

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
 Data File : VY015988.D
 Acq On : 16 Oct 2023 16:40
 Operator : SY/MD
 Sample : 04892-01
 Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 EO-02-10132023-A

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\82Y101223S.M

Title : SW846 8260

Signal : TIC: VY015988.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.728	467	477	485	rBV2	5066	15317	4.26%	0.447%
2	5.758	637	646	653	rBV5	3017	8839	2.46%	0.258%
3	7.728	959	969	974	rBV	53697	125191	34.85%	3.657%
4	7.795	974	980	992	rVB	90095	203539	56.66%	5.946%
5	8.148	1030	1038	1052	rBV	55061	122989	34.24%	3.593%
6	8.697	1120	1128	1140	rBV	133859	260463	72.51%	7.608%
7	10.185	1364	1372	1383	rBV	206656	359231	100.00%	10.494%
8	10.587	1433	1438	1443	rBV2	7156	12390	3.45%	0.362%
9	11.044	1510	1513	1516	rVV4	3027	4332	1.21%	0.127%
10	11.093	1516	1521	1527	rVB4	5359	9223	2.57%	0.269%
11	11.209	1535	1540	1545	rBV7	2335	3963	1.10%	0.116%
12	11.288	1545	1553	1564	rVB9	4644	15414	4.29%	0.450%
13	11.386	1564	1569	1573	rBV2	2982	4776	1.33%	0.140%
14	11.495	1580	1587	1597	rBV	196678	319252	88.87%	9.326%
15	11.678	1613	1617	1622	rVV5	3329	7325	2.04%	0.214%
16	11.739	1622	1627	1632	rVV4	5664	13160	3.66%	0.384%
17	11.782	1632	1634	1640	rVB3	3876	5450	1.52%	0.159%
18	12.013	1664	1672	1683	rBV9	6624	24794	6.90%	0.724%
19	12.129	1683	1691	1694	rVB3	5741	9568	2.66%	0.279%
20	12.209	1699	1704	1706	rBV2	5597	8347	2.32%	0.244%
21	12.245	1706	1710	1716	rVB5	6027	12070	3.36%	0.353%
22	12.489	1744	1750	1760	rBV	158064	258399	71.93%	7.548%
23	12.647	1772	1776	1783	rVB5	7426	16357	4.55%	0.478%
24	12.733	1783	1790	1794	rBV8	5786	14549	4.05%	0.425%
25	12.818	1798	1804	1809	rVV4	20119	46272	12.88%	1.352%
26	12.867	1809	1812	1817	rVB2	7434	11072	3.08%	0.323%
27	12.977	1822	1830	1835	rVV3	13409	31757	8.84%	0.928%
28	13.044	1838	1841	1847	rVV4	13802	21900	6.10%	0.640%
29	13.135	1848	1856	1859	rVV8	4038	10334	2.88%	0.302%
30	13.178	1859	1863	1866	rVV5	4454	7602	2.12%	0.222%
31	13.227	1867	1871	1873	rVV4	4429	7393	2.06%	0.216%
32	13.251	1873	1875	1879	rVV2	5978	7349	2.05%	0.215%
33	13.324	1879	1887	1891	rVV5	17545	36692	10.21%	1.072%
34	13.367	1891	1894	1898	rVV4	5594	9975	2.78%	0.291%
35	13.428	1898	1904	1914	rVV2	204627	352099	98.01%	10.285%
36	13.599	1929	1932	1933	rBV2	5187	5036	1.40%	0.147%

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Integrator: RTE

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Start Thrs: 0.2

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Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101223S.M
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37	13.629	1933	1937	1940	rVB	15127	19901	5.54%	0.581%
38	13.672	1940	1944	1952	rVB3	19885	37517	10.44%	1.096%
39	13.757	1952	1958	1962	rBV3	28356	49676	13.83%	1.451%
40	13.830	1962	1970	1973	rBV	18639	35928	10.00%	1.049%
41	13.922	1981	1985	1988	rVB3	10183	13246	3.69%	0.387%
42	13.977	1989	1994	1999	rVB3	31766	52949	14.74%	1.547%
43	14.019	1999	2001	2006	rVB3	5329	9161	2.55%	0.268%
44	14.080	2006	2011	2017	rBV2	29339	47673	13.27%	1.393%
45	14.147	2017	2022	2028	rBV8	7118	18870	5.25%	0.551%
46	14.214	2028	2033	2037	rBV2	11623	20792	5.79%	0.607%
47	14.263	2037	2041	2045	rBV5	13318	23077	6.42%	0.674%
48	14.300	2045	2047	2050	rVB2	5602	5603	1.56%	0.164%
49	14.336	2050	2053	2057	rVB	13431	18641	5.19%	0.545%
50	14.385	2057	2061	2064	rBV3	14040	20200	5.62%	0.590%
51	14.428	2064	2068	2071	rBV3	15231	21920	6.10%	0.640%
52	14.489	2074	2078	2079	rBV4	5648	6925	1.93%	0.202%
53	14.568	2087	2091	2096	rBV5	8055	15393	4.28%	0.450%
54	14.617	2096	2099	2103	rVB	19998	25951	7.22%	0.758%
55	14.684	2103	2110	2116	rBV4	49235	116249	32.36%	3.396%
56	14.745	2116	2120	2125	rVB4	13611	25887	7.21%	0.756%
57	14.836	2130	2135	2142	rBV	26565	44465	12.38%	1.299%
58	14.946	2149	2153	2157	rBV2	18329	29019	8.08%	0.848%
59	15.062	2166	2172	2177	rVB2	20751	43729	12.17%	1.277%
60	15.220	2192	2198	2201	rBV7	7498	17007	4.73%	0.497%
61	15.299	2206	2211	2217	rBV4	16976	40545	11.29%	1.184%
62	15.354	2217	2220	2222	rBV3	4589	7741	2.15%	0.226%
63	15.440	2231	2234	2242	rVB9	9487	16834	4.69%	0.492%
64	15.531	2243	2249	2256	rVV5	9231	21961	6.11%	0.642%
65	15.690	2267	2275	2278	rBV7	12465	31715	8.83%	0.926%
66	15.732	2278	2282	2289	rVV2	23782	46416	12.92%	1.356%
67	15.824	2293	2297	2302	rVV7	6489	15265	4.25%	0.446%
68	15.891	2305	2308	2322	rVB6	10831	30920	8.61%	0.903%
69	16.043	2326	2333	2341	rBV3	19633	41579	11.57%	1.215%
70	16.232	2358	2364	2371	rVB4	14406	30455	8.48%	0.890%
71	16.299	2371	2375	2380	rBV3	5681	10578	2.94%	0.309%
72	16.488	2402	2406	2407	rBV4	6388	8575	2.39%	0.250%
73	16.549	2414	2416	2420	rVB5	4830	5708	1.59%	0.167%
74	16.635	2423	2430	2437	rVB8	4844	12874	3.58%	0.376%

Sum of corrected areas: 3423364

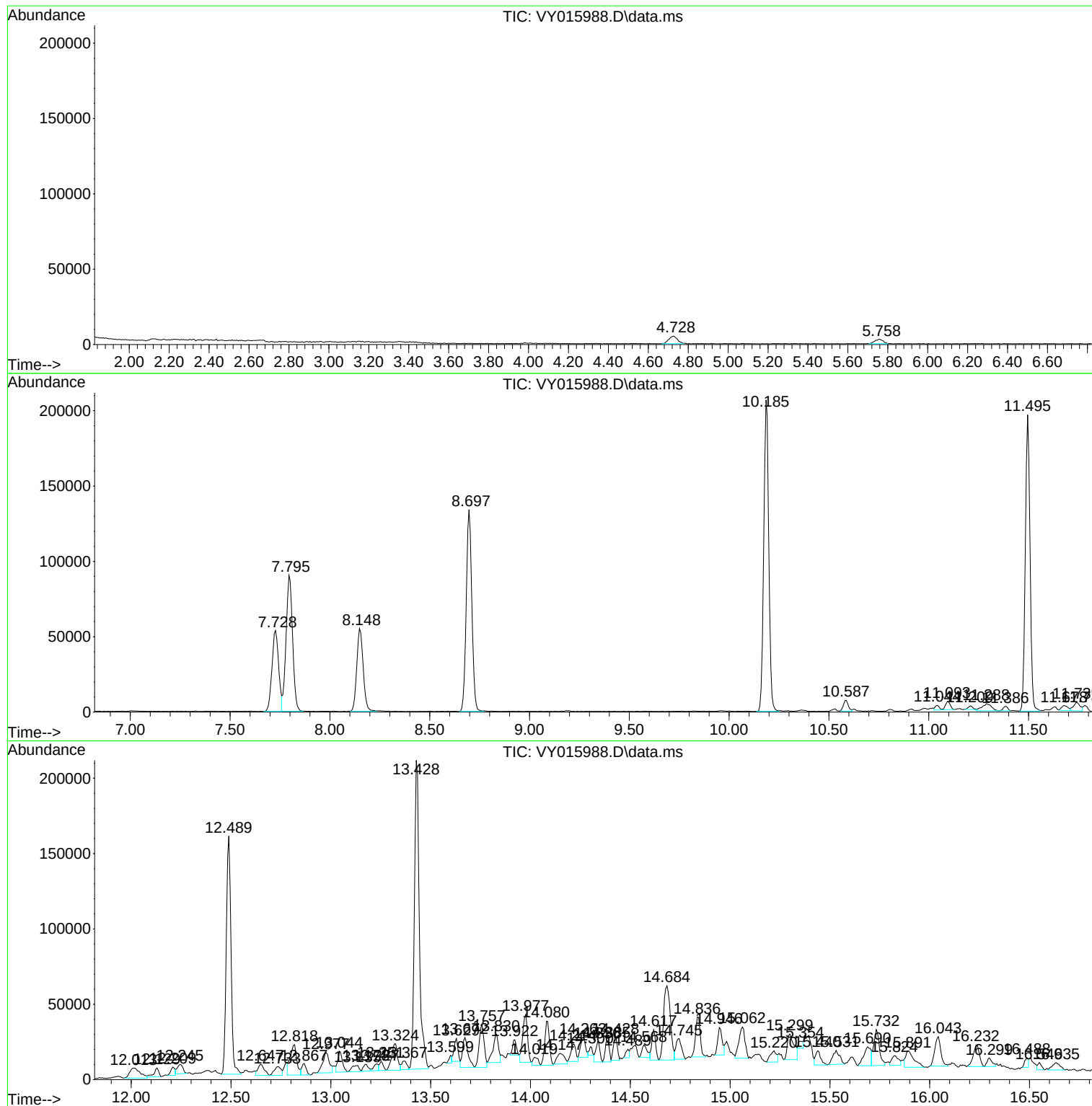
Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

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Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101223S.M
Quant Title : SW846 8260

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



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ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

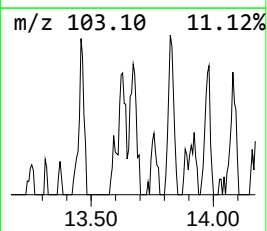
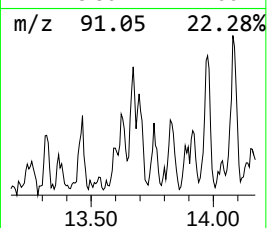
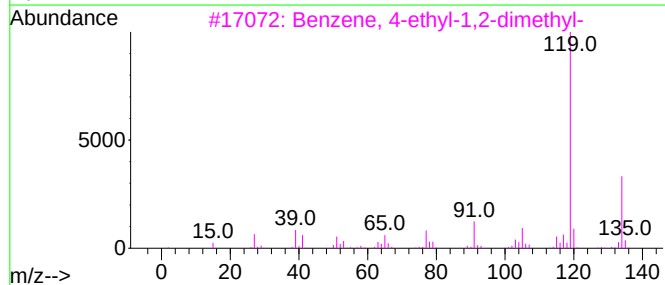
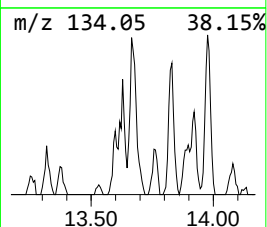
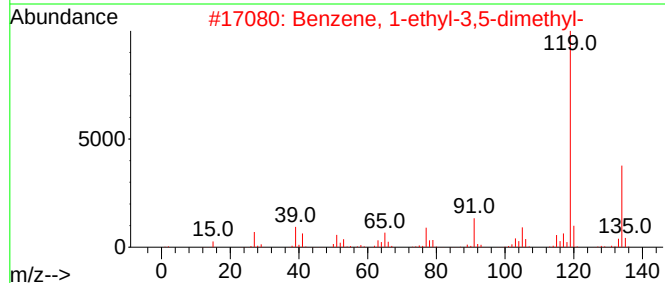
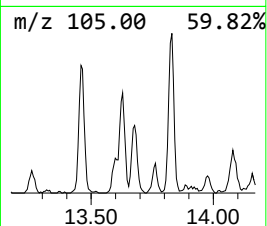
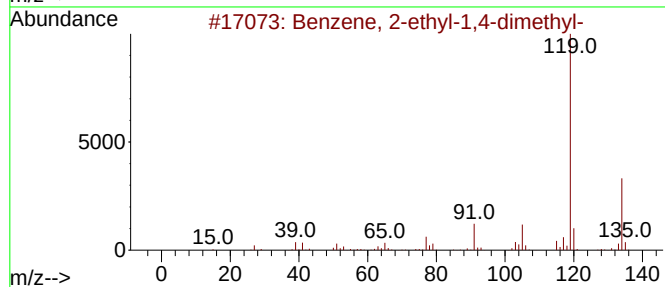
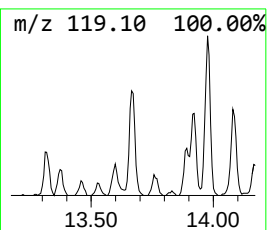
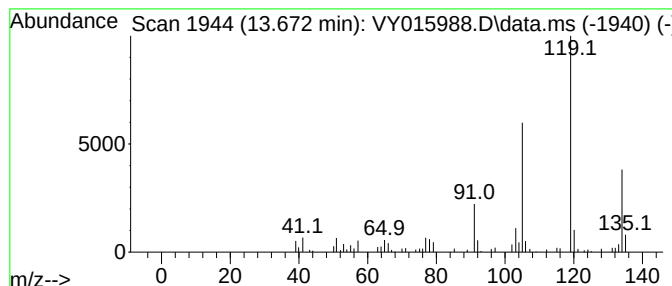
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101223S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.672	5.33 ug/l	37517	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
2			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	83



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Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

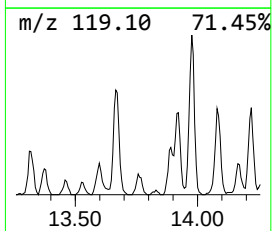
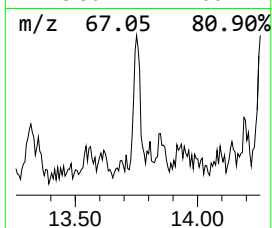
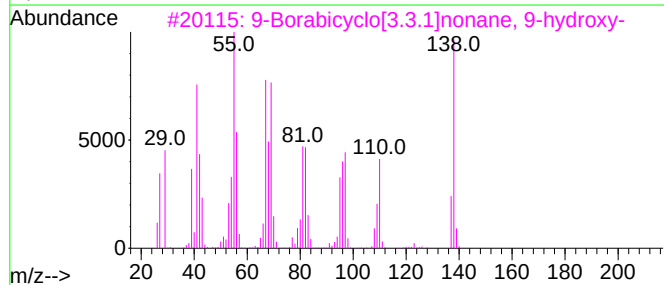
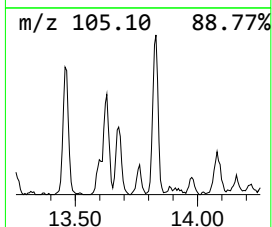
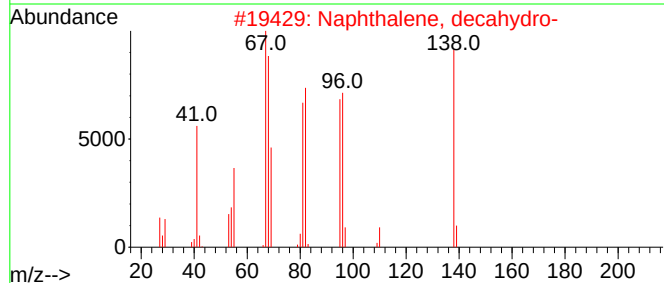
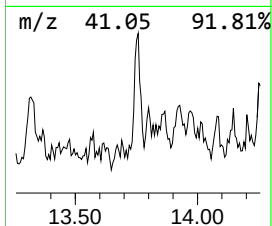
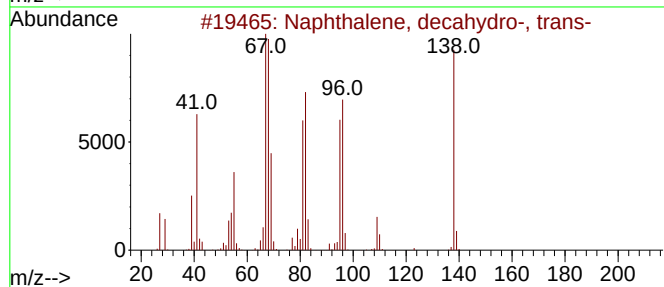
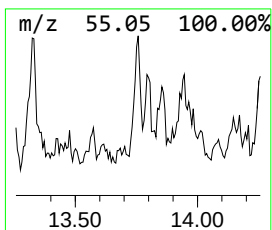
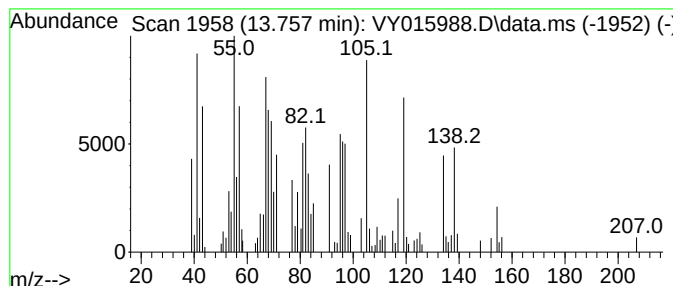
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101223S.M
Quant Title : SW846 8260

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

Peak Number 2 Naphthalene, decahydro-, tr... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.757	7.05 ug/l	49676	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	64
2			Naphthalene, decahydro-	138	C10H18	000091-17-8	50
3			9-Borabicyclo[3.3.1]nonane, 9-hy...	138	C8H15BO	063366-65-4	46
4			1H-Inden-1-one, octahydro-, trans-	138	C9H14O	016783-22-5	41
5			Cyclohexane, 1-methyl-3-(1-methy...	138	C10H18	024399-15-3	38



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ClientSampleId :
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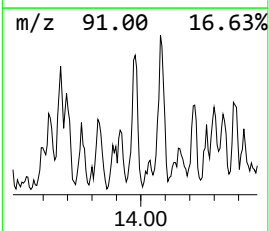
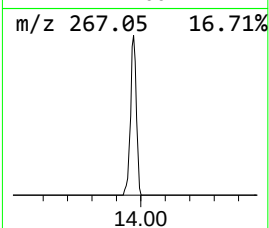
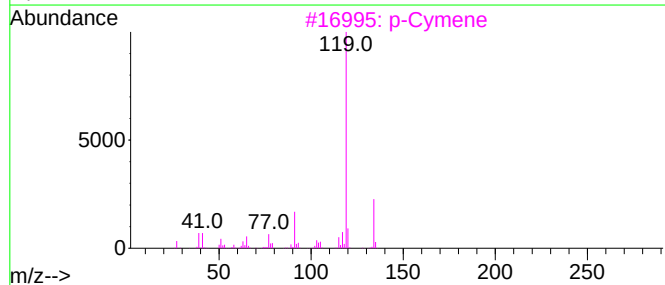
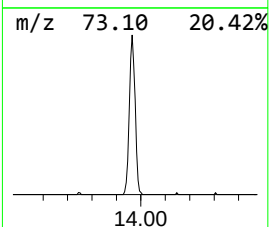
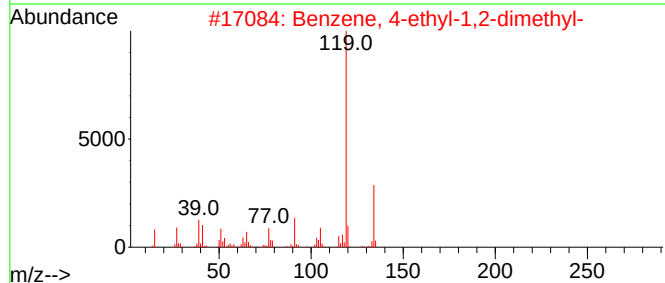
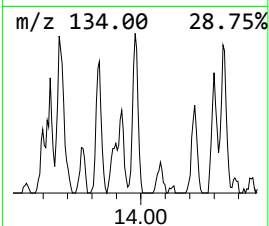
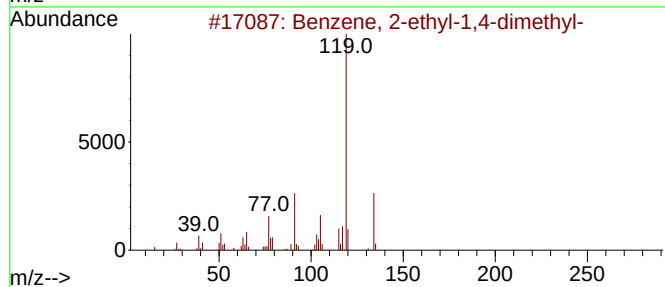
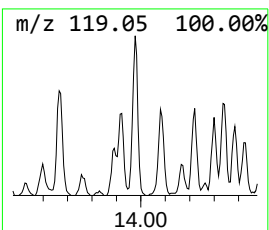
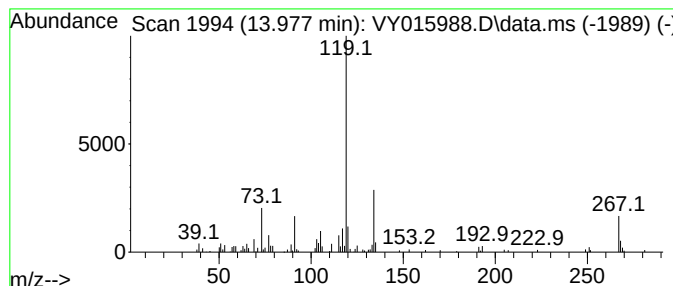
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.977	7.52 ug/l	52949	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
3			p-Cymene	134	C10H14	000099-87-6	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



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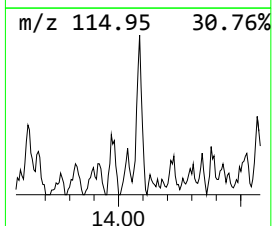
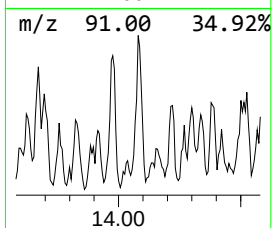
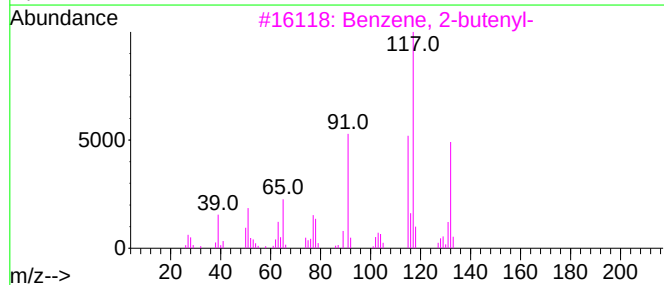
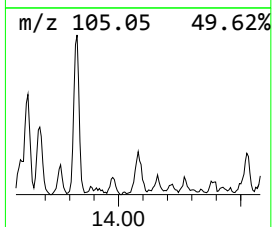
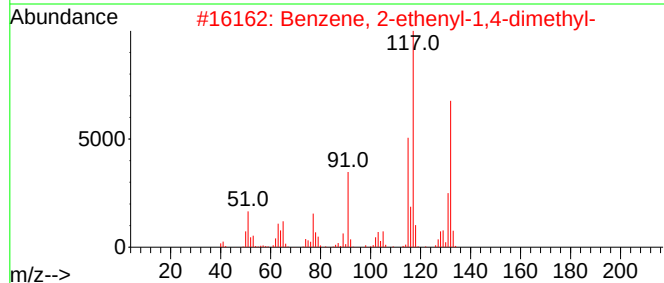
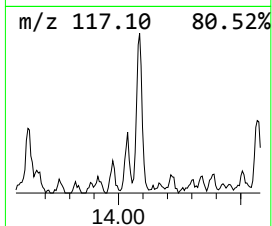
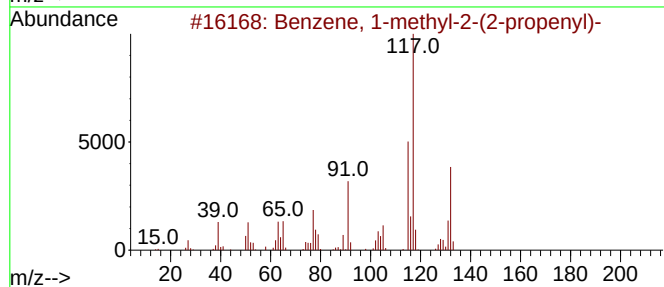
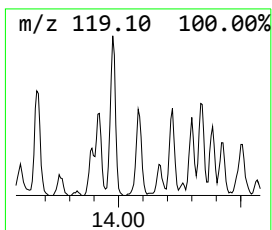
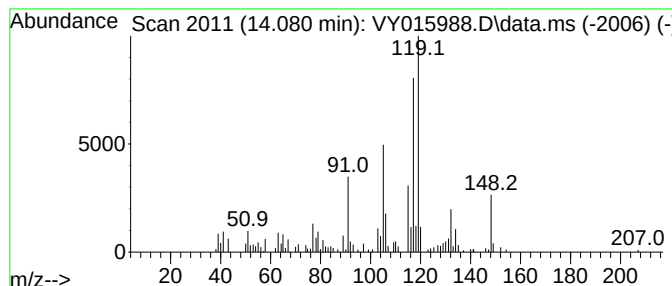
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101223S.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-methyl-2-(2-prop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	6.77 ug/l	47673	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	64
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
4			2,4-Dimethylstyrene	132	C10H12	002234-20-0	64
5			(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

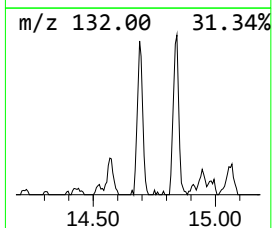
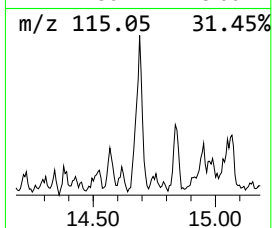
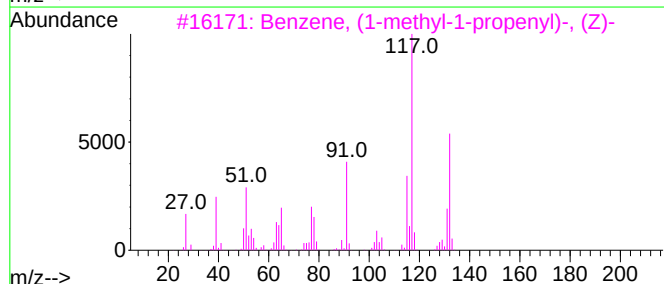
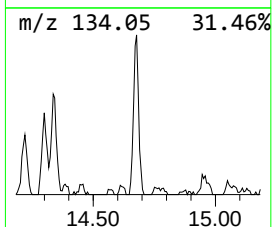
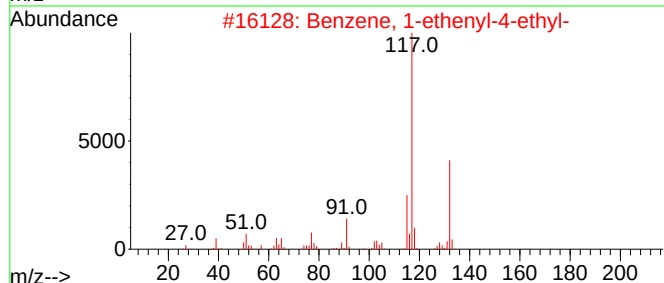
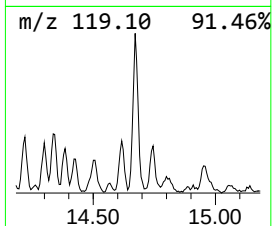
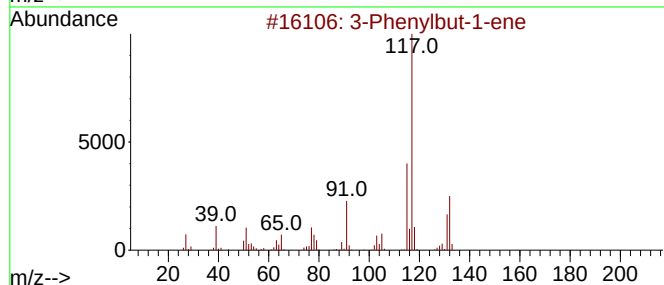
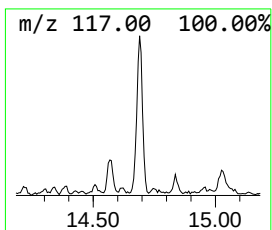
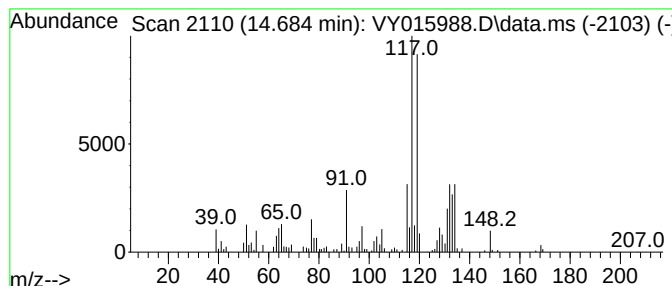
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Quant Title : SW846 8260

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

Peak Number 5 3-Phenylbut-1-ene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	16.51 ug/l	116249	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	46
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	46
4			Benzene, 2-butenyl-	132	C10H12	001560-06-1	46
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/soil
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

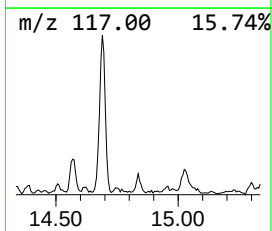
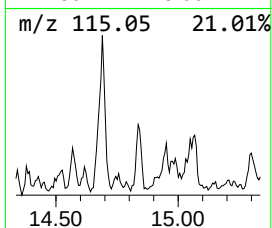
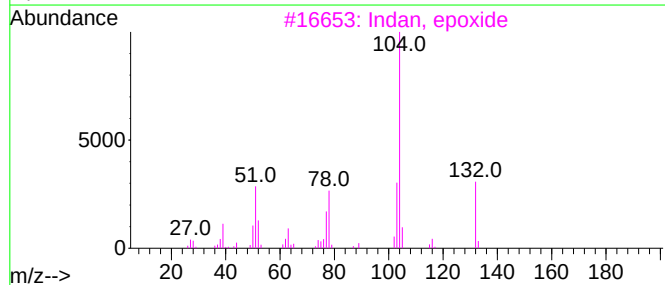
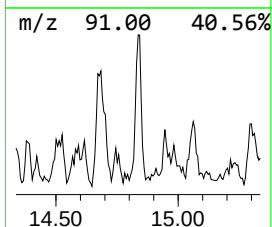
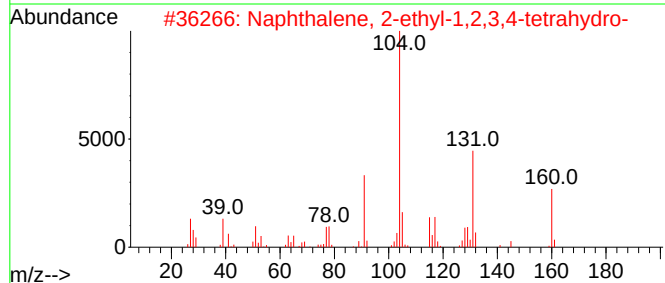
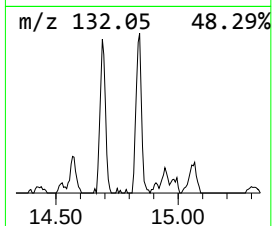
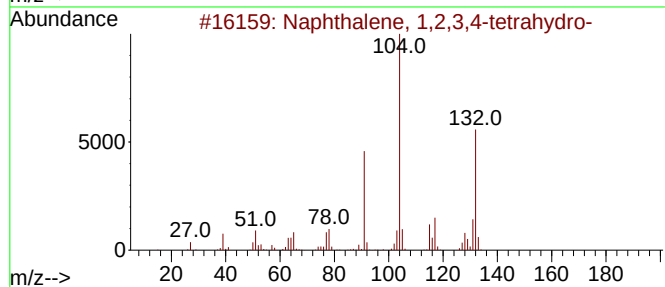
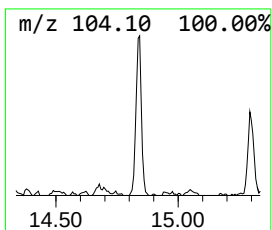
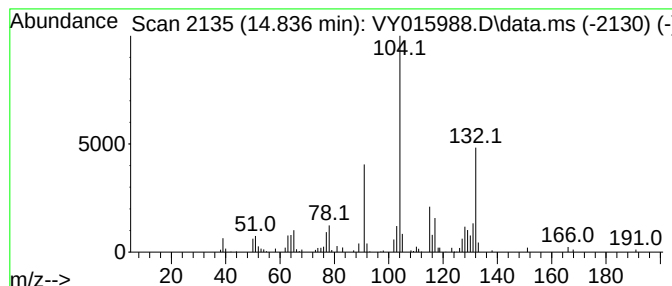
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Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.836	6.31 ug/l	44465	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	50
3			Indan, epoxide	132	C9H8O	000768-22-9	49
4			4-Ethylbenzoic acid, 2-phenyleth...	254	C17H18O2	1000293-50-1	35
5			Isoquinoline, 1,2,3,4-tetrahydro-	133	C9H11N	000091-21-4	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

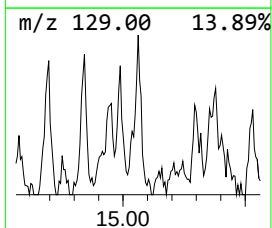
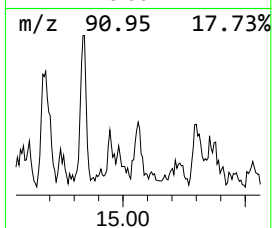
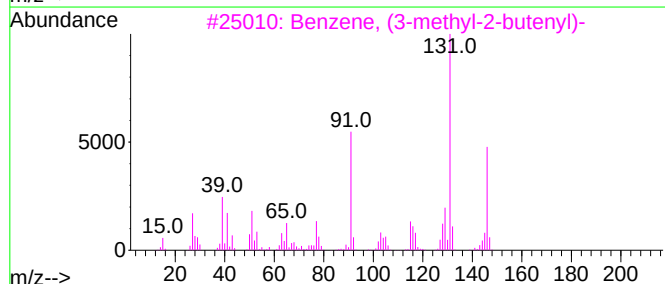
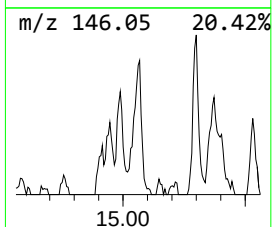
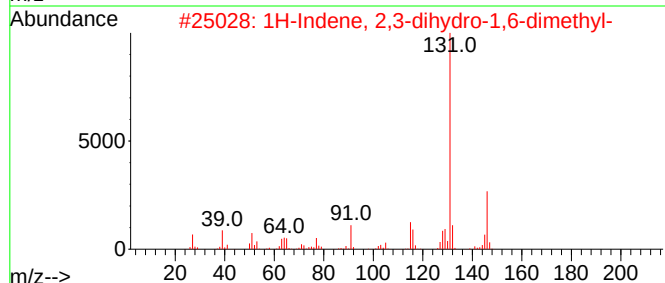
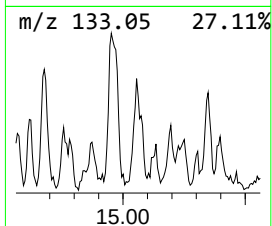
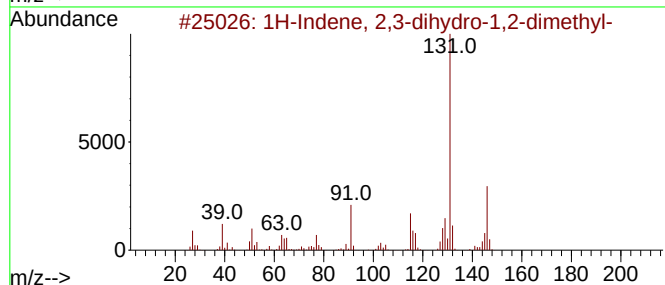
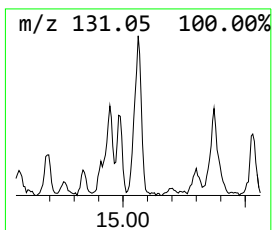
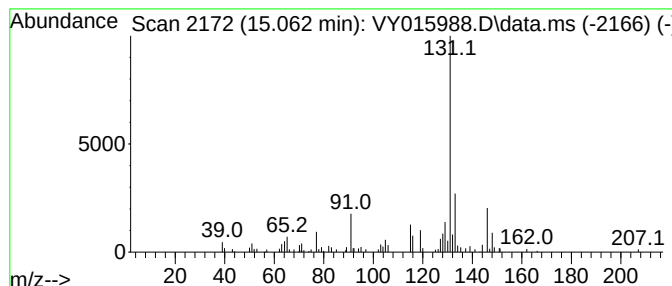
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.062	6.21 ug/l	43729	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	76
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

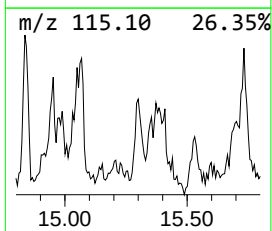
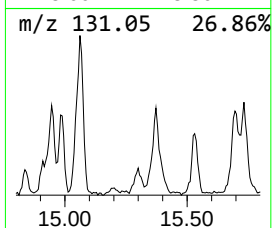
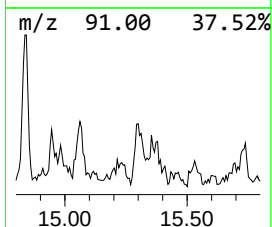
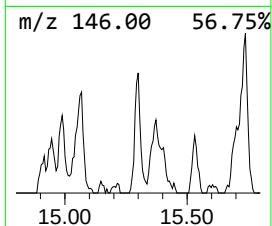
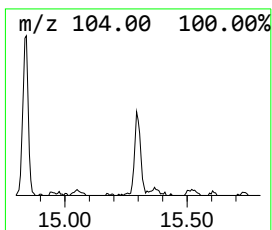
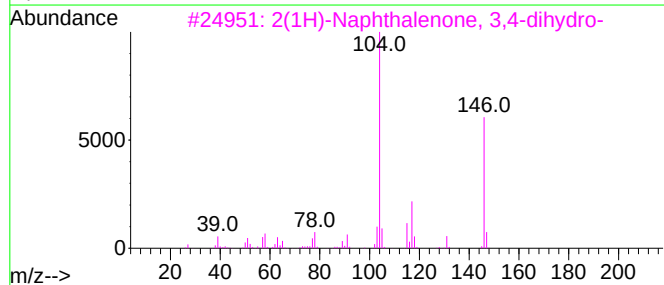
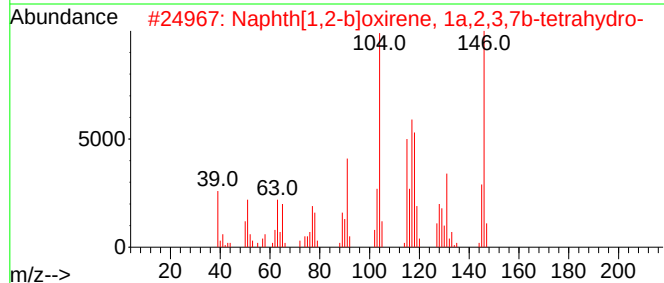
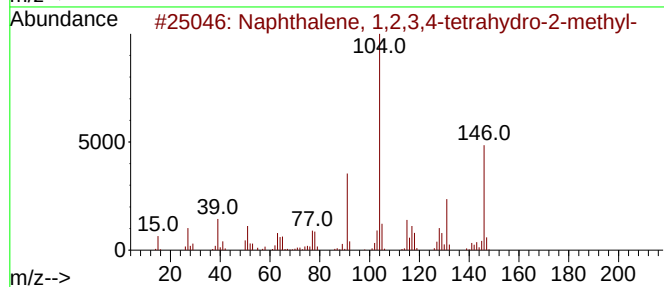
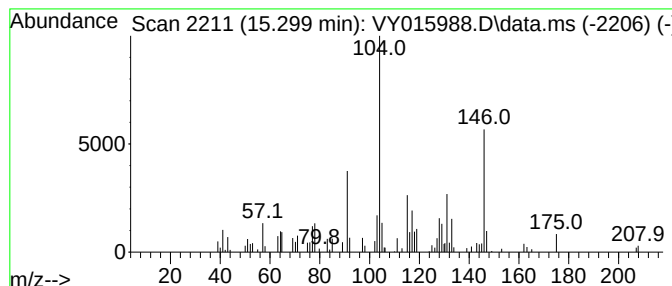
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.300	5.76 ug/l	40545	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	95
2			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	78
3			2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	68
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46
5			Benzeneethanol, .alpha.,.alpha.-...	162	C11H14O	025982-72-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

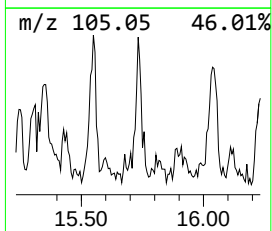
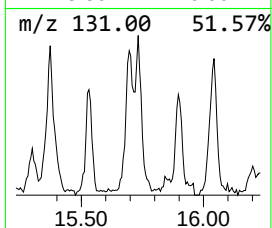
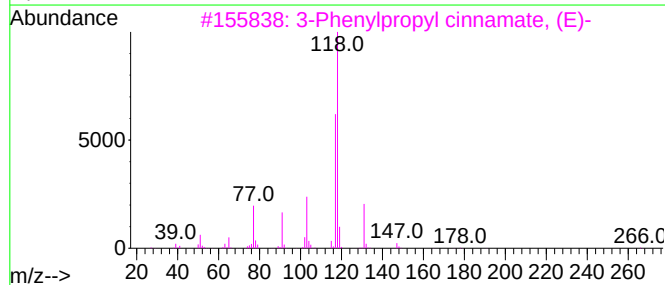
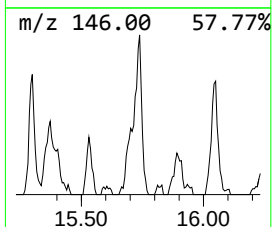
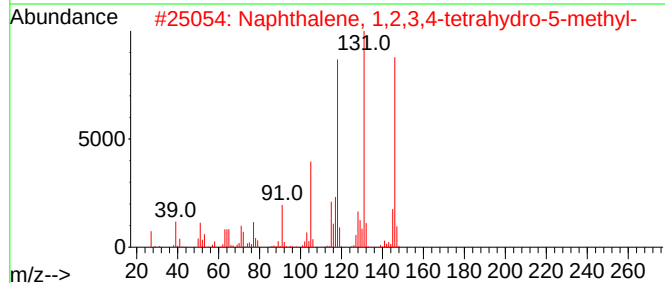
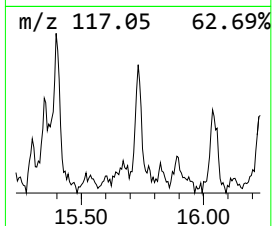
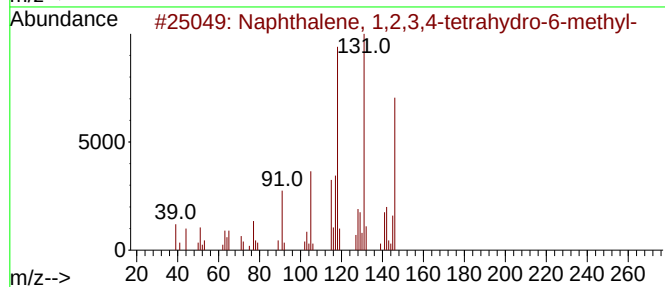
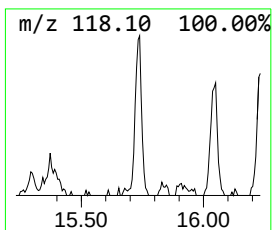
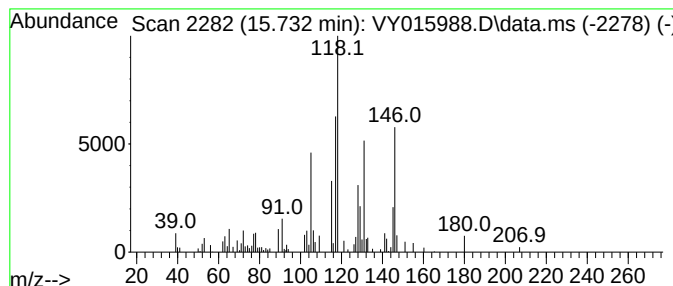
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.732	6.59 ug/l	46416	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	76
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	68
3			3-Phenylpropyl cinnamate, (E)-	266	C18H18O2	290311-72-7	38
4			Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	38
5			Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-02-10132023-A

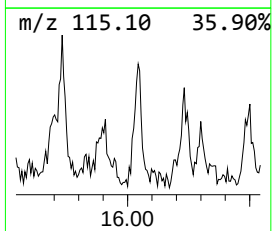
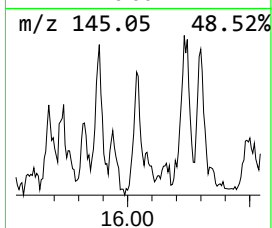
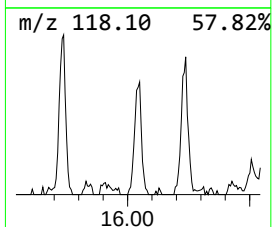
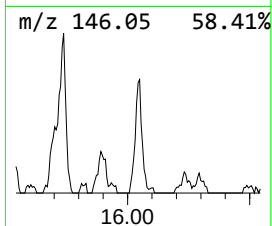
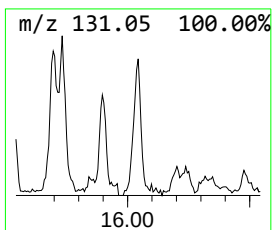
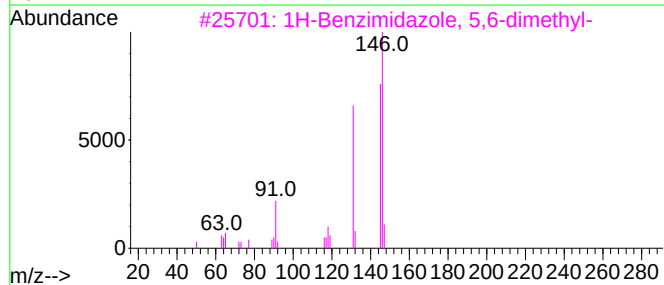
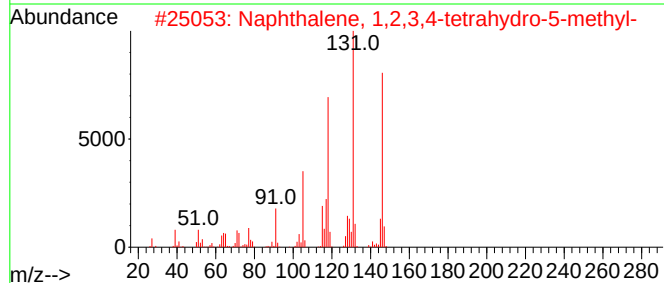
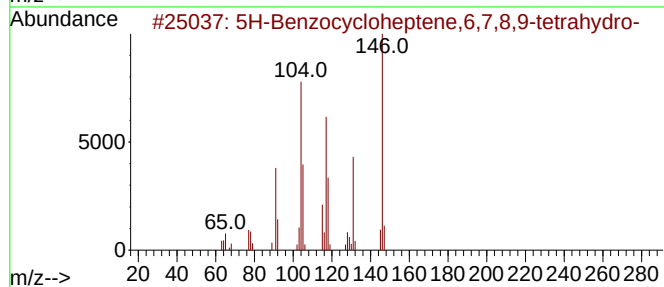
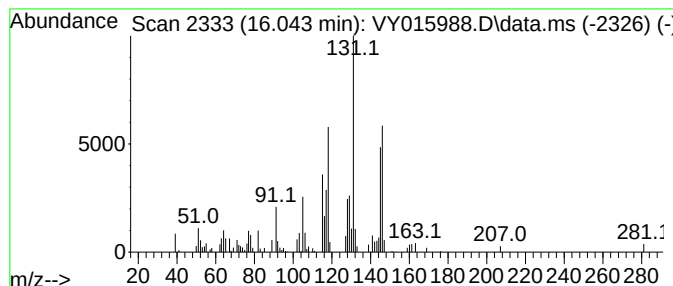
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Quant Title : SW846 8260

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

Peak Number 10 5H-Benzocycloheptene,6,7,8,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.043	5.90 ug/l	41579	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5H-Benzocycloheptene,6,7,8,9-tet...	146	C11H14	001075-16-7	62
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58
3			1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	58
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	53
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY101623\
Data File : VY015988.D
Acq On : 16 Oct 2023 16:40
Operator : SY/MD
Sample : 04892-01
Misc : 5.01g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
EO-02-10132023-A

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101223S.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
Benzene, 2-ethy...	13.672	5.3	ug/l	37517	4	13.428	352099	50.0
Naphthalene, de...	13.757	7.0	ug/l	49676	4	13.428	352099	50.0
Benzene, 4-ethy...	13.977	7.5	ug/l	52949	4	13.428	352099	50.0
Benzene, 1-meth...	14.080	6.8	ug/l	47673	4	13.428	352099	50.0
3-Phenylbut-1-ene	14.684	16.5	ug/l	116249	4	13.428	352099	50.0
Naphthalene, 1,...	14.836	6.3	ug/l	44465	4	13.428	352099	50.0
1H-Indene, 2,3-...	15.062	6.2	ug/l	43729	4	13.428	352099	50.0
Naphthalene, 1,...	15.300	5.8	ug/l	40545	4	13.428	352099	50.0
Naphthalene, 1,...	15.732	6.6	ug/l	46416	4	13.428	352099	50.0
5H-Benzocyclohe...	16.043	5.9	ug/l	41579	4	13.428	352099	50.0