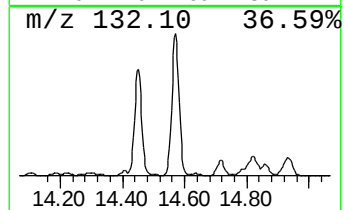
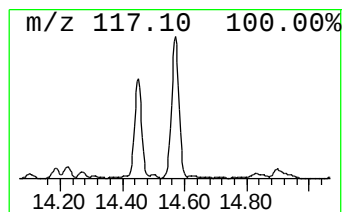
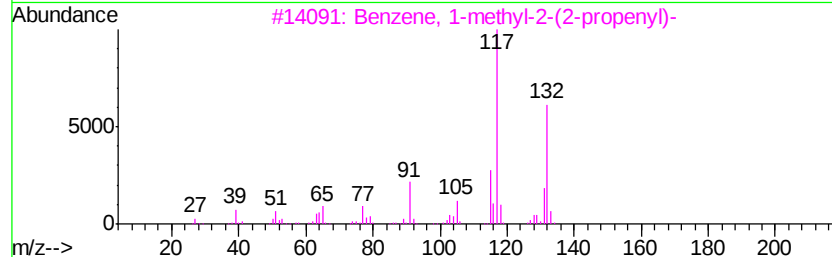
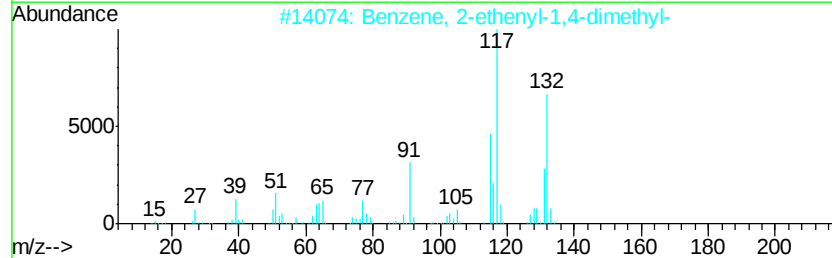
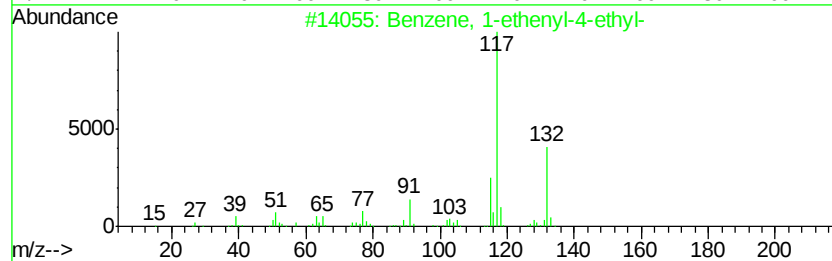
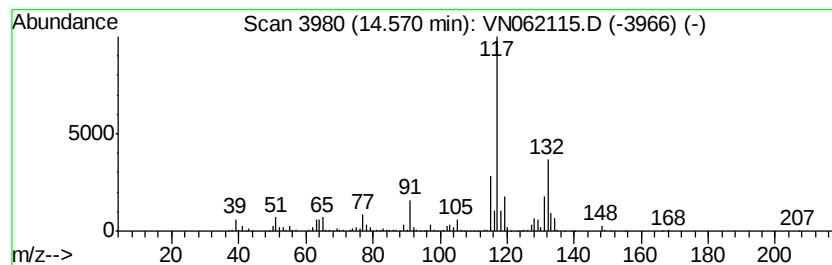
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	36.77 ug/l	933256	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	81
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062115.D
Acq On : 25 Jun 2020 16:53
Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleID :
MW-27(16.5-18.5)DL

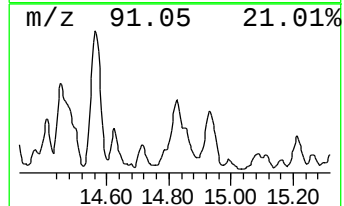
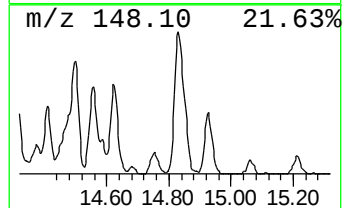
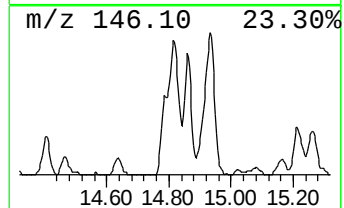
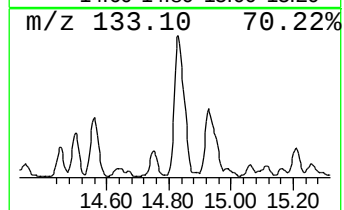
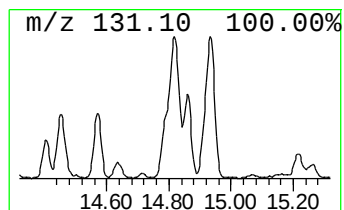
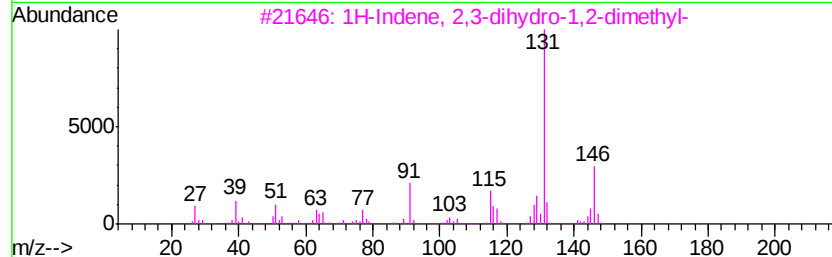
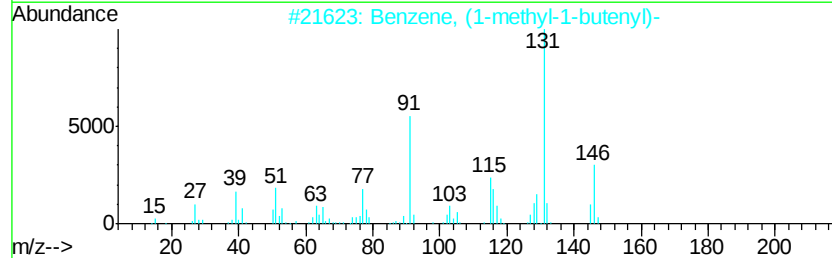
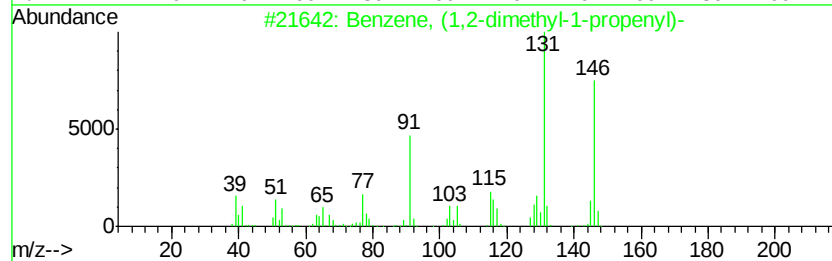
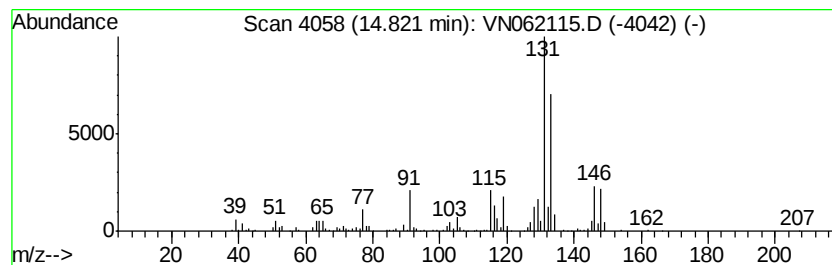
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	18.41 ug/l	467160	1,4-Dichlorobenzene-d4	13.32

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	90
3	1H-Indene, 2,3-dihydro-1,2-dimethyl-	146	C11H14	017057-82-8	89
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062520\
Data File : VN062115.D
Acq On : 25 Jun 2020 16:53
Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

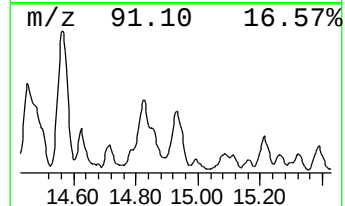
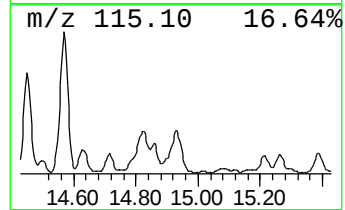
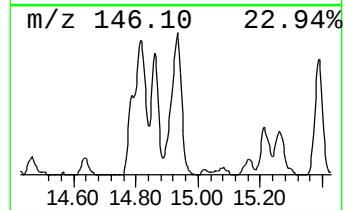
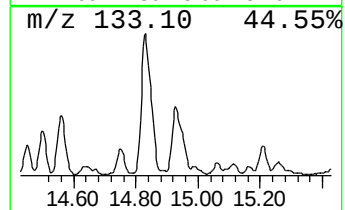
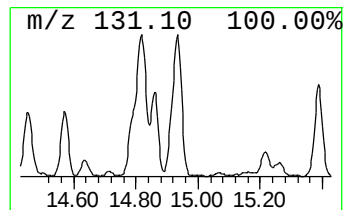
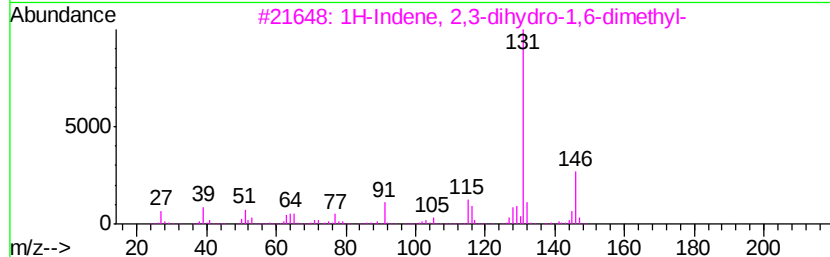
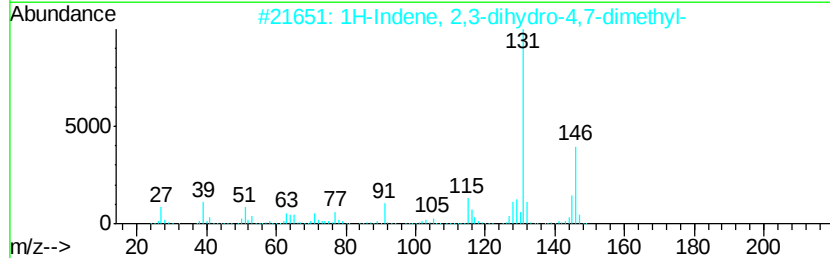
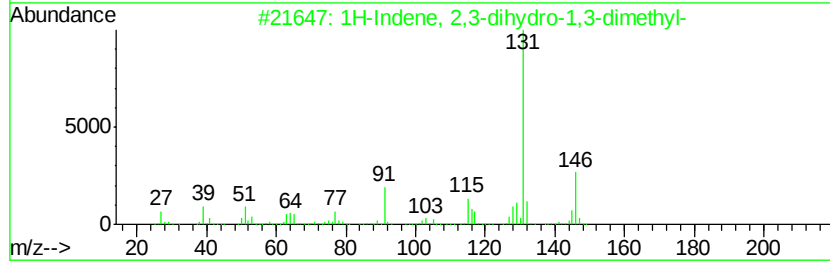
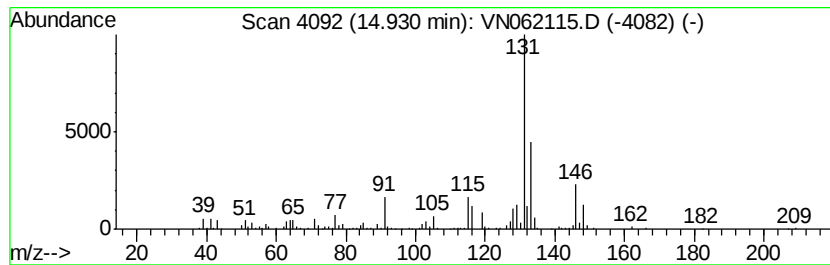
Instrument :
MSVOA_N
ClientSampleId :
MW-27(16.5-18.5)DL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.93	15.10 ug/l	383161	1,4-Dichlorobenzene-d4	13.32
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	70
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	70
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	70
4	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	70
5	Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN062520\
Data File : VN062115.D
Acq On : 25 Jun 2020 16:53
Operator : JC/MD
Sample : L3020-17 20X
Misc : 5.34g/5.0mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-27(16.5-18.5)DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.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
Octane, 2-methyl-	11.17	15.1	ug/l	409860	3	11.38	1358280	50.0
Indane	13.52	24.5	ug/l	621765	4	13.32	1269000	50.0
Benzene, 2-ethyl-...	13.78	15.1	ug/l	383057	4	13.32	1269000	50.0
Benzene, (1-methy...	13.97	20.5	ug/l	519763	4	13.32	1269000	50.0
Benzene, 1,2,3,4-...	14.22	17.3	ug/l	438112	4	13.32	1269000	50.0
1H-Indene, 2,3-di...	14.45	17.8	ug/l	452014	4	13.32	1269000	50.0
Benzene, 1-etheny...	14.57	36.8	ug/l	933256	4	13.32	1269000	50.0
Benzene, (1,2-dim...	14.82	18.4	ug/l	467160	4	13.32	1269000	50.0
1H-Indene, 2,3-di...	14.93	15.1	ug/l	383161	4	13.32	1269000	50.0