

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN062420\
 Data File : VN062075.D
 Acq On : 24 Jun 2020 20:13
 Operator : JC/MD
 Sample : L3063-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-13

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	1.793	3	6	24	rVB3	69825	84685	5.61%	0.463%
2	2.134	102	112	132	rBV2	46153	105155	6.97%	0.575%
3	2.288	152	160	171	rVB2	17170	25581	1.70%	0.140%
4	2.420	190	201	211	rVB8	18548	34928	2.31%	0.191%
5	2.770	295	310	329	rBV	146745	300342	19.90%	1.642%
6	3.021	381	388	400	rVB3	8917	17063	1.13%	0.093%
7	3.101	401	413	432	rVB6	12771	34129	2.26%	0.187%
8	3.761	597	618	644	rBV4	51957	178394	11.82%	0.975%
9	4.015	683	697	714	rBV3	12683	35555	2.36%	0.194%
10	4.484	823	843	854	rBV3	60856	205439	13.61%	1.123%
11	4.998	979	1003	1027	rBV3	62555	228145	15.12%	1.247%
12	6.474	1442	1462	1479	rBV3	28818	95973	6.36%	0.525%
13	6.580	1479	1495	1514	rVV3	56289	173613	11.50%	0.949%
14	6.661	1514	1520	1533	rVV4	6967	16121	1.07%	0.088%
15	6.776	1537	1556	1581	rVB9	8818	31472	2.09%	0.172%
16	7.349	1721	1734	1749	rVB5	10303	29598	1.96%	0.162%
17	7.548	1774	1796	1807	rBV	179281	458401	30.37%	2.505%
18	7.625	1808	1820	1834	rVV	317181	791242	52.43%	4.325%
19	7.702	1834	1844	1862	rVB3	65501	168547	11.17%	0.921%
20	7.838	1869	1886	1899	rBV3	24940	62353	4.13%	0.341%
21	7.995	1914	1935	1957	rBV3	414059	1099486	72.85%	6.009%
22	8.169	1957	1989	2009	rVV	247912	779100	51.63%	4.258%
23	8.262	2010	2018	2034	rVB4	23238	53659	3.56%	0.293%
24	8.407	2052	2063	2078	rVB4	8704	20031	1.33%	0.109%
25	8.548	2090	2107	2130	rBV	498891	1053808	69.83%	5.760%
26	8.786	2168	2181	2192	rBV3	35794	79943	5.30%	0.437%
27	9.014	2234	2252	2255	rBV3	36942	67754	4.49%	0.370%
28	9.108	2273	2281	2293	rVB4	21708	41939	2.78%	0.229%
29	9.188	2293	2306	2315	rBV2	28425	58646	3.89%	0.321%
30	9.243	2315	2323	2336	rVB6	23481	51814	3.43%	0.283%
31	9.336	2340	2352	2363	rBV6	7107	17843	1.18%	0.098%
32	9.403	2365	2373	2385	rVB5	11593	23460	1.55%	0.128%
33	9.487	2385	2399	2415	rVB	218584	433324	28.71%	2.368%
34	9.622	2418	2441	2456	rBV2	411964	950091	62.96%	5.193%

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Integrator: RTE
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35	9.728	2469	2474	2488	rVB6	17739	33761	2.24%	0.185%
36	9.815	2488	2501	2513	rBV8	16852	41801	2.77%	0.228%
37	9.879	2513	2521	2531	rVB5	9405	19112	1.27%	0.104%
38	9.995	2544	2557	2565	rVV4	93786	170635	11.31%	0.933%
39	10.059	2565	2577	2590	rVV	814319	1509147	100.00%	8.248%
40	10.124	2592	2597	2607	rVB2	34278	54081	3.58%	0.296%
41	10.223	2621	2628	2634	rBV9	10237	18471	1.22%	0.101%
42	10.355	2660	2669	2677	rBV2	14410	23659	1.57%	0.129%
43	10.407	2677	2685	2694	rVB5	10527	17300	1.15%	0.095%
44	10.490	2700	2711	2730	rVV6	38726	90241	5.98%	0.493%
45	10.577	2730	2738	2753	rVB5	12135	26765	1.77%	0.146%
46	10.767	2781	2797	2812	rBV2	35696	79711	5.28%	0.436%
47	10.889	2827	2835	2842	rBV8	16452	27364	1.81%	0.150%
48	11.381	2972	2988	3005	rBV	745561	1313317	87.02%	7.178%
49	11.480	3012	3019	3038	rVB4	17896	39662	2.63%	0.217%
50	11.751	3088	3103	3104	rBV4	17187	28960	1.92%	0.158%
51	12.223	3237	3250	3269	rBV	420206	690237	45.74%	3.773%
52	12.378	3278	3298	3312	rBV	579023	1109522	73.52%	6.064%
53	12.567	3345	3357	3369	rBV	349929	591673	39.21%	3.234%
54	12.866	3437	3450	3461	rBV	91671	149935	9.94%	0.819%
55	13.014	3485	3496	3507	rBV	88326	139376	9.24%	0.762%
56	13.146	3519	3537	3548	rBV4	40327	108054	7.16%	0.591%
57	13.320	3579	3591	3613	rVB	700778	1222752	81.02%	6.683%
58	13.426	3613	3624	3633	rBV7	12663	24996	1.66%	0.137%
59	13.519	3633	3653	3662	rVV2	441236	890256	58.99%	4.866%
60	13.567	3662	3668	3684	rVV3	90741	176721	11.71%	0.966%
61	13.648	3684	3693	3703	rVV4	30209	56347	3.73%	0.308%
62	13.715	3705	3714	3725	rVB	57770	100966	6.69%	0.552%
63	13.776	3725	3733	3750	rVB2	34679	62842	4.16%	0.343%
64	13.866	3750	3761	3770	rBV	50986	79120	5.24%	0.432%
65	13.921	3770	3778	3784	rBV2	26079	43284	2.87%	0.237%
66	13.969	3784	3793	3805	rVB2	120123	198660	13.16%	1.086%
67	14.104	3827	3835	3844	rBV7	11520	18362	1.22%	0.100%
68	14.185	3847	3860	3866	rVV	104703	166742	11.05%	0.911%
69	14.223	3867	3872	3880	rVV	62587	93219	6.18%	0.510%
70	14.268	3880	3886	3893	rVB3	12676	17426	1.15%	0.095%
71	14.342	3902	3909	3918	rVB2	32585	44798	2.97%	0.245%

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Integration Parameters: RTEINT.P

Integrator: RTE
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72	14.451	3934	3943	3953	rBV	79956	126764	8.40%	0.693%
73	14.500	3953	3958	3965	rVB2	14710	18181	1.20%	0.099%
74	14.570	3965	3980	3992	rBV3	187135	377906	25.04%	2.066%
75	14.828	4049	4060	4079	rVB5	44913	123671	8.19%	0.676%
76	14.934	4080	4093	4111	rVB3	40288	98834	6.55%	0.540%
77	15.107	4135	4147	4156	rVB	22830	41320	2.74%	0.226%
78	15.213	4172	4180	4187	rBV8	12928	23694	1.57%	0.130%
79	15.258	4189	4194	4212	rVB4	11796	25307	1.68%	0.138%
80	15.387	4223	4234	4246	rBV2	24175	47755	3.16%	0.261%
81	15.548	4271	4284	4301	rBV5	14301	39934	2.65%	0.218%
82	15.664	4311	4320	4330	rBV2	14056	24925	1.65%	0.136%
83	16.127	4453	4464	4475	rVB2	14829	27504	1.82%	0.150%
84	16.313	4509	4522	4538	rVB5	15336	33396	2.21%	0.183%

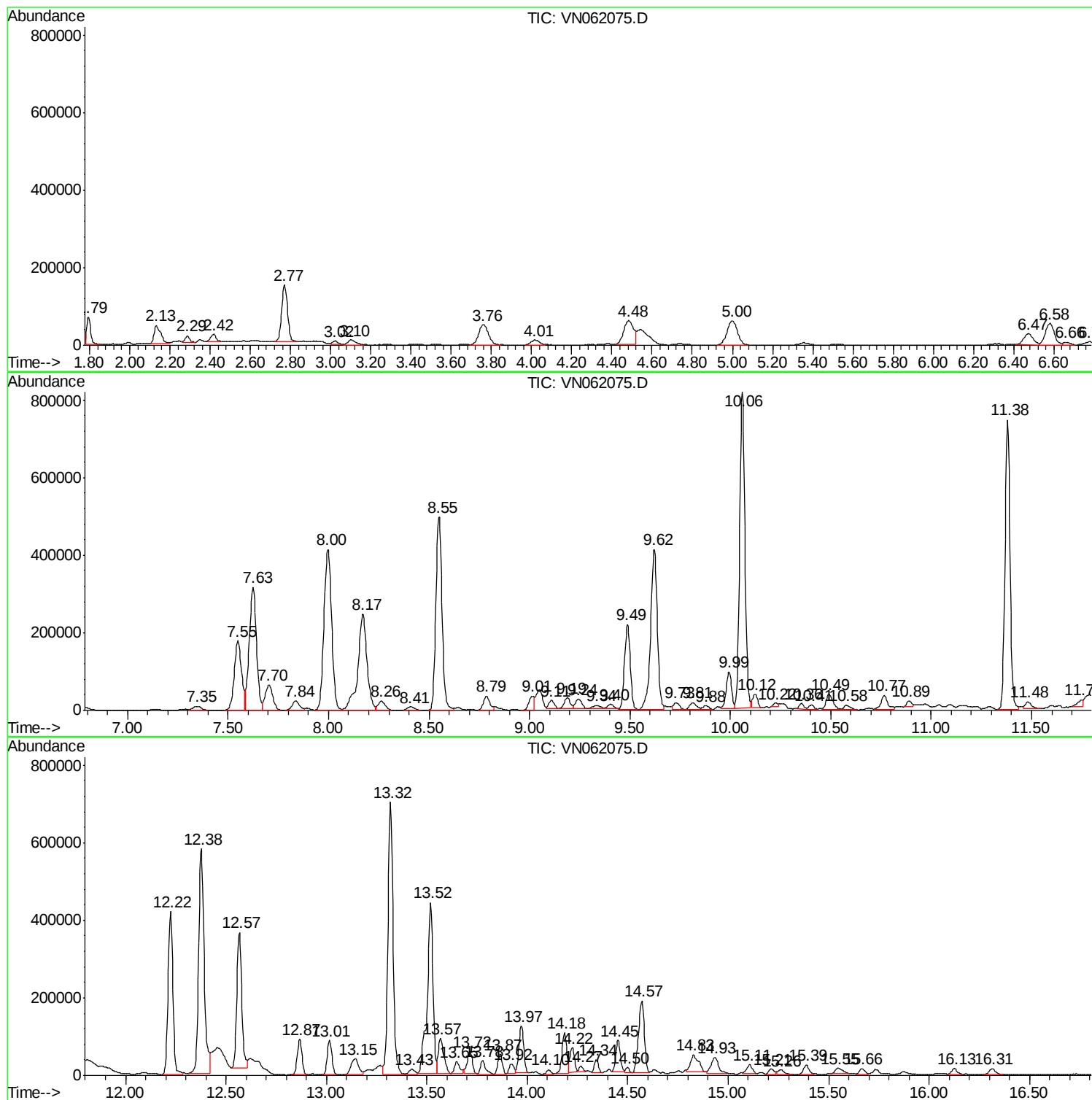
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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



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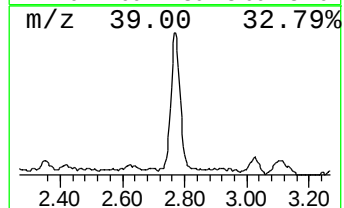
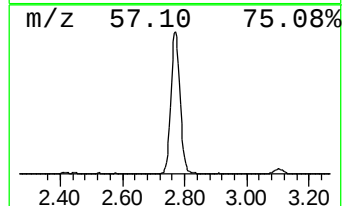
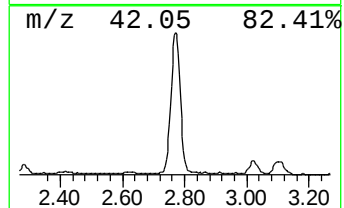
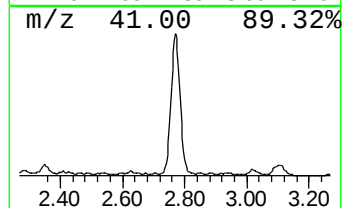
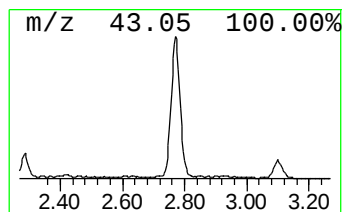
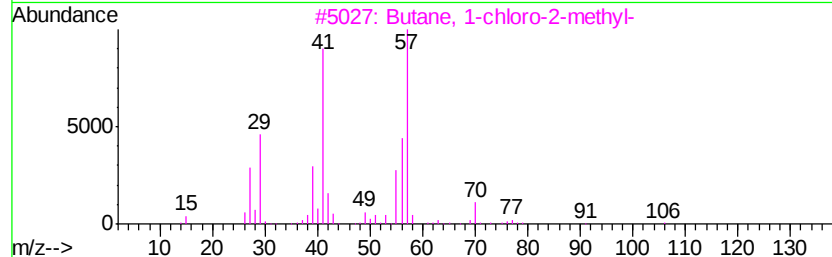
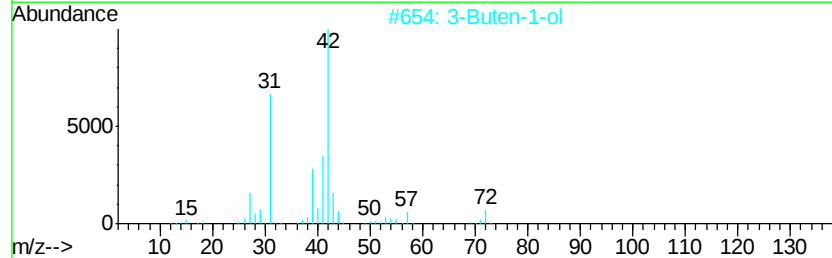
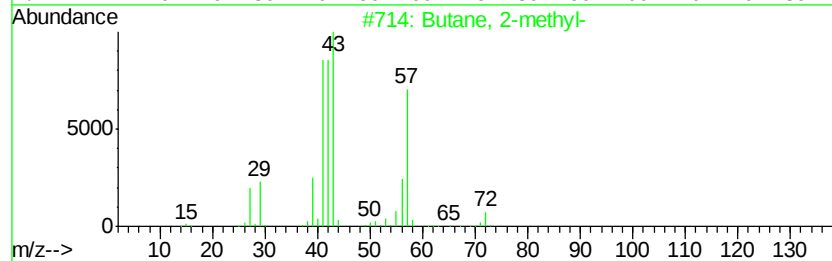
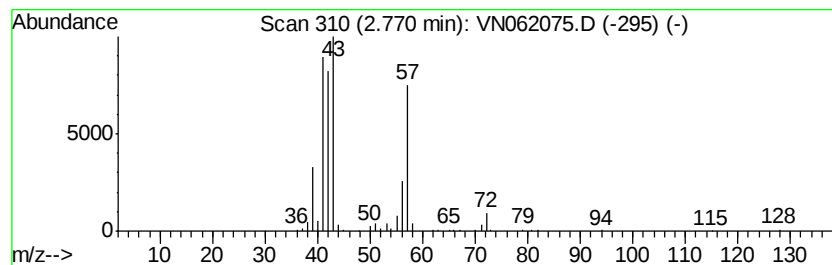
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.77	18.98 ug/l	300342	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		3-Buten-1-ol	72	C4H8O	000627-27-0	43
3		Butane, 1-chloro-2-methyl-	106	C5H11Cl	000616-13-7	35
4		Oxirane, trimethyl-	86	C5H10O	005076-19-7	33
5		Pentane	72	C5H12	000109-66-0	28



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Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

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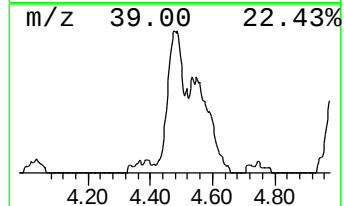
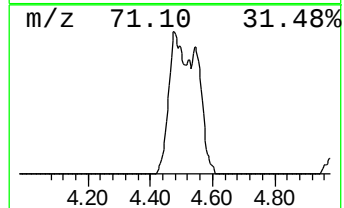
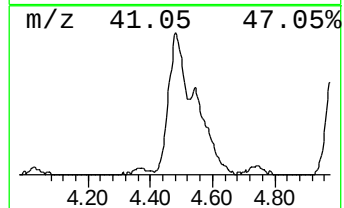
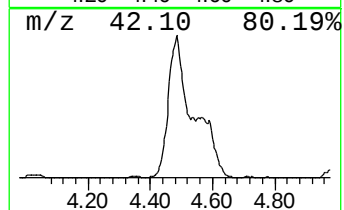
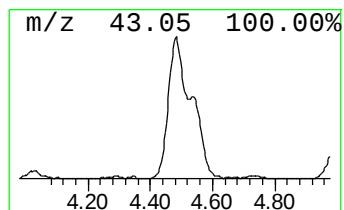
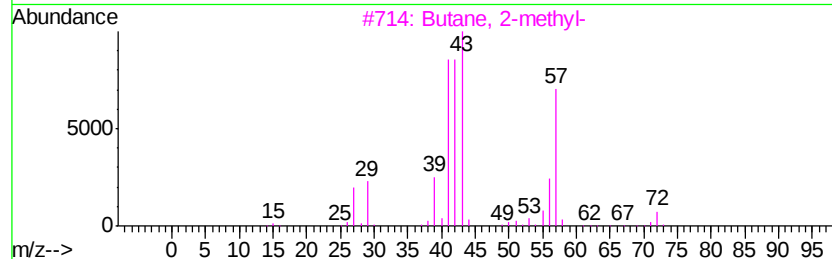
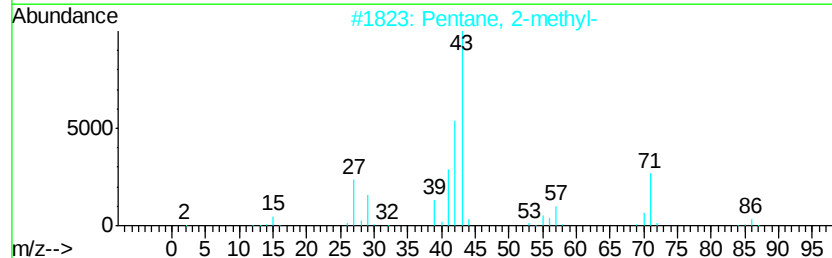
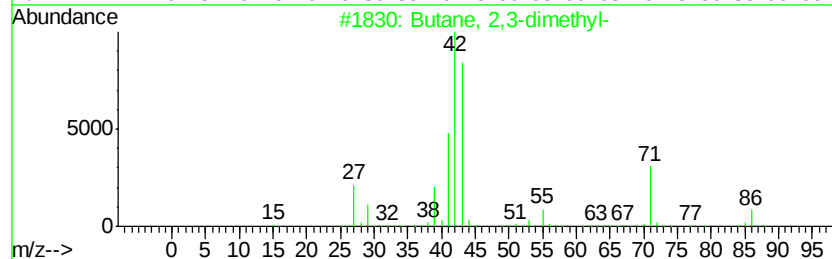
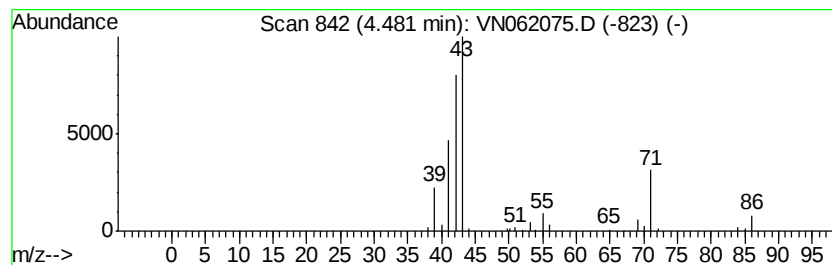
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.48	12.98 ug/l	205439	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2		Pentane, 2-methyl-	86	C6H14	000107-83-5	38
3		Butane, 2-methyl-	72	C5H12	000078-78-4	36
4		Pentane	72	C5H12	000109-66-0	33
5		1-Pentene	70	C5H10	000109-67-1	33



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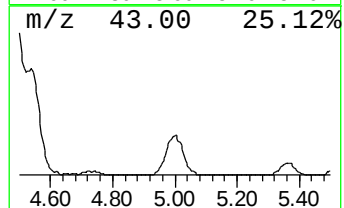
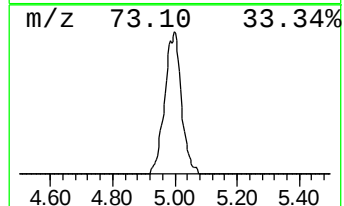
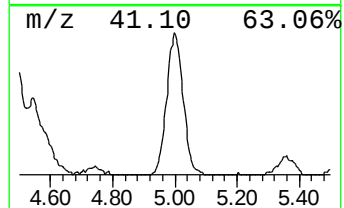
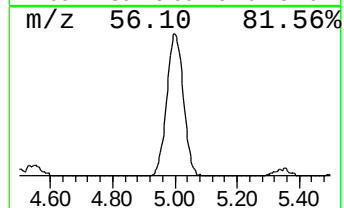
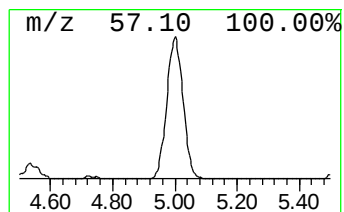
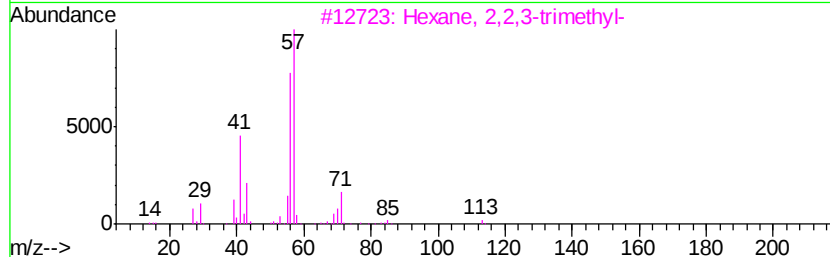
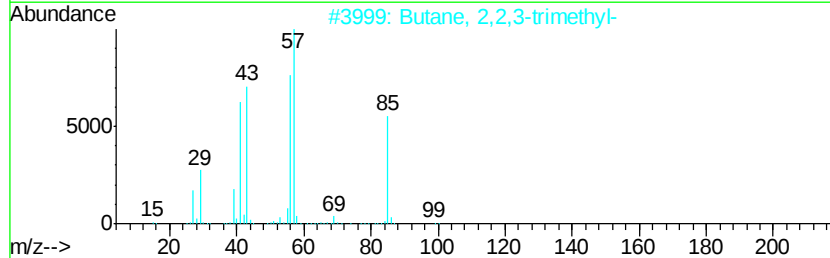
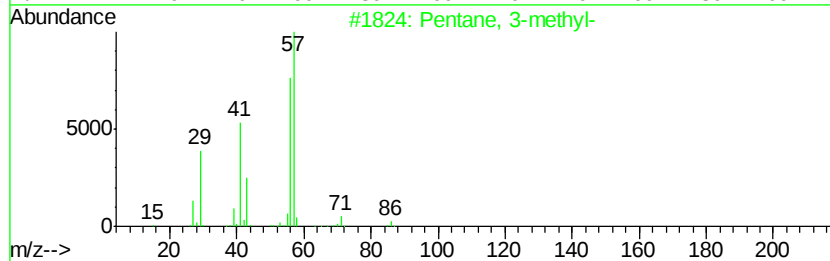
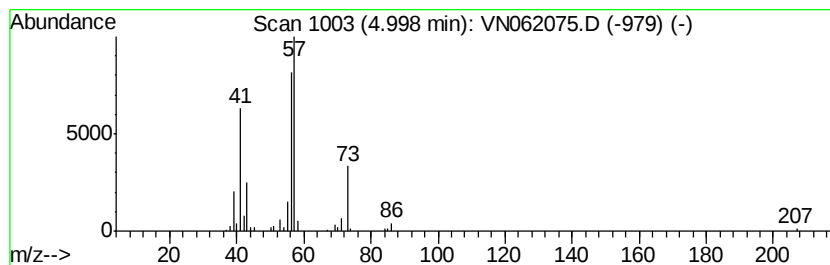
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N062420W.M
Quant Title : SW846 8260

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

Peak Number 3 Pentane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	14.42 ug/l	228145	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 3-methyl-	86	C6H14	000096-14-0	72
2		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	59
4		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	59
5		Heptane, 2,2-dimethyl-	128	C9H20	001071-26-7	47



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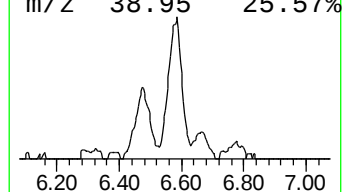
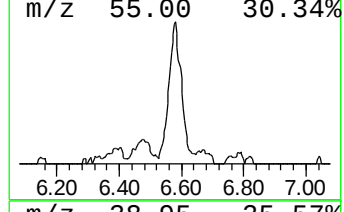
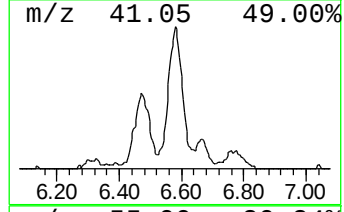
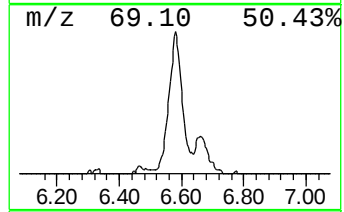
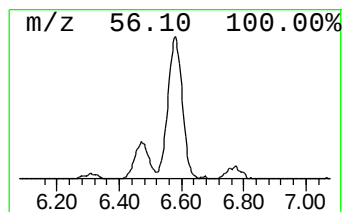
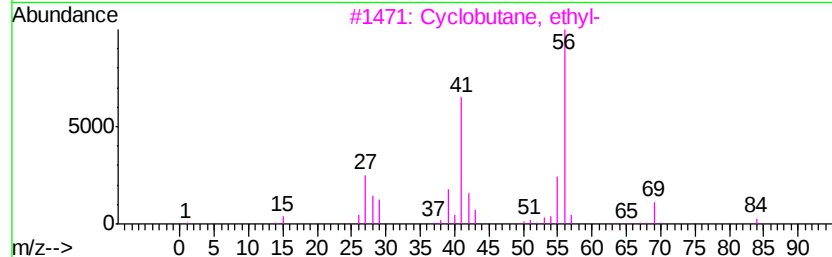
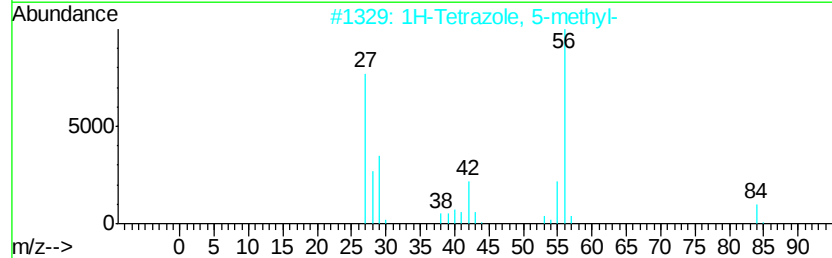
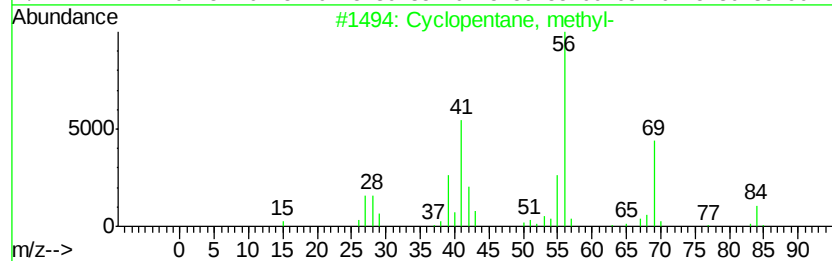
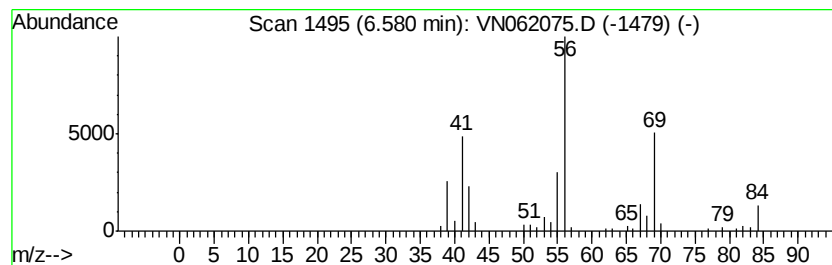
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Quant Title : SW846 8260

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

Peak Number 4 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	10.97 ug/l	173613	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	59
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	58
4		2H-Pyran-2,6(3H)-dione, dihydro-...	142	C7H10O3	004160-82-1	42
5		1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TW-13

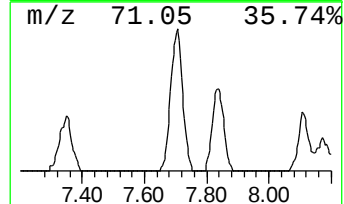
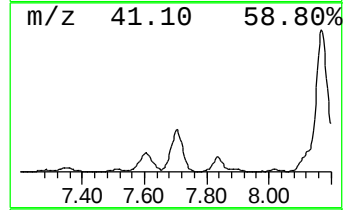
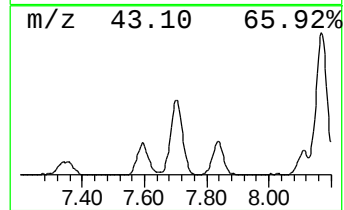
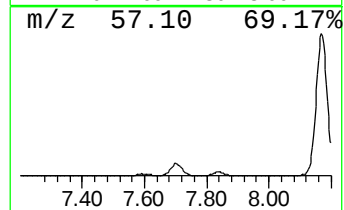
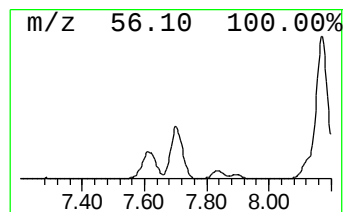
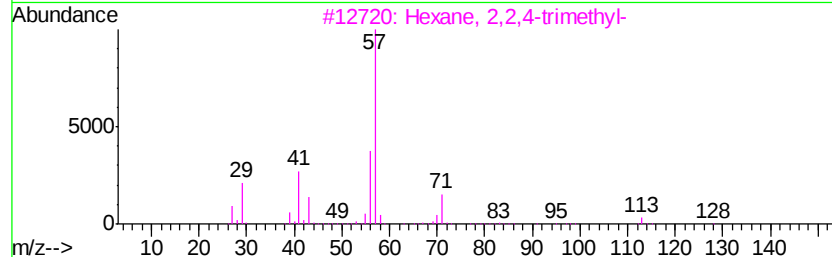
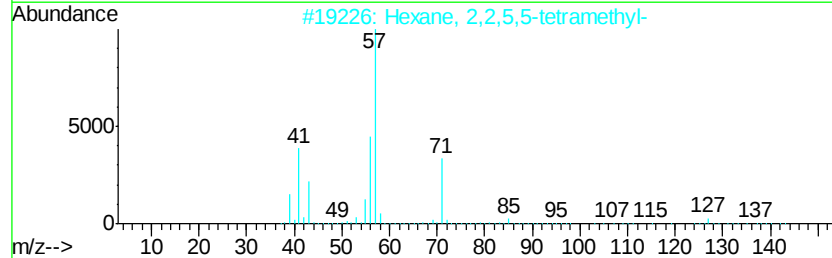
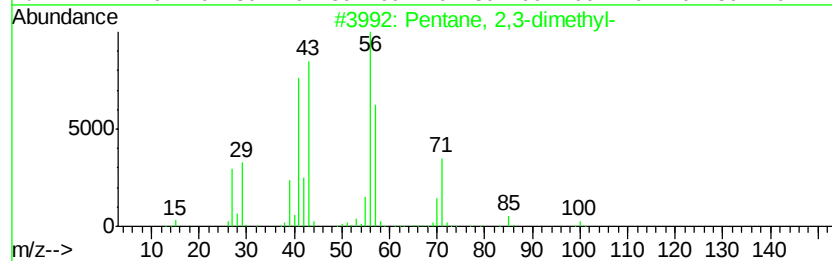
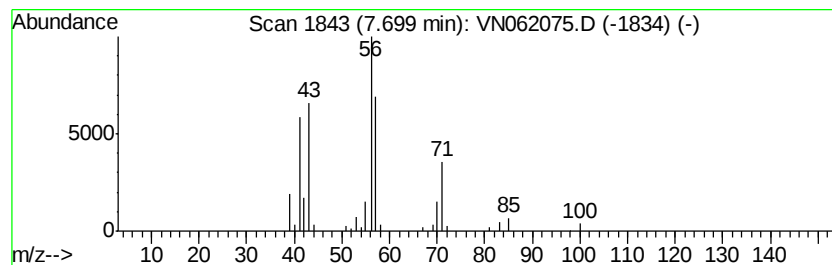
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Quant Title : SW846 8260

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

Peak Number 5 Pentane, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.70	10.65 ug/l	168547	Pentafluorobenzene	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	90
2		Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	64
3		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	50
4		1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	50
5		Oxalic acid, isobutyl octyl ester	258	C14H26O4	1000309-37-3	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TW-13

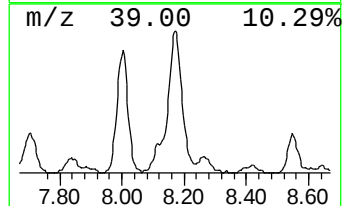
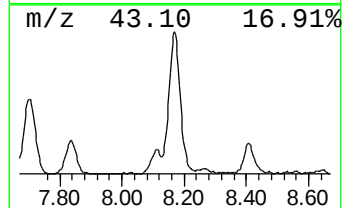
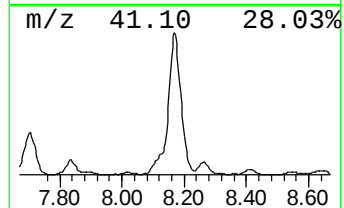
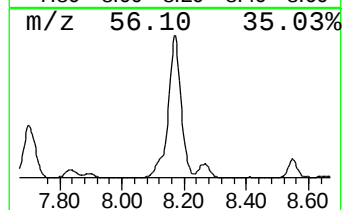
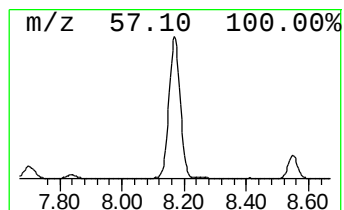
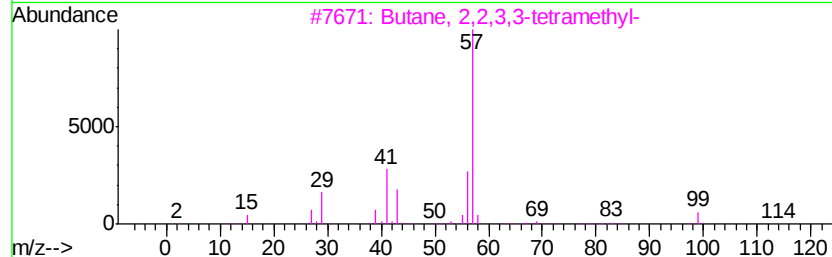
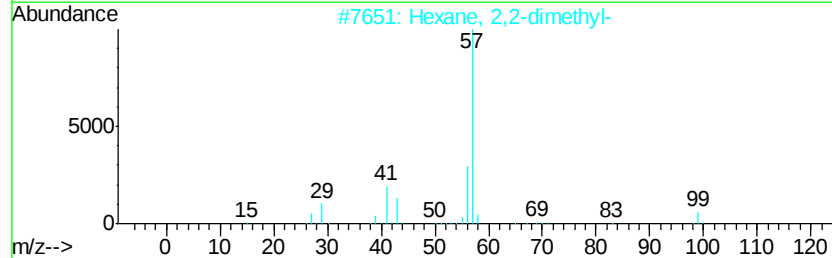
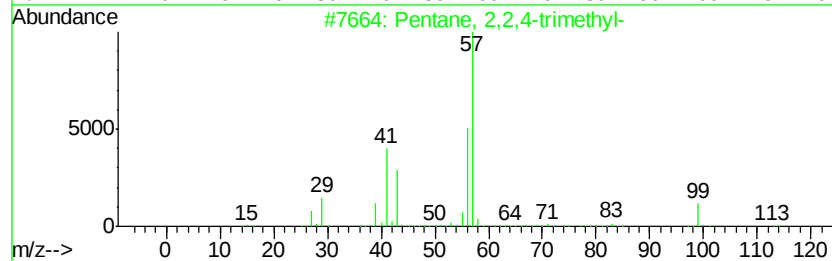
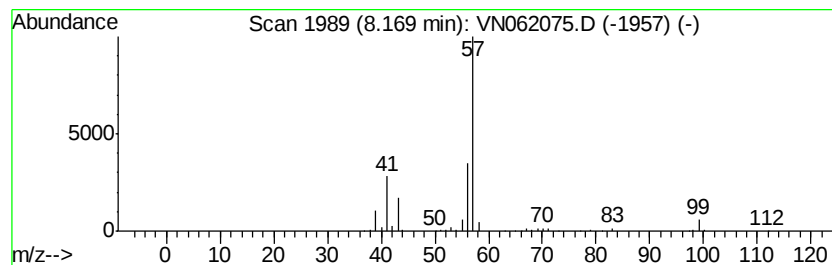
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Quant Title : SW846 8260

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

Peak Number 6 Pentane, 2,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.17	36.97 ug/l	779100	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
3		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
4		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-13

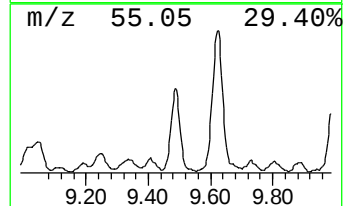
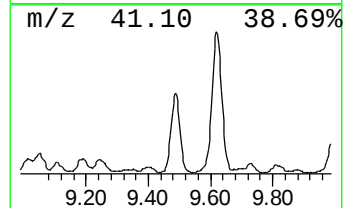
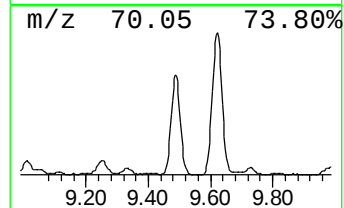
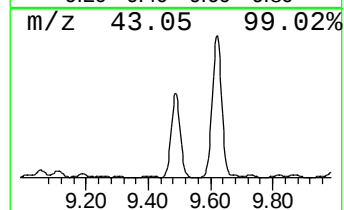
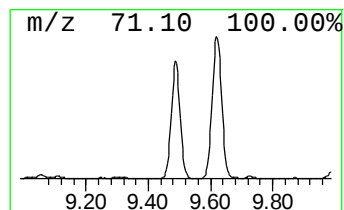
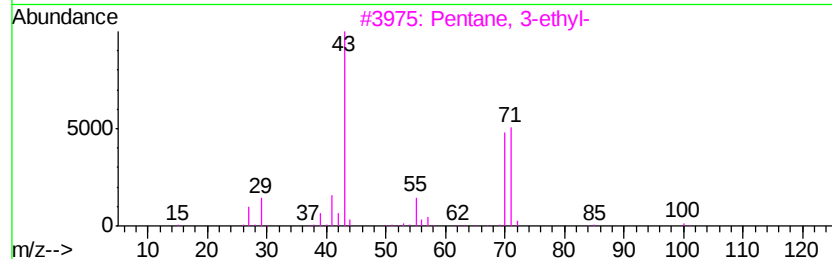
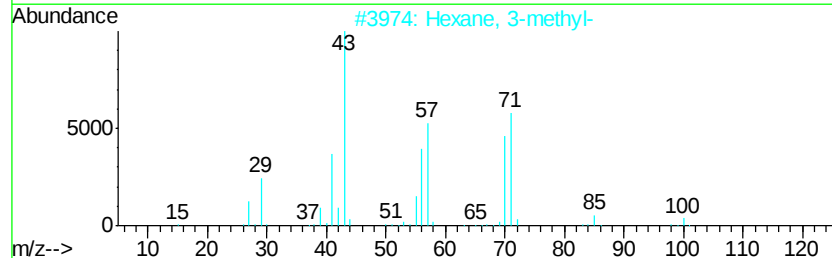
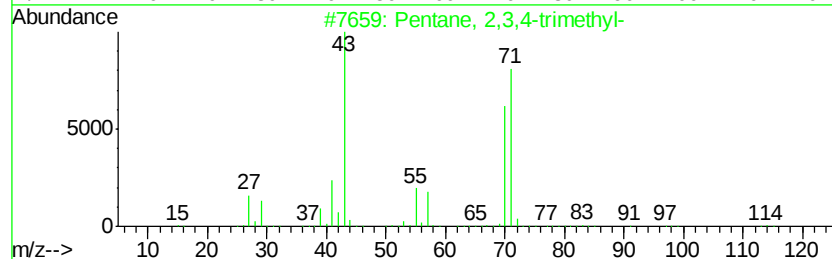
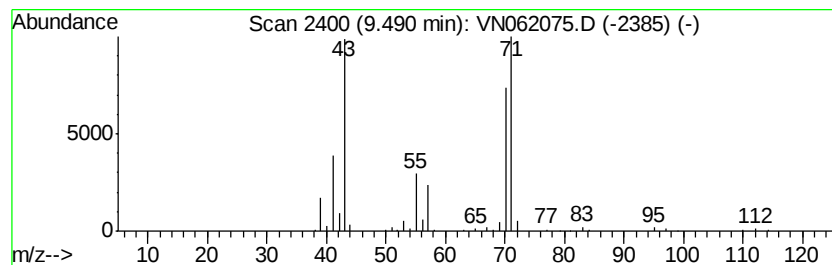
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Quant Title : SW846 8260

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

Peak Number 7 Pentane, 2,3,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	20.56 ug/l	433324	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	90
2		Hexane, 3-methyl-	100	C7H16	000589-34-4	83
3		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
4		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	74
5		Heptane, 4-methyl-	114	C8H18	000589-53-7	72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-13

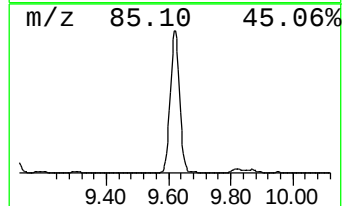
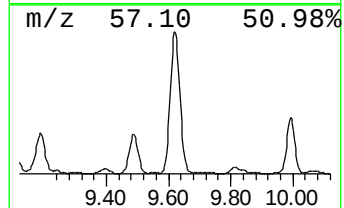
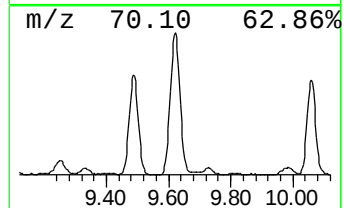
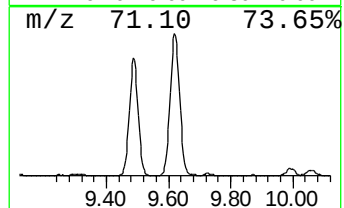
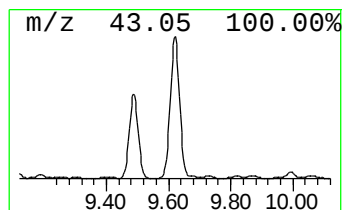
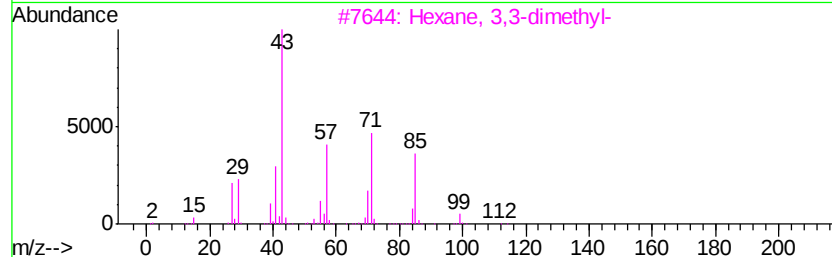
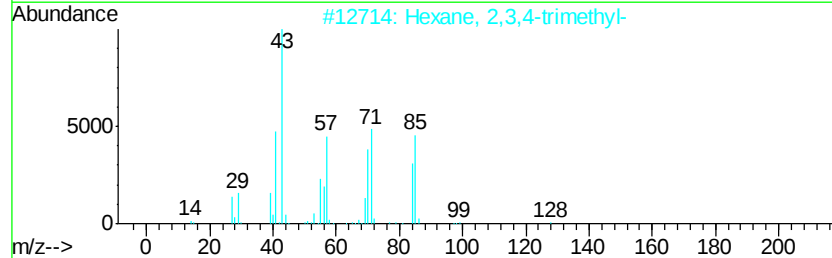
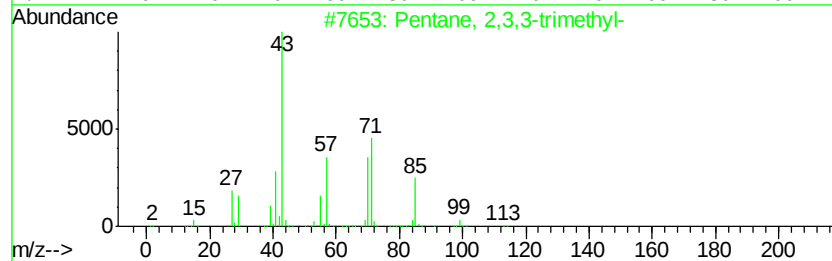
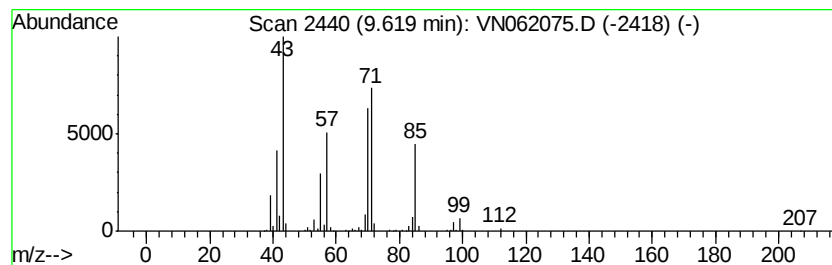
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Quant Title : SW846 8260

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

Peak Number 8 Pentane, 2,3,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	45.08 ug/l	950091	1,4-Difluorobenzene	8.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4		Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	59
5		1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
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ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-13

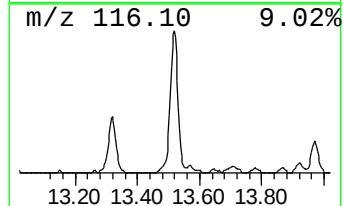
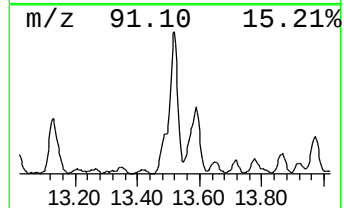
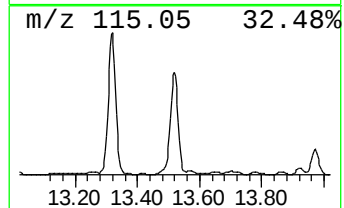
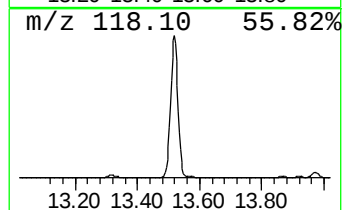
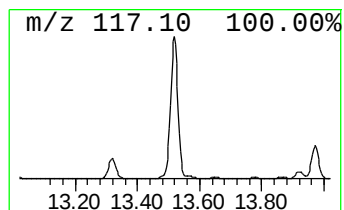
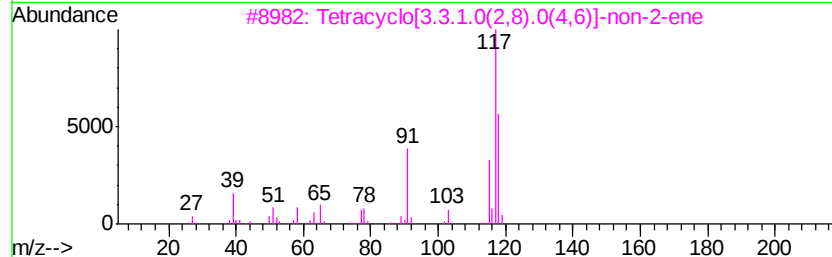
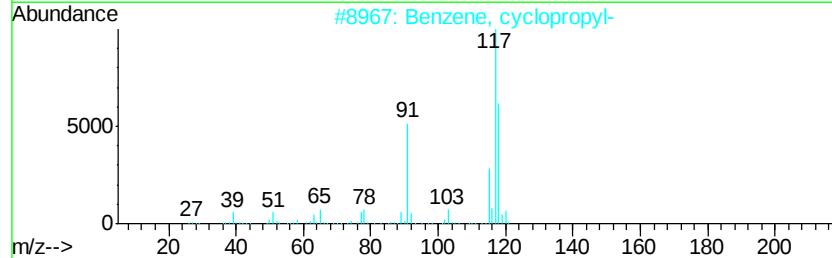
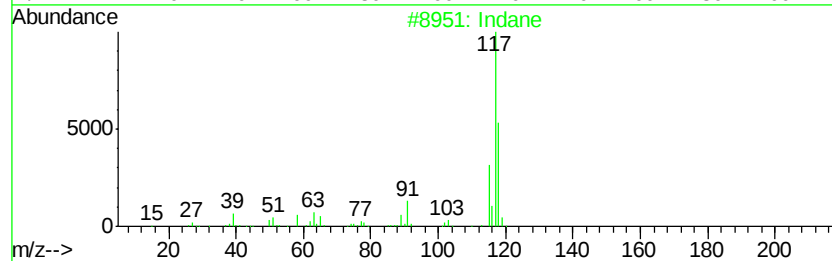
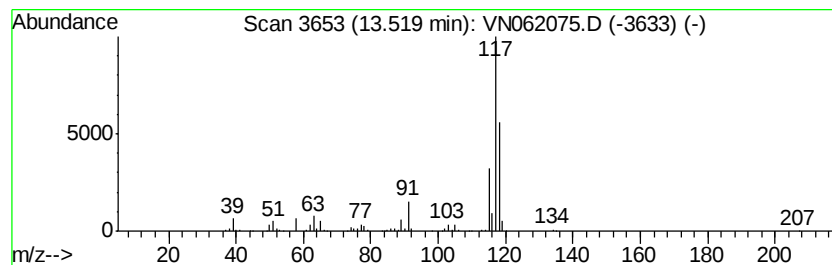
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Quant Title : SW846 8260

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

Peak Number 9 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	36.40 ug/l	890256	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	95
2	Benzene, cyclopropyl-		118	C9H10	000873-49-4	83
3	Tetracyclo[3.3.1.0(2,8).0(4,6)]-	...	118	C9H10	1000191-13-7	80
4	Benzene, 1-ethenyl-4-methyl-		118	C9H10	000622-97-9	74
5	Deltacyclene		118	C9H10	007785-10-6	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TW-13

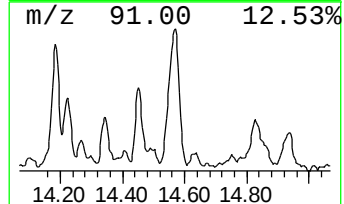
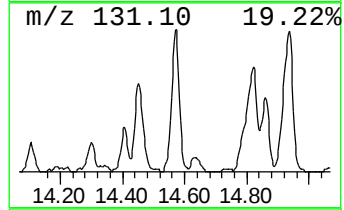
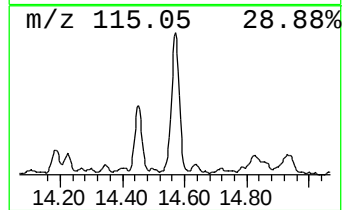
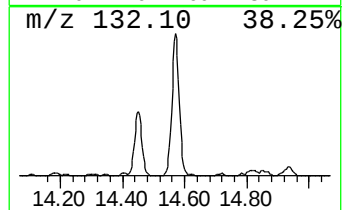
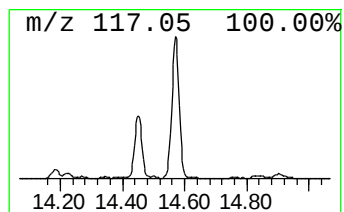
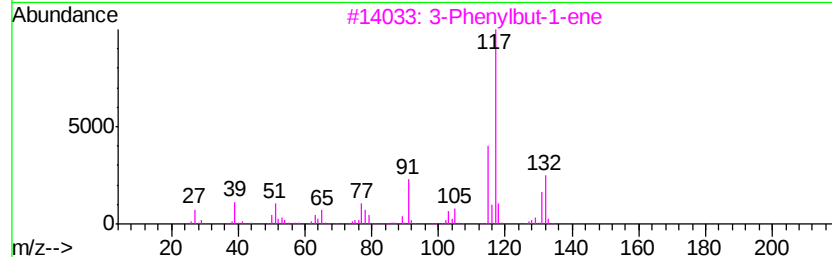
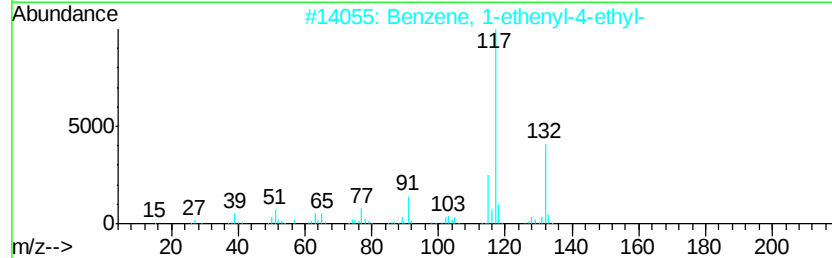
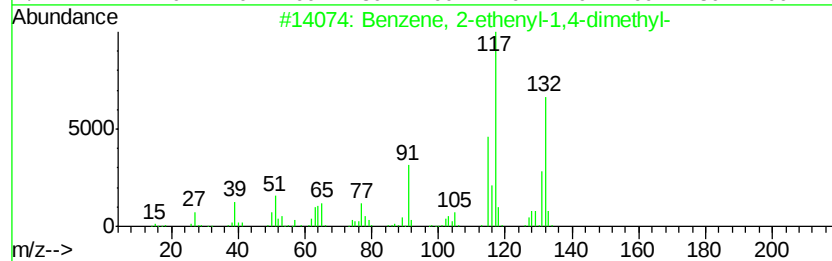
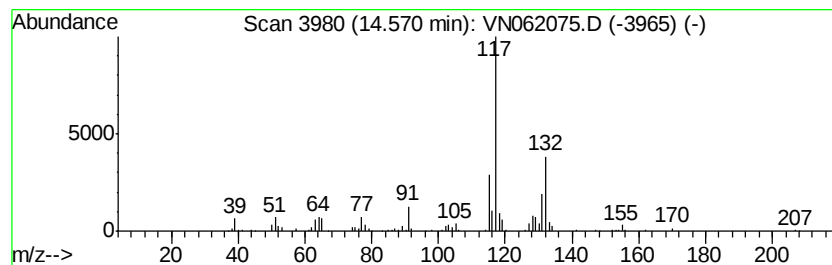
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	15.45 ug/l	377906	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN062420\
Data File : VN062075.D
Acq On : 24 Jun 2020 20:13
Operator : JC/MD
Sample : L3063-12
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TW-13

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
Butane, 2-methyl-	2.77	19.0	ug/l	300342	1	7.63	791242	50.0
Butane, 2,3-dimet...	4.48	13.0	ug/l	205439	1	7.63	791242	50.0
Pentane, 3-methyl-	5.00	14.4	ug/l	228145	1	7.63	791242	50.0
Cyclopentane, met...	6.58	11.0	ug/l	173613	1	7.63	791242	50.0
Pentane, 2,3-dime...	7.70	10.7	ug/l	168547	1	7.63	791242	50.0
Pentane, 2,2,4-tr...	8.17	37.0	ug/l	779100	2	8.55	1053810	50.0
Pentane, 2,3,4-tr...	9.49	20.6	ug/l	433324	2	8.55	1053810	50.0
Pentane, 2,3,3-tr...	9.62	45.1	ug/l	950091	2	8.55	1053810	50.0
Indane	13.52	36.4	ug/l	890256	4	13.32	1222750	50.0
Benzene, 2-etheny...	14.57	15.4	ug/l	377906	4	13.32	1222750	50.0