

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
 Data File : VY006031.D
 Acq On : 15 Sep 2021 18:00
 Operator : SY/MD
 Sample : M3733-09
 Misc : 7.25G/5ML/MSVOA_Y/SOIL
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 NAPL-01-(8-10)-END-POINT

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_Y\methods\82Y091421S.M

Title : SW846 8260

Signal : TIC: VY006031.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.996	835	849	863	rBV	126762	353596	22.27%	3.188%
2	7.795	974	980	990	rVB2	9235	20703	1.30%	0.187%
3	8.149	1031	1038	1045	rVB2	8267	17807	1.12%	0.161%
4	8.691	1121	1127	1135	rVB	16385	33433	2.11%	0.301%
5	8.941	1160	1168	1178	rBV	111385	217552	13.70%	1.962%
6	10.185	1363	1372	1380	rBV2	29289	51719	3.26%	0.466%
7	10.721	1452	1460	1470	rBV	670115	1100664	69.33%	9.925%
8	11.490	1579	1586	1594	rBV	29901	48996	3.09%	0.442%
9	12.008	1663	1671	1677	rBV5	16965	42586	2.68%	0.384%
10	12.203	1694	1703	1706	rBV4	17826	38882	2.45%	0.351%
11	12.246	1706	1710	1714	rVB4	14933	27030	1.70%	0.244%
12	12.489	1745	1750	1754	rBV3	73638	127250	8.01%	1.147%
13	12.648	1768	1776	1783	rVB6	31169	78548	4.95%	0.708%
14	12.739	1783	1791	1795	rBV6	13945	40207	2.53%	0.363%
15	12.806	1795	1802	1807	rBV	424980	626812	39.48%	5.652%
16	12.861	1807	1811	1817	rVB4	31859	55340	3.49%	0.499%
17	12.953	1821	1826	1829	rBV2	86691	134815	8.49%	1.216%
18	13.001	1829	1834	1837	rBV3	104691	160399	10.10%	1.446%
19	13.038	1837	1840	1846	rVB	144690	191095	12.04%	1.723%
20	13.099	1846	1850	1857	rVB3	99647	198193	12.48%	1.787%
21	13.166	1857	1861	1866	rBV	142253	216208	13.62%	1.950%
22	13.221	1866	1870	1873	rBV2	82286	130805	8.24%	1.179%
23	13.318	1878	1886	1889	rBV4	131357	243842	15.36%	2.199%
24	13.361	1889	1893	1896	rVV	276584	396604	24.98%	3.576%
25	13.398	1896	1899	1901	rVV	229046	294127	18.53%	2.652%
26	13.428	1901	1904	1909	rVB	325597	424918	26.76%	3.831%
27	13.471	1909	1911	1912	rBV	23320	21253	1.34%	0.192%
28	13.501	1912	1916	1923	rVB	298445	422928	26.64%	3.814%
29	13.568	1923	1927	1928	rBV3	25984	35767	2.25%	0.323%
30	13.678	1941	1945	1946	rBV2	36551	48697	3.07%	0.439%
31	13.751	1951	1957	1962	rVV	1151868	1587683	100.00%	14.316%
32	13.794	1962	1964	1971	rVV6	72661	191880	12.09%	1.730%
33	13.867	1972	1976	1979	rVV3	77364	176028	11.09%	1.587%
34	13.916	1979	1984	1989	rVV3	183451	418322	26.35%	3.772%
35	13.971	1989	1993	1997	rVV2	141611	244397	15.39%	2.204%
36	14.013	1997	2000	2006	rVB2	93959	138191	8.70%	1.246%

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37	14.074	2006	2010	2015	rVB3	164998	243816	15.36%	2.198%
38	14.142	2015	2021	2027	rVB6	99693	228442	14.39%	2.060%
39	14.196	2027	2030	2032	rBV3	48207	68526	4.32%	0.618%
40	14.239	2032	2037	2039	rVV4	120815	206827	13.03%	1.865%
41	14.276	2040	2043	2047	rVV3	180398	296938	18.70%	2.677%
42	14.312	2047	2049	2056	rVB3	148751	216465	13.63%	1.952%
43	14.379	2056	2060	2064	rBV2	175933	259207	16.33%	2.337%
44	14.459	2070	2073	2078	rVB3	31707	56217	3.54%	0.507%
45	14.611	2093	2098	2103	rVB	339311	452313	28.49%	4.078%
46	14.672	2103	2108	2115	rVV4	59488	148000	9.32%	1.335%
47	14.733	2115	2118	2123	rVV5	55813	101120	6.37%	0.912%
48	14.788	2124	2127	2134	rVB3	45412	86680	5.46%	0.782%
49	14.873	2137	2141	2149	rVB5	24769	48689	3.07%	0.439%
50	14.946	2149	2153	2160	rBV4	30578	58492	3.68%	0.527%
51	15.050	2166	2170	2176	rVB6	16989	35205	2.22%	0.317%
52	15.141	2182	2185	2191	rVB6	15185	26070	1.64%	0.235%

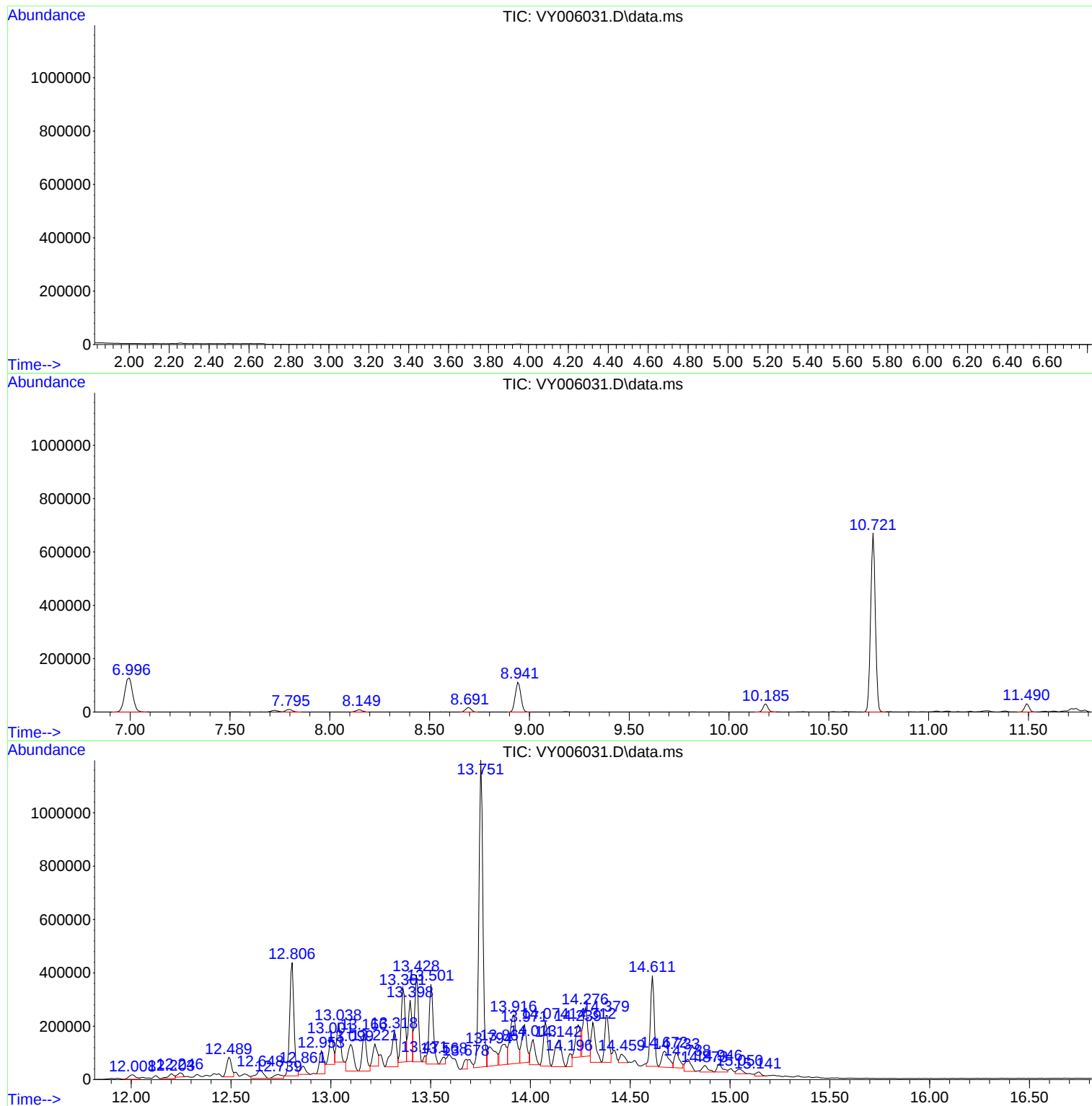
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091421S.M
Quant Title : SW846 8260

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



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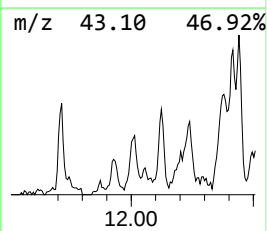
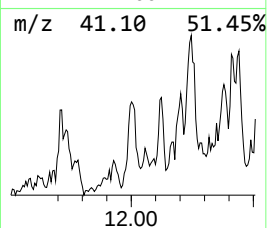
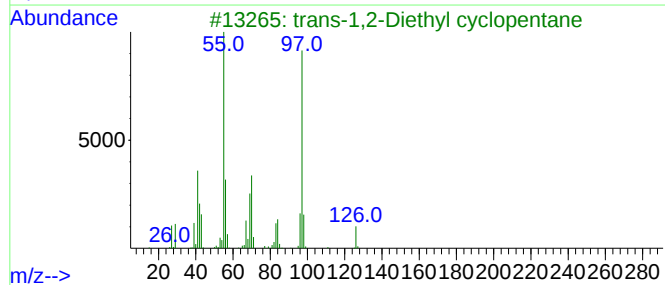
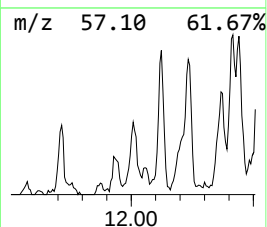
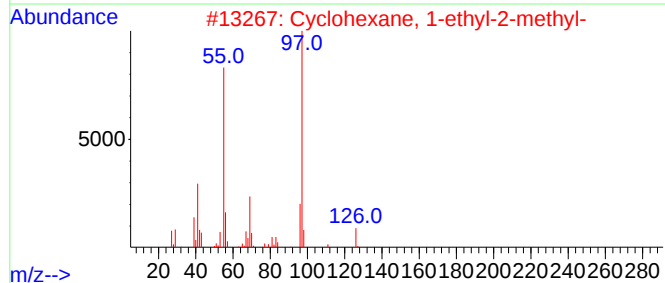
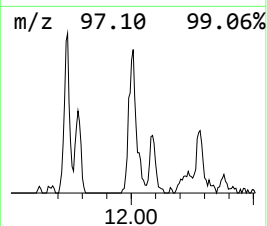
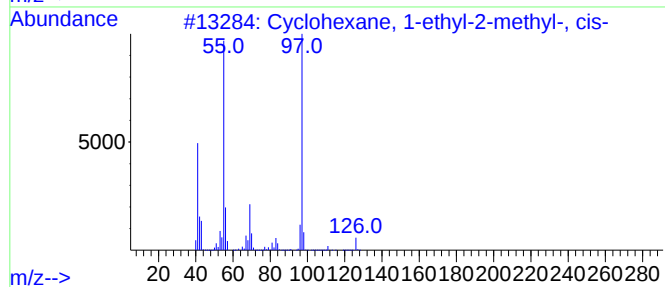
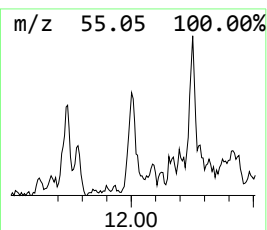
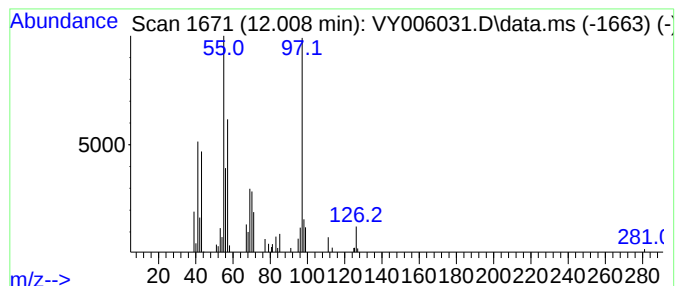
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091421S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.008	43.46 ug/l	42586	Chlorobenzene-d5	11.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	64
2			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	53
3			trans-1,2-Diethyl cyclopentane	126	C9H18	000932-40-1	49
4			1H-Pyrazol-4-amine, 3-methyl-	97	C4H7N3	1010338-28-2	49
5			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	46



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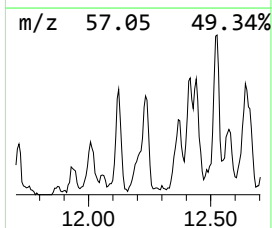
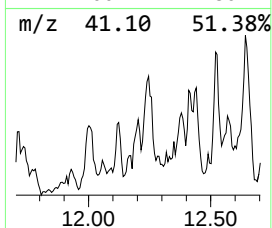
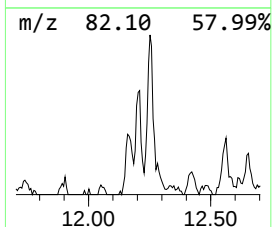
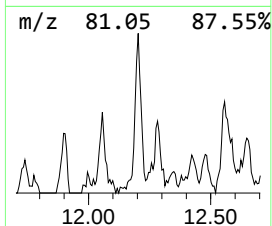
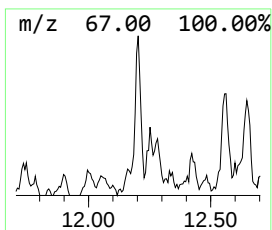
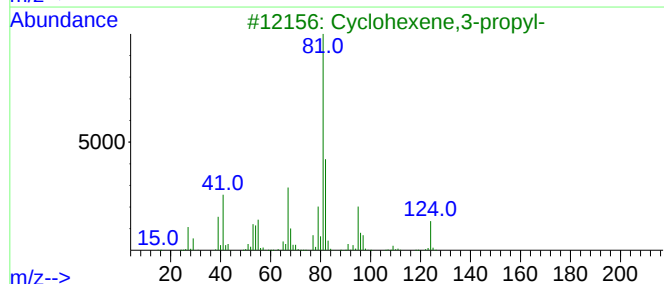
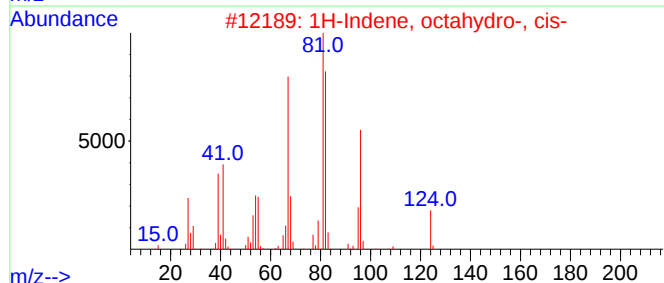
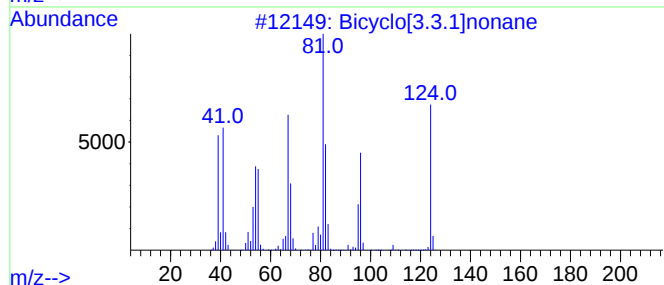
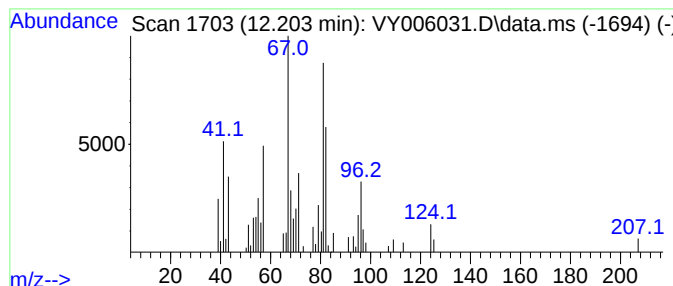
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091421S.M
Quant Title : SW846 8260

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

Peak Number 2 Bicyclo[3.3.1]nonane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.203	39.68 ug/l	38882	Chlorobenzene-d5	11.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.3.1]nonane	124	C9H16	000280-65-9	68
2			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	58
3			Cyclohexene,3-propyl-	124	C9H16	003983-06-0	49
4			Cyclohexene, 4-propyl-	124	C9H16	013487-64-4	49
5			Pentalene, octahydro-1-methyl-	124	C9H16	032273-77-1	46



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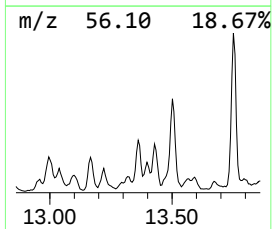
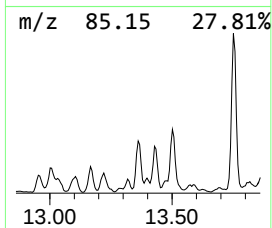
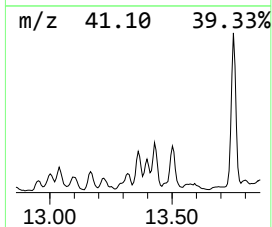
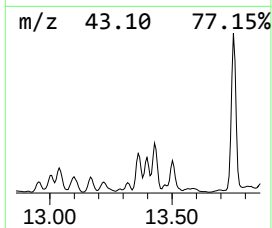
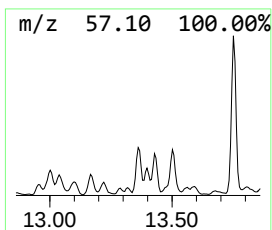
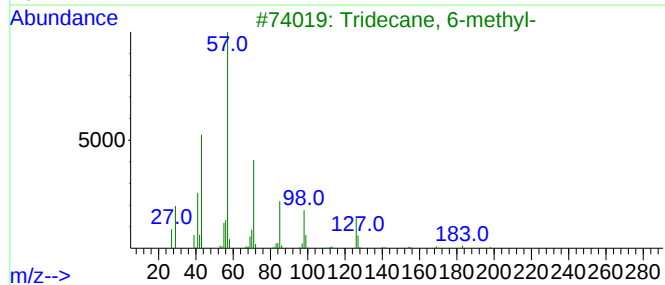
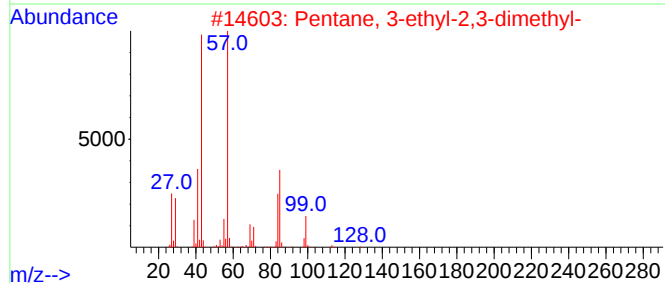
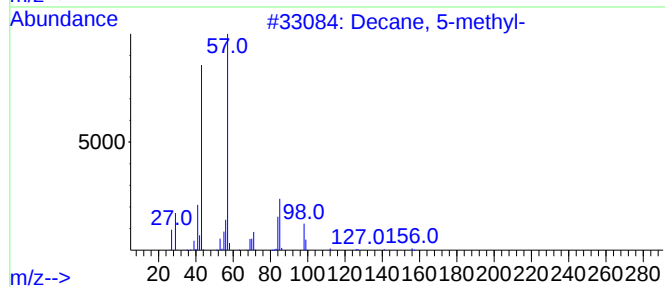
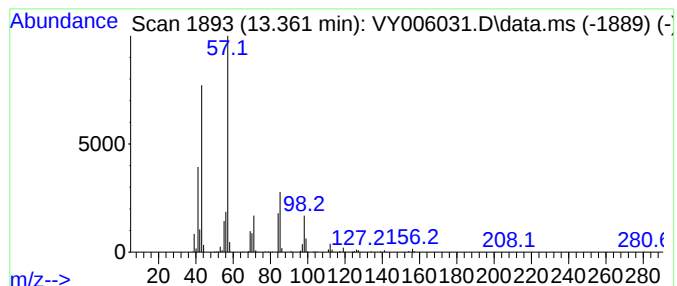
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091421S.M
Quant Title : SW846 8260

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

Peak Number 3 Decane, 5-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.361	46.67 ug/l	396604	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5-methyl-	156	C11H24	013151-35-4	87
2			Pentane, 3-ethyl-2,3-dimethyl-	128	C9H20	016747-33-4	59
3			Tridecane, 6-methyl-	198	C14H30	013287-21-3	59
4			4,4-Dimethyl octane	142	C10H22	015869-95-1	50
5			Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	50



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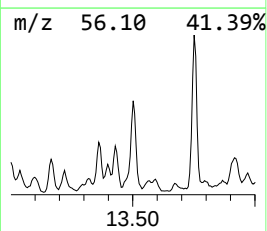
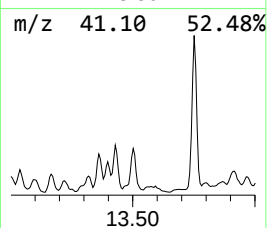
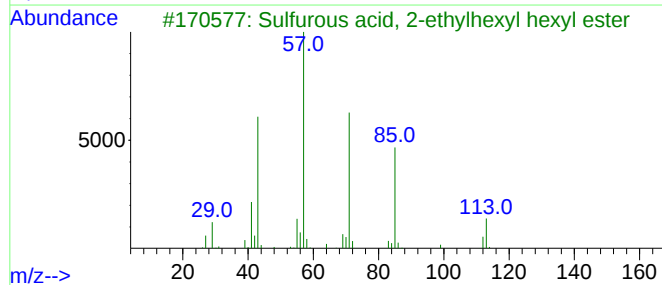
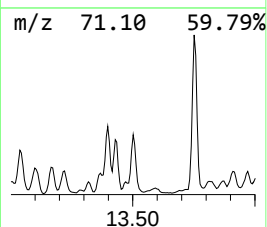
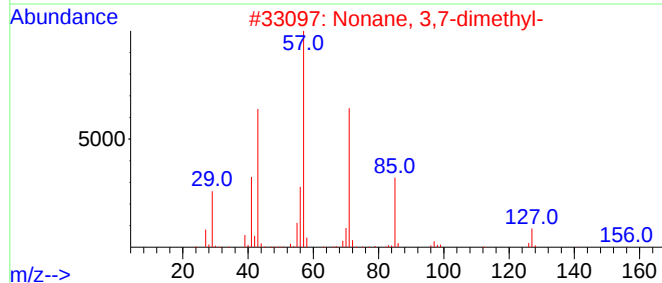
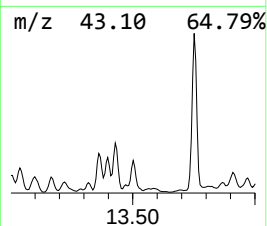
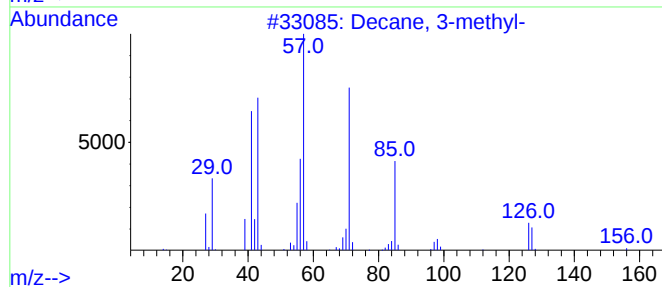
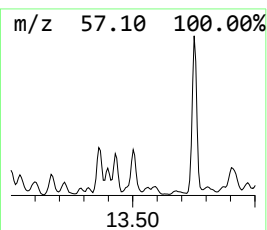
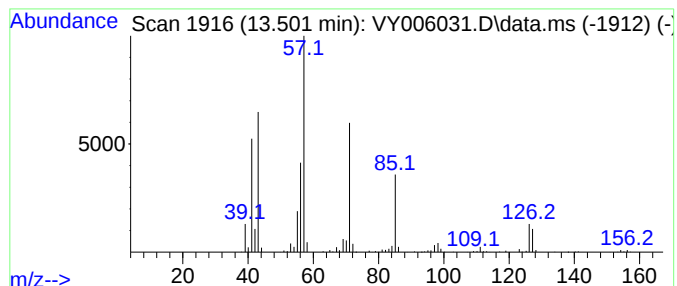
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Peak Number 4 Decane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.501	49.77 ug/l	422928	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	74
2			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	53
3			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	47
4			Nonane, 5-butyl-	184	C13H28	017312-63-9	43
5			Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	38



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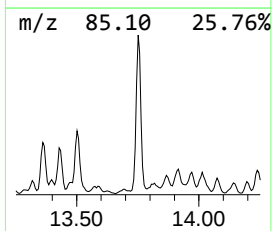
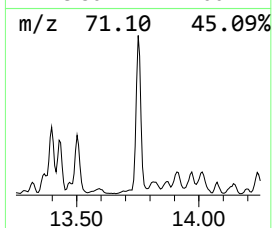
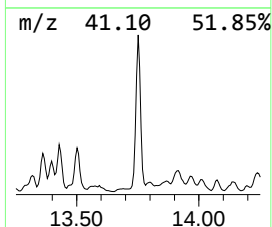
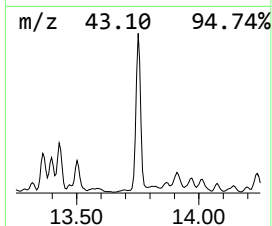
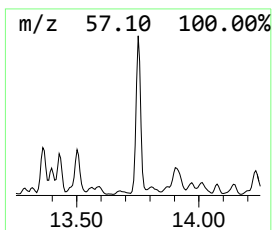
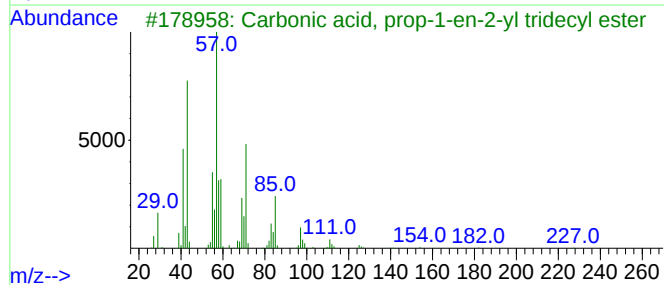
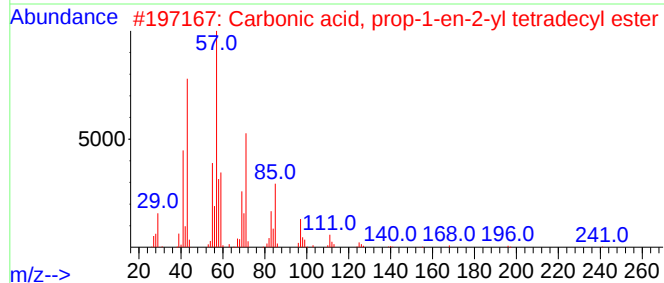
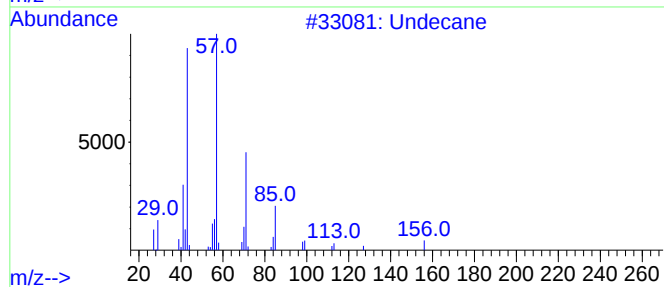
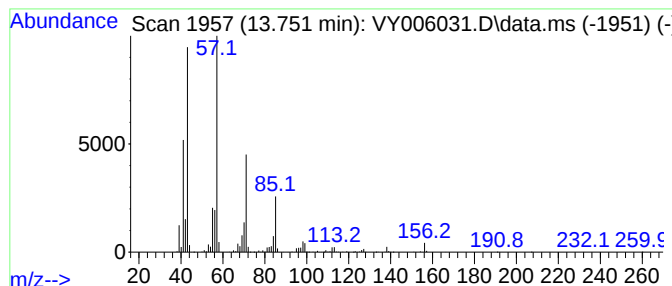
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Quant Title : SW846 8260

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

Peak Number 5 Undecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	186.82 ug/l	1587680	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	94
2			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	90
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	78
4			Nonane	128	C9H20	000111-84-2	72
5			Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
Data File : VY006031.D
Acq On : 15 Sep 2021 18:00
Operator : SY/MD
Sample : M3733-09
Misc : 7.25G/5ML/MSVOA_Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NAPL-01-(8-10)-END-POINT

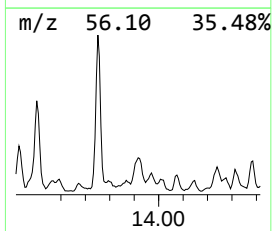
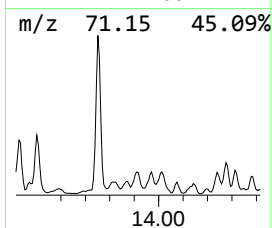
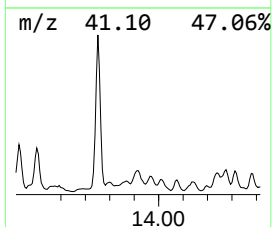
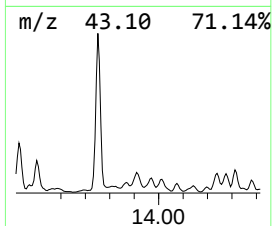
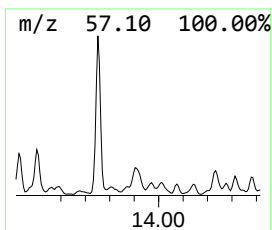
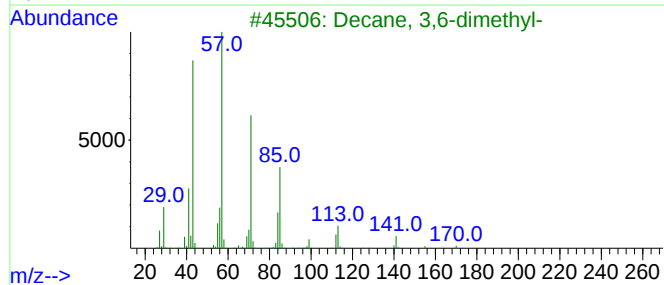
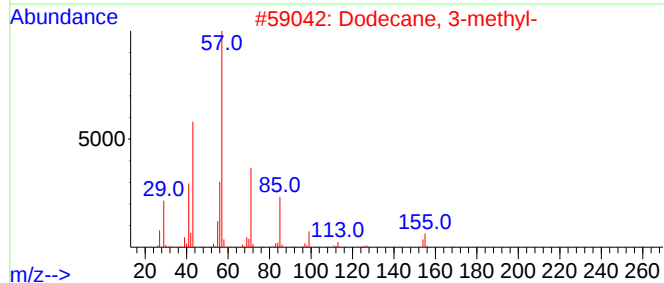
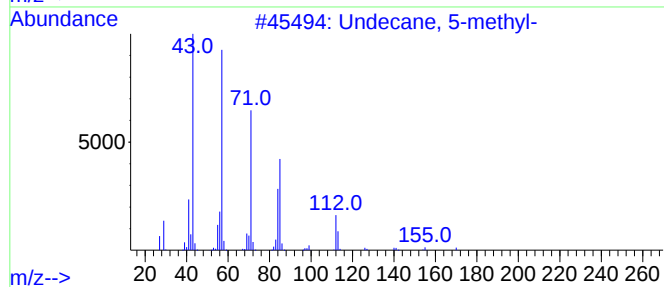
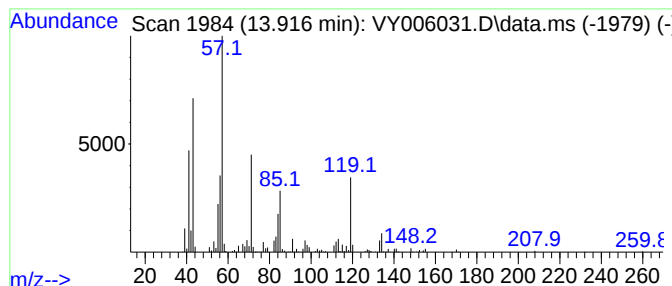
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Quant Title : SW846 8260

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

Peak Number 6 Undecane, 5-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.916	49.22 ug/l	418322	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 5-methyl-	170	C12H26	001632-70-8	53
2			Dodecane, 3-methyl-	184	C13H28	017312-57-1	53
3			Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	53
4			2-Methyl-1-pentanol, pentafluoro...	248	C9H13F5O2	1000352-35-9	50
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
Data File : VY006031.D
Acq On : 15 Sep 2021 18:00
Operator : SY/MD
Sample : M3733-09
Misc : 7.25G/5ML/MSVOA_Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NAPL-01-(8-10)-END-POINT

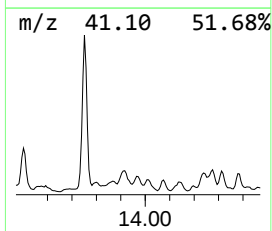
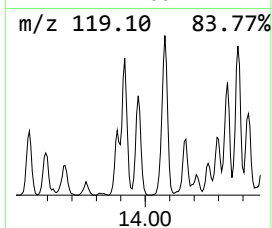
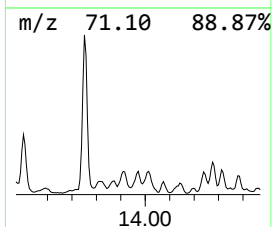
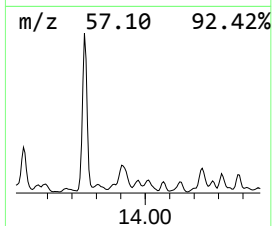
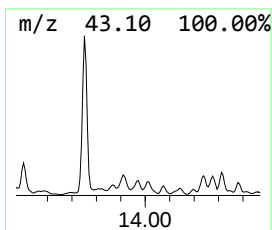
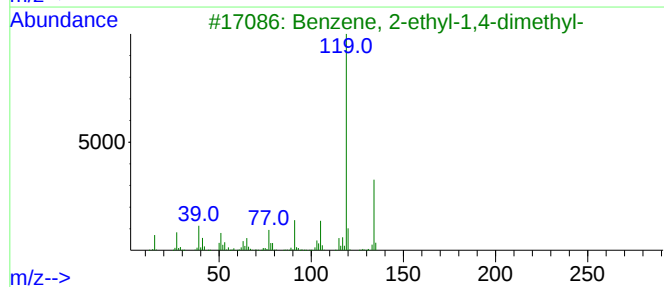
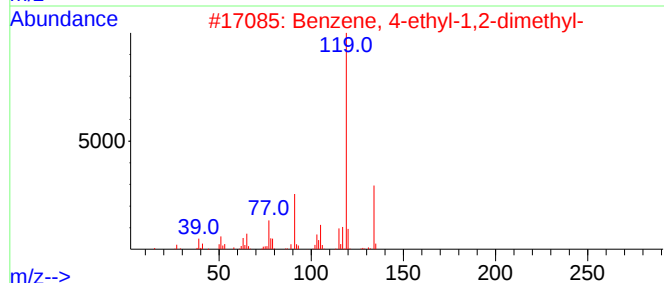
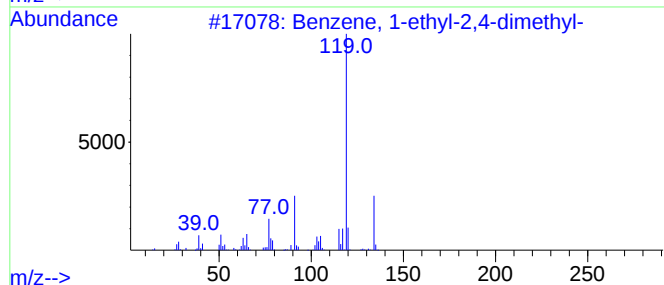
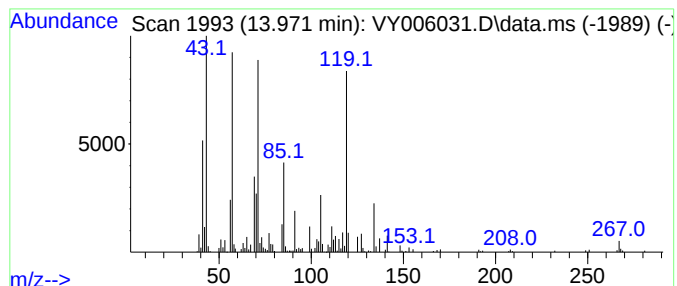
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	28.76 ug/l	244397	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	70
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	70
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	42
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	38
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
Data File : VY006031.D
Acq On : 15 Sep 2021 18:00
Operator : SY/MD
Sample : M3733-09
Misc : 7.25G/5ML/MSVOA_Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NAPL-01-(8-10)-END-POINT

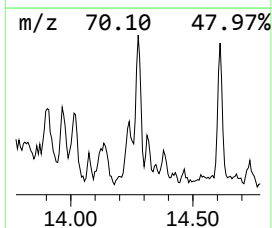
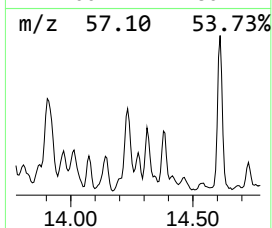
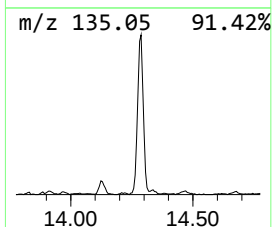
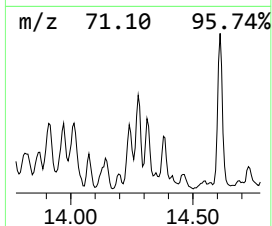
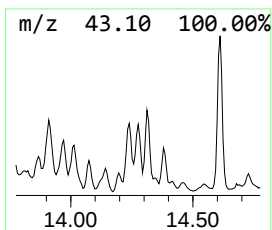
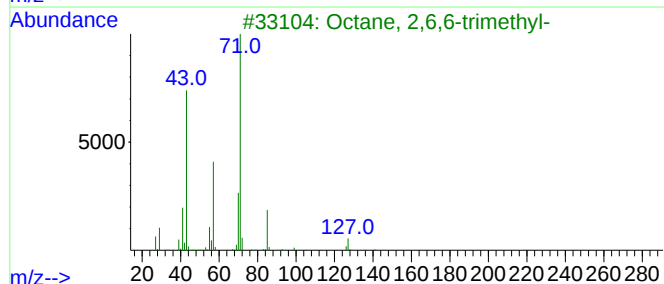
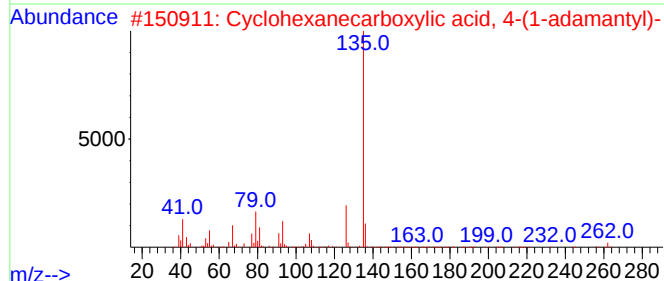
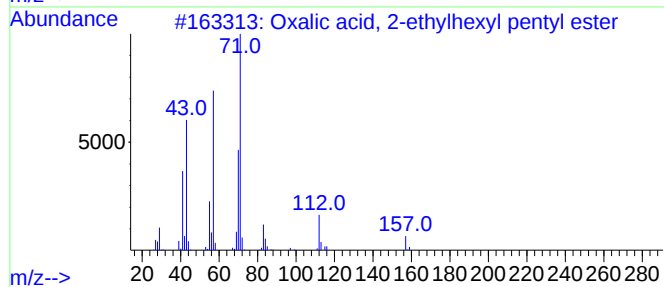
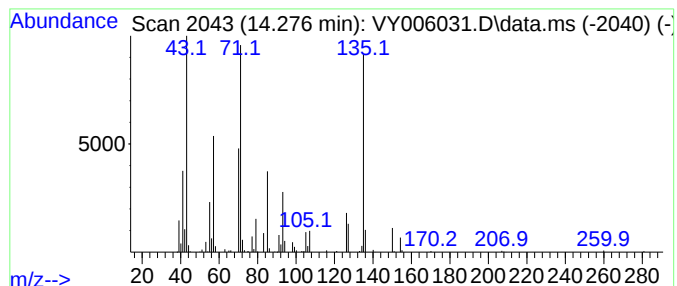
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Quant Title : SW846 8260

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

Peak Number 8 unknown14.276 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.276	34.94 ug/l	296938	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	27
2			Cyclohexanecarboxylic acid, 4-(1...	262	C17H26O2	175881-24-0	27
3			Octane, 2,6,6-trimethyl-	156	C11H24	054166-32-4	27
4			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	27
5			Ethanedioic acid, bis(3-methylbu...	230	C12H22O4	002051-00-5	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
Data File : VY006031.D
Acq On : 15 Sep 2021 18:00
Operator : SY/MD
Sample : M3733-09
Misc : 7.25G/5ML/MSVOA_Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NAPL-01-(8-10)-END-POINT

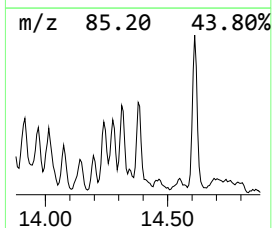
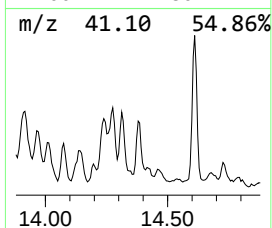
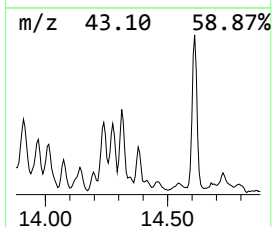
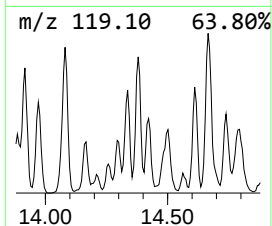
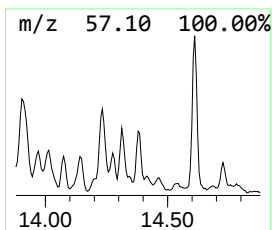
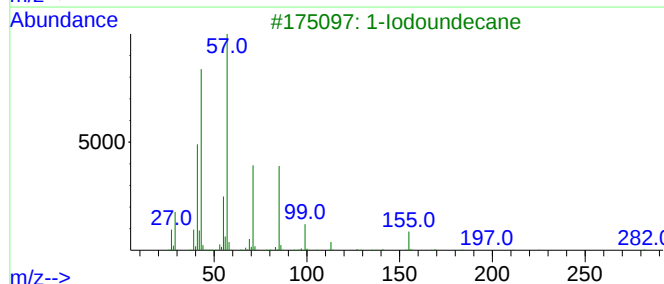
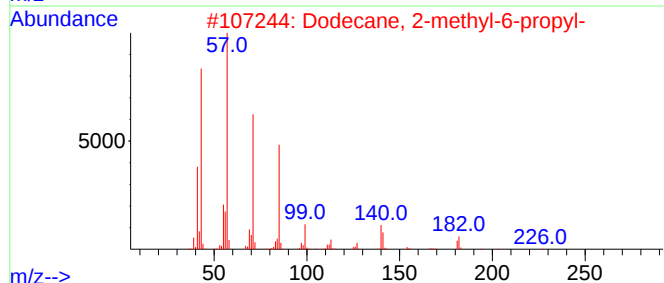
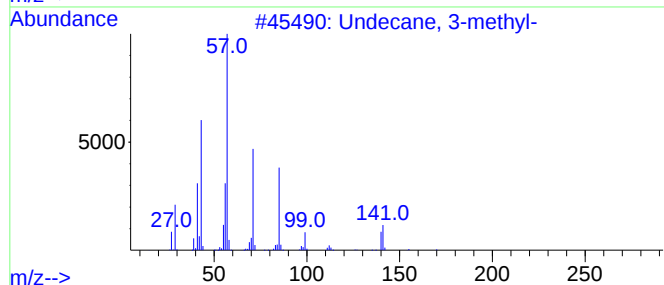
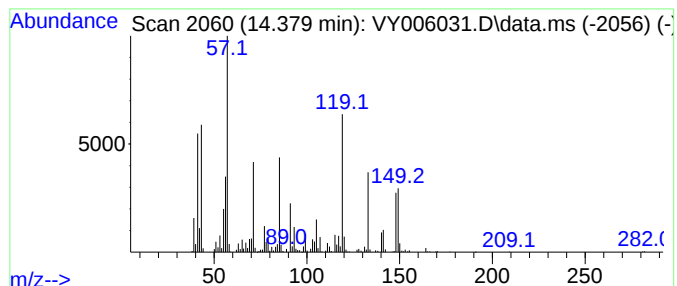
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Quant Title : SW846 8260

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.
14.379	30.50 ug/l	259207	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 3-methyl-	170	C12H26	001002-43-3	55
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	46
3			1-Iodoundecane	282	C11H23I	004282-44-4	30
4			Sulfurous acid, decyl 2-propyl e...	264	C13H28O3S	1000309-12-1	27
5			Octane, 4-methyl-	128	C9H20	002216-34-4	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
Data File : VY006031.D
Acq On : 15 Sep 2021 18:00
Operator : SY/MD
Sample : M3733-09
Misc : 7.25G/5ML/MSVOA_Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NAPL-01-(8-10)-END-POINT

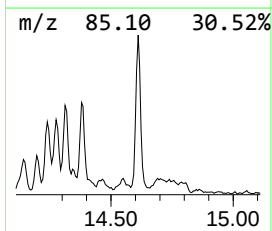
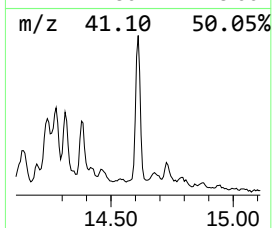
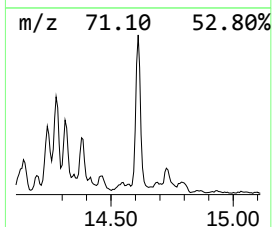
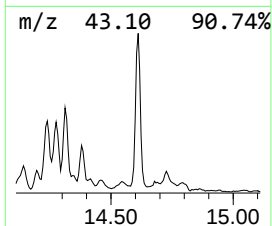
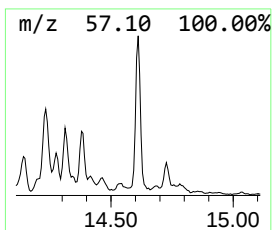
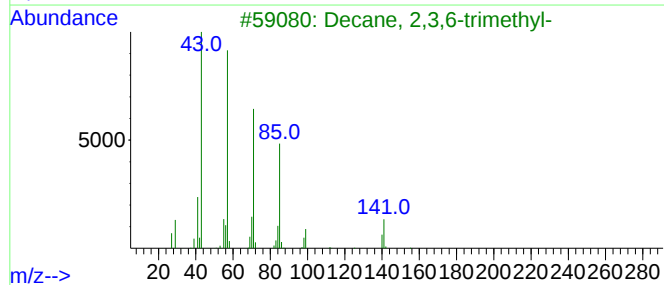
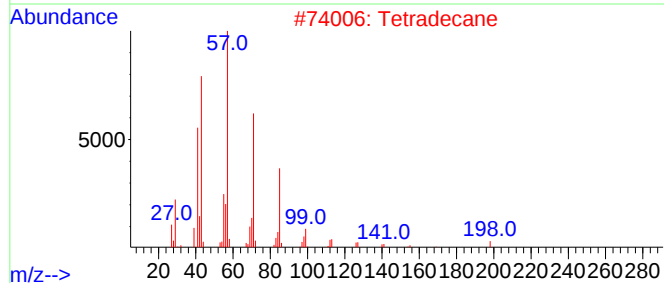
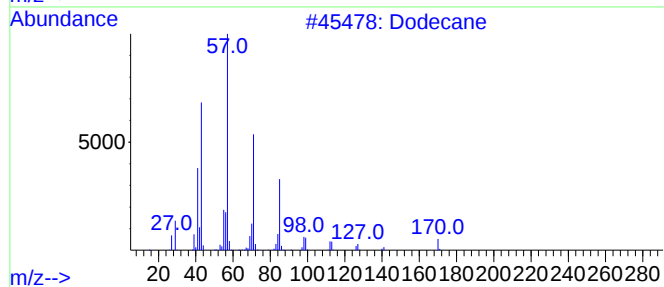
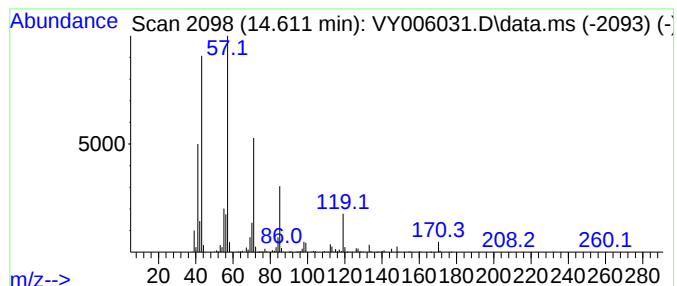
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Quant Title : SW846 8260

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

Peak Number 10 Dodecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.611	53.22 ug/l	452313	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	93
2			Tetradecane	198	C14H30	000629-59-4	72
3			Decane, 2,3,6-trimethyl-	184	C13H28	062238-12-4	72
4			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	72
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY091521\
Data File : VY006031.D
Acq On : 15 Sep 2021 18:00
Operator : SY/MD
Sample : M3733-09
Misc : 7.25G/5ML/MSVOA_Y/SOIL
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
NAPL-01-(8-10)-END-POINT

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y091421S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Bicyclo[3.3.1]n...	12.203	39.7	ug/l	38882	3	11.490	48996	50.0
Decane, 5-methyl-	13.361	46.7	ug/l	396604	4	13.428	424918	50.0
Decane, 3-methyl-	13.501	49.8	ug/l	422928	4	13.428	424918	50.0
Undecane	13.751	186.8	ug/l	1587680	4	13.428	424918	50.0
Undecane, 5-met...	13.916	49.2	ug/l	418322	4	13.428	424918	50.0
Benzene, 1-ethy...	13.971	28.8	ug/l	244397	4	13.428	424918	50.0
unknown14.276	14.276	34.9	ug/l	296938	4	13.428	424918	50.0
Undecane, 3-met...	14.379	30.5	ug/l	259207	4	13.428	424918	50.0
Dodecane	14.611	53.2	ug/l	452313	4	13.428	424918	50.0