

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
 Data File : VX043116.D
 Acq On : 24 Sep 2024 12:21
 Operator : JC/MD
 Sample : P4123-02
 Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 613-D

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_X\Method\82X092324W.M
 Title : SW846 8260

Signal : TIC: VX043116.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.391	695	706	712	rBV	55278	168718	1.17%	0.162%
2	5.483	712	721	734	rVV2	555116	1906537	13.25%	1.833%
3	5.599	734	740	755	rVB	200191	638862	4.44%	0.614%
4	5.806	760	774	789	rVV	1106016	3195234	22.20%	3.072%
5	5.952	790	798	817	rVV	63278	181277	1.26%	0.174%
6	6.172	819	834	841	rVV3	163727	598508	4.16%	0.575%
7	6.245	841	846	853	rVV2	69318	183748	1.28%	0.177%
8	6.342	853	862	874	rVB2	88016	245579	1.71%	0.236%
9	6.598	891	904	922	rBV	2548473	6317513	43.89%	6.074%
10	6.757	922	930	947	rVB	148798	374574	2.60%	0.360%
11	7.232	999	1008	1018	rBV	102432	236751	1.64%	0.228%
12	7.367	1018	1030	1042	rVV3	401550	1039855	7.22%	1.000%
13	7.488	1042	1050	1055	rVV	128698	262857	1.83%	0.253%
14	7.549	1055	1060	1069	rVV	185998	390029	2.71%	0.375%
15	7.647	1069	1076	1085	rVB	201454	397800	2.76%	0.382%
16	8.647	1234	1240	1248	rBV	320308	500701	3.48%	0.481%
17	10.055	1466	1471	1485	rVB	336044	471662	3.28%	0.454%
18	11.085	1634	1640	1650	rVB	303437	449145	3.12%	0.432%
19	11.476	1700	1704	1710	rVB	429901	505705	3.51%	0.486%
20	11.713	1740	1743	1746	rBV	289052	295109	2.05%	0.284%
21	11.835	1759	1763	1766	rBV	228188	309817	2.15%	0.298%
22	11.890	1766	1772	1776	rVV	224401	353033	2.45%	0.339%
23	11.939	1776	1780	1783	rVV2	345564	488994	3.40%	0.470%
24	11.975	1783	1786	1788	rVV2	141097	164791	1.14%	0.158%
25	12.043	1790	1797	1799	rVV2	855264	1488099	10.34%	1.431%
26	12.073	1799	1802	1804	rVV	899648	976653	6.79%	0.939%
27	12.103	1804	1807	1811	rVV	1372510	1564770	10.87%	1.505%
28	12.146	1811	1814	1816	rVV	395485	540682	3.76%	0.520%
29	12.177	1816	1819	1823	rVB	1192923	1376914	9.57%	1.324%
30	12.219	1823	1826	1830	rBV	262898	374068	2.60%	0.360%
31	12.262	1830	1833	1836	rVB2	268197	321483	2.23%	0.309%
32	12.305	1836	1840	1847	rBV5	605742	1366028	9.49%	1.313%
33	12.384	1847	1853	1855	rBV3	474354	819892	5.70%	0.788%
34	12.427	1855	1860	1864	rVV	10671625	14393987	100.00%	13.840%
35	12.469	1864	1867	1874	rVB3	845702	1687021	11.72%	1.622%
36	12.555	1875	1881	1883	rBV3	940030	1894404	13.16%	1.821%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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37	12.585	1883	1886	1891	rVV3	1684114	2924151	20.32%	2.812%
38	12.640	1891	1895	1899	rVV2	711493	1086231	7.55%	1.044%
39	12.683	1899	1902	1906	rVV2	1262104	1591958	11.06%	1.531%
40	12.738	1906	1911	1913	rVV3	455659	798320	5.55%	0.768%
41	12.762	1913	1915	1920	rVV3	772714	1351001	9.39%	1.299%
42	12.817	1920	1924	1930	rVB3	1682461	2568806	17.85%	2.470%
43	12.890	1930	1936	1938	rBV3	2512478	4223809	29.34%	4.061%
44	12.945	1942	1945	1947	rVV	1717456	1913296	13.29%	1.840%
45	12.981	1947	1951	1955	rVV2	2759849	3965245	27.55%	3.813%
46	13.042	1958	1961	1966	rVB	2387177	2698256	18.75%	2.594%
47	13.164	1977	1981	1982	rBV2	311404	421591	2.93%	0.405%
48	13.213	1986	1989	1991	rBV3	289228	342140	2.38%	0.329%
49	13.274	1994	1999	2002	rBV	9107772	11271374	78.31%	10.838%
50	13.347	2008	2011	2014	rBV2	608866	779010	5.41%	0.749%
51	13.384	2014	2017	2020	rBV	1288601	1558425	10.83%	1.498%
52	13.414	2020	2022	2026	rVB4	618222	833976	5.79%	0.802%
53	13.500	2032	2036	2041	rBV4	444246	699875	4.86%	0.673%
54	13.548	2042	2044	2047	rVB2	183620	193028	1.34%	0.186%
55	13.591	2047	2051	2055	rBV5	305405	535853	3.72%	0.515%
56	13.682	2061	2066	2067	rBV3	543499	803780	5.58%	0.773%
57	13.737	2071	2075	2078	rVV2	856614	1205639	8.38%	1.159%
58	13.768	2078	2080	2083	rVV2	525149	480735	3.34%	0.462%
59	13.798	2083	2085	2088	rVV3	310447	343420	2.39%	0.330%
60	13.841	2088	2092	2099	rVB4	655524	1390233	9.66%	1.337%
61	13.926	2099	2106	2111	rVB7	346657	862624	5.99%	0.829%
62	14.036	2120	2124	2128	rVB	1385183	1448769	10.07%	1.393%
63	14.073	2128	2130	2133	rBV2	252134	247166	1.72%	0.238%
64	14.225	2150	2155	2161	rVB3	750049	1309446	9.10%	1.259%
65	14.323	2167	2171	2176	rBV4	135651	197322	1.37%	0.190%
66	14.371	2176	2179	2182	rVV	273977	292300	2.03%	0.281%
67	14.414	2183	2186	2188	rVV3	113827	164761	1.14%	0.158%
68	14.481	2192	2197	2206	rVB6	490468	1047330	7.28%	1.007%
69	14.560	2206	2210	2215	rVB6	182468	357633	2.48%	0.344%
70	14.615	2215	2219	2225	rBV2	818934	1320302	9.17%	1.269%
71	14.670	2225	2228	2233	rVB2	379366	516781	3.59%	0.497%
72	14.731	2233	2238	2239	rBV2	216478	295935	2.06%	0.285%
73	14.755	2239	2242	2245	rVB	582537	572565	3.98%	0.551%
74	14.847	2254	2257	2260	rVB3	243390	265973	1.85%	0.256%
75	14.902	2263	2266	2276	rVB3	168711	364282	2.53%	0.350%
76	15.042	2284	2289	2293	rBV4	132473	236843	1.65%	0.228%

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77	15.103	2294	2299	2302	rBV4	114326	215092	1.49%	0.207%
78	15.182	2307	2312	2314	rVV2	414415	630626	4.38%	0.606%
79	15.207	2314	2316	2320	rVV	306694	383099	2.66%	0.368%
80	15.274	2324	2327	2331	rVB3	166049	218806	1.52%	0.210%
81	15.371	2340	2343	2348	rVB3	124105	163663	1.14%	0.157%
82	15.487	2356	2362	2367	rBV2	336300	551771	3.83%	0.531%
83	15.560	2369	2374	2378	rVV	1062934	1536287	10.67%	1.477%
84	15.609	2378	2382	2386	rVB2	119864	176385	1.23%	0.170%
85	15.877	2423	2426	2433	rVB5	85383	166437	1.16%	0.160%

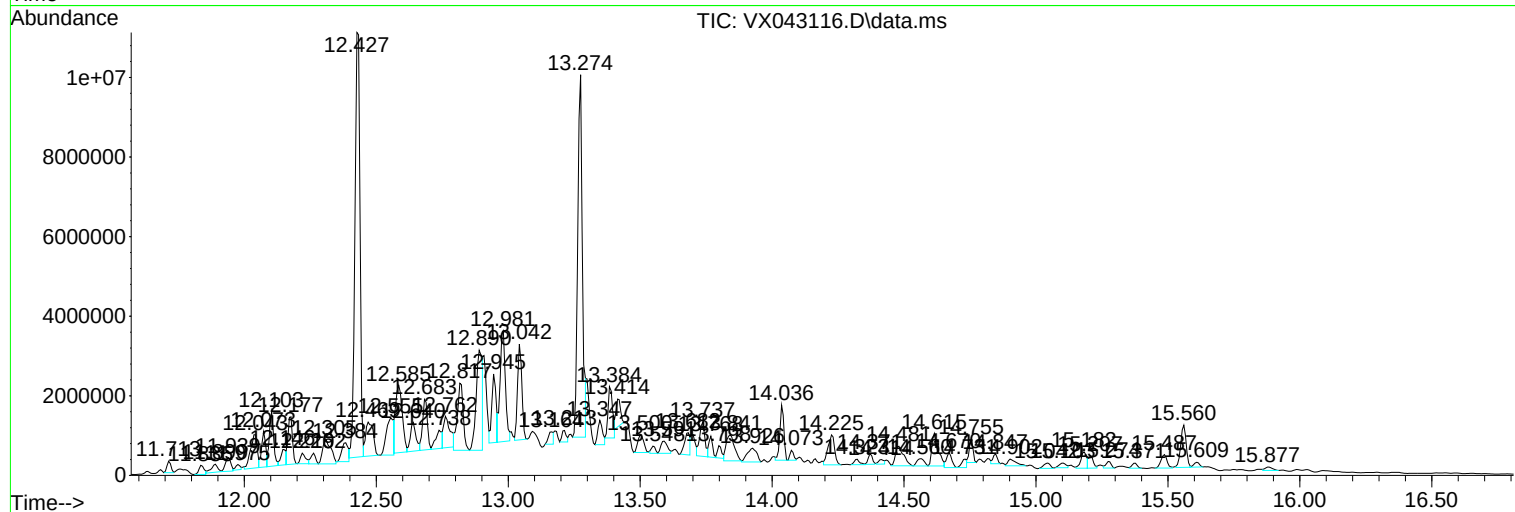
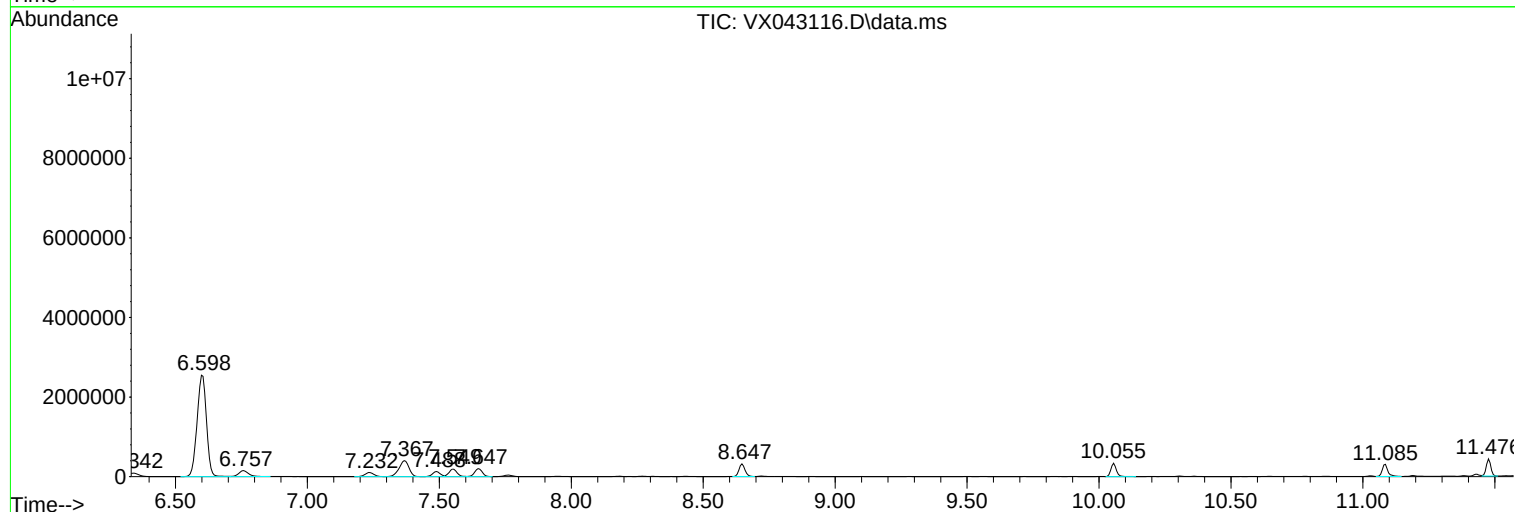
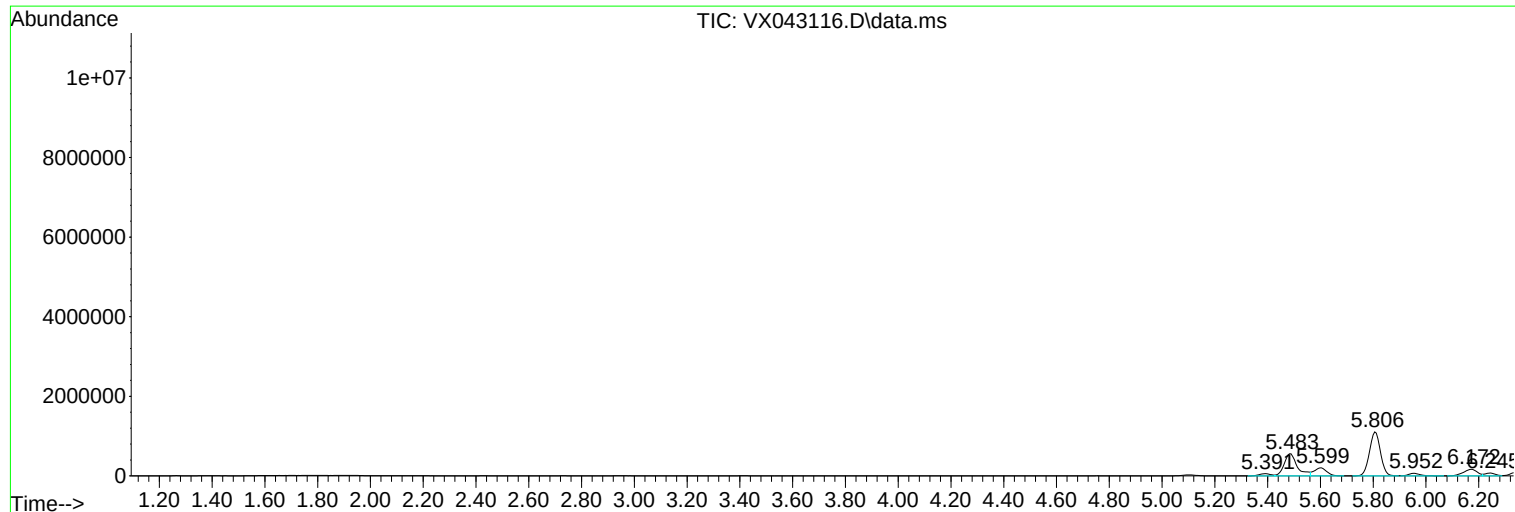
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
Quant Title : SW846 8260

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



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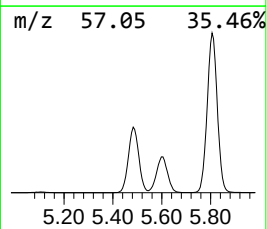
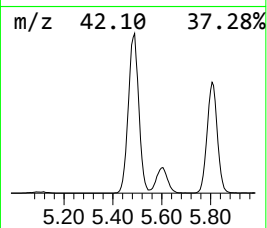
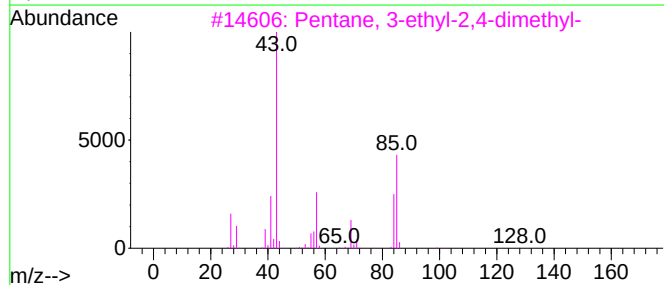
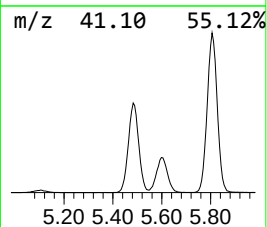
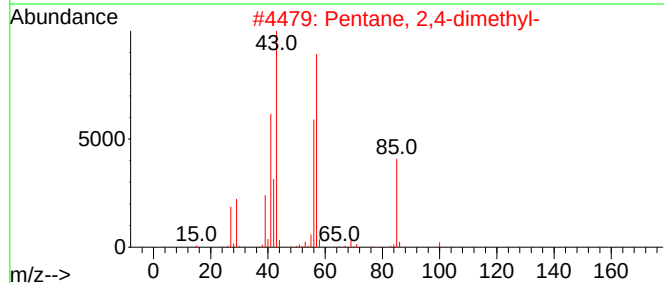
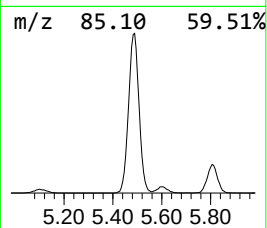
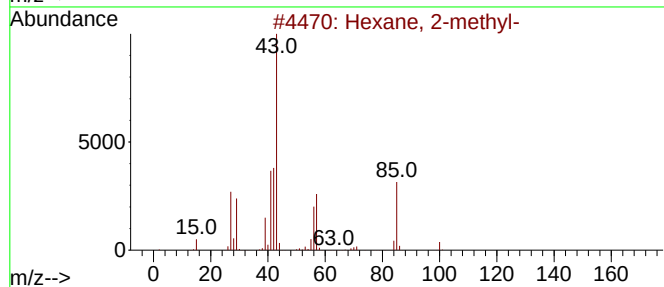
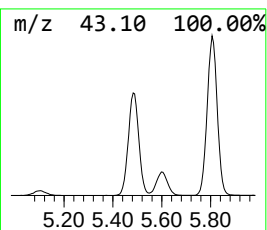
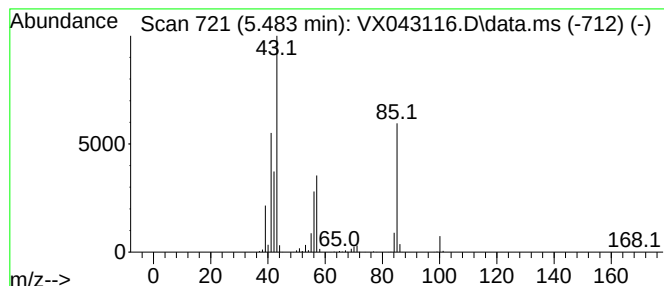
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.483	149.21 ug/l	1906540	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2-methyl-	100	C7H16	000591-76-4	90
2			Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
3			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	50
4			Heptane, 4,4-dimethyl-	128	C9H20	001068-19-5	50
5			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	50



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Instrument :
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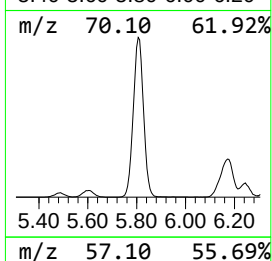
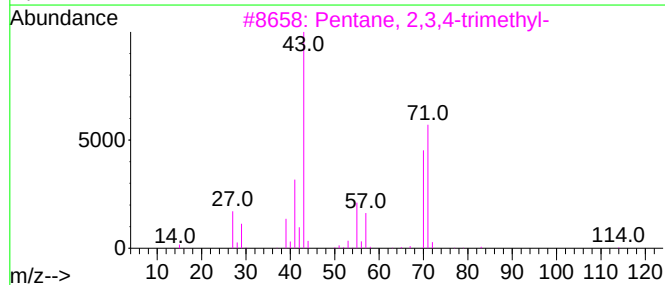
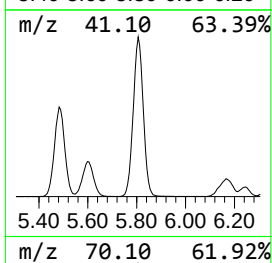
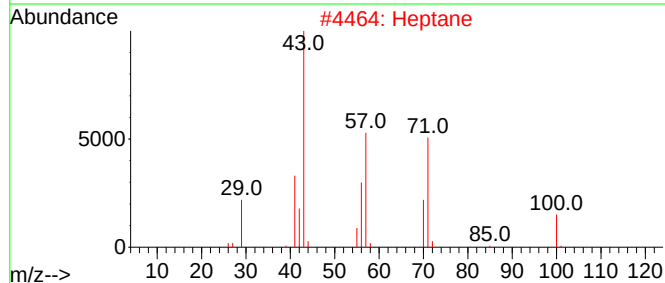
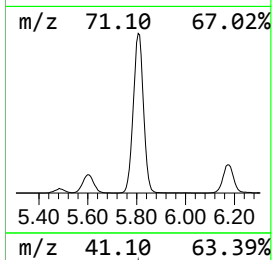
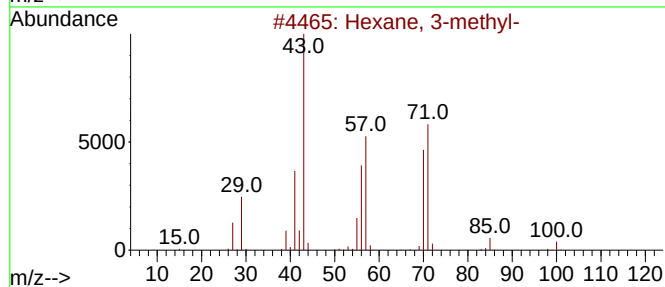
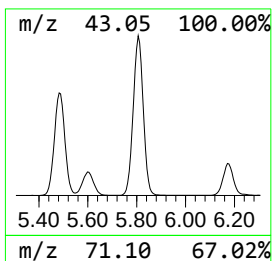
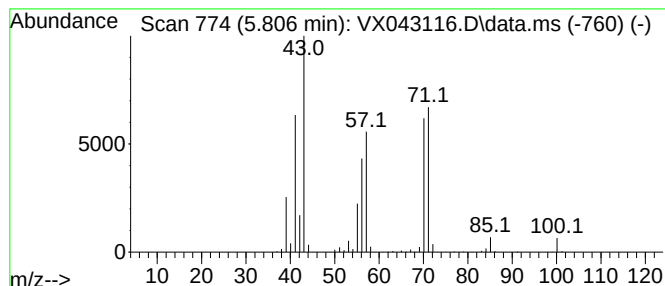
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
Quant Title : SW846 8260

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

Peak Number 2 Hexane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.806	250.07 ug/l	3195230	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Heptane	100	C7H16	000142-82-5	80
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
4			Heptane, 4-methyl-	114	C8H18	000589-53-7	53
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	53



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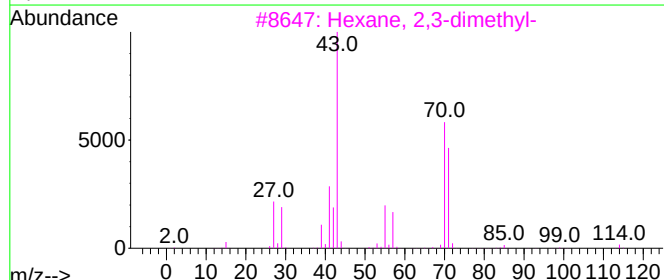
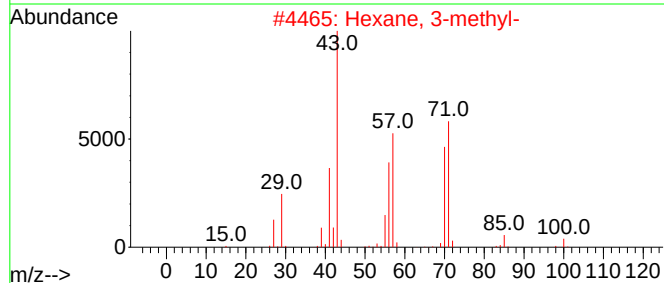
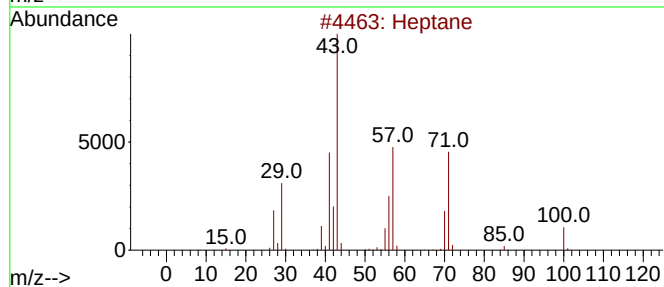
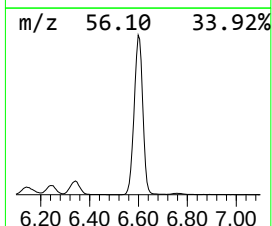
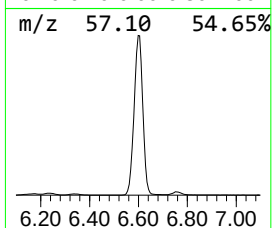
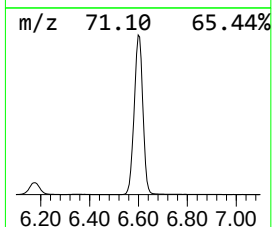
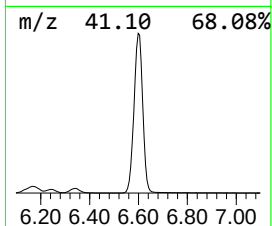
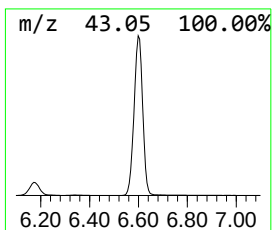
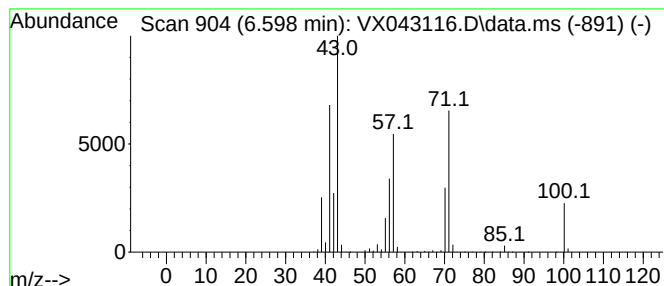
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.M
Quant Title : SW846 8260

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

Peak Number 3 Heptane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.598	843.29 ug/l	6317510	1,4-Difluorobenzene	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	40
3			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
4			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	38
5			Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	37



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613-D

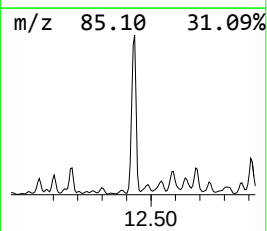
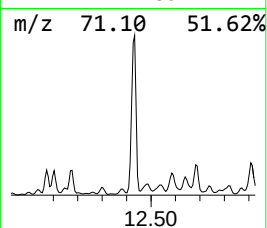
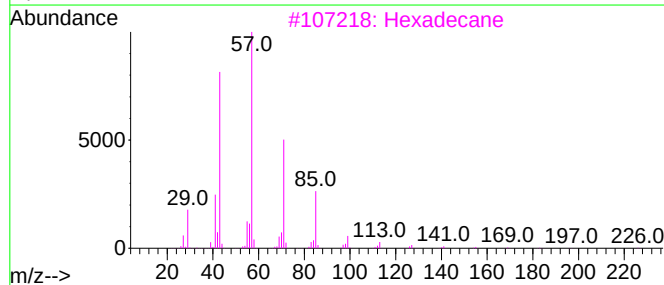
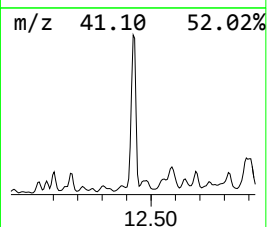
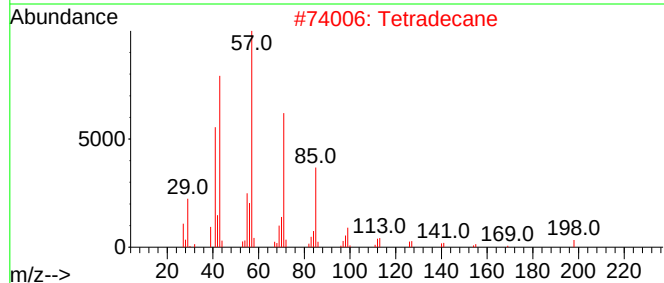
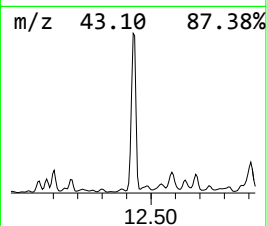
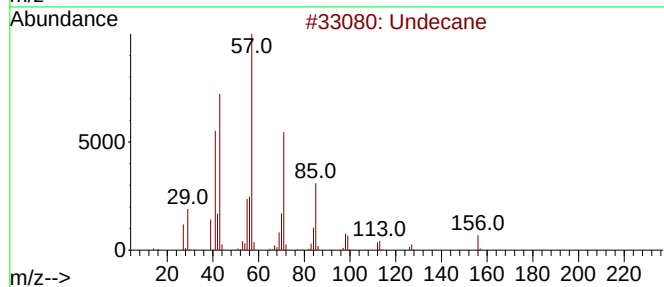
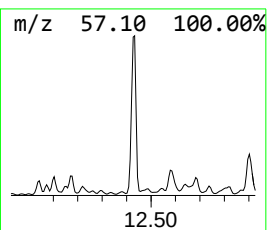
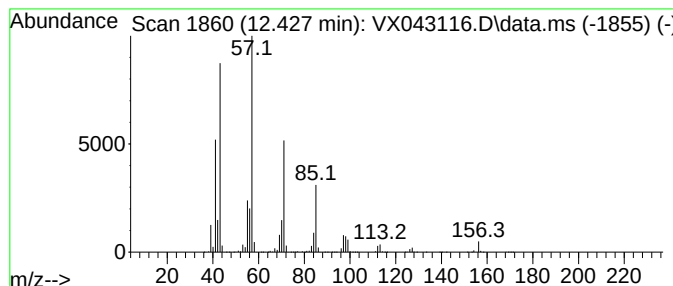
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Quant Title : SW846 8260

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

Peak Number 4 Undecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.427	483.64 ug/l	14394000	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	96
2	Tetradecane			198	C14H30	000629-59-4	86
3	Hexadecane			226	C16H34	000544-76-3	80
4	Tridecane			184	C13H28	000629-50-5	78
5	Carbonic acid, nonyl vinyl ester			214	C12H22O3	1000383-25-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
613-D

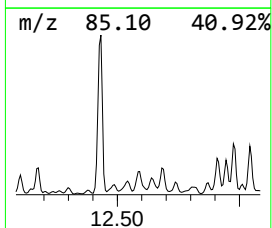
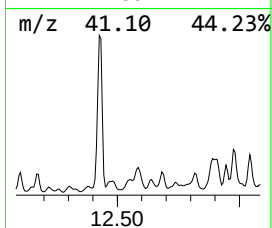
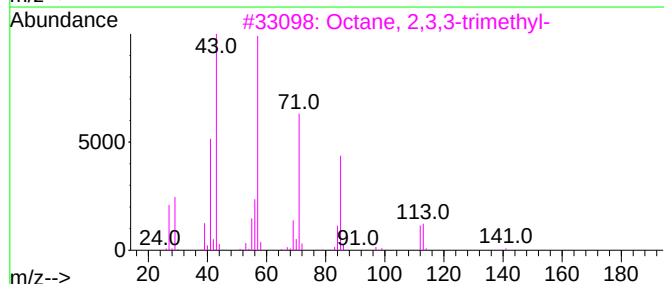
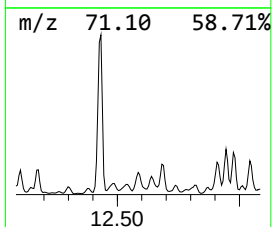
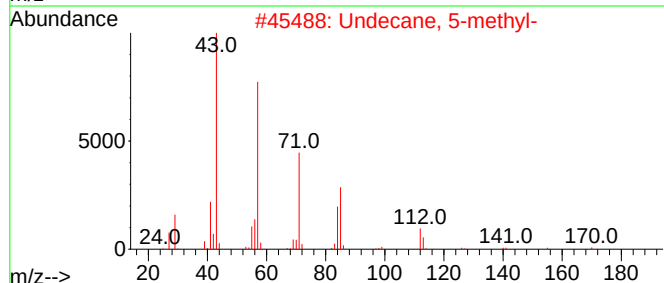
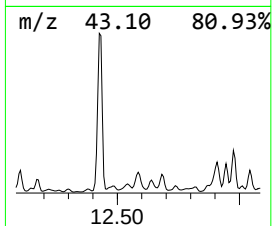
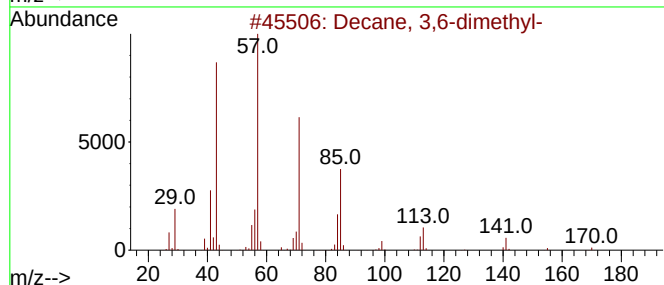
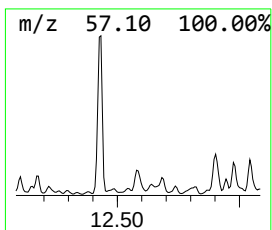
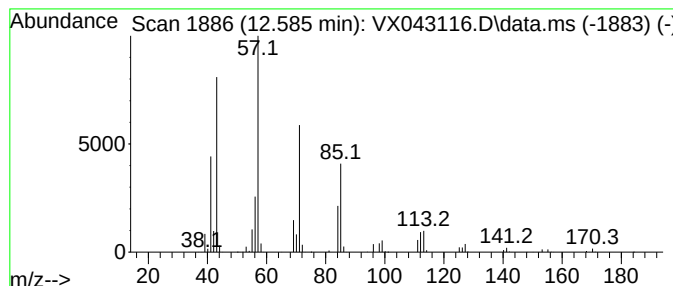
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Quant Title : SW846 8260

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

Peak Number 5 Decane, 3,6-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.585	98.25 ug/l	2924150	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	95
2			Undecane, 5-methyl-	170	C12H26	001632-70-8	87
3			Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	80
4			Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	72
5			2,6-Dimethyldecane	170	C12H26	013150-81-7	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
613-D

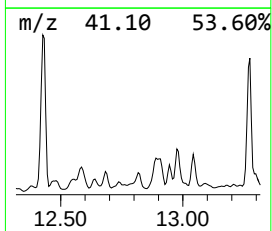
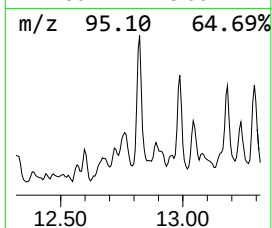
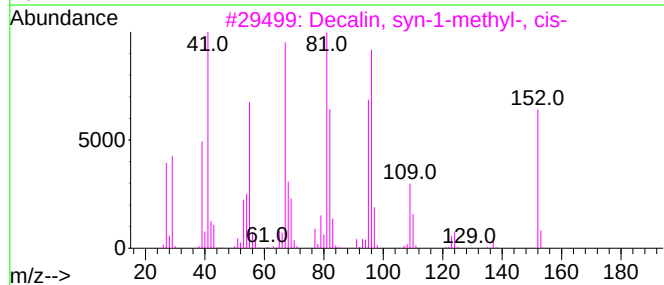
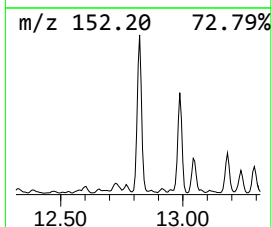
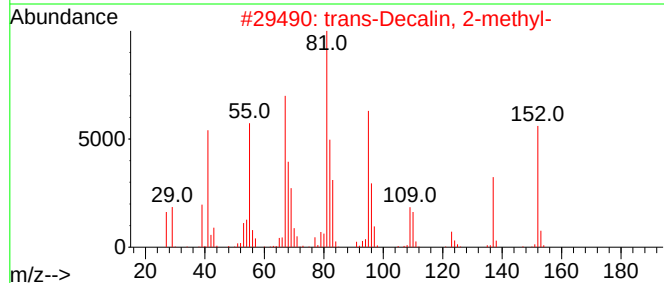
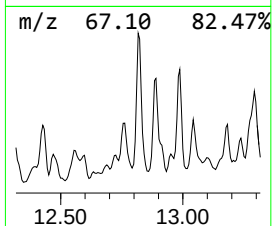
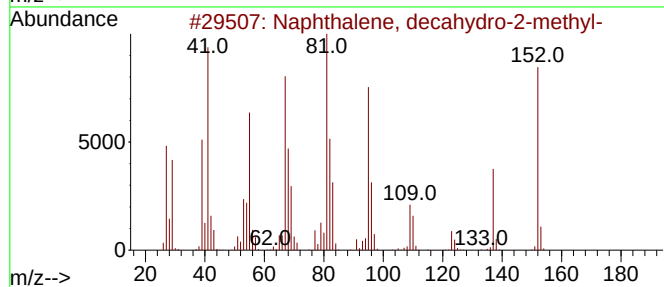
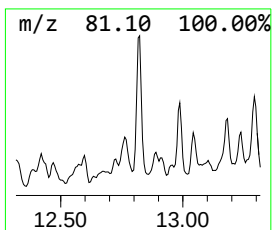
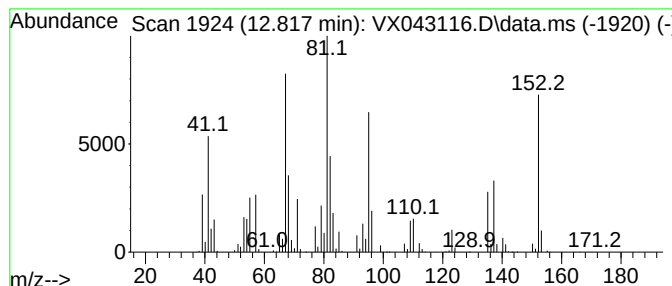
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, decahydro-2-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.817	86.31 ug/l	2568810	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	91
2			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	87
3			Decalin, syn-1-methyl-, cis-	152	C11H20	1000158-89-1	76
4			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	74
5			Pulegone	152	C10H16O	000089-82-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
613-D

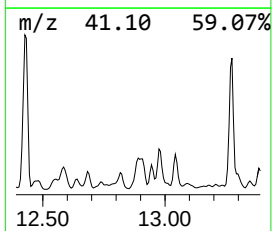
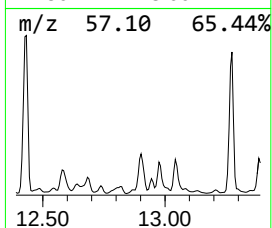
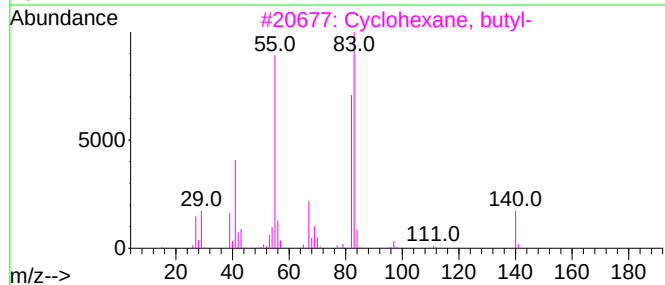
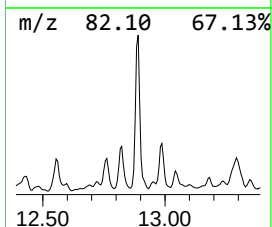
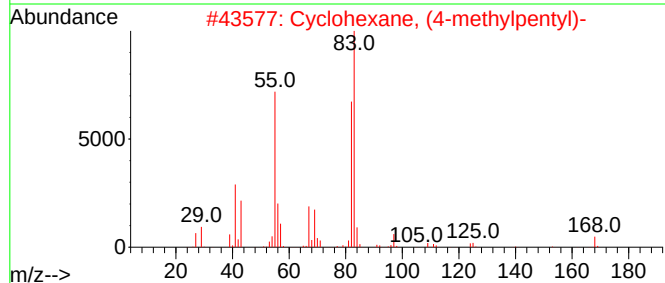
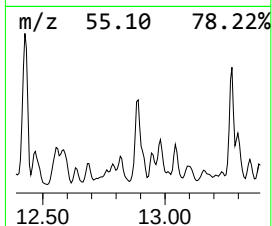
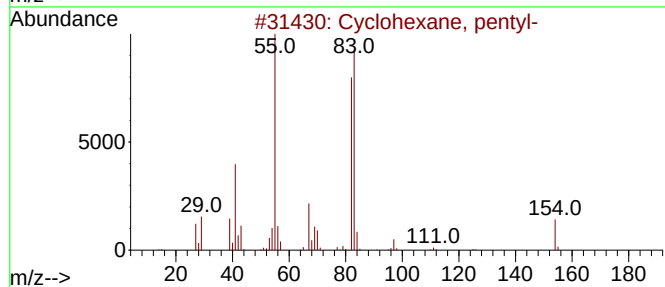
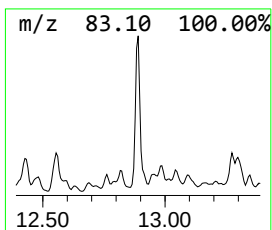
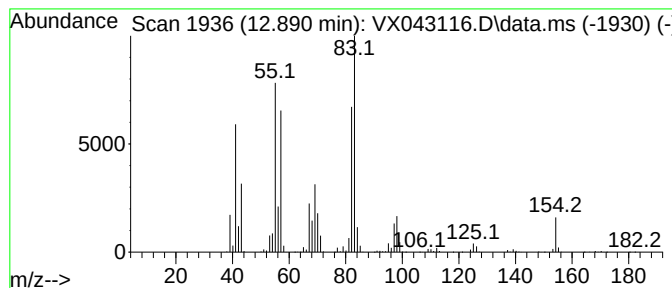
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Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, pentyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	141.92 ug/l	4223810	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	83
2			Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	72
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	59
4			Cyclohexane, methyl-	98	C7H14	000108-87-2	52
5			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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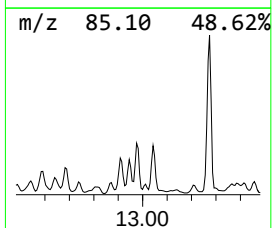
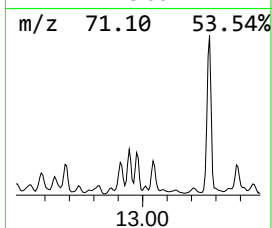
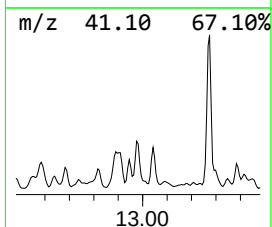
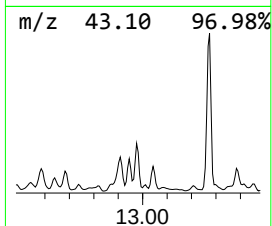
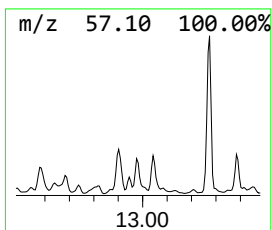
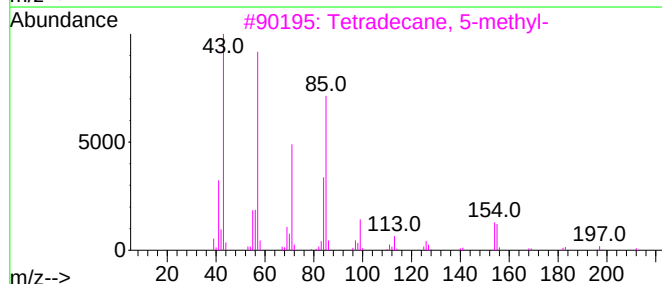
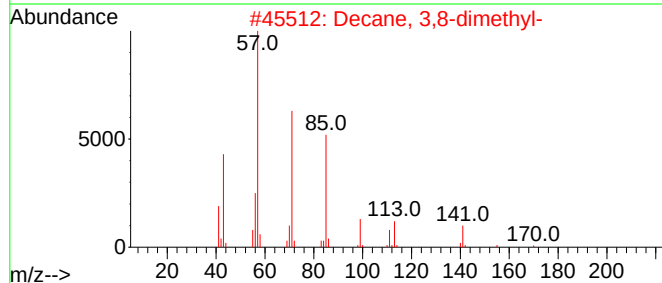
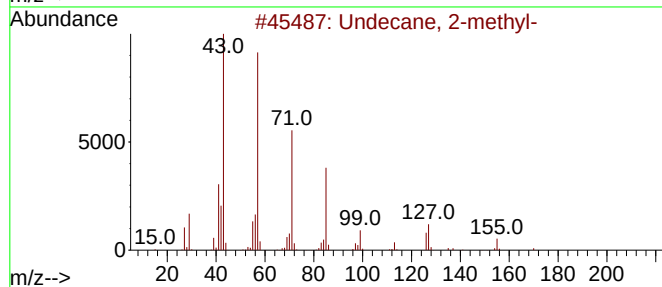
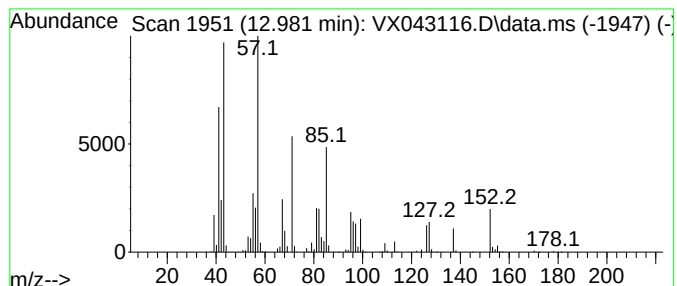
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Quant Title : SW846 8260

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

Peak Number 8 Undecane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.981	133.23 ug/l	3965250	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2-methyl-	170	C12H26	007045-71-8	64
2			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	43
3			Tetradecane, 5-methyl-	212	C15H32	025117-32-2	43
4			Undecane, 3-methyl-	170	C12H26	001002-43-3	43
5			Decane, 1-iodo-	268	C10H21I	002050-77-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
613-D

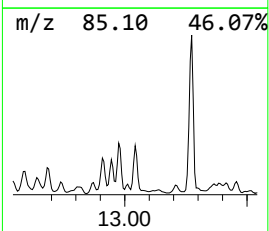
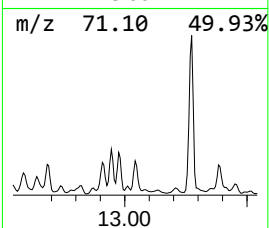
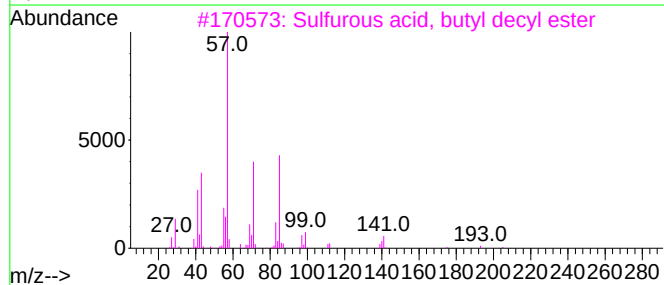
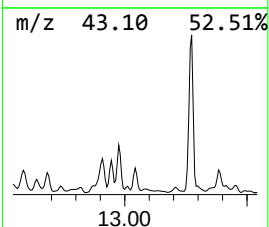
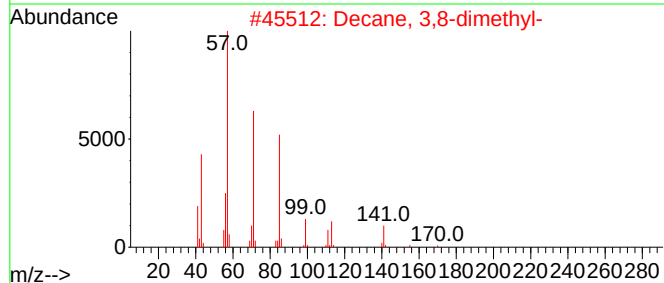
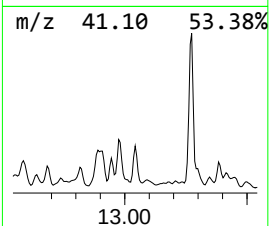
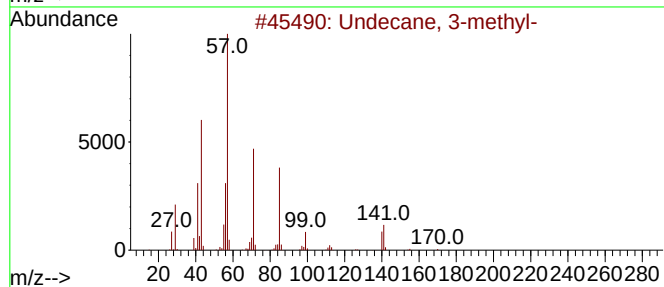
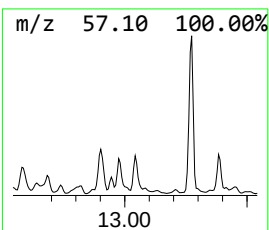
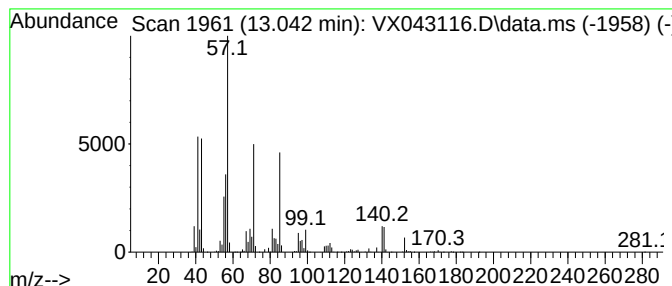
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Quant Title : SW846 8260

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

Peak Number 9 Undecane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.042	90.66 ug/l	2698260	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 3-methyl-	170	C12H26	001002-43-3	94
2			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	81
3			Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	72
4			Decane, 1-iodo-	268	C10H21I	002050-77-3	59
5			Nonane, 3-methyl-	142	C10H22	005911-04-6	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
613-D

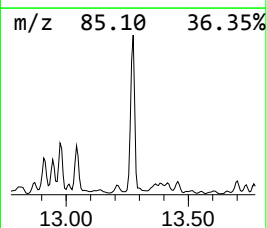
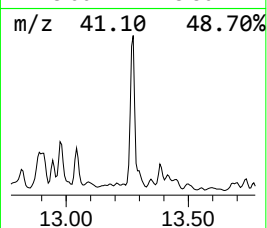
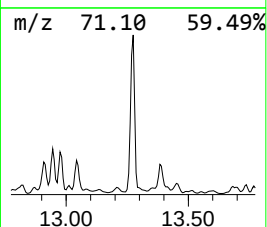
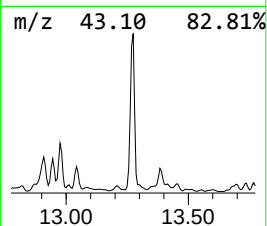
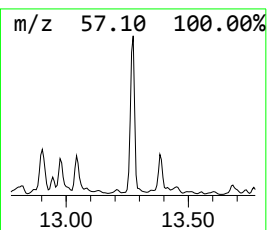
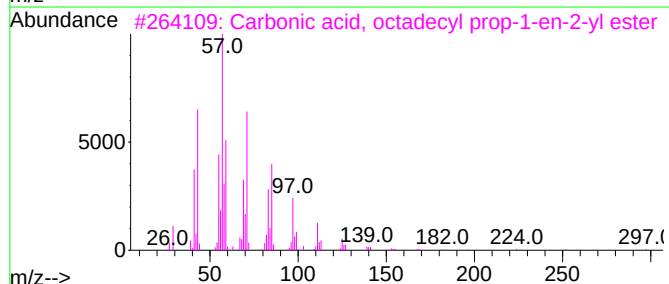
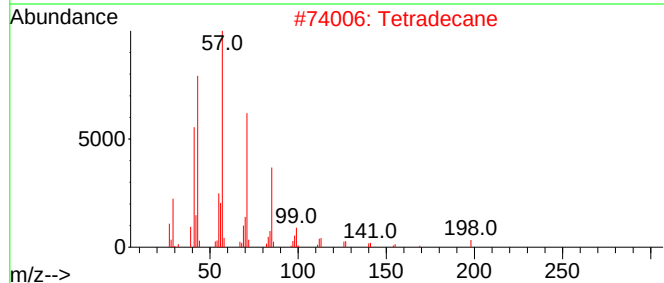
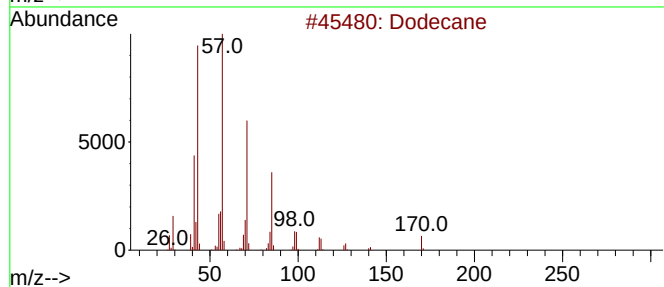
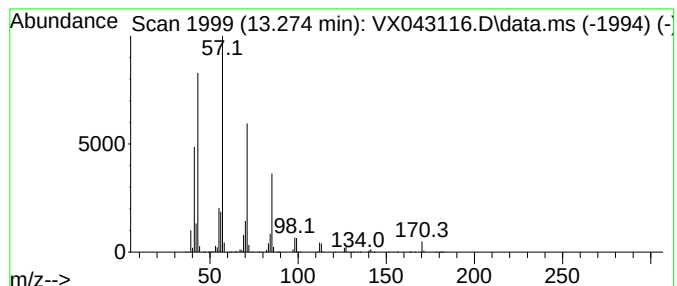
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Quant Title : SW846 8260

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

Peak Number 10 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.274	378.72 ug/l	11271400	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	96
2			Tetradecane	198	C14H30	000629-59-4	91
3			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90
4			Carbonic acid, tetradecyl vinyl ...	284	C17H32O3	1000382-54-5	90
5			Tridecane	184	C13H28	000629-50-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092424\
Data File : VX043116.D
Acq On : 24 Sep 2024 12:21
Operator : JC/MD
Sample : P4123-02
Misc : 1.15g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
613-D

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092324W.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
Hexane, 2-methyl-	5.483	149.2	ug/l	1906540	1	5.544	638862	50.0
Hexane, 3-methyl-	5.806	250.1	ug/l	3195230	1	5.544	638862	50.0
Heptane	6.598	843.3	ug/l	6317510	2	6.757	374574	50.0
Undecane	12.427	483.6	ug/l	14394000	4	12.024	1488100	50.0
Decane, 3,6-dim...	12.585	98.3	ug/l	2924150	4	12.024	1488100	50.0
Naphthalene, de...	12.817	86.3	ug/l	2568810	4	12.024	1488100	50.0
Cyclohexane, pe...	12.890	141.9	ug/l	4223810	4	12.024	1488100	50.0
Undecane, 2-met...	12.981	133.2	ug/l	3965250	4	12.024	1488100	50.0
Undecane, 3-met...	13.042	90.7	ug/l	2698260	4	12.024	1488100	50.0
Dodecane	13.274	378.7	ug/l	11271400	4	12.024	1488100	50.0