

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
 Data File : VN087347.D
 Acq On : 17 Jul 2025 16:09
 Operator : JC\MD
 Sample : Q2620-01
 Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GAS-BUR-25-0010

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\82N071625W.M
 Title : SW846 8260

Signal : TIC: VN087347.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.312	55	61	66	rBV	25378	41690	1.92%	0.149%
2	2.895	142	160	161	rBV6	24347	78649	3.63%	0.281%
3	4.436	415	422	433	rBV2	9423	29613	1.37%	0.106%
4	6.271	723	734	746	rBV5	23014	72813	3.36%	0.260%
5	8.147	1043	1053	1057	rBV	209304	505733	23.34%	1.809%
6	8.200	1057	1062	1075	rVB	345919	804501	37.13%	2.878%
7	8.424	1091	1100	1107	rBV4	15866	34158	1.58%	0.122%
8	8.559	1113	1123	1134	rBV	248406	550777	25.42%	1.970%
9	8.941	1183	1188	1203	rVB3	14388	29500	1.36%	0.106%
10	9.083	1203	1212	1223	rBV	572382	1170209	54.01%	4.186%
11	10.547	1451	1461	1467	rBV	872161	1624542	74.98%	5.812%
12	10.612	1467	1472	1480	rVB	270268	497346	22.96%	1.779%
13	10.724	1485	1491	1498	rVB6	13441	29835	1.38%	0.107%
14	11.624	1635	1644	1648	rBV4	23705	52418	2.42%	0.188%
15	11.724	1656	1661	1668	rVB2	27713	46340	2.14%	0.166%
16	11.847	1674	1682	1693	rBV	715295	1321158	60.98%	4.727%
17	12.053	1708	1717	1721	rBV	139259	252423	11.65%	0.903%
18	12.129	1727	1730	1736	rVB3	24549	43505	2.01%	0.156%
19	12.265	1748	1753	1761	rVB5	12825	25025	1.16%	0.090%
20	12.353	1761	1768	1771	rBV5	55180	123863	5.72%	0.443%
21	12.459	1780	1786	1791	rVB2	68381	110955	5.12%	0.397%
22	12.577	1791	1806	1818	rBV9	70533	349223	16.12%	1.249%
23	12.682	1818	1824	1826	rVV5	33572	76823	3.55%	0.275%
24	12.724	1827	1831	1832	rVV3	55784	88668	4.09%	0.317%
25	12.747	1832	1835	1838	rVV	101924	181826	8.39%	0.650%
26	12.776	1838	1840	1844	rVV	101606	121476	5.61%	0.435%
27	12.829	1844	1849	1859	rVV2	550948	1091185	50.37%	3.904%
28	12.918	1860	1864	1867	rVB5	25222	37539	1.73%	0.134%
29	13.006	1874	1879	1885	rVB5	44500	90692	4.19%	0.324%
30	13.076	1885	1891	1895	rBV2	104699	206020	9.51%	0.737%
31	13.135	1895	1901	1909	rVV2	702628	1293481	59.70%	4.627%
32	13.206	1909	1913	1918	rVB3	76377	125859	5.81%	0.450%
33	13.294	1918	1928	1930	rBV4	74699	163487	7.55%	0.585%
34	13.371	1937	1941	1948	rVB2	196680	319274	14.74%	1.142%
35	13.459	1951	1956	1961	rBV	139460	232806	10.75%	0.833%
36	13.500	1961	1963	1968	rVB3	54291	63540	2.93%	0.227%

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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\82N071625W.M
Title : SW846 8260

37	13.553	1968	1972	1976	rBV4	67993	112901	5.21%	0.404%
38	13.594	1976	1979	1982	rVB4	41546	53372	2.46%	0.191%
39	13.653	1983	1989	1993	rBV5	198076	406732	18.77%	1.455%
40	13.694	1993	1996	1999	rVV	138669	181526	8.38%	0.649%
41	13.723	1999	2001	2004	rVV	70316	81011	3.74%	0.290%
42	13.771	2004	2009	2016	rVB2	753820	1350097	62.32%	4.830%
43	13.835	2016	2020	2026	rVB	204854	322992	14.91%	1.156%
44	13.923	2027	2035	2037	rBV7	97147	230758	10.65%	0.826%
45	13.959	2037	2041	2045	rBV2	119176	192987	8.91%	0.690%
46	14.082	2057	2062	2068	rBV	1352613	2166492	100.00%	7.751%
47	14.247	2084	2090	2094	rBV6	109389	221367	10.22%	0.792%
48	14.300	2096	2099	2104	rVB	131267	188558	8.70%	0.675%
49	14.341	2104	2106	2113	rVB4	146912	237698	10.97%	0.850%
50	14.418	2113	2119	2124	rBV5	184825	373503	17.24%	1.336%
51	14.476	2124	2129	2136	rBV8	89767	200289	9.24%	0.717%
52	14.559	2138	2143	2146	rBV4	213494	416202	19.21%	1.489%
53	14.606	2146	2151	2154	rVV3	459394	695660	32.11%	2.489%
54	14.641	2154	2157	2160	rVB2	176319	197812	9.13%	0.708%
55	14.712	2165	2169	2173	rBV4	227735	320562	14.80%	1.147%
56	14.770	2173	2179	2183	rVB6	148145	301360	13.91%	1.078%
57	14.823	2184	2188	2191	rBV4	107737	179063	8.27%	0.641%
58	14.935	2198	2207	2215	rBV	1146737	1926536	88.92%	6.892%
59	15.023	2215	2222	2225	rBV6	289488	714246	32.97%	2.555%
60	15.188	2247	2250	2256	rVB3	148313	278687	12.86%	0.997%
61	15.241	2257	2259	2263	rVB4	42986	50196	2.32%	0.180%
62	15.294	2263	2268	2273	rBV5	133497	232959	10.75%	0.833%
63	15.482	2296	2300	2309	rVB5	173543	418906	19.34%	1.499%
64	15.565	2309	2314	2320	rBV	277560	535910	24.74%	1.917%
65	15.806	2349	2355	2365	rVB	763941	1472842	67.98%	5.269%
66	15.917	2365	2374	2381	rBV5	156395	495226	22.86%	1.772%
67	16.141	2407	2412	2423	rVB4	195817	407254	18.80%	1.457%
68	16.459	2456	2466	2472	rBV6	153246	570148	26.32%	2.040%
69	16.582	2483	2487	2496	rVB2	101906	185681	8.57%	0.664%
70	16.676	2498	2503	2510	rVB4	114931	242975	11.22%	0.869%
71	16.759	2511	2517	2518	rBV4	55429	102646	4.74%	0.367%

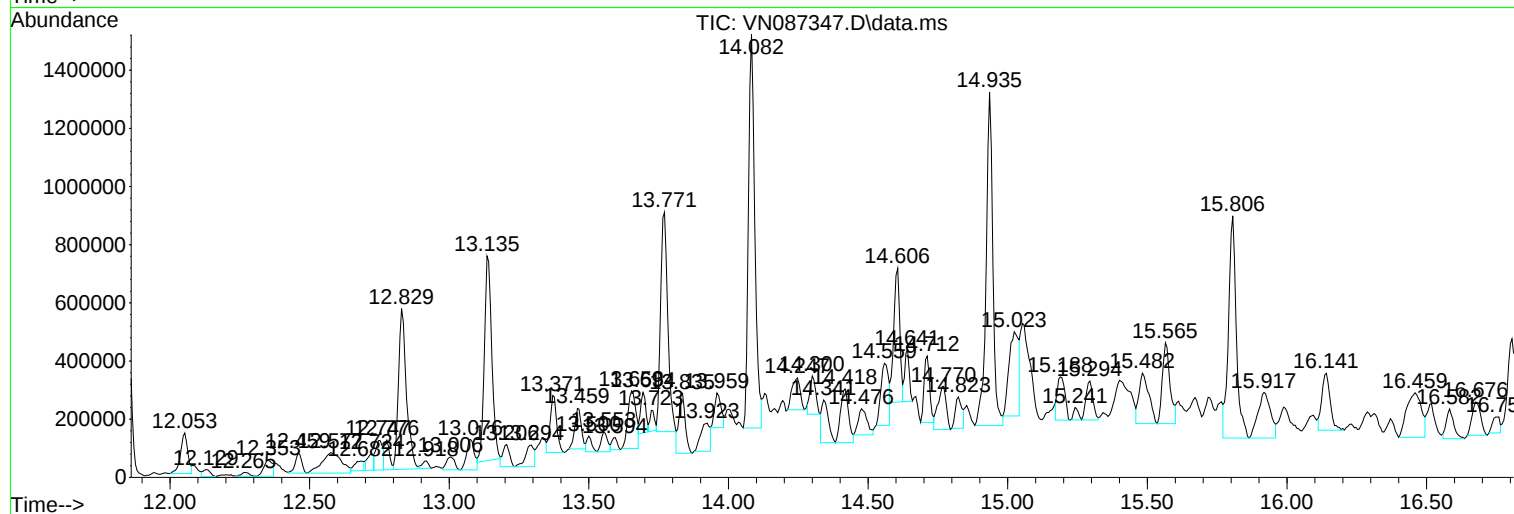
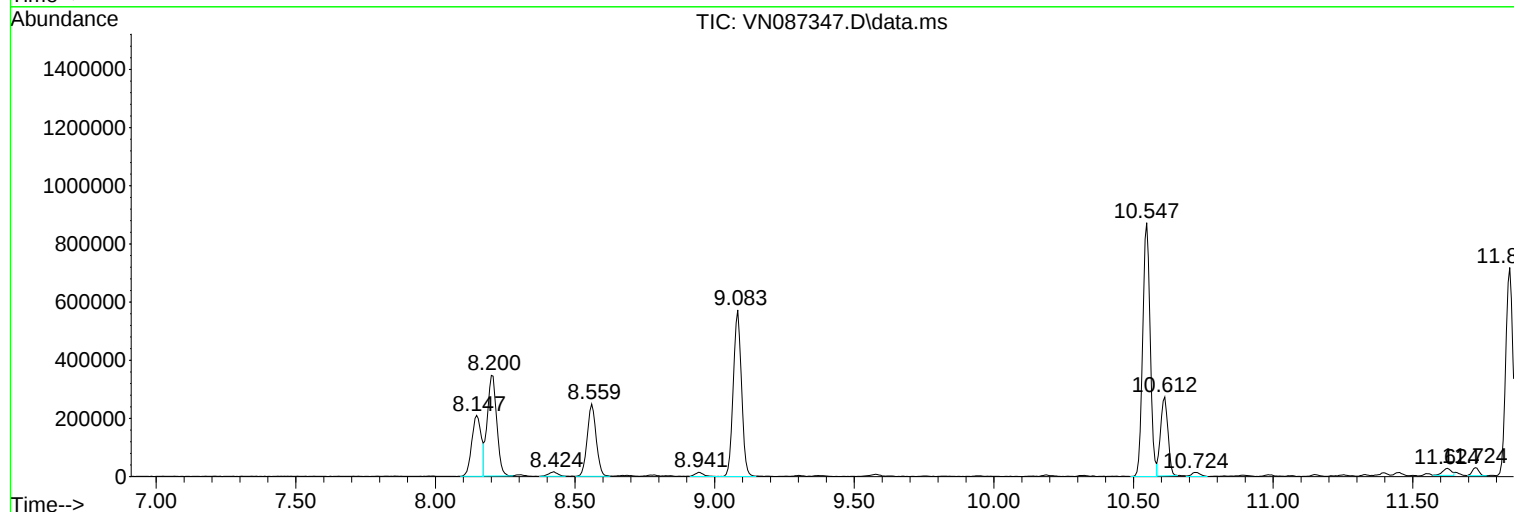
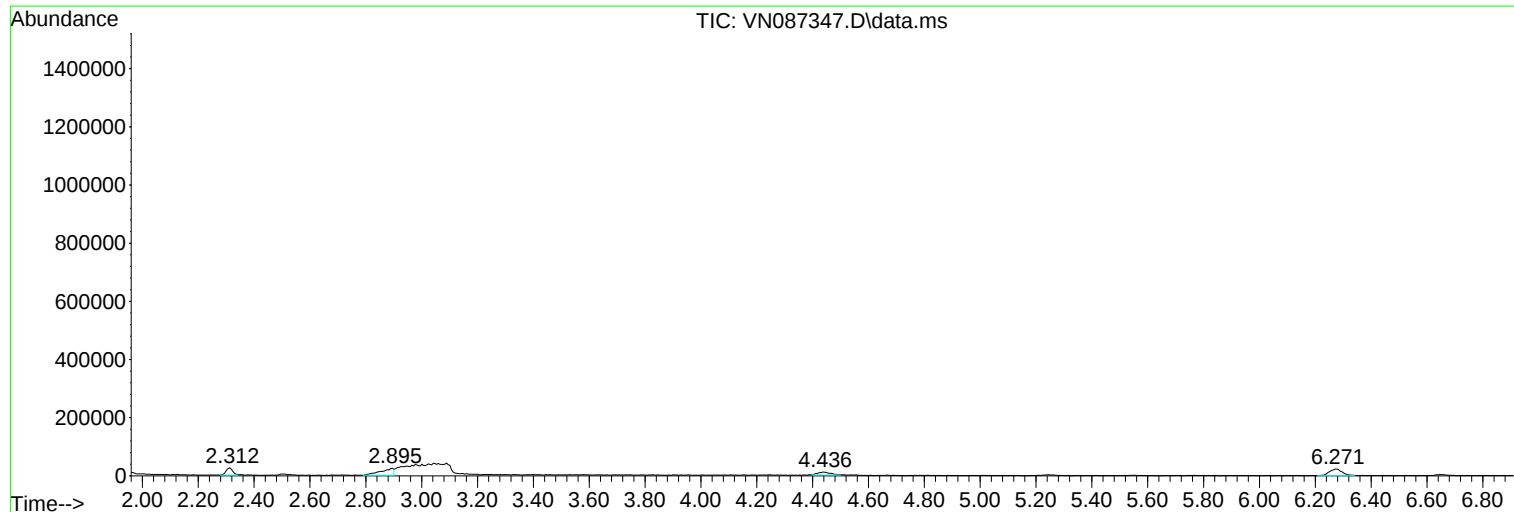
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Instrument :
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ClientSampleId :
GAS-BUR-25-0010

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.M
Quant Title : SW846 8260

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



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Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

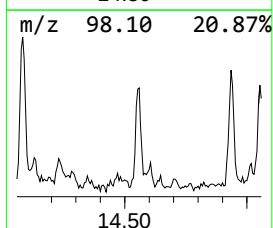
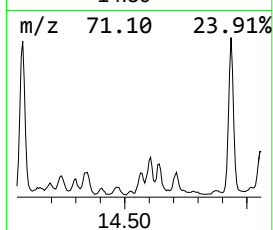
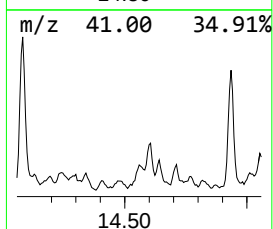
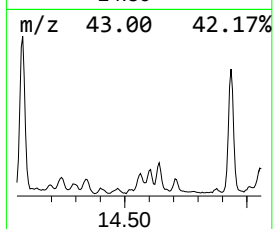
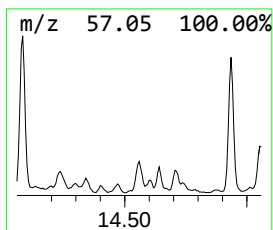
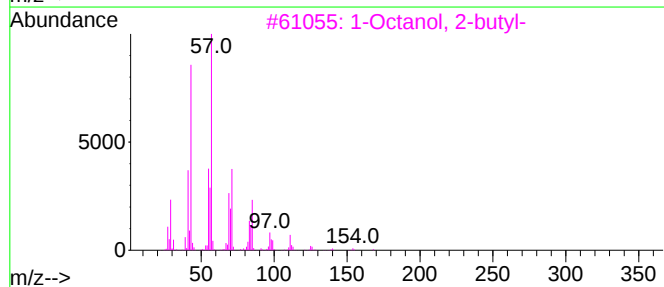
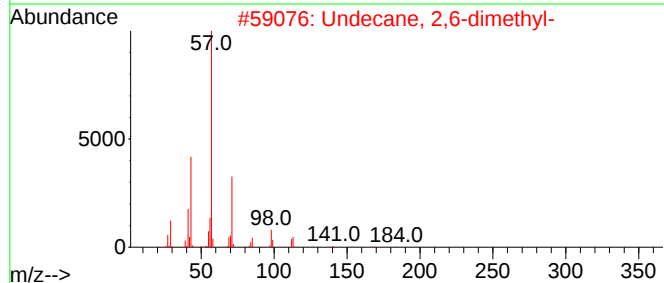
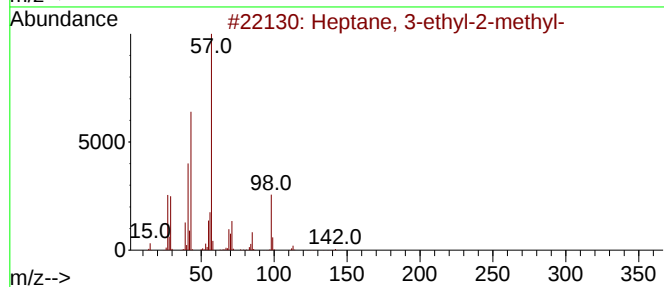
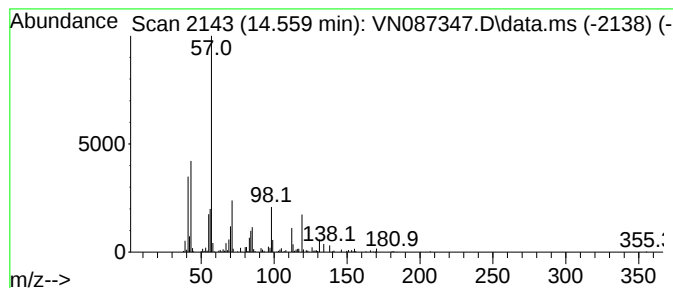
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.559	15.41 ug/l	416202	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50
2			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	43
3			1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	43
4			Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	43
5			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	38



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Instrument :
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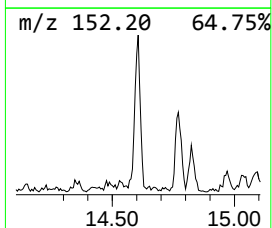
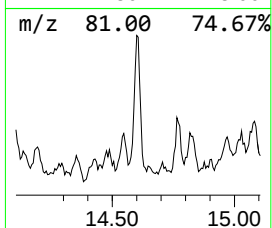
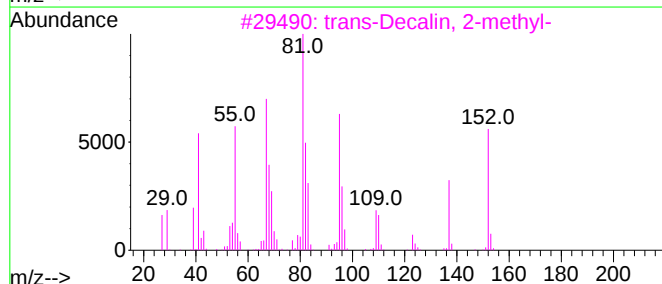
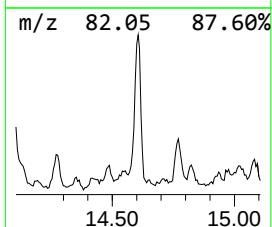
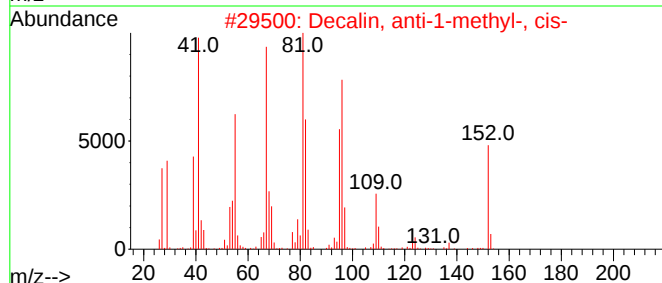
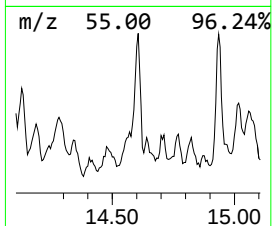
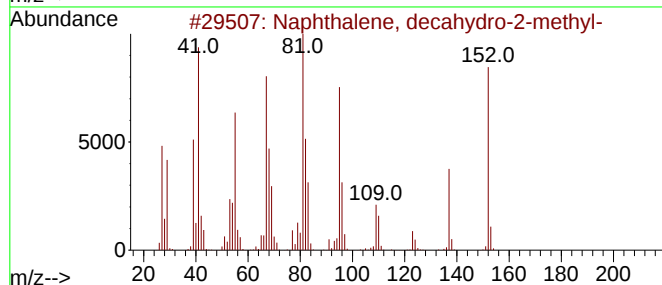
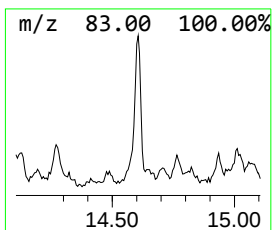
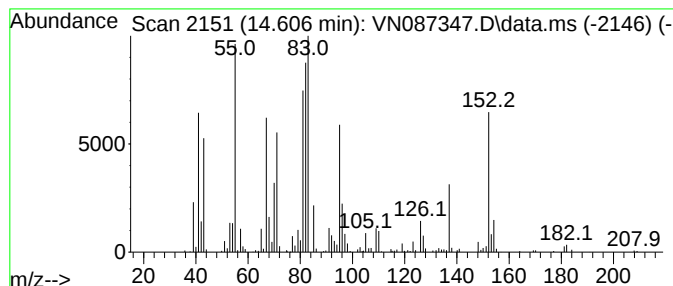
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.M
Quant Title : SW846 8260

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

Peak Number 2 Naphthalene, decahydro-2-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.606	25.76 ug/l	695660	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	70
2			Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	50
3			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	47
4			Pulegone	152	C10H16O	000089-82-7	46
5			cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	43



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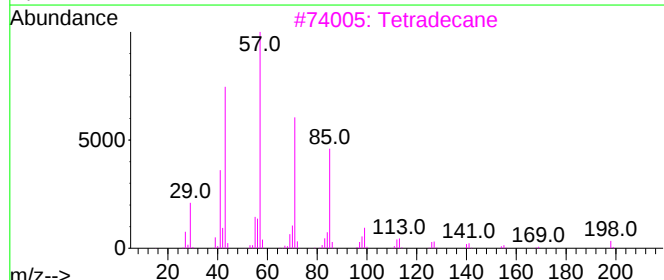
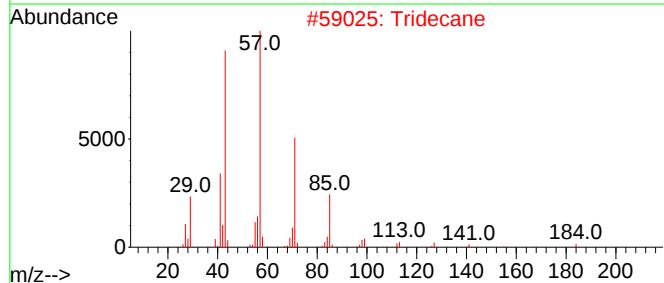
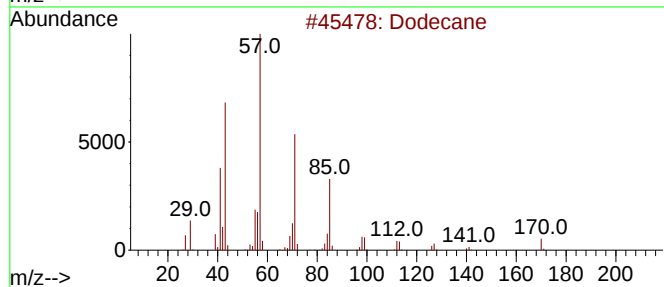
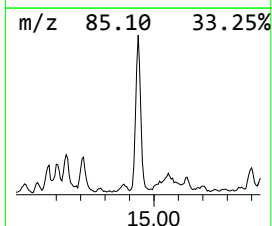
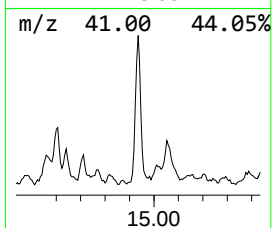
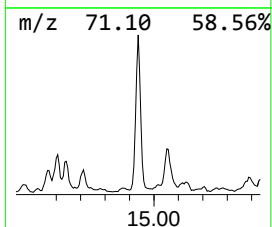
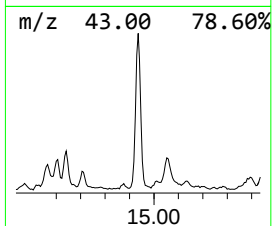
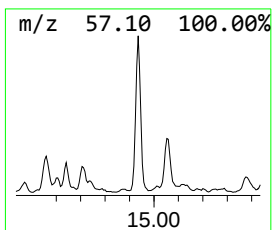
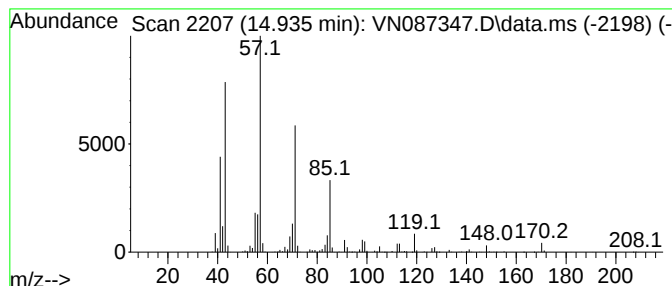
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.M
Quant Title : SW846 8260

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

Peak Number 3 Dodecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.935	71.35 ug/l	1926540	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	96
2			Tridecane	184	C13H28	000629-50-5	86
3			Tetradecane	198	C14H30	000629-59-4	86
4			Undecane	156	C11H24	001120-21-4	80
5			Pentadecane	212	C15H32	000629-62-9	72



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GAS-BUR-25-0010

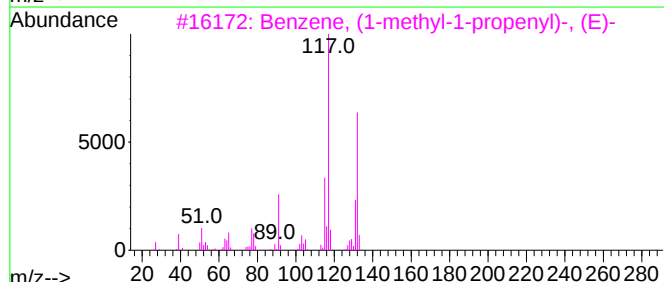
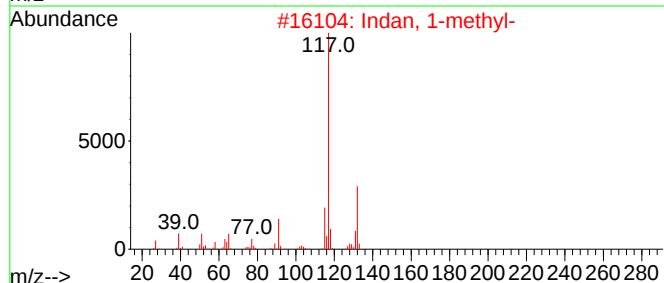
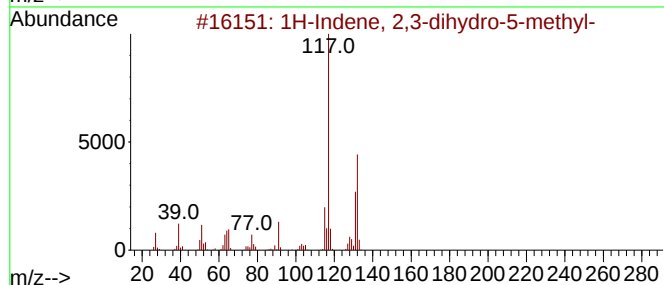
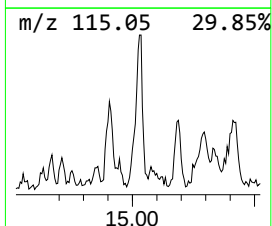
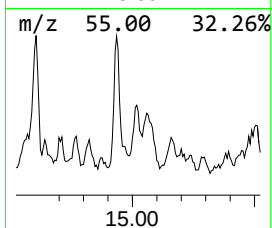
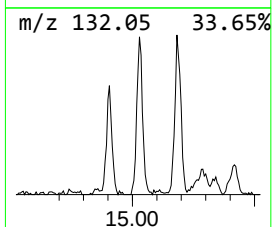
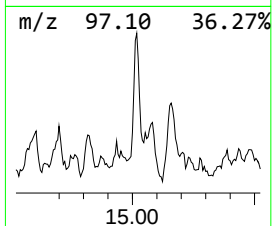
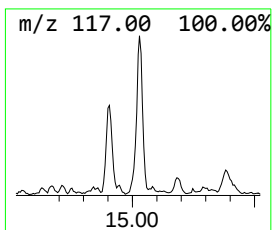
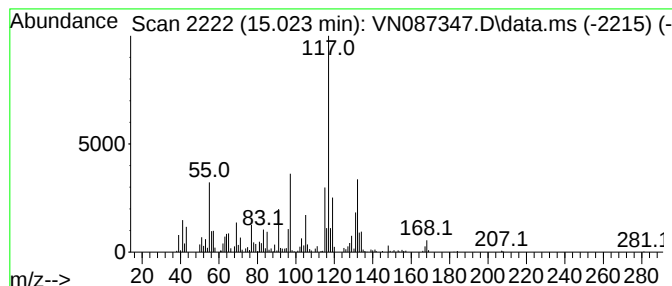
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Quant Title : SW846 8260

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

Peak Number 4 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.023	26.45 ug/l	714246	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2			Indan, 1-methyl-	132	C10H12	000767-58-8	70
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	62
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
Acq On : 17 Jul 2025 16:09
Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

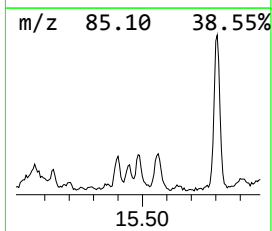
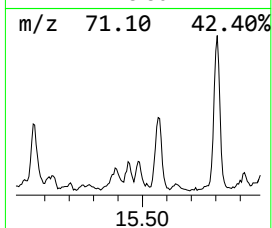
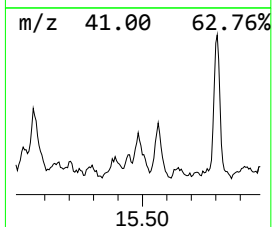
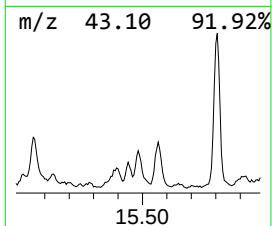
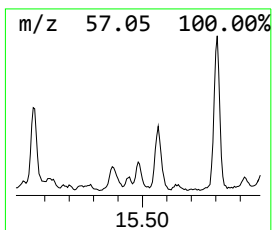
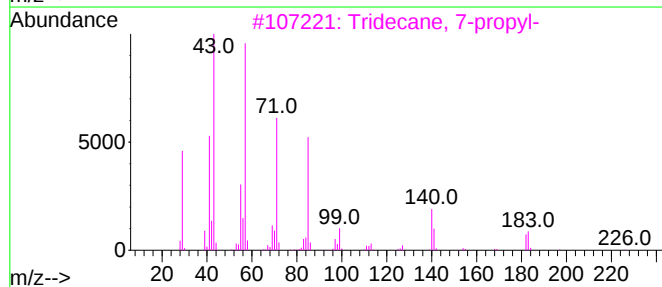
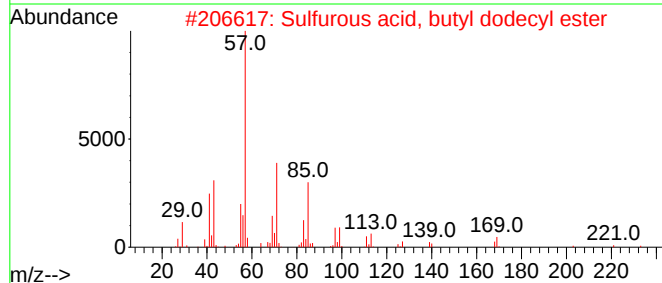
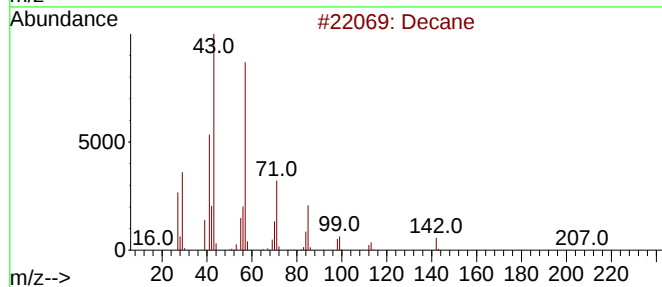
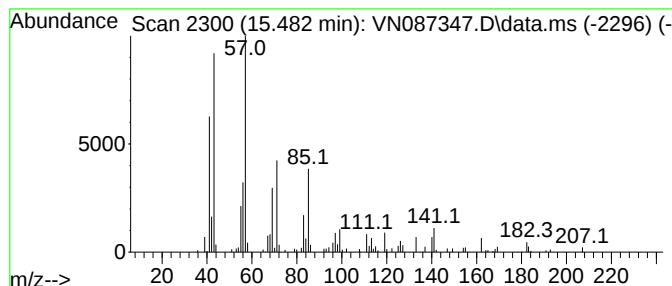
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Quant Title : SW846 8260

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

Peak Number 5 Decane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.482	15.51 ug/l	418906	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	68
2			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	64
3			Tridecane, 7-propyl-	226	C16H34	055045-09-5	59
4			Octacosane	394	C28H58	000630-02-4	59
5			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
Acq On : 17 Jul 2025 16:09
Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

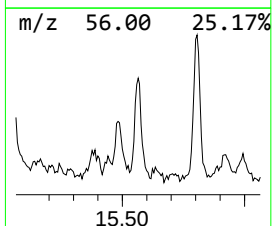
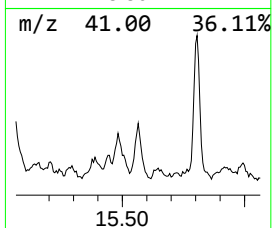
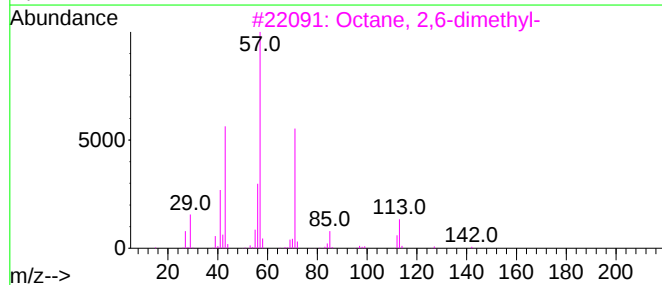
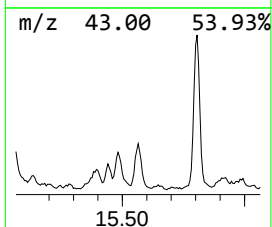
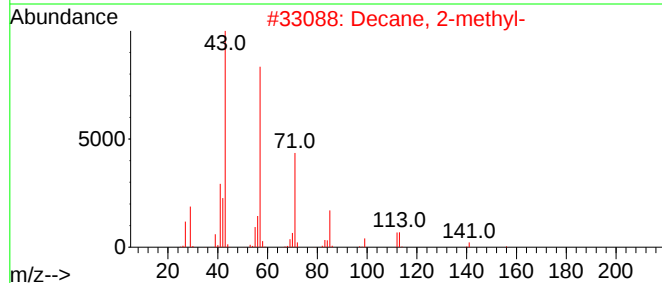
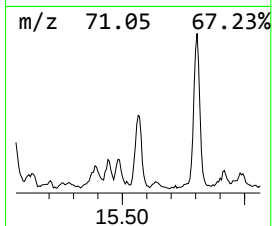
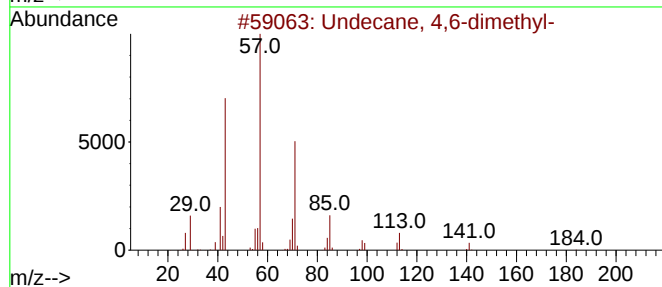
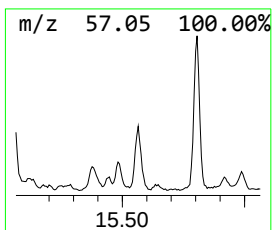
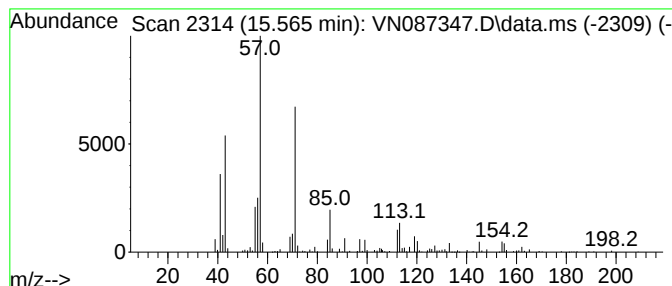
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Quant Title : SW846 8260

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

Peak Number 6 Undecane, 4,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.565	19.85 ug/l	535910	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	76
2			Decane, 2-methyl-	156	C11H24	006975-98-0	68
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	68
4			Undecane, 6-ethyl-	184	C13H28	017312-60-6	64
5			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
Acq On : 17 Jul 2025 16:09
Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

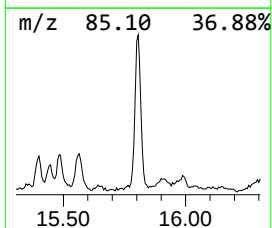
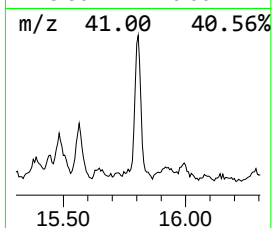
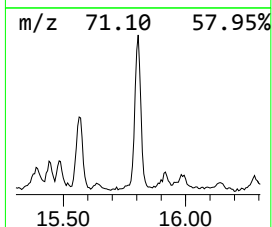
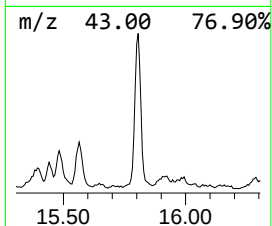
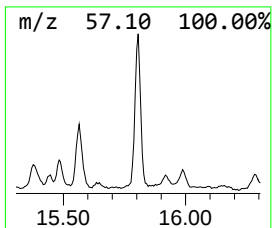
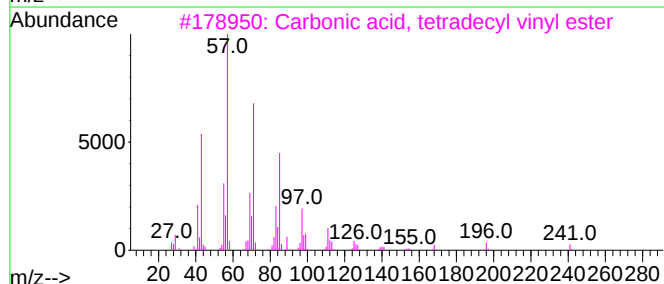
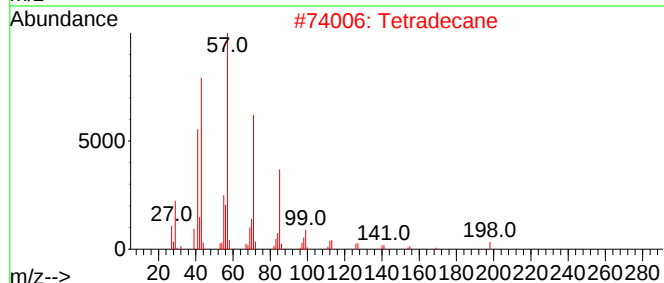
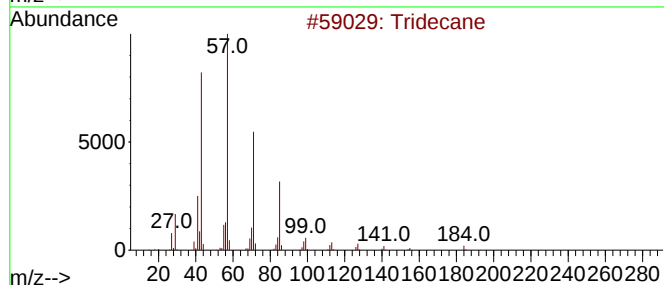
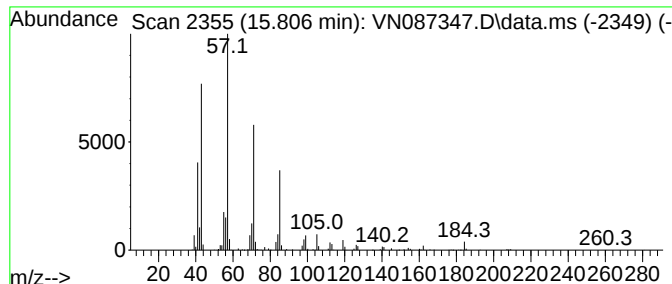
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Quant Title : SW846 8260

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

Peak Number 7 Tridecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.806	54.55 ug/l	1472840	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Tetradecane	198	C14H30	000629-59-4	87
3			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	86
4			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	86
5			Hexadecane	226	C16H34	000544-76-3	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
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Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

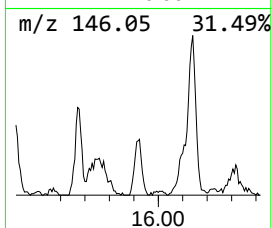
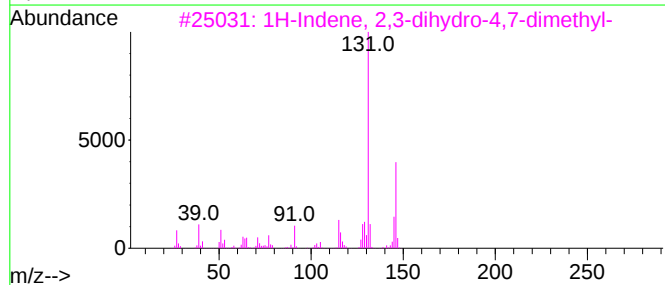
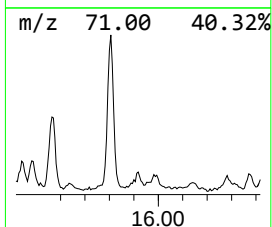
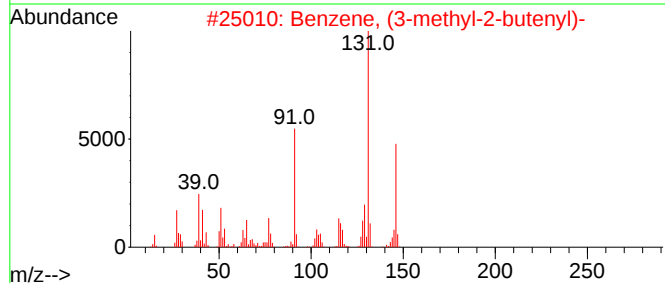
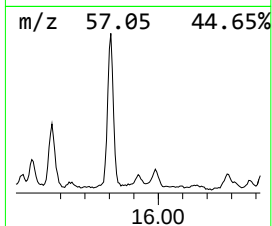
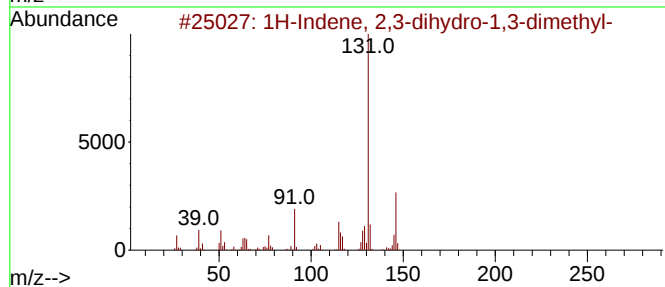
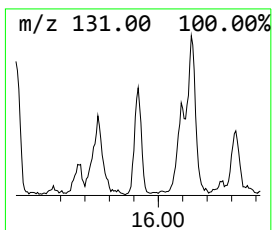
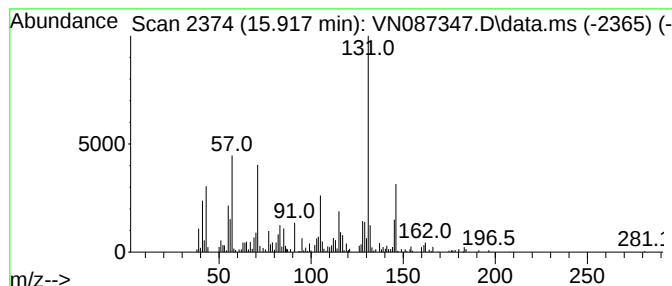
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.918	18.34 ug/l	495226	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	83
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	83
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
Acq On : 17 Jul 2025 16:09
Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

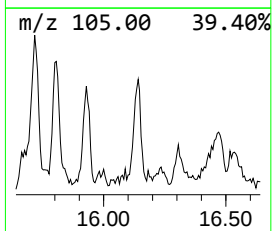
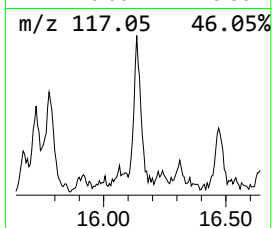
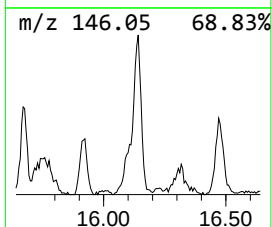
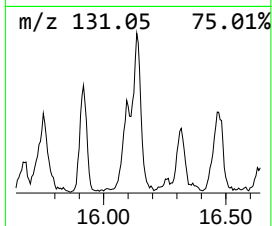
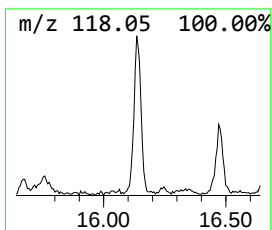
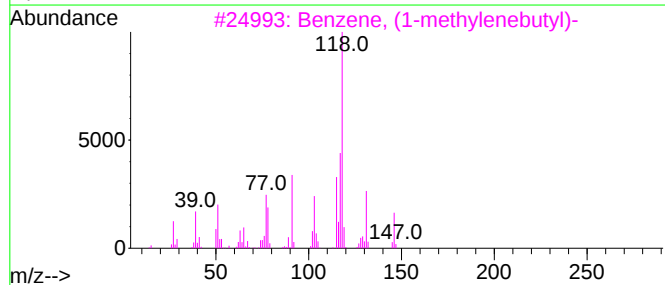
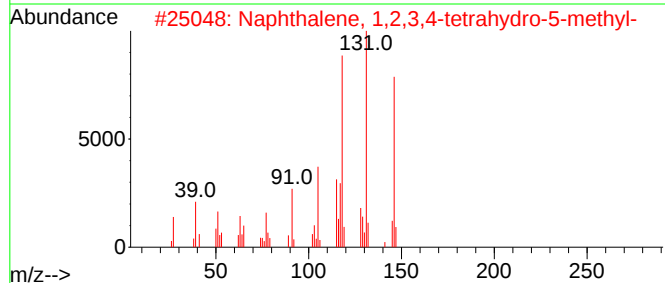
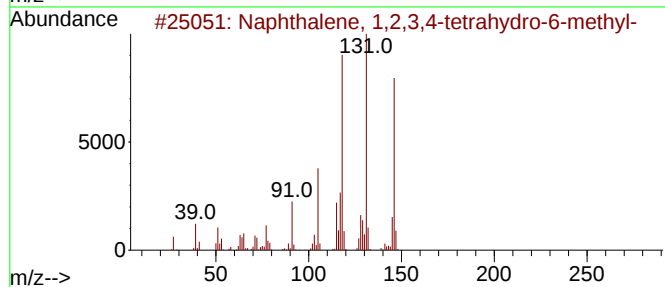
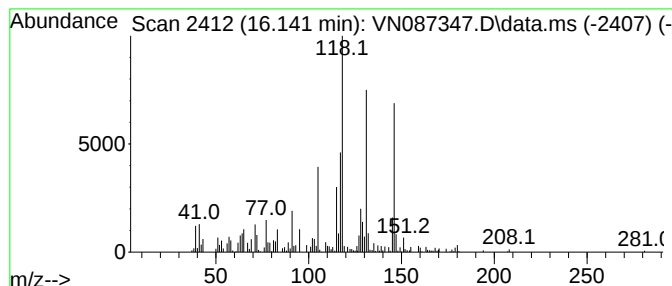
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.141	15.08 ug/l	407254	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	87
3			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	80
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	49
5			1H-Inden-1-one, 2,3-dihydro-3-me...	146	C10H10O	006072-57-7	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
Acq On : 17 Jul 2025 16:09
Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

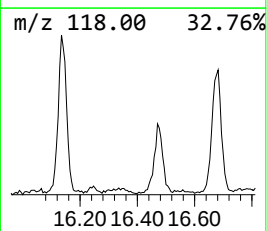
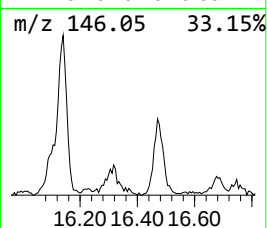
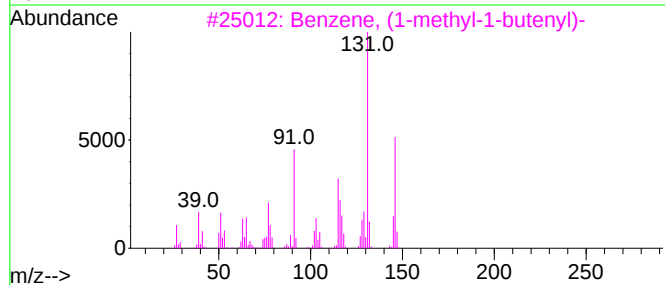
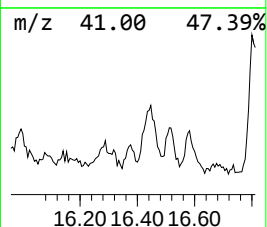
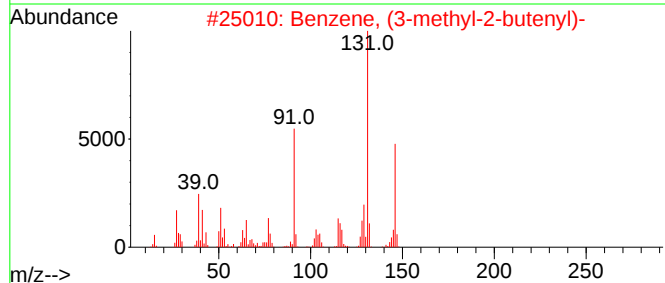
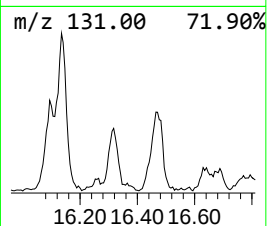
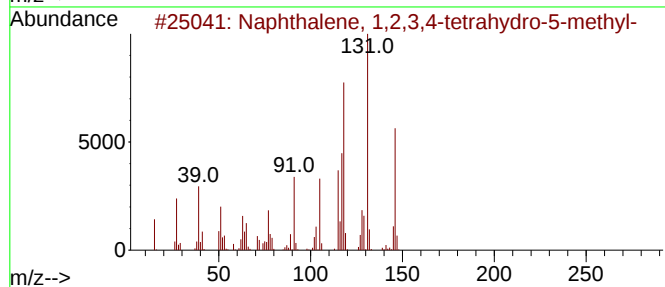
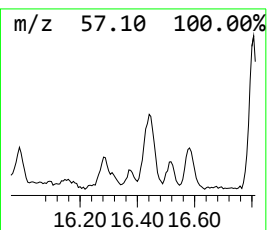
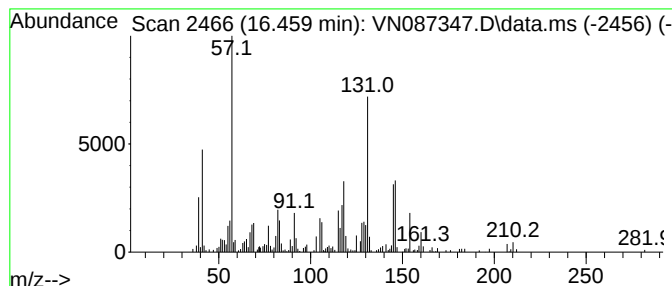
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.459	21.12 ug/l	570148	1,4-Dichlorobenzene-d4	13.771

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	70
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	50
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	50
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	49
5			Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071725\
Data File : VN087347.D
Acq On : 17 Jul 2025 16:09
Operator : JC\MD
Sample : Q2620-01
Misc : 1.00g/10.0mL/100uL/5.0mL/MSVOA_N/MEOH
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
GAS-BUR-25-0010

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071625W.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
Heptane, 3-ethy...	14.559	15.4	ug/l	416202	4	13.771	1350100	50.0
Naphthalene, de...	14.606	25.8	ug/l	695660	4	13.771	1350100	50.0
Dodecane	14.935	71.3	ug/l	1926540	4	13.771	1350100	50.0
1H-Indene, 2,3-...	15.023	26.4	ug/l	714246	4	13.771	1350100	50.0
Decane	15.482	15.5	ug/l	418906	4	13.771	1350100	50.0
Undecane, 4,6-d...	15.565	19.9	ug/l	535910	4	13.771	1350100	50.0
Tridecane	15.806	54.5	ug/l	1472840	4	13.771	1350100	50.0
1H-Indene, 2,3-...	15.918	18.3	ug/l	495226	4	13.771	1350100	50.0
Naphthalene, 1,...	16.141	15.1	ug/l	407254	4	13.771	1350100	50.0
Naphthalene, 1,...	16.459	21.1	ug/l	570148	4	13.771	1350100	50.0