

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
 Data File : VY009408.D
 Acq On : 30 Jun 2022 21:04
 Operator : KP/MD
 Sample : N3550-03
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 HR-03-06292022

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\82Y062822S.M

Title : SW846 8260

Signal : TIC: VY009408.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.704	465	473	485	rBV3	6986	22422	2.89%	0.252%
2	5.728	633	641	652	rBV4	4175	13164	1.69%	0.148%
3	6.972	836	845	864	rBV	27918	87001	11.20%	0.979%
4	7.710	957	966	972	rBV2	62089	147912	19.04%	1.664%
5	7.783	972	978	991	rVB	167027	370073	47.63%	4.164%
6	8.136	1024	1036	1046	rVB2	79099	201244	25.90%	2.264%
7	8.685	1118	1126	1140	rVB	261060	510688	65.73%	5.746%
8	10.173	1362	1370	1386	rBV	312524	543631	69.97%	6.116%
9	10.362	1394	1401	1406	rVB3	6273	11166	1.44%	0.126%
10	10.514	1421	1426	1431	rVB4	5668	9508	1.22%	0.107%
11	10.795	1463	1472	1481	rBV4	4975	10139	1.30%	0.114%
12	10.929	1481	1494	1504	rBV8	19809	61989	7.98%	0.697%
13	11.032	1507	1511	1515	rVV	12091	19570	2.52%	0.220%
14	11.081	1515	1519	1525	rVB3	12195	22761	2.93%	0.256%
15	11.197	1533	1538	1542	rBV3	7863	13823	1.78%	0.156%
16	11.282	1542	1552	1562	rVB5	30378	89180	11.48%	1.003%
17	11.374	1562	1567	1577	rVB2	19340	33399	4.30%	0.376%
18	11.483	1577	1585	1595	rBV	401518	657679	84.64%	7.399%
19	11.624	1601	1608	1611	rBV4	12794	19285	2.48%	0.217%
20	11.672	1611	1616	1617	rVV4	14242	20664	2.66%	0.232%
21	11.709	1617	1622	1630	rVV3	72714	162368	20.90%	1.827%
22	11.776	1630	1633	1638	rVB2	18599	28930	3.72%	0.325%
23	11.922	1654	1657	1662	rVB3	9041	14245	1.83%	0.160%
24	12.002	1662	1670	1682	rBV7	42363	131436	16.92%	1.479%
25	12.117	1685	1689	1693	rVB	29474	43687	5.62%	0.492%
26	12.197	1697	1702	1705	rVV4	22436	45137	5.81%	0.508%
27	12.239	1705	1709	1714	rVV5	36357	79932	10.29%	0.899%
28	12.325	1718	1723	1726	rVV4	14925	28951	3.73%	0.326%
29	12.380	1728	1732	1734	rVV3	18940	37379	4.81%	0.421%
30	12.410	1734	1737	1739	rVV	29301	41937	5.40%	0.472%
31	12.434	1739	1741	1743	rVV	27446	29854	3.84%	0.336%
32	12.477	1743	1748	1753	rVV	251877	390831	50.30%	4.397%
33	12.520	1753	1755	1759	rVV2	27069	32166	4.14%	0.362%
34	12.666	1771	1779	1783	rVB	37578	82141	10.57%	0.924%
35	12.733	1783	1790	1793	rBV	263980	441710	56.85%	4.970%
36	12.764	1793	1795	1798	rVV	137595	185777	23.91%	2.090%

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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\82Y062822S.M
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37	12.806	1798	1802	1808	rVV2	473065	751012	96.66%	8.450%
38	12.855	1808	1810	1815	rVB3	28138	38852	5.00%	0.437%
39	12.916	1815	1820	1821	rBV4	6070	9410	1.21%	0.106%
40	12.971	1821	1829	1836	rBV	236388	426018	54.83%	4.793%
41	13.038	1836	1840	1846	rVB	58919	87788	11.30%	0.988%
42	13.117	1846	1853	1858	rBV	396892	578212	74.42%	6.505%
43	13.166	1858	1861	1866	rVB3	16386	22276	2.87%	0.251%
44	13.221	1866	1870	1872	rBV2	15923	22732	2.93%	0.256%
45	13.312	1878	1885	1889	rBV3	57026	108847	14.01%	1.225%
46	13.361	1889	1893	1896	rVV3	39247	61194	7.88%	0.688%
47	13.422	1896	1903	1906	rVV	461358	776988	100.00%	8.742%
48	13.453	1906	1908	1913	rVV	262153	330332	42.51%	3.717%
49	13.495	1913	1915	1922	rVV	29732	42086	5.42%	0.474%
50	13.562	1922	1926	1927	rVV4	6211	9259	1.19%	0.104%
51	13.587	1927	1930	1931	rVV	23012	25626	3.30%	0.288%
52	13.623	1931	1936	1939	rVV2	103294	171090	22.02%	1.925%
53	13.660	1939	1942	1951	rVV2	53267	90346	11.63%	1.016%
54	13.751	1951	1957	1961	rVV	149343	221111	28.46%	2.488%
55	13.788	1961	1963	1964	rVV2	14197	14022	1.80%	0.158%
56	13.818	1966	1968	1972	rVV	21915	34670	4.46%	0.390%
57	13.885	1975	1979	1980	rVV2	18019	28588	3.68%	0.322%
58	13.916	1980	1984	1988	rVV5	28969	55241	7.11%	0.622%
59	13.965	1988	1992	1997	rVV2	35785	60080	7.73%	0.676%
60	14.013	1997	2000	2005	rVB6	7808	12340	1.59%	0.139%
61	14.074	2005	2010	2015	rBV5	16380	25232	3.25%	0.284%
62	14.141	2015	2021	2026	rBV10	6804	13861	1.78%	0.156%
63	14.209	2027	2032	2034	rBV6	8452	15805	2.03%	0.178%
64	14.251	2035	2039	2047	rVV9	16537	53187	6.85%	0.598%
65	14.330	2050	2052	2056	rVB	12749	16374	2.11%	0.184%
66	14.379	2056	2060	2063	rBV6	8653	11069	1.42%	0.125%
67	14.416	2063	2066	2070	rVB5	8025	11308	1.46%	0.127%
68	14.611	2094	2098	2102	rVB	18603	26127	3.36%	0.294%
69	14.678	2102	2109	2114	rBV5	16647	36406	4.69%	0.410%
70	14.727	2114	2117	2123	rVB5	7507	12010	1.55%	0.135%
71	14.830	2125	2134	2140	rBV8	8275	16064	2.07%	0.181%
72	14.946	2148	2153	2156	rBV7	4277	8116	1.04%	0.091%
73	15.233	2191	2200	2206	rBV5	5252	13418	1.73%	0.151%
74	15.312	2206	2213	2216	rBV7	3409	9293	1.20%	0.105%

Sum of corrected areas: 8888169

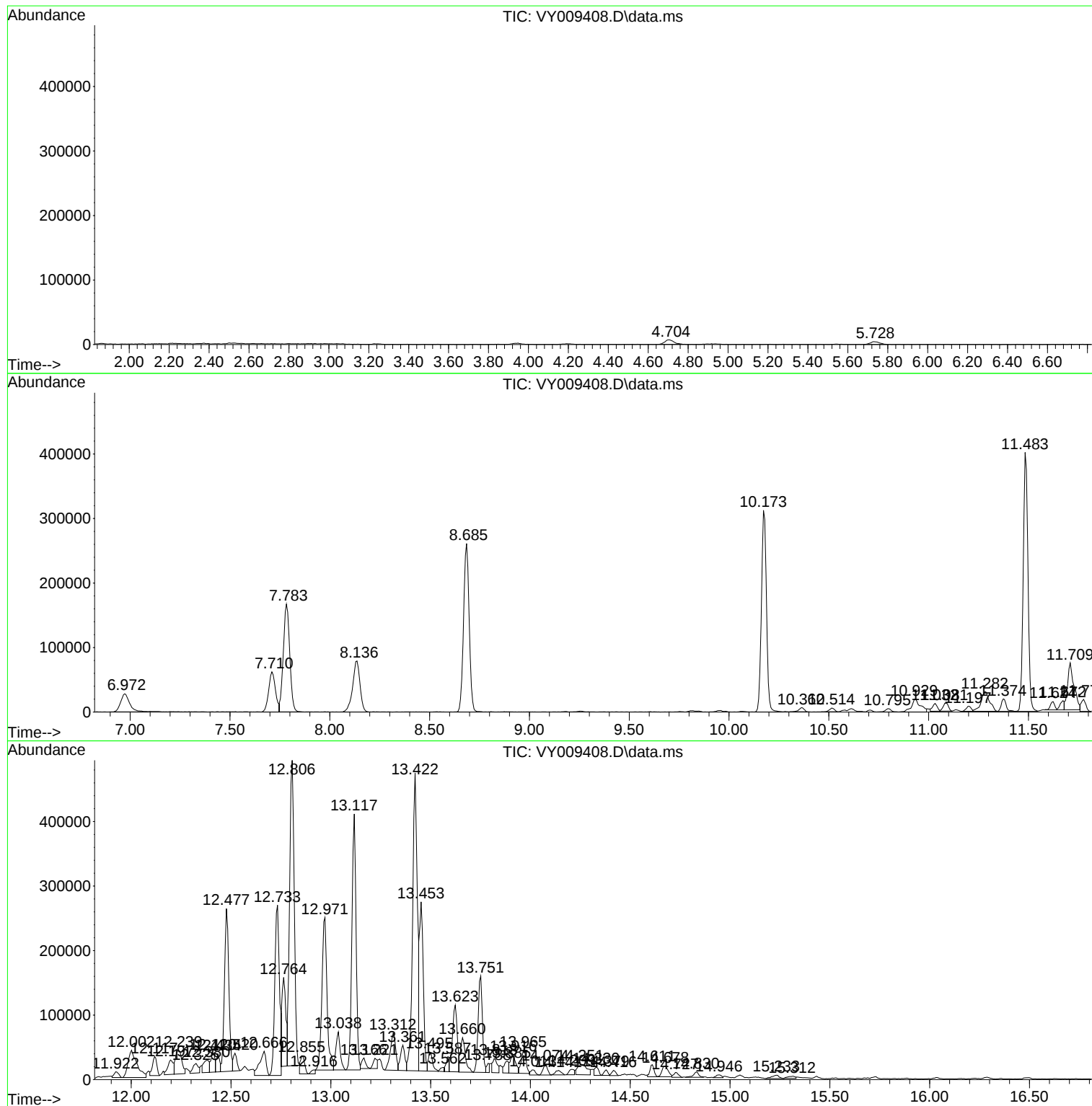
Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062822S.M
Quant Title : SW846 8260

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



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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

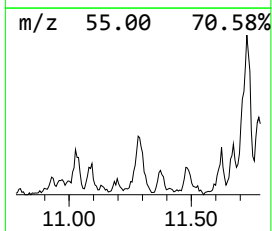
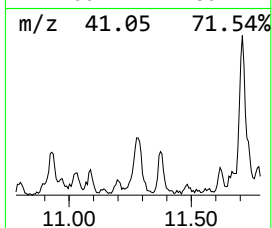
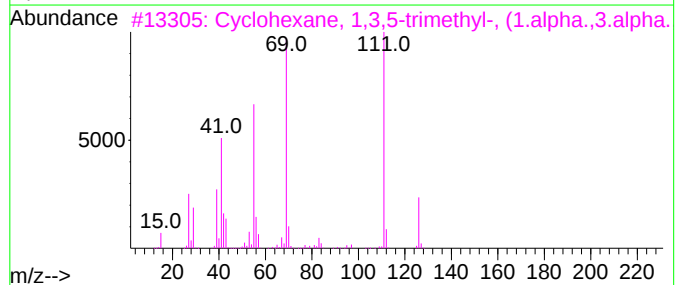
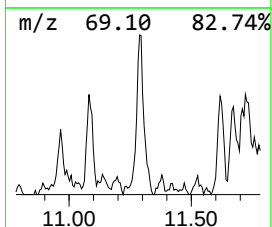
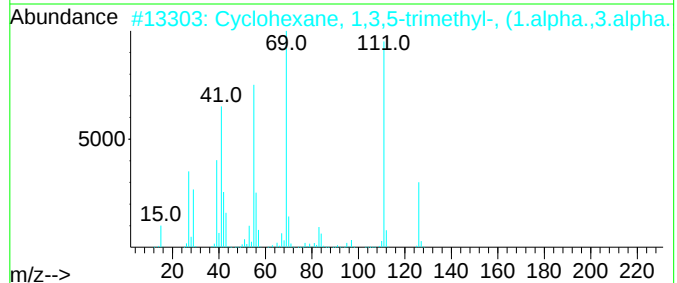
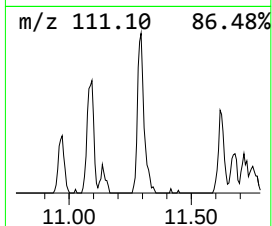
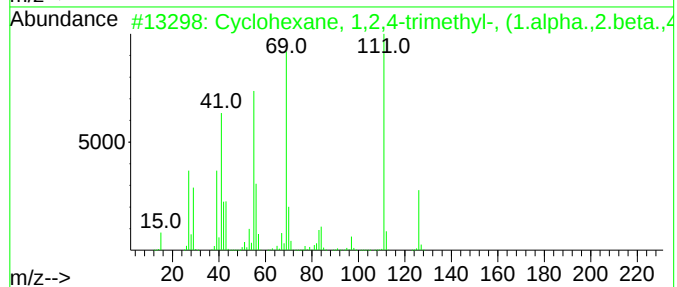
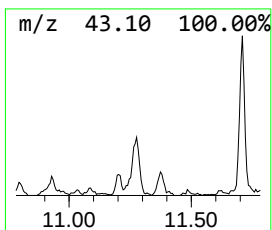
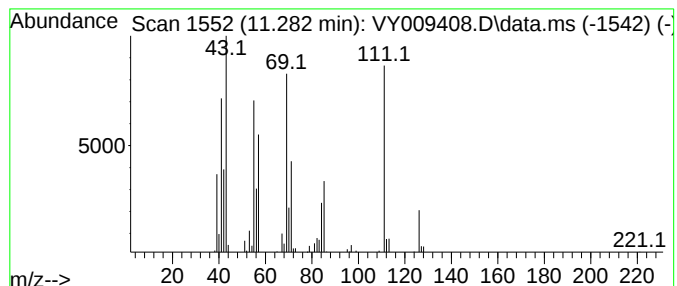
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062822S.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclohexane, 1,2,4-trimethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.282	6.78 ug/l	89180	Chlorobenzene-d5	11.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	91
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	60
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	49
4			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	49
5			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	43



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Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

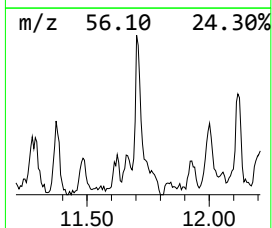
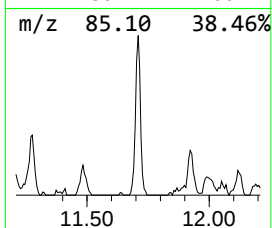
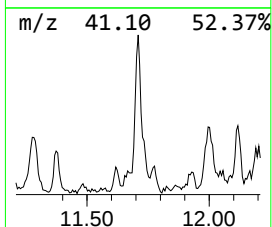
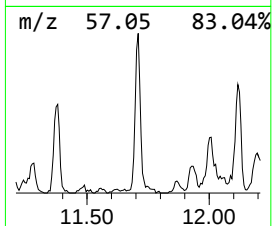
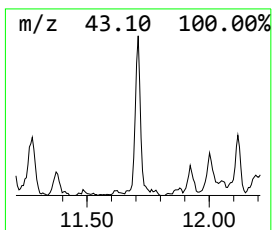
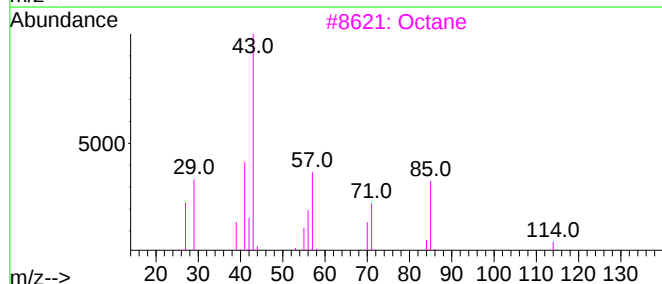
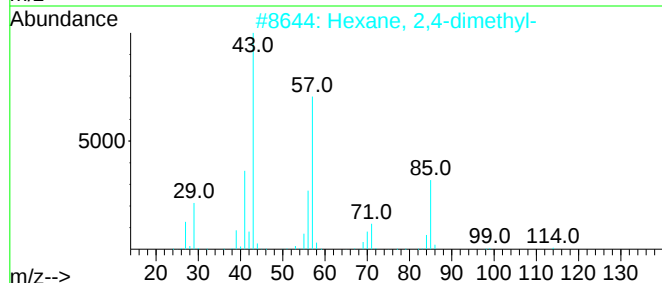
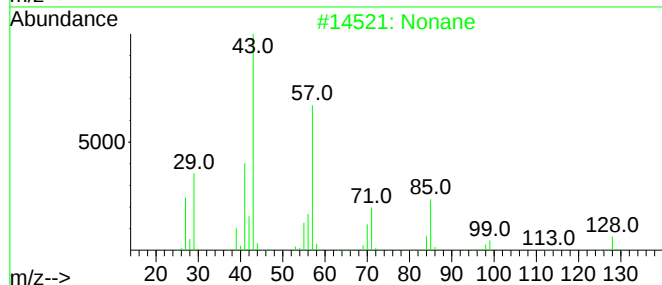
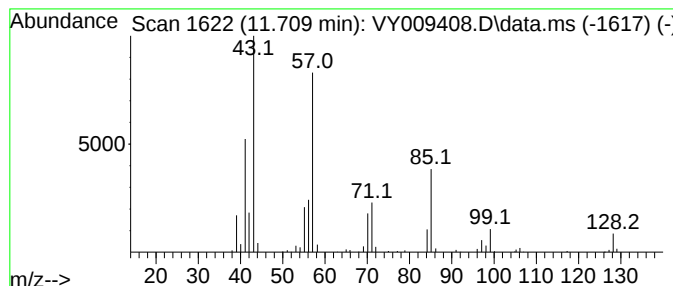
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062822S.M
Quant Title : SW846 8260

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

Peak Number 3 Nonane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.709	12.34 ug/l	162368	Chlorobenzene-d5	11.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	91
2			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59
3			Octane	114	C8H18	000111-65-9	59
4			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	53
5			3(2H)-Furanone, dihydro-5-isopro...	128	C7H12O2	034004-69-8	47



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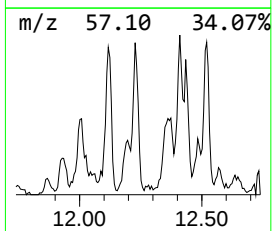
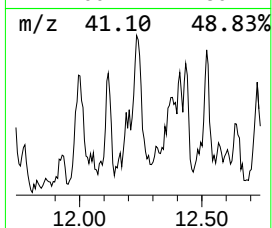
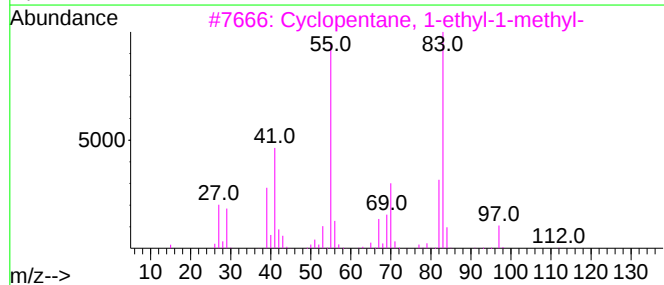
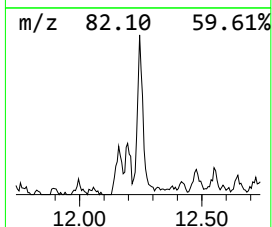
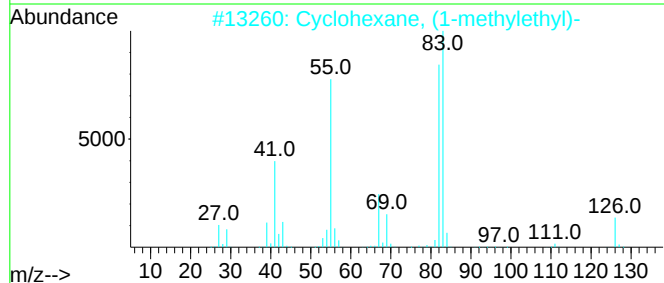
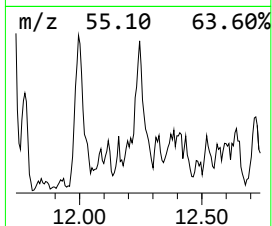
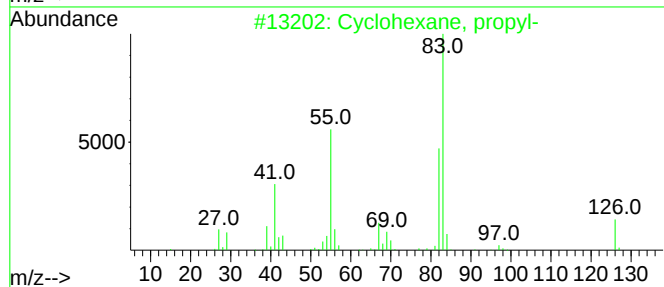
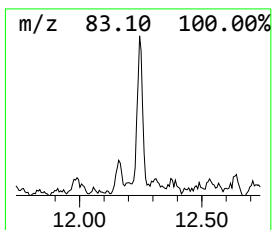
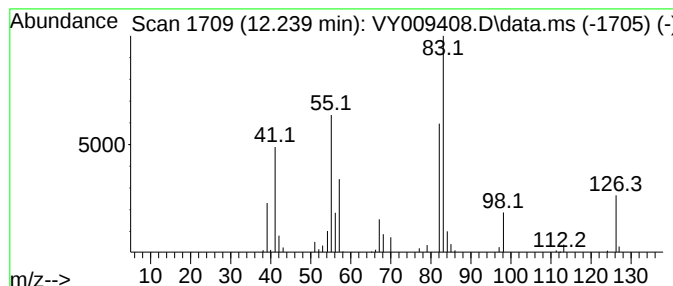
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.239	6.08 ug/l	79932	Chlorobenzene-d5	11.483

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	87
2			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	87
3			Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	53
4			1-Butene, 2,3,3-trimethyl-	98	C7H14	000594-56-9	49
5			2-Pyrazoline, 1-methyl-4-propyl-	126	C7H14N2	033063-77-3	47



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ClientSampleId :
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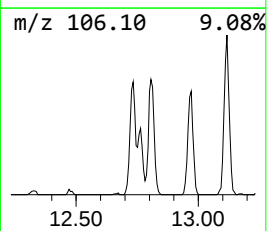
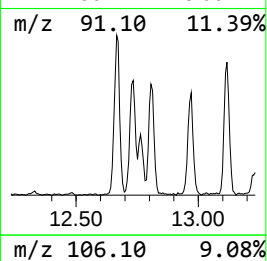
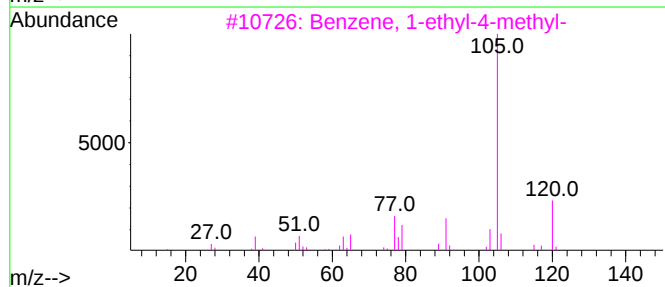
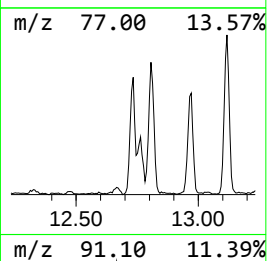
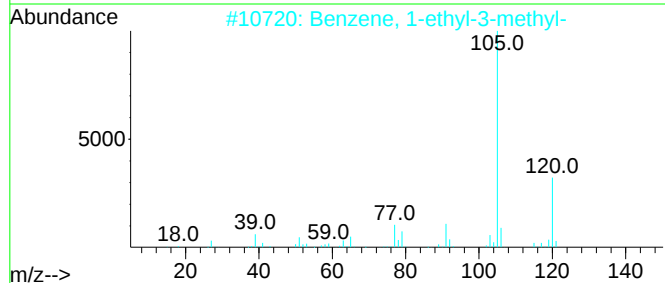
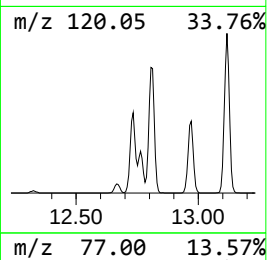
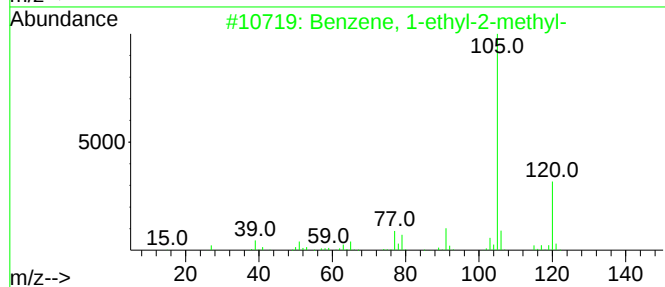
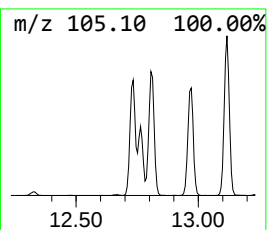
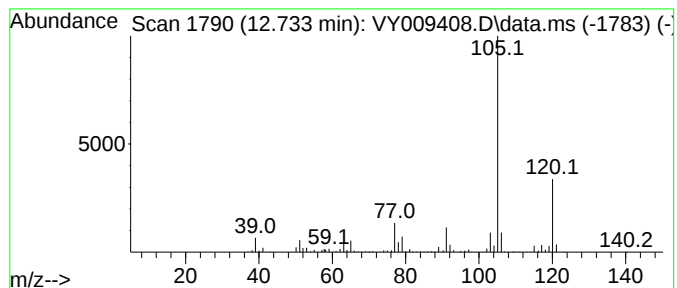
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	28.42 ug/l	441710	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	95
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
Data File : VY009408.D
Acq On : 30 Jun 2022 21:04
Operator : KP/MD
Sample : N3550-03
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

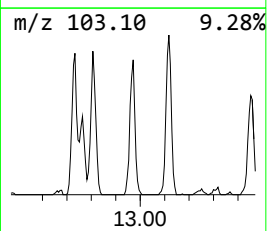
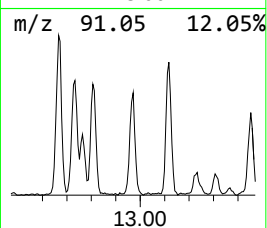
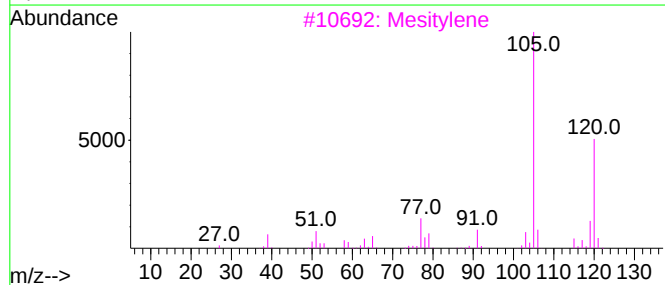
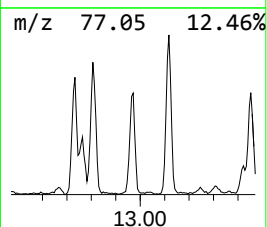
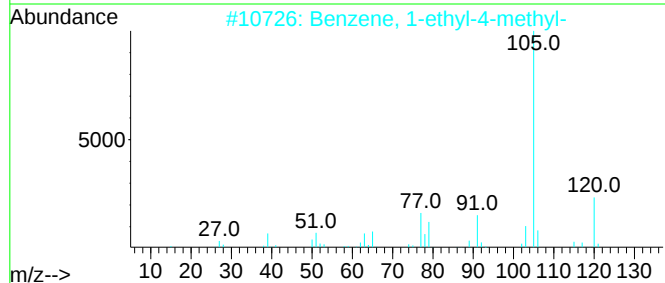
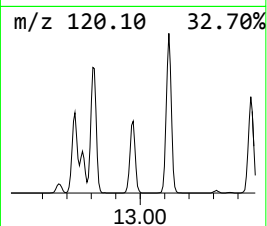
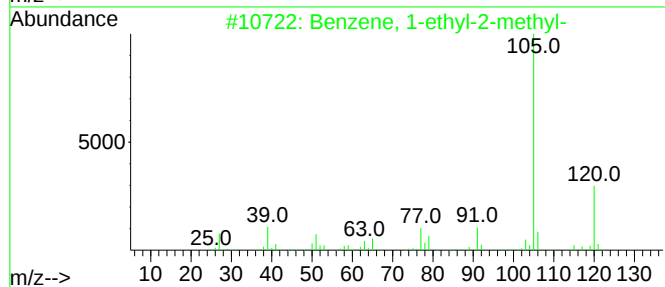
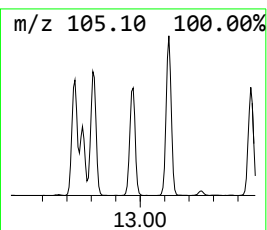
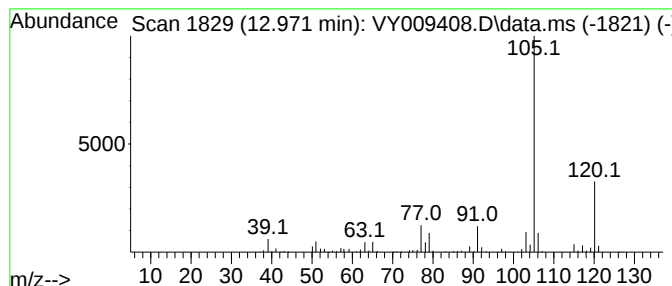
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	27.41 ug/l	426018	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
Data File : VY009408.D
Acq On : 30 Jun 2022 21:04
Operator : KP/MD
Sample : N3550-03
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

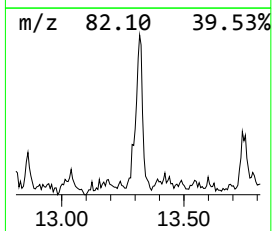
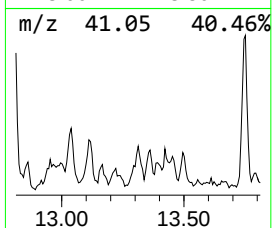
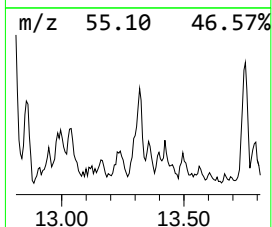
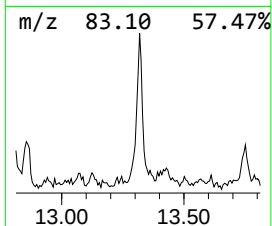
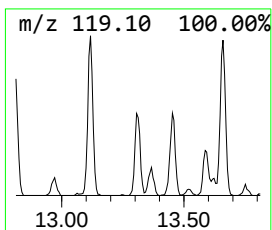
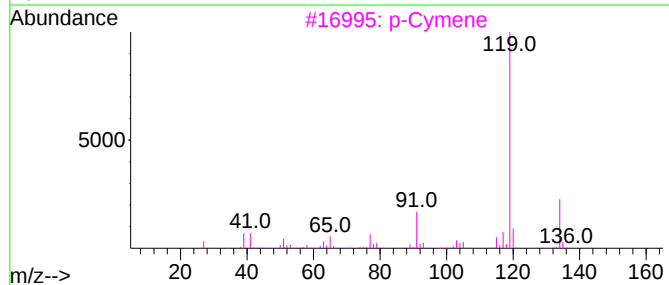
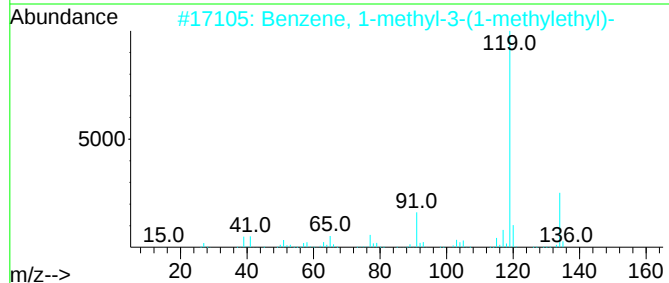
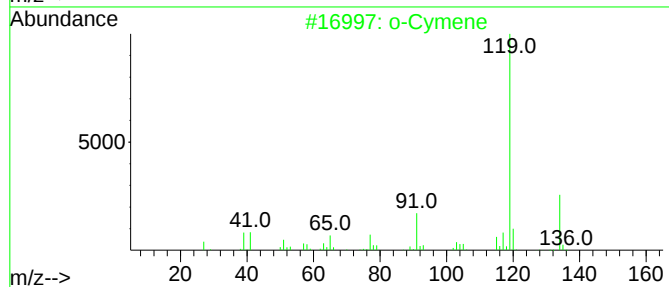
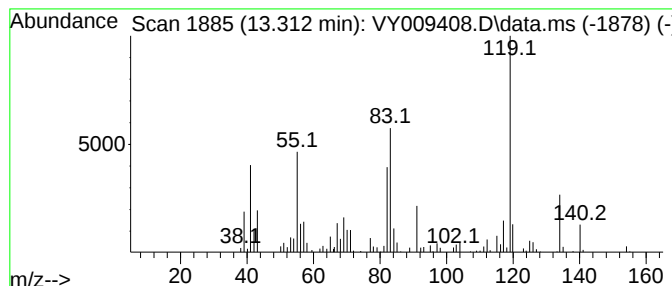
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Quant Title : SW846 8260

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

Peak Number 7 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.312	7.00 ug/l	108847	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	89
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	86
3			p-Cymene	134	C10H14	000099-87-6	86
4			1,3,5-Cycloheptatriene, 3,7,7-trimethyl-	134	C10H14	003479-89-8	50
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
Data File : VY009408.D
Acq On : 30 Jun 2022 21:04
Operator : KP/MD
Sample : N3550-03
Misc : 5.00g/5.0mL/MSVOA_Y/soil
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

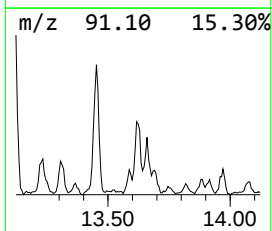
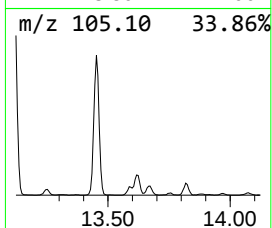
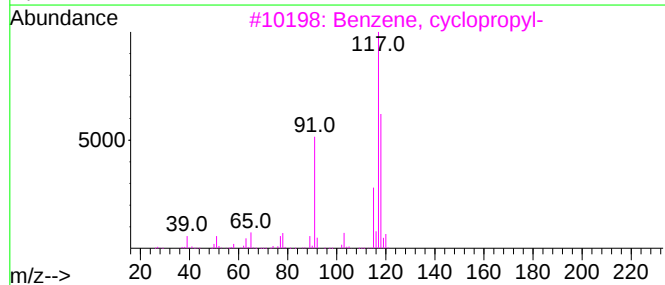
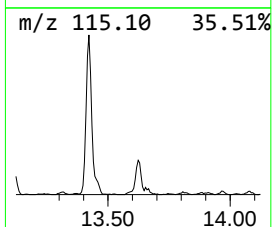
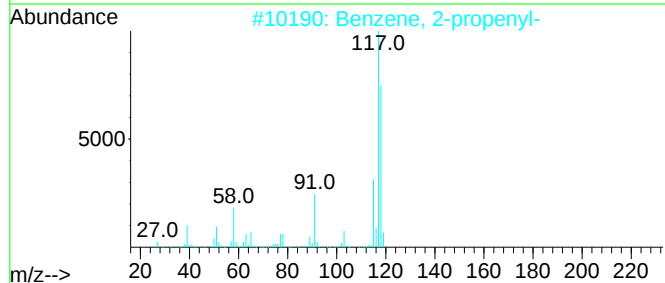
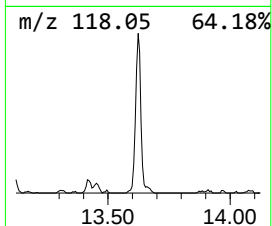
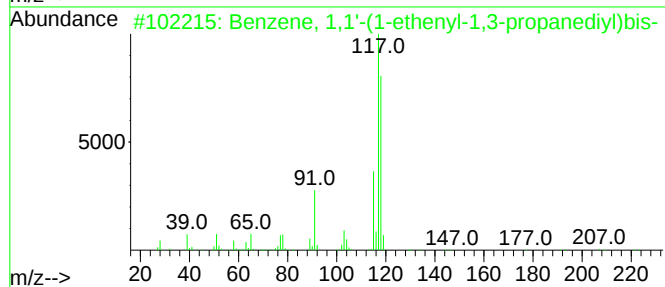
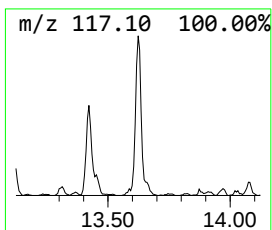
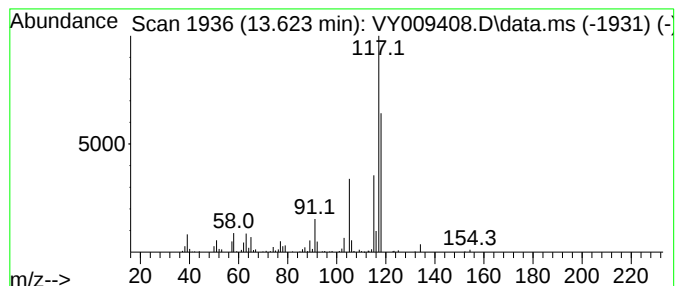
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,1'-(1-ethenyl-1,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	11.01 ug/l	171090	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	72
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4			Indane	118	C9H10	000496-11-7	70
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
Data File : VY009408.D
Acq On : 30 Jun 2022 21:04
Operator : KP/MD
Sample : N3550-03
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

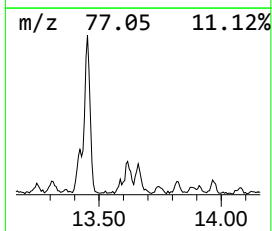
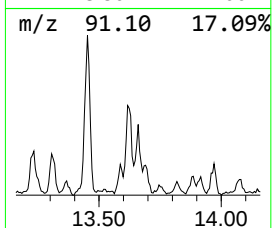
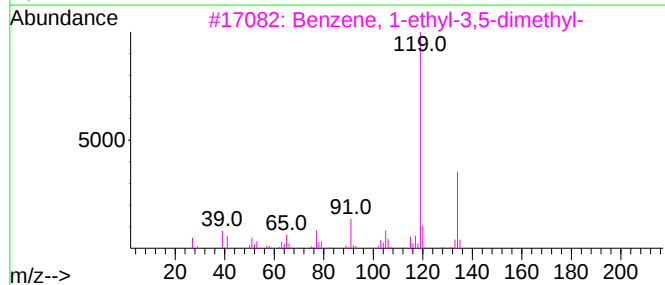
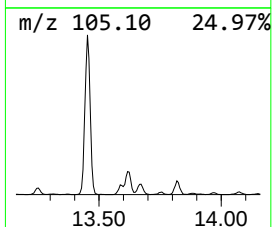
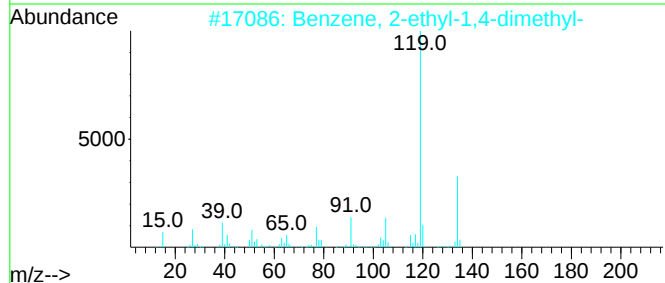
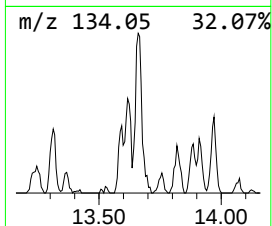
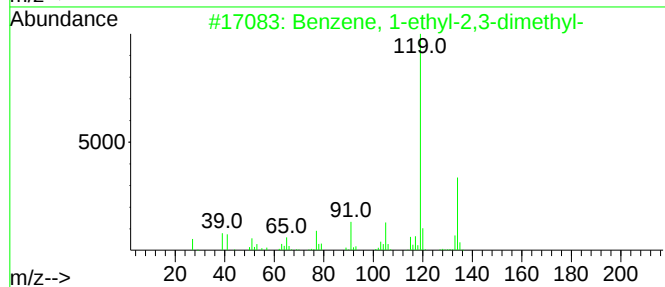
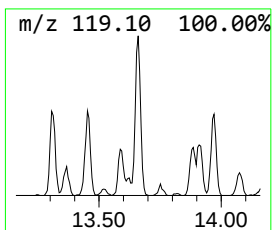
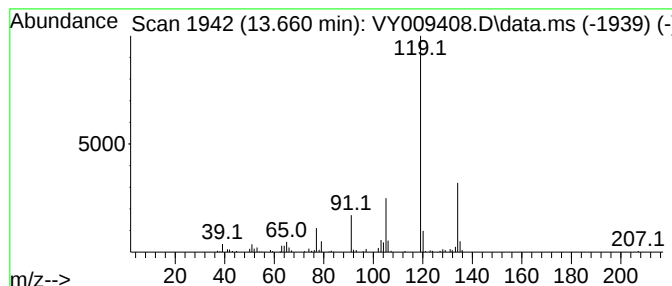
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.660	5.81 ug/l	90346	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87
3			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	83
4			o-Cymene	134	C10H14	000527-84-4	83
5			p-Cymene	134	C10H14	000099-87-6	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
Data File : VY009408.D
Acq On : 30 Jun 2022 21:04
Operator : KP/MD
Sample : N3550-03
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

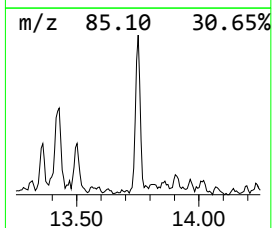
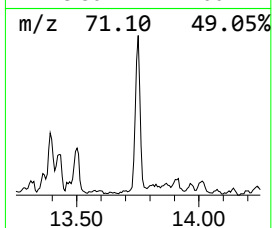
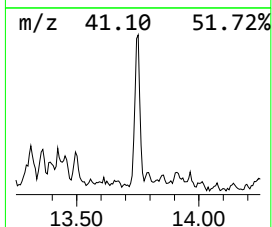
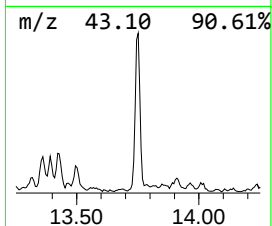
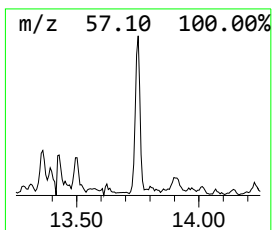
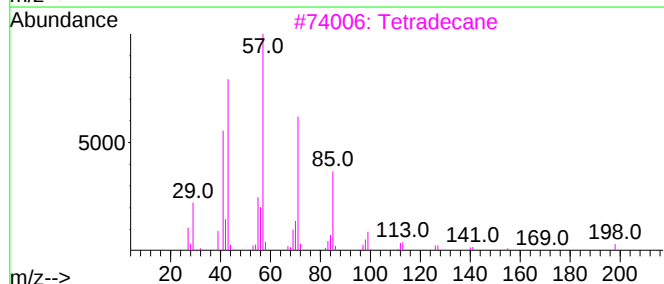
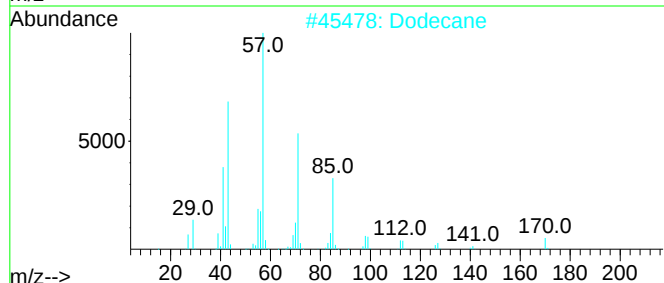
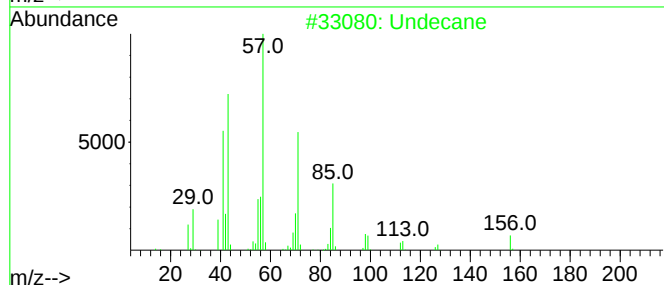
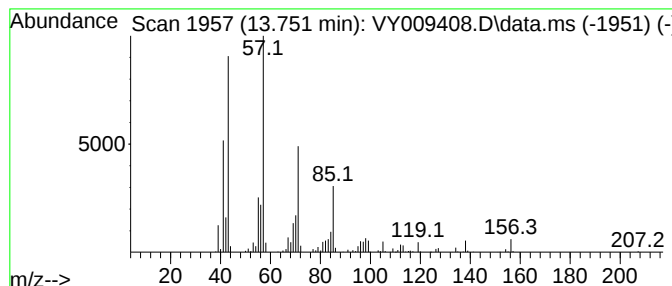
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Quant Title : SW846 8260

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

Peak Number 10 Undecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	14.23 ug/l	221111	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	96
2			Dodecane	170	C12H26	000112-40-3	80
3			Tetradecane	198	C14H30	000629-59-4	80
4			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
5			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY063022\
Data File : VY009408.D
Acq On : 30 Jun 2022 21:04
Operator : KP/MD
Sample : N3550-03
Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-03-06292022

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062822S.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
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Nonane	11.709	12.3	ug/l	162368	3	11.483	657679	50.0
Cyclohexane, pr...	12.239	6.1	ug/l	79932	3	11.483	657679	50.0
Benzene, 1-ethy...	12.733	28.4	ug/l	441710	4	13.422	776988	50.0
Benzene, 1-ethy...	12.971	27.4	ug/l	426018	4	13.422	776988	50.0
o-Cymene	13.312	7.0	ug/l	108847	4	13.422	776988	50.0
Benzene, 1,1'-(...	13.623	11.0	ug/l	171090	4	13.422	776988	50.0
Benzene, 1-ethy...	13.660	5.8	ug/l	90346	4	13.422	776988	50.0
Undecane	13.751	14.2	ug/l	221111	4	13.422	776988	50.0