

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
 Data File : VN070407.D
 Acq On : 30 Dec 2021 18:32
 Operator : JC/MD
 Sample : M5212-06
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DUP122121

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\82N122721W.M
 Title : SW846 8260

Signal : TIC: VN070407.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.263	93	102	128	rVB3	356094	684241	7.09%	0.577%
2	2.440	161	168	194	rVB	350079	606042	6.28%	0.511%
3	3.135	408	427	466	rVB	4148860	9649874	100.00%	8.131%
4	3.508	553	566	586	rVB2	85735	208232	2.16%	0.175%
5	3.843	672	691	713	rVB4	41200	130915	1.36%	0.110%
6	3.982	725	743	767	rBV2	71017	195225	2.02%	0.165%
7	4.248	813	842	884	rBV2	755892	2693225	27.91%	2.269%
8	5.074	1115	1150	1173	rBV3	1682596	6810194	70.57%	5.739%
9	5.616	1317	1352	1389	rBV3	559254	1967377	20.39%	1.658%
10	6.474	1651	1672	1696	rBV8	36733	131465	1.36%	0.111%
11	6.611	1701	1723	1748	rVV6	104449	325060	3.37%	0.274%
12	6.892	1801	1828	1851	rBV4	141618	490249	5.08%	0.413%
13	7.045	1863	1885	1904	rBV2	538126	1610196	16.69%	1.357%
14	7.147	1905	1923	1938	rVV	619658	1807952	18.74%	1.523%
15	7.222	1940	1951	1973	rVV3	388433	1053856	10.92%	0.888%
16	7.324	1977	1989	2020	rVB8	64966	222267	2.30%	0.187%
17	7.654	2095	2112	2130	rBV5	48855	140069	1.45%	0.118%
18	7.791	2143	2163	2164	rBV5	55379	105256	1.09%	0.089%
19	7.844	2170	2183	2209	rVB5	284578	779827	8.08%	0.657%
20	8.024	2225	2250	2260	rBV	634355	1576670	16.34%	1.329%
21	8.088	2261	2274	2290	rVV2	1407549	3362062	34.84%	2.833%
22	8.174	2290	2306	2332	rVB	1593858	4062932	42.10%	3.424%
23	8.295	2338	2351	2362	rBV3	59865	133609	1.38%	0.113%
24	8.357	2363	2374	2387	rVV4	98300	224215	2.32%	0.189%
25	8.458	2388	2412	2433	rVV3	2572810	6986430	72.40%	5.887%
26	8.617	2436	2471	2492	rVV	1853666	5544354	57.46%	4.672%
27	8.705	2493	2504	2525	rVB2	395495	911336	9.44%	0.768%
28	8.971	2579	2603	2627	rBV	1997958	4528916	46.93%	3.816%
29	9.062	2628	2637	2652	rVV2	199868	419846	4.35%	0.354%
30	9.199	2672	2688	2698	rBV3	258960	553589	5.74%	0.466%
31	9.247	2699	2706	2732	rVB	227055	446917	4.63%	0.377%
32	9.424	2753	2772	2796	rBV3	603074	1925840	19.96%	1.623%
33	9.515	2797	2806	2822	rVB5	187477	396274	4.11%	0.334%
34	9.593	2824	2835	2845	rBV3	80371	154854	1.60%	0.130%
35	9.732	2868	2887	2902	rBV6	163267	390157	4.04%	0.329%
36	9.883	2930	2943	2962	rVB	834692	1728152	17.91%	1.456%

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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

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37	9.976	2964	2978	2981	rBV3	288914	456610	4.73%	0.385%
38	10.014	2983	2992	3021	rVB	871391	1871432	19.39%	1.577%
39	10.124	3023	3033	3046	rVB7	84296	195607	2.03%	0.165%
40	10.199	3047	3061	3075	rBV5	185274	365690	3.79%	0.308%
41	10.322	3095	3107	3116	rBV	361845	652659	6.76%	0.550%
42	10.443	3136	3152	3197	rBV	2816420	5635147	58.40%	4.748%
43	10.644	3199	3227	3241	rVV3	1006300	2171653	22.50%	1.830%
44	10.717	3242	3254	3269	rVV2	875043	1581498	16.39%	1.333%
45	10.784	3269	3279	3291	rVV5	94000	167995	1.74%	0.142%
46	10.859	3292	3307	3327	rVB4	737710	1627656	16.87%	1.372%
47	10.947	3330	3340	3358	rVB8	83722	202441	2.10%	0.171%
48	11.132	3398	3409	3418	rBV5	127666	223857	2.32%	0.189%
49	11.258	3448	3456	3467	rVB5	143670	232466	2.41%	0.196%
50	11.452	3518	3528	3546	rVB4	128509	230410	2.39%	0.194%
51	11.559	3558	3568	3586	rVB10	47231	122069	1.26%	0.103%
52	11.747	3622	3638	3653	rBV	2513126	4555545	47.21%	3.839%
53	11.846	3662	3675	3699	rVB	3193175	5640782	58.45%	4.753%
54	12.581	3935	3949	3969	rBV	1636458	2742199	28.42%	2.311%
55	12.731	3991	4005	4027	rVB	1959124	3441489	35.66%	2.900%
56	12.921	4062	4076	4096	rVB	1069167	1832033	18.99%	1.544%
57	13.222	4176	4188	4203	rVB	268250	453868	4.70%	0.382%
58	13.369	4235	4243	4254	rVB	98844	159420	1.65%	0.134%
59	13.484	4273	4286	4304	rBV3	282002	653923	6.78%	0.551%
60	13.675	4343	4357	4369	rBV	1850703	3339706	34.61%	2.814%
61	13.771	4385	4393	4404	rVB	128118	200033	2.07%	0.169%
62	13.838	4405	4418	4422	rBV	696217	1005129	10.42%	0.847%
63	13.876	4423	4432	4445	rVB	3493591	5682955	58.89%	4.789%
64	14.005	4467	4480	4492	rVB	568371	921677	9.55%	0.777%
65	14.069	4493	4504	4517	rVB	416977	670021	6.94%	0.565%
66	14.150	4518	4534	4547	rBV2	369105	691441	7.17%	0.583%
67	14.217	4548	4559	4571	rBV	307065	471973	4.89%	0.398%
68	14.278	4572	4582	4589	rBV	177059	293500	3.04%	0.247%
69	14.329	4589	4601	4615	rVB	1650034	2758087	28.58%	2.324%
70	14.407	4618	4630	4642	rVB5	173666	332680	3.45%	0.280%
71	14.466	4642	4652	4662	rBV	156737	229923	2.38%	0.194%
72	14.538	4666	4679	4690	rBV	452955	741163	7.68%	0.625%
73	14.656	4717	4723	4733	rVB2	92778	127356	1.32%	0.107%
74	14.764	4752	4763	4771	rBV2	149709	237332	2.46%	0.200%
75	14.812	4771	4781	4806	rVB5	295125	618695	6.41%	0.521%
76	14.930	4809	4825	4838	rBV3	255472	561923	5.82%	0.474%

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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\82N122721W.M
Title : SW846 8260

77	15.000	4839	4851	4863	rVB4	135206	236257	2.45%	0.199%
78	15.083	4872	4882	4893	rBV2	107887	188120	1.95%	0.159%
79	15.196	4911	4924	4936	rBV3	178144	356150	3.69%	0.300%
80	15.306	4956	4965	4984	rVB2	232612	526780	5.46%	0.444%
81	15.804	5140	5151	5170	rBV4	54070	100620	1.04%	0.085%
82	16.110	5254	5265	5282	rBV	97232	207946	2.15%	0.175%

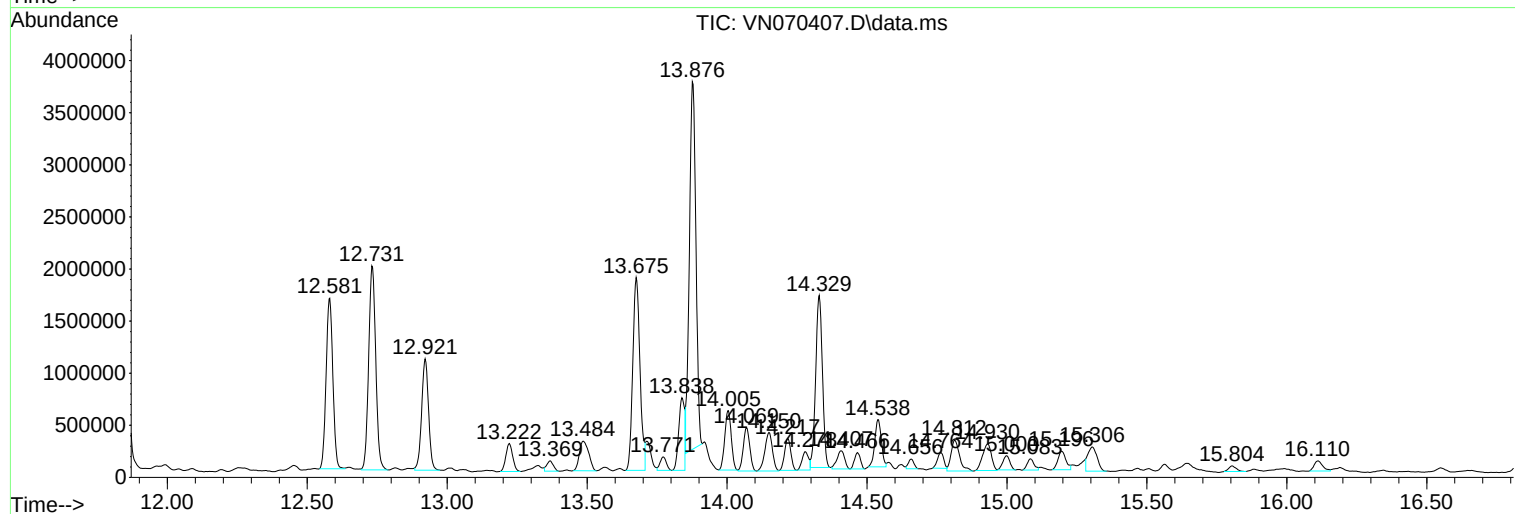
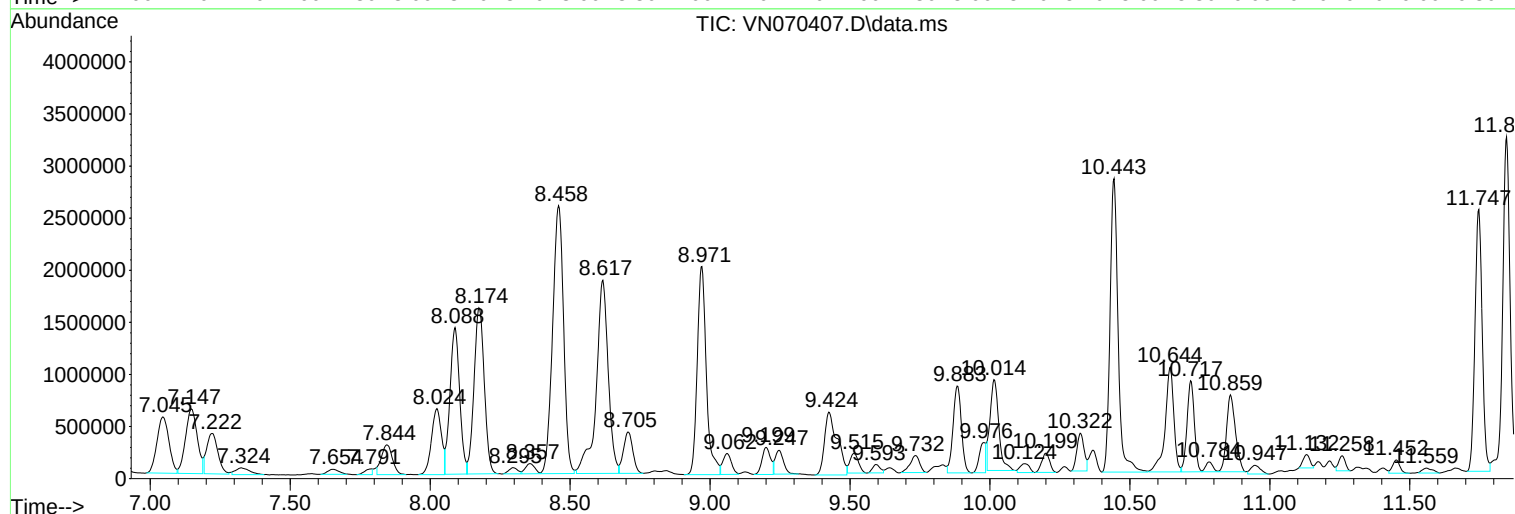
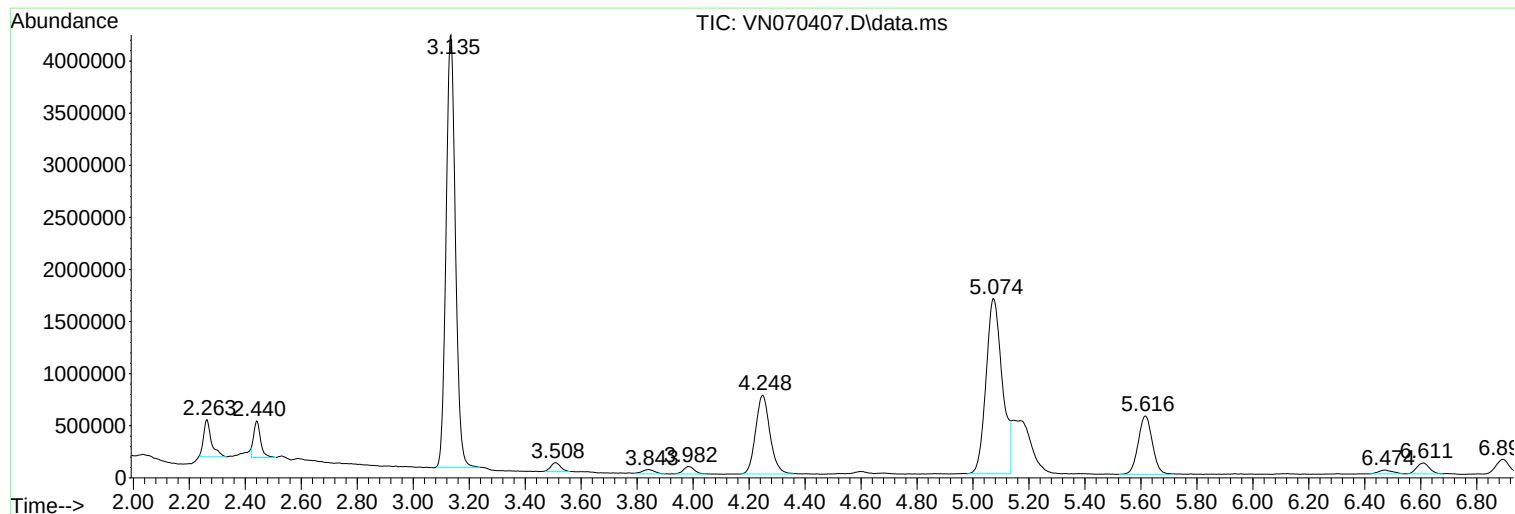
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

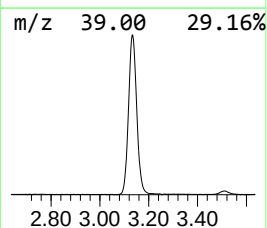
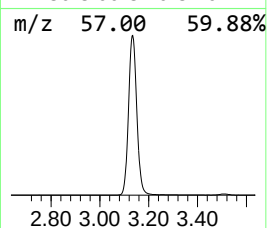
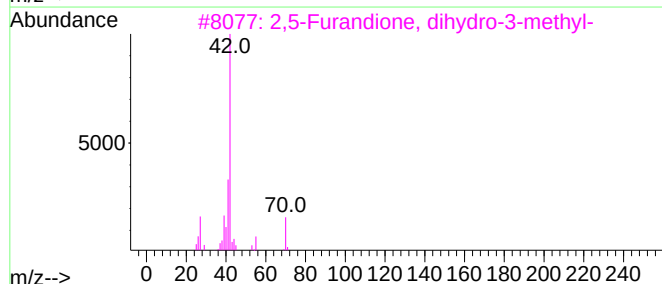
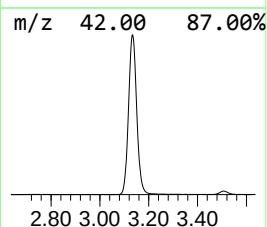
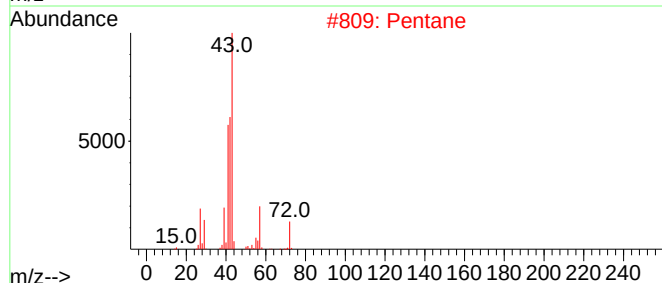
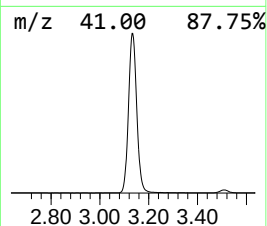
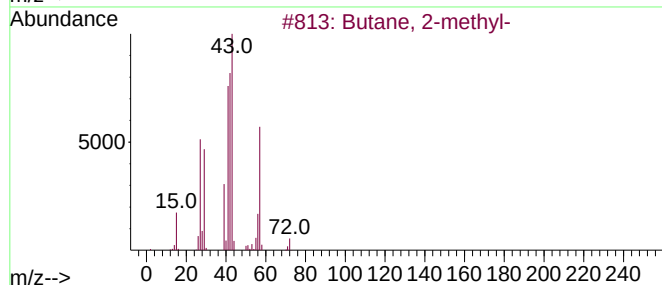
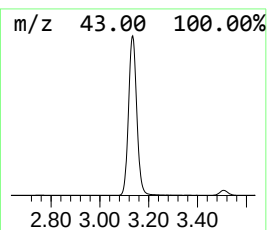
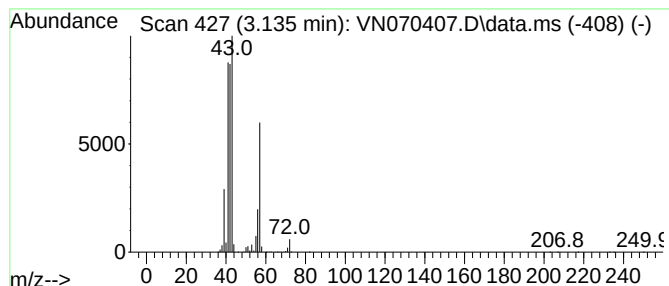
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.135	143.51 ug/l	9649870	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methyl-	72	C5H12	000078-78-4	91
2			Pentane	72	C5H12	000109-66-0	42
3			2,5-Furandione, dihydro-3-methyl-	114	C5H6O3	004100-80-5	25
4			3-Buten-1-ol	72	C4H8O	000627-27-0	12
5			1-Propene, 2-methyl-	56	C4H8	000115-11-7	9



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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

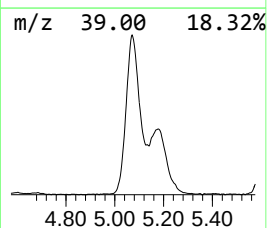
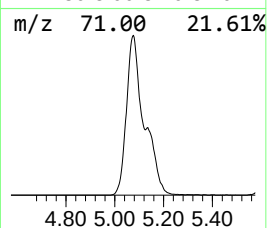
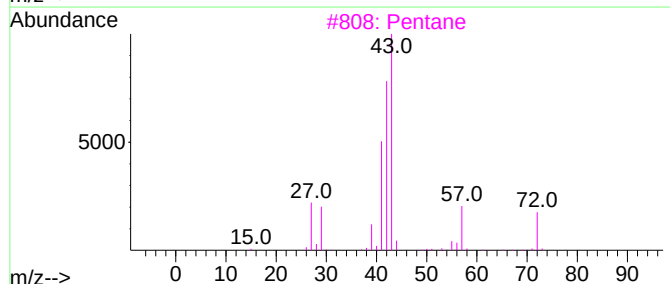
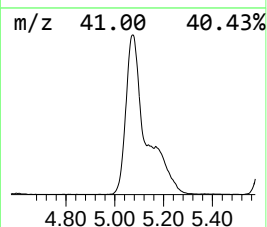
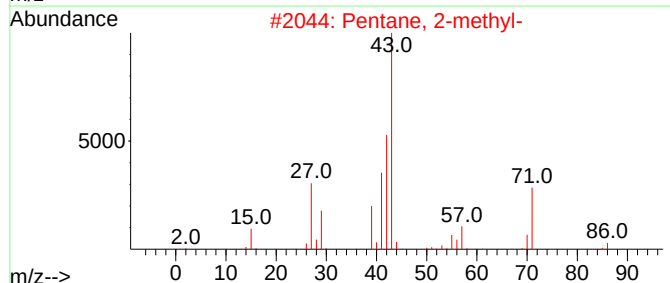
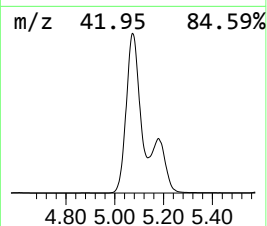
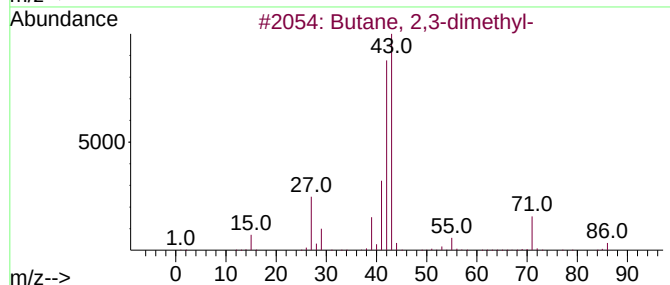
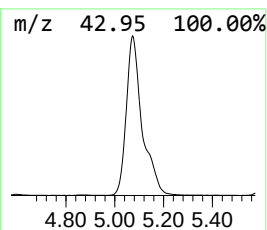
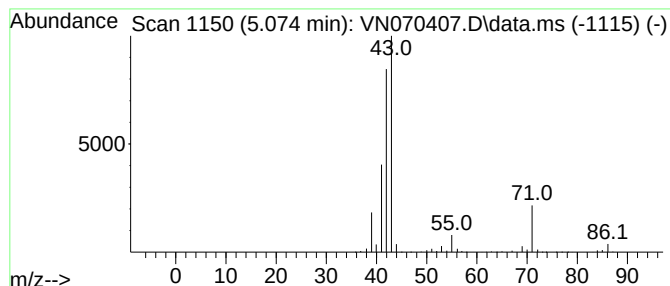
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.074	101.28 ug/l	6810190	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	91
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3			Pentane	72	C5H12	000109-66-0	38
4			Pyrrolidine	71	C4H9N	000123-75-1	9
5			Butanal, 2-ethyl-	100	C6H12O	000097-96-1	9



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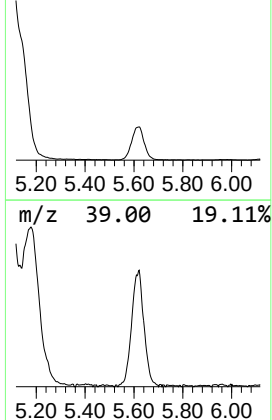
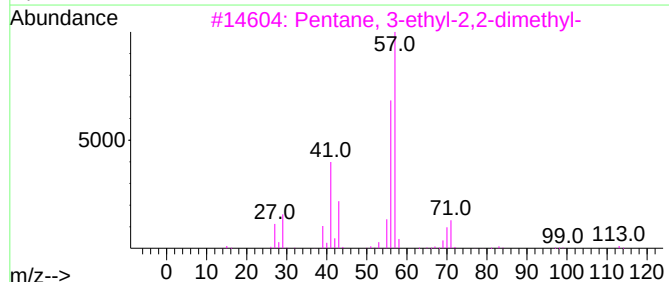
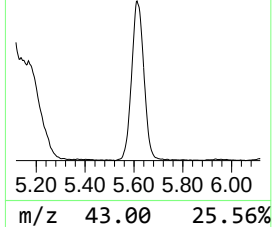
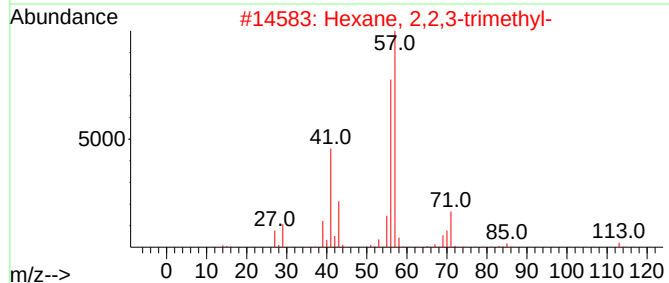
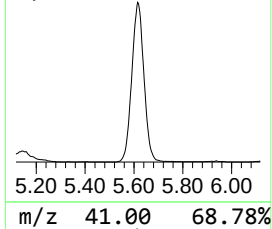
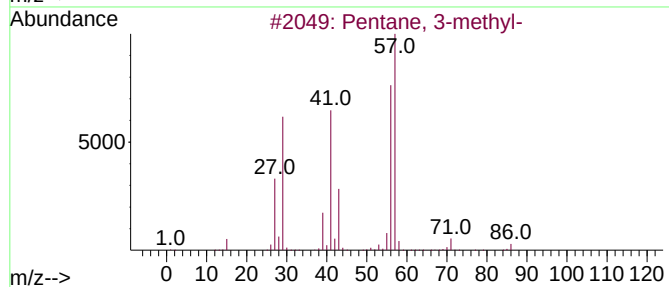
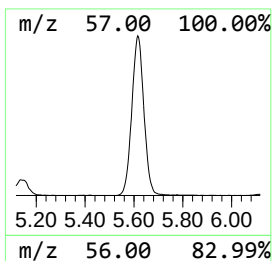
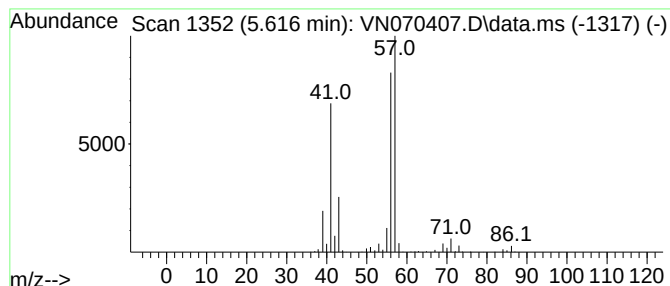
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.616	29.26 ug/l	1967380	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	87
2			Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	74
3			Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	64
5			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	42



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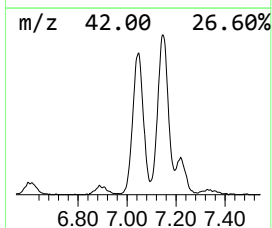
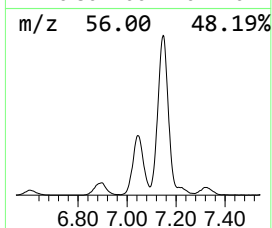
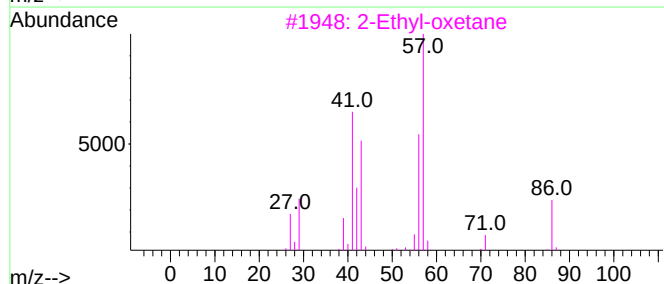
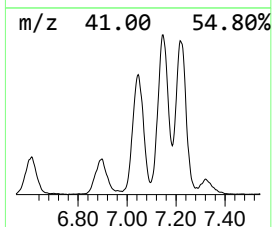
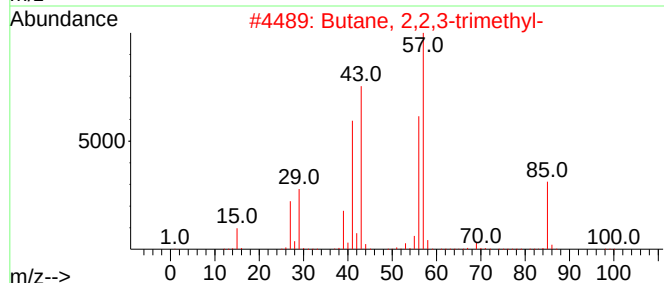
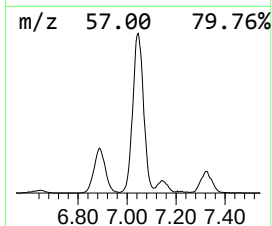
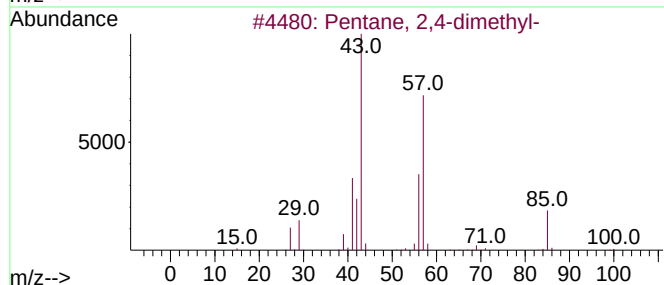
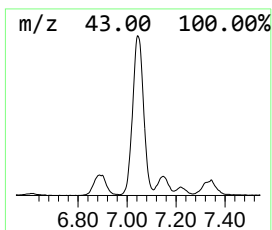
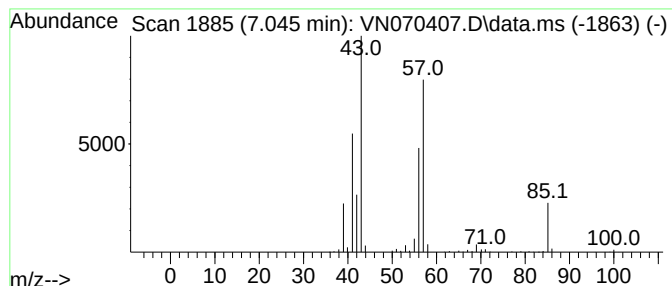
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Pentane, 2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.045	23.95 ug/l	1610200	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3			2-Ethyl-oxetane	86	C5H10O	1010386-40-2	50
4			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
5			2,2,6,6-Tetramethylheptane	156	C11H24	040117-45-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

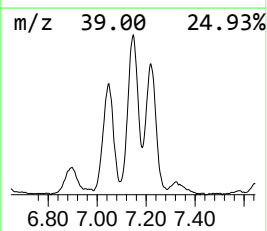
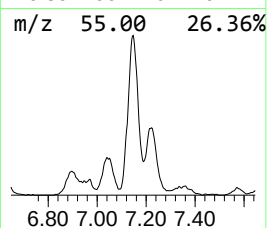
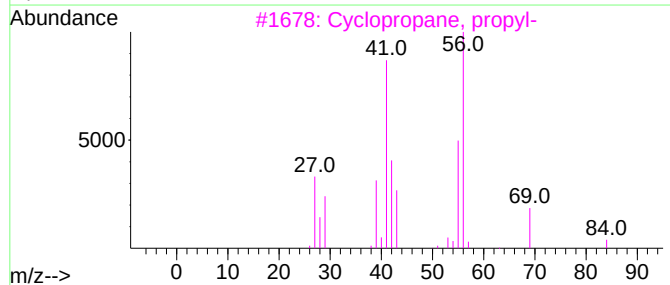
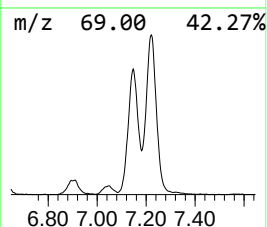
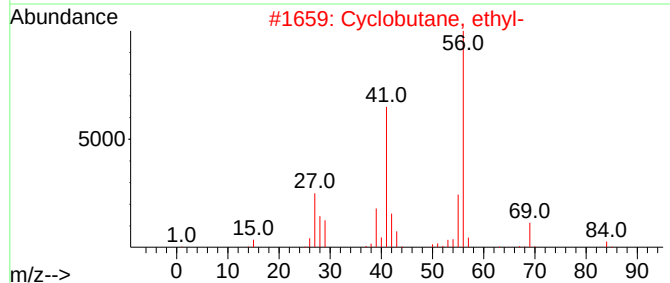
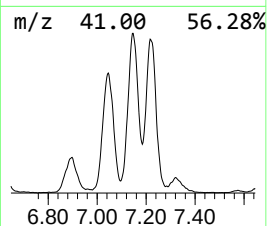
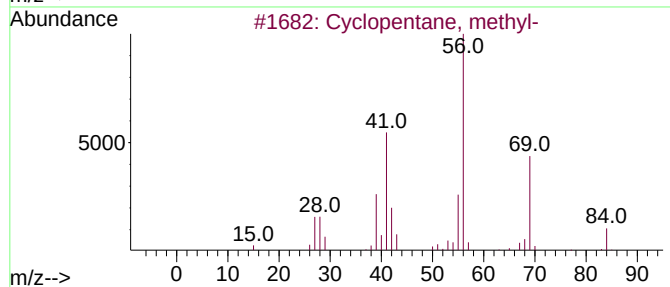
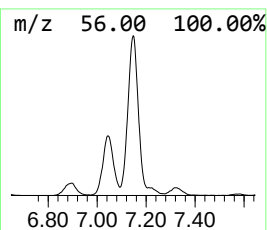
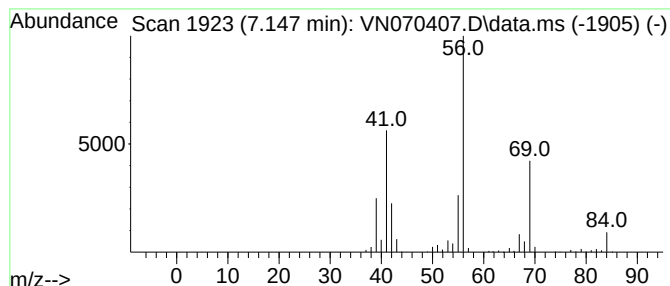
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Cyclopentane, methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.147	26.89 ug/l	1807950	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3			Cyclopropane, propyl-	84	C6H12	002415-72-7	72
4			Cyclobutane	56	C4H8	000287-23-0	50
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

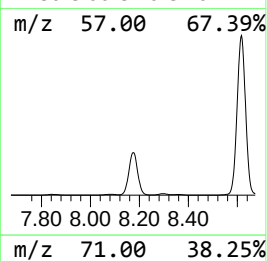
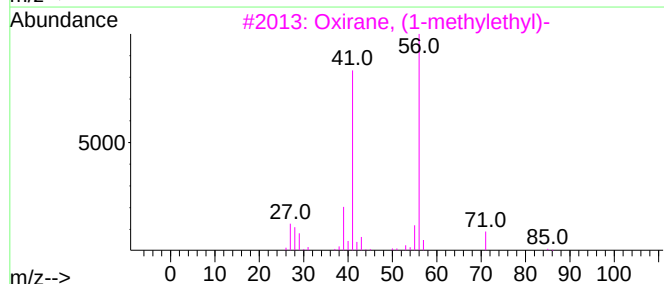
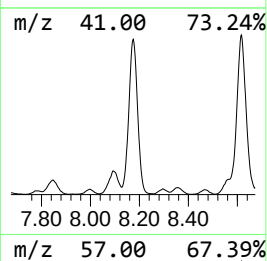
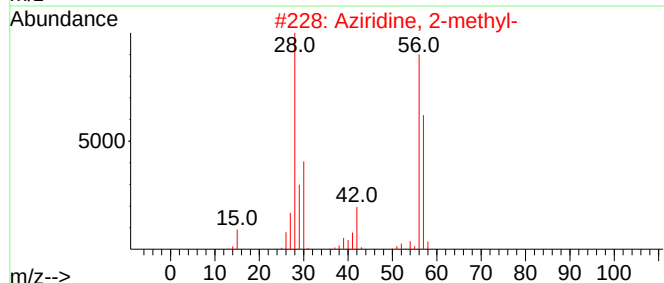
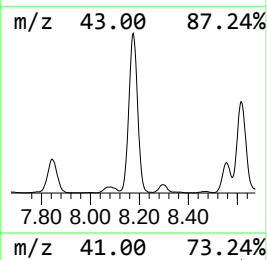
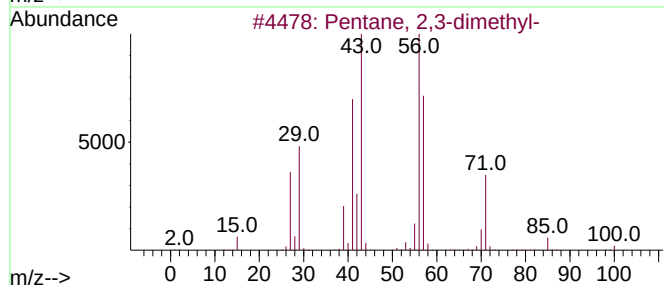
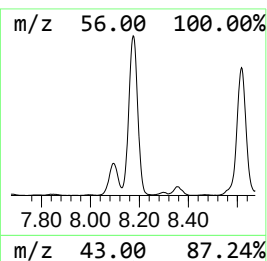
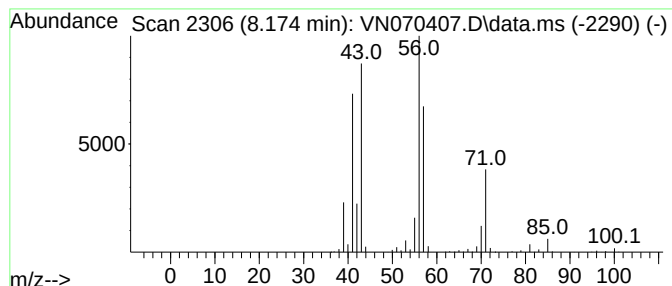
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Pentane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.174	60.42 ug/l	4062930	Pentafluorobenzene	8.088

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2			Aziridine, 2-methyl-	57	C3H7N	000075-55-8	43
3			Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	37
4			Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	37
5			1-Hexene, 2-methyl-	98	C7H14	006094-02-6	32



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

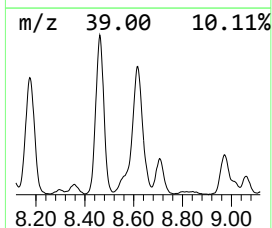
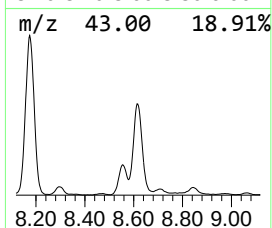
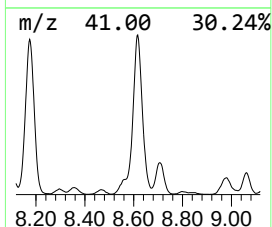
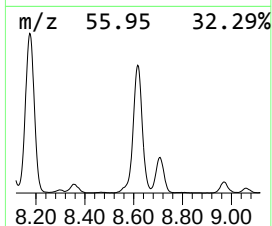
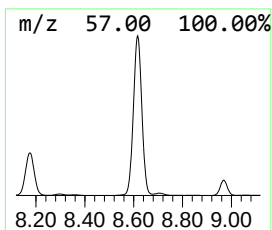
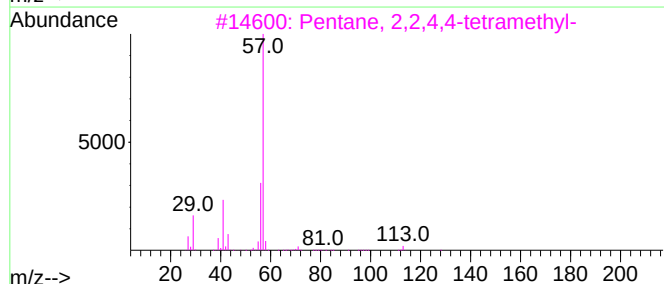
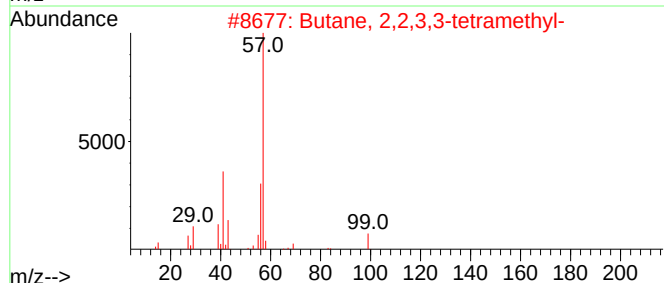
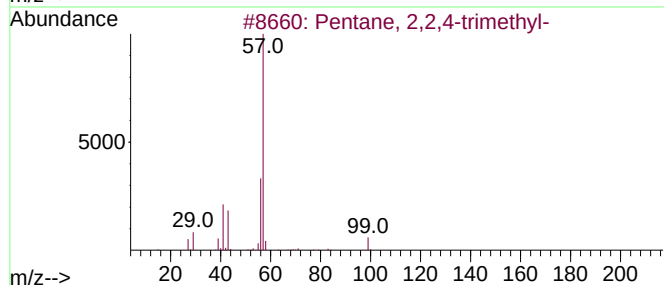
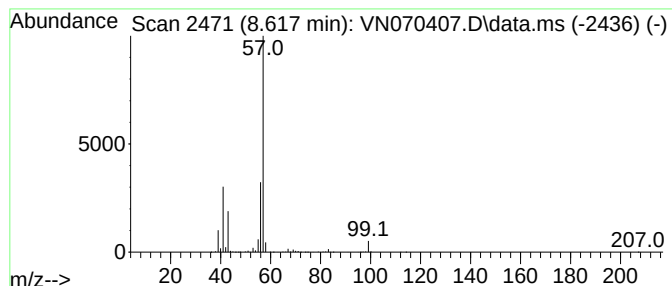
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.617	61.21 ug/l	5544350	1,4-Difluorobenzene	8.971

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
4			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5			Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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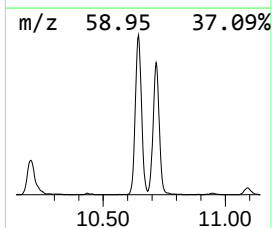
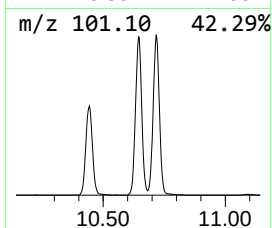
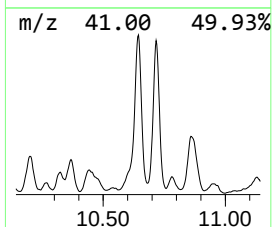
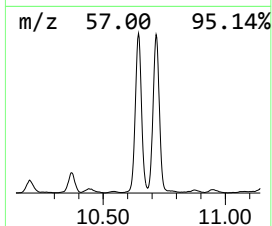
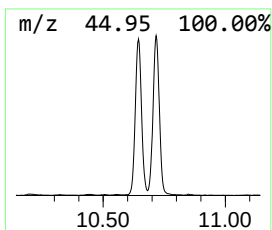
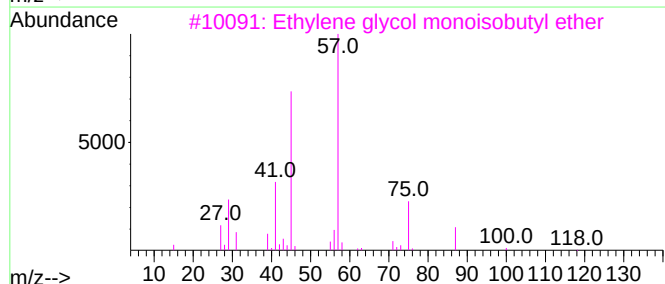
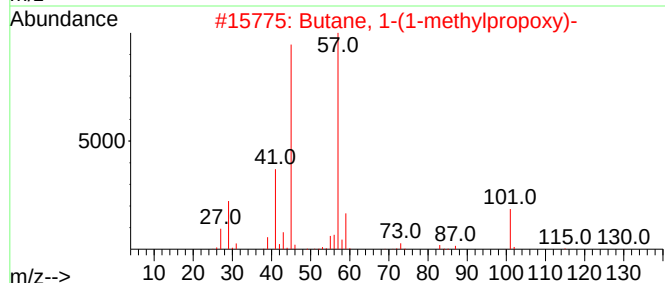
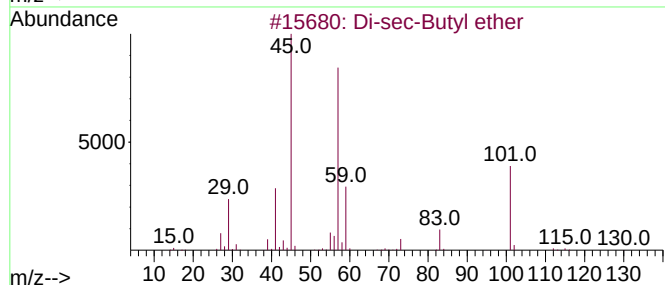
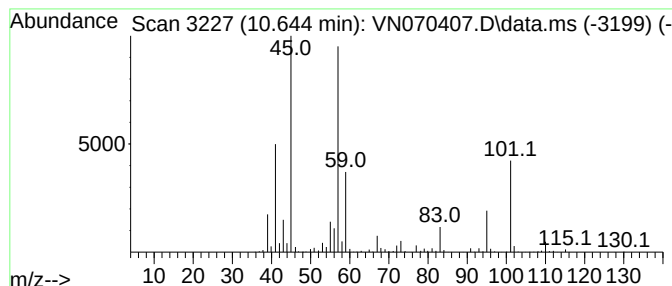
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Di-sec-Butyl ether Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.644	23.84 ug/l	2171650	Chlorobenzene-d5	11.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Di-sec-Butyl ether	130	C8H18O	006863-58-7	87
2			Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	43
4			Pentane, 2-butoxy-	144	C9H20O	062238-02-2	42
5			Fructofuranose, 2,6-anhydro-1,3,...	204	C9H16O5	004451-11-0	28



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

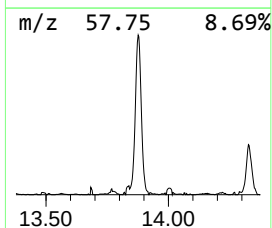
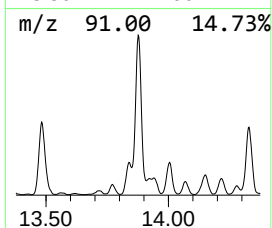
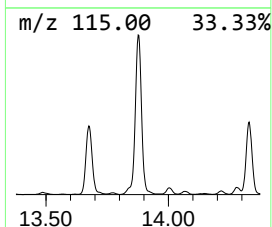
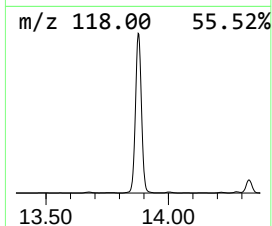
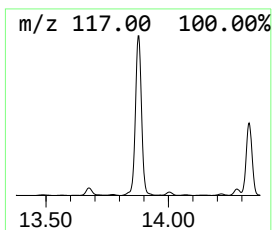
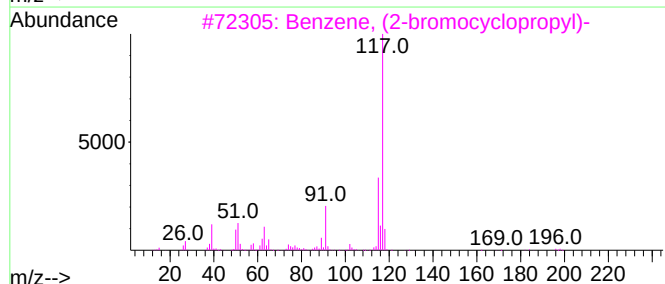
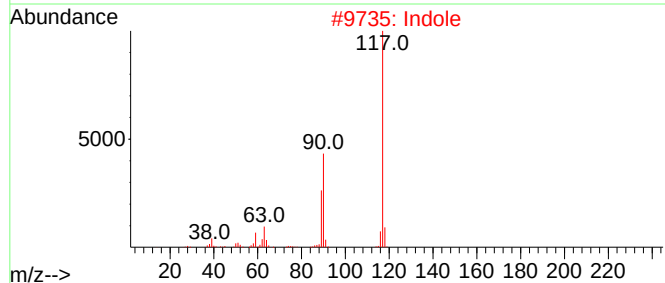
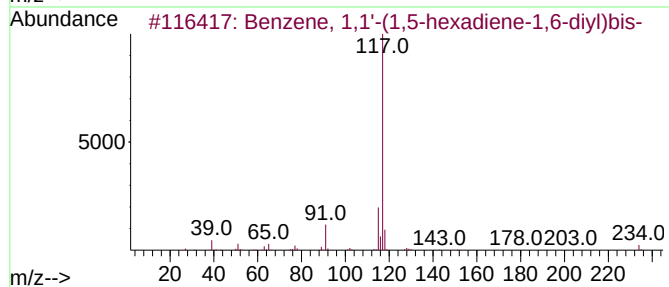
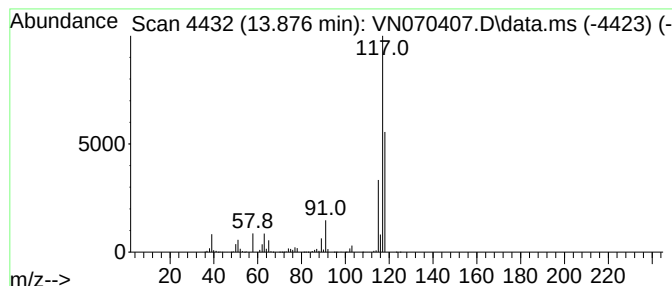
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.876	85.08 ug/l	5682960	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2			Indole	117	C8H7N	000120-72-9	49
3			Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
4			5H-1-Pyridine	117	C8H7N	000270-91-7	43
5			4-Ethynylaniline	117	C8H7N	014235-81-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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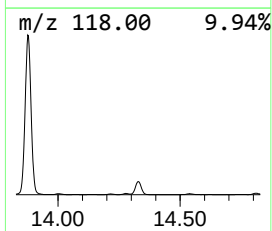
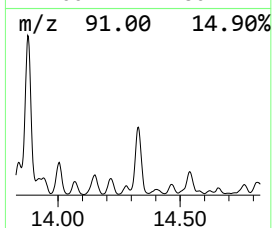
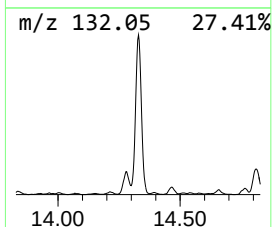
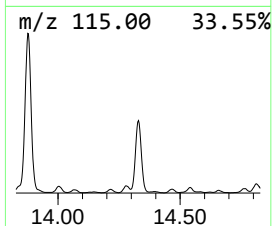
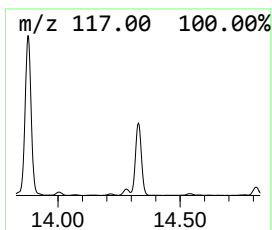
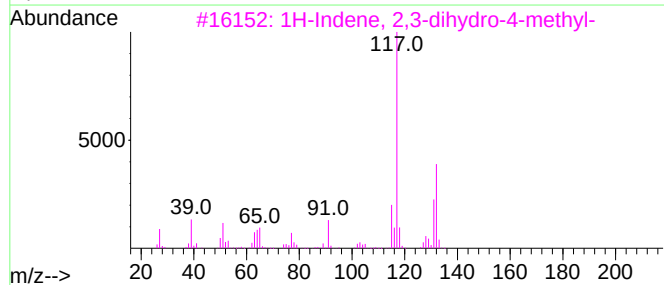
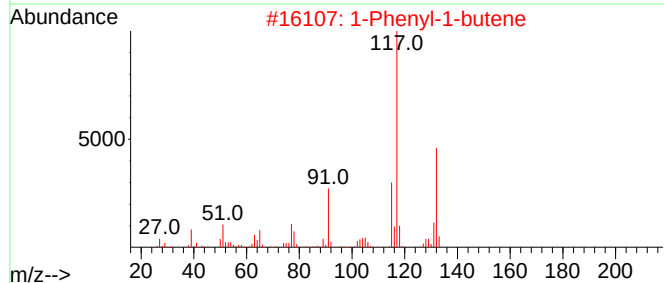
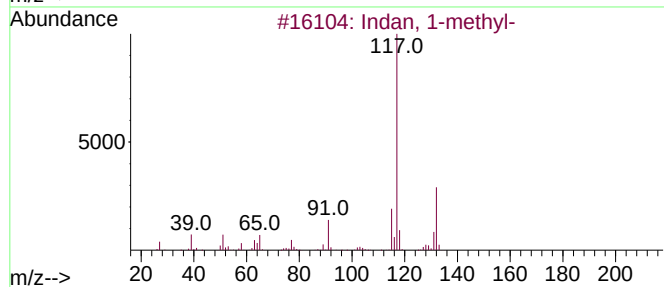
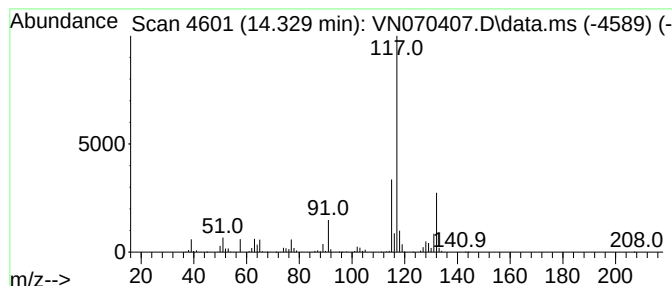
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Indan, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.329	41.29 ug/l	2758090	1,4-Dichlorobenzene-d4	13.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	80
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	80
4			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	80
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN123021\
Data File : VN070407.D
Acq On : 30 Dec 2021 18:32
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
DUP122121

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methyl-	3.135	143.5	ug/l	9649870	1	8.088	3362060	50.0
Butane, 2,3-dim...	5.074	101.3	ug/l	6810190	1	8.088	3362060	50.0
Pentane, 3-methyl-	5.616	29.3	ug/l	1967380	1	8.088	3362060	50.0
Pentane, 2,4-di...	7.045	23.9	ug/l	1610200	1	8.088	3362060	50.0
Cyclopentane, m...	7.147	26.9	ug/l	1807950	1	8.088	3362060	50.0
Pentane, 2,3-di...	8.174	60.4	ug/l	4062930	1	8.088	3362060	50.0
Pentane, 2,2,4-...	8.617	61.2	ug/l	5544350	2	8.971	4528920	50.0
Di-sec-Butyl ether	10.644	23.8	ug/l	2171650	3	11.746	4555550	50.0
Benzene, 1,1'-(...	13.876	85.1	ug/l	5682960	4	13.675	3339710	50.0
Indan, 1-methyl-	14.329	41.3	ug/l	2758090	4	13.675	3339710	50.0