

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
 Data File : VY017259.D
 Acq On : 06 Feb 2024 13:05
 Operator : SY/MD
 Sample : P1334-07
 Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 HR-04-020124

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\82Y020124S.M
 Title : SW846 8260

Signal : TIC: VY017259.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.734	959	970	976	rBV	67205	155971	1.16%	0.242%
2	7.807	976	982	995	rVB	98497	220108	1.63%	0.341%
3	8.161	1027	1040	1052	rBV	58524	135344	1.00%	0.210%
4	8.703	1121	1129	1140	rBV	157853	316478	2.35%	0.491%
5	10.197	1366	1374	1380	rBV	274553	468115	3.47%	0.726%
6	11.294	1545	1554	1565	rVB4	120985	305511	2.27%	0.474%
7	11.392	1565	1570	1575	rBV	103287	166254	1.23%	0.258%
8	11.508	1583	1589	1599	rBV	239190	391472	2.91%	0.607%
9	11.721	1614	1624	1633	rVV4	622164	1498252	11.12%	2.323%
10	11.794	1633	1636	1641	rVB2	93682	147270	1.09%	0.228%
11	12.044	1665	1677	1685	rBV2	487113	1172315	8.70%	1.818%
12	12.136	1688	1692	1696	rVB	144014	210627	1.56%	0.327%
13	12.215	1701	1705	1708	rBV2	105179	181033	1.34%	0.281%
14	12.258	1708	1712	1717	rVB3	144111	273709	2.03%	0.424%
15	12.343	1722	1726	1729	rVV2	84757	157524	1.17%	0.244%
16	12.398	1731	1735	1737	rVV4	91544	183715	1.36%	0.285%
17	12.428	1737	1740	1742	rVV	187206	273757	2.03%	0.424%
18	12.453	1742	1744	1748	rVV	180185	212457	1.58%	0.329%
19	12.495	1748	1751	1755	rVV	176599	266083	1.97%	0.413%
20	12.538	1755	1758	1762	rVB	148900	194560	1.44%	0.302%
21	12.684	1774	1782	1786	rVB	145014	318350	2.36%	0.494%
22	12.751	1786	1793	1796	rBV	695709	1178333	8.75%	1.827%
23	12.782	1796	1798	1800	rVV	362204	451681	3.35%	0.700%
24	12.818	1800	1804	1810	rVV2	1412511	2263417	16.80%	3.509%
25	12.873	1810	1813	1818	rVB2	148039	216051	1.60%	0.335%
26	12.983	1818	1831	1836	rBV	410642	919296	6.82%	1.425%
27	13.056	1839	1843	1849	rVB2	307908	489176	3.63%	0.758%
28	13.135	1849	1856	1861	rBV	851098	1356504	10.07%	2.103%
29	13.263	1869	1877	1881	rVB4	135611	292362	2.17%	0.453%
30	13.331	1881	1888	1892	rBV4	330254	641549	4.76%	0.995%
31	13.379	1892	1896	1899	rVV2	199064	302031	2.24%	0.468%
32	13.440	1903	1906	1908	rVV2	338499	494471	3.67%	0.767%
33	13.471	1908	1911	1915	rVV	433029	616129	4.57%	0.955%
34	13.513	1915	1918	1922	rVB2	145201	172002	1.28%	0.267%
35	13.635	1930	1938	1942	rBV2	713997	1220083	9.06%	1.892%
36	13.678	1942	1945	1954	rVB3	405541	852218	6.32%	1.321%

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Peak Location: TOP

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37	13.763	1954	1959	1964	rBV	1087503	1604743	11.91%	2.488%
38	13.837	1968	1971	1977	rVB	163385	231649	1.72%	0.359%
39	13.898	1977	1981	1983	rBV	174805	228839	1.70%	0.355%
40	13.928	1983	1986	1991	rVB2	278964	377908	2.80%	0.586%
41	13.983	1991	1995	2000	rVB	378033	518277	3.85%	0.804%
42	14.032	2000	2003	2008	rVB2	114021	206943	1.54%	0.321%
43	14.093	2008	2013	2018	rVB	492885	759754	5.64%	1.178%
44	14.154	2018	2023	2029	rBV6	92752	231796	1.72%	0.359%
45	14.269	2029	2042	2047	rBV8	314225	1040821	7.72%	1.614%
46	14.349	2052	2055	2059	rVV	250315	334038	2.48%	0.518%
47	14.391	2059	2062	2066	rVV	254730	323207	2.40%	0.501%
48	14.434	2066	2069	2074	rVB2	199869	257336	1.91%	0.399%
49	14.532	2082	2085	2089	rVV	179564	244794	1.82%	0.380%
50	14.580	2089	2093	2096	rVV	382707	641883	4.76%	0.995%
51	14.623	2096	2100	2105	rVB	779191	1130329	8.39%	1.752%
52	14.696	2105	2112	2117	rBV3	761529	1593871	11.83%	2.471%
53	14.745	2117	2120	2127	rVB4	305257	521767	3.87%	0.809%
54	14.849	2132	2137	2145	rVB	473989	766483	5.69%	1.188%
55	14.958	2150	2155	2158	rVV3	265623	453192	3.36%	0.703%
56	14.989	2158	2160	2166	rVB	129076	169912	1.26%	0.263%
57	15.068	2166	2173	2178	rBV4	326400	667267	4.95%	1.035%
58	15.147	2183	2186	2192	rVB6	131092	231080	1.72%	0.358%
59	15.251	2193	2203	2209	rBV	2219120	4031242	29.92%	6.250%
60	15.306	2209	2212	2216	rVB	179876	235088	1.74%	0.364%
61	15.361	2216	2221	2226	rBV3	455978	825890	6.13%	1.280%
62	15.409	2226	2229	2232	rVB2	123875	157180	1.17%	0.244%
63	15.446	2232	2235	2244	rVB3	418446	718662	5.33%	1.114%
64	15.544	2244	2251	2258	rBV	866020	1702173	12.63%	2.639%
65	15.617	2260	2263	2267	rVB2	122844	171526	1.27%	0.266%
66	15.708	2267	2278	2281	rBV2	542702	1475085	10.95%	2.287%
67	15.830	2290	2298	2304	rBV	876771	1679597	12.47%	2.604%
68	15.909	2304	2311	2326	rVB2	491388	1146207	8.51%	1.777%
69	16.050	2327	2334	2339	rBV3	202437	422446	3.14%	0.655%
70	16.245	2357	2366	2370	rBV3	145055	356261	2.64%	0.552%
71	16.312	2370	2377	2385	rVB	7198441	13473813	100.00%	20.889%
72	16.507	2401	2409	2423	rVB	3287338	6855295	50.88%	10.628%

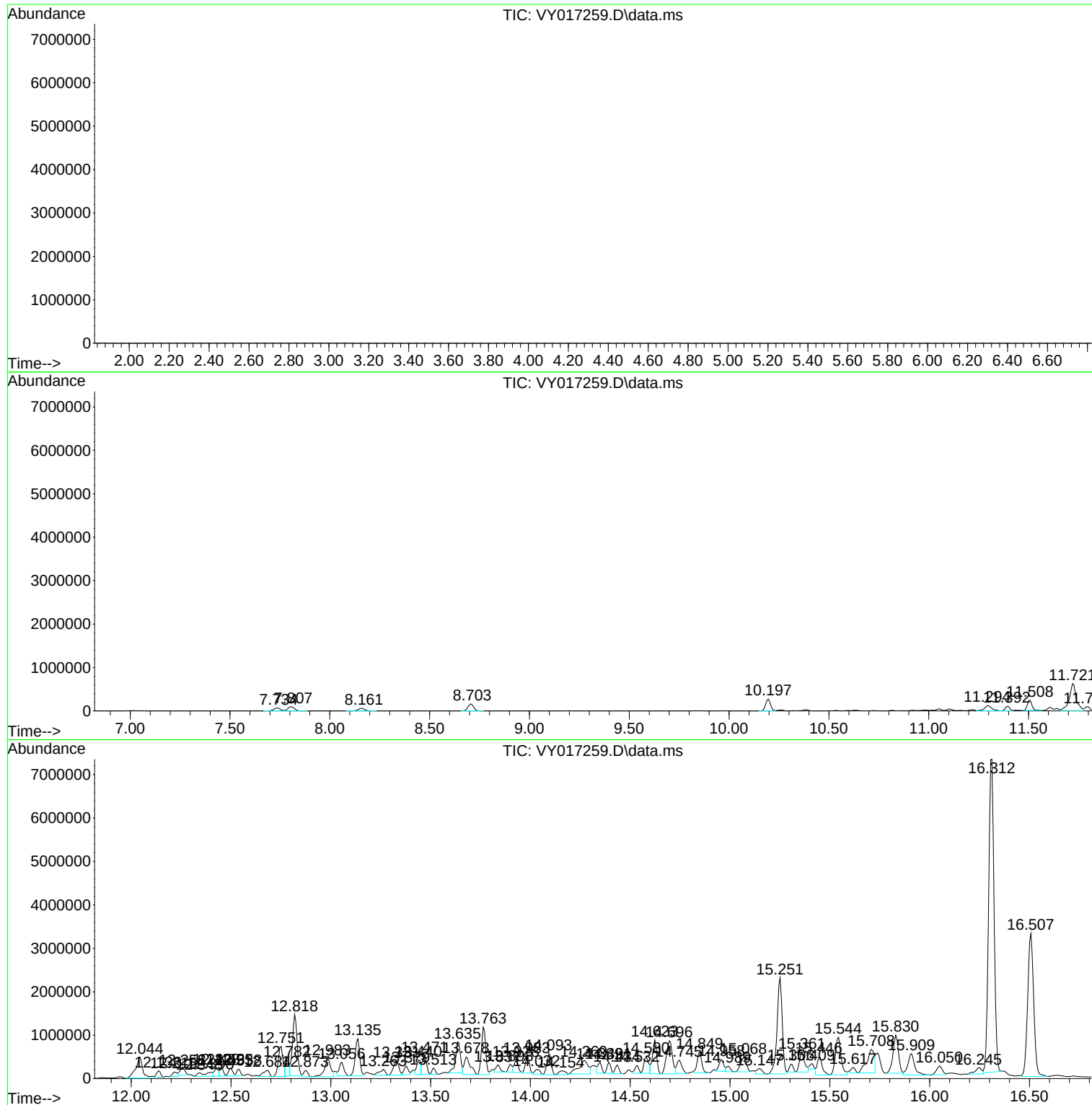
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ClientSampleId :
HR-04-020124

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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



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ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

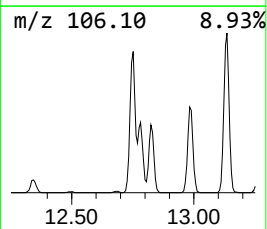
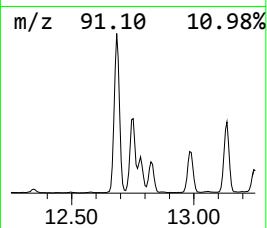
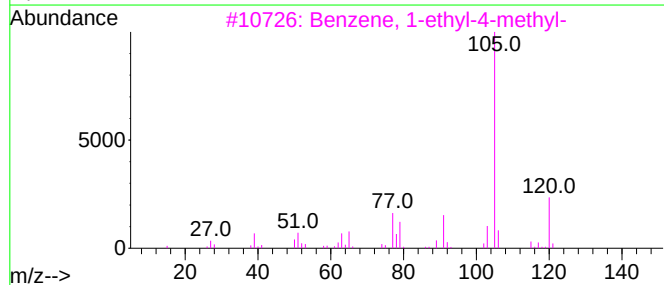
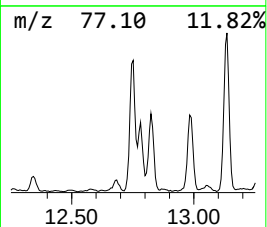
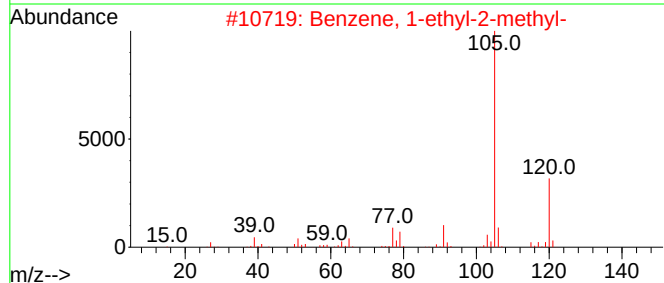
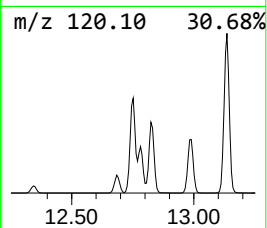
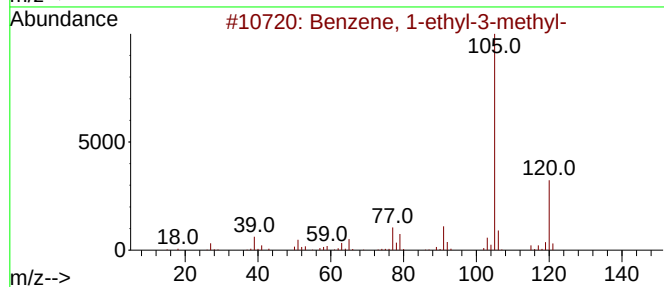
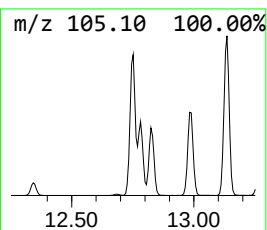
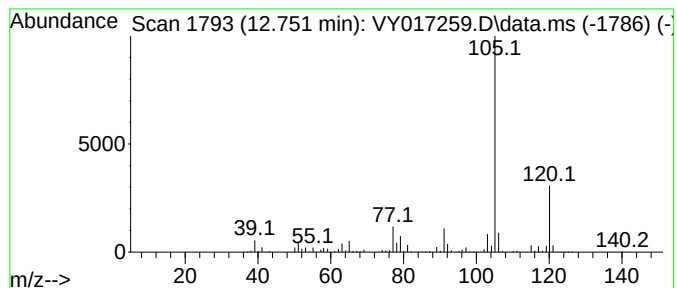
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.751	119.15 ug/l	1178330	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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MSVOA_Y
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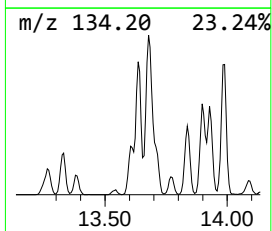
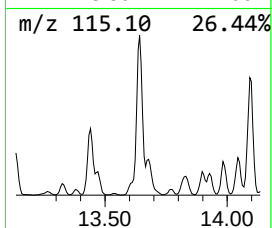
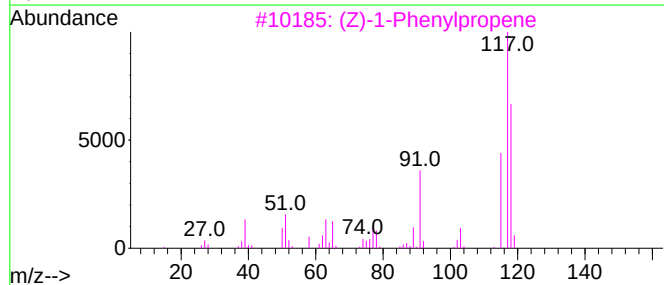
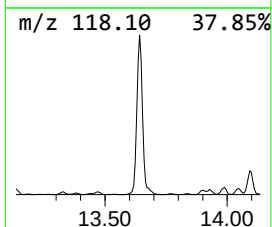
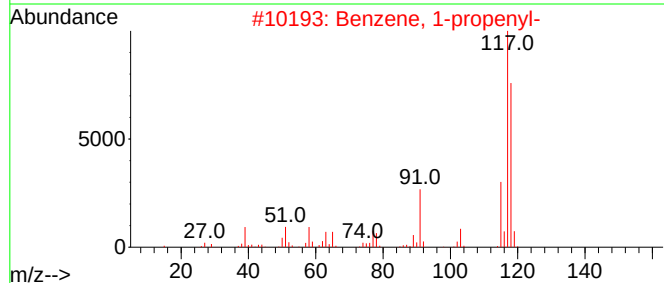
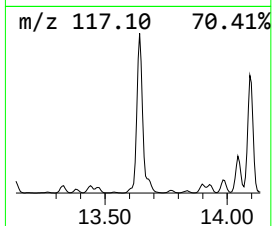
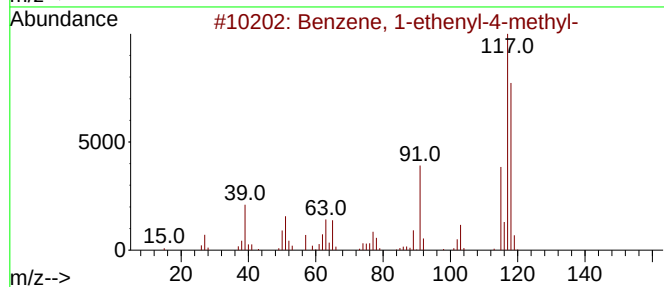
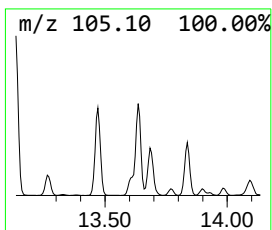
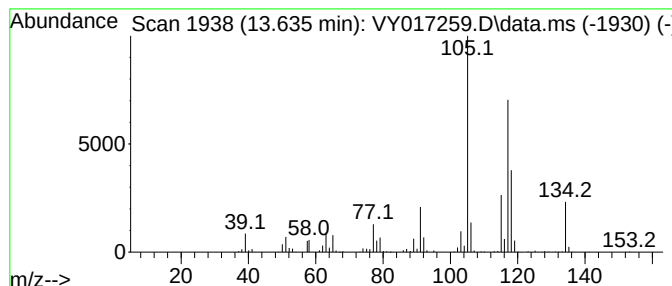
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethenyl-4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.635	123.37 ug/l	1220080	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	50
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	50
3			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	50
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	50
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	50



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MSVOA_Y
ClientSampleId :
HR-04-020124

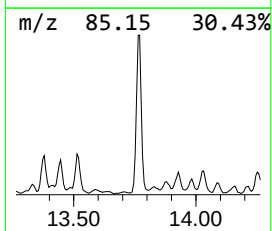
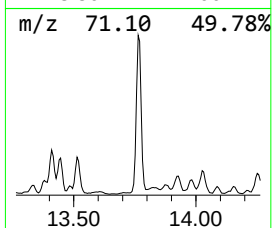
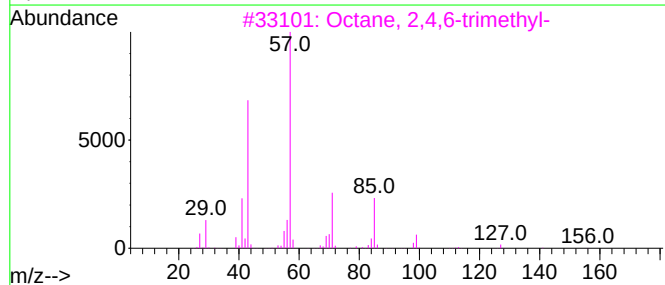
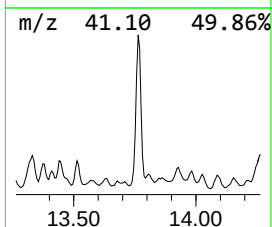
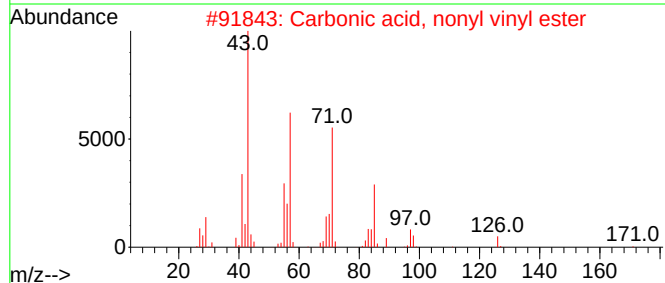
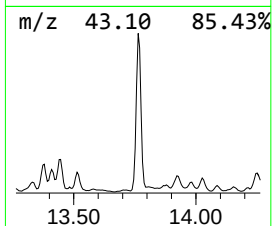
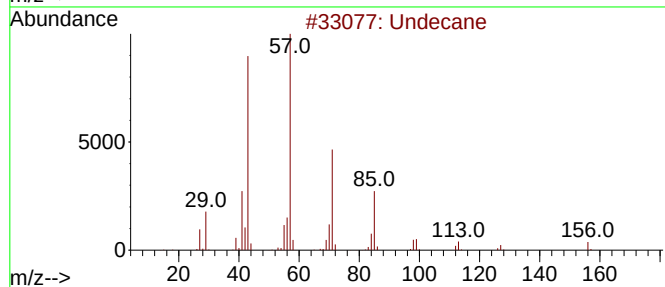
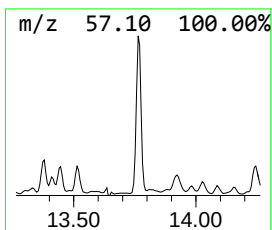
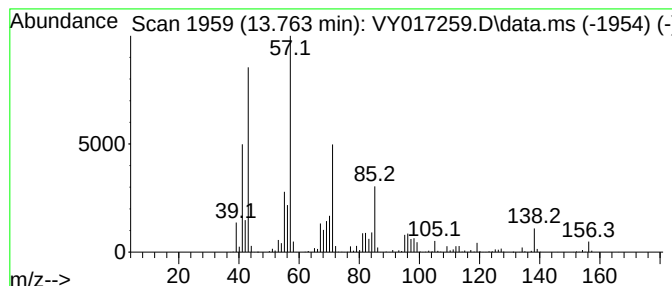
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Quant Title : SW846 8260

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

Peak Number 3 Undecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.763	162.27 ug/l	1604740	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	81
2			Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	64
3			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	64
4			Tridecane	184	C13H28	000629-50-5	59
5			Nonane, 5-(1-methylpropyl)-	184	C13H28	062185-54-0	53



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MSVOA_Y
ClientSampleId :
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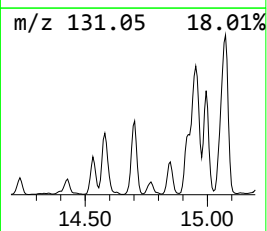
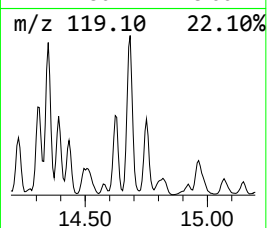
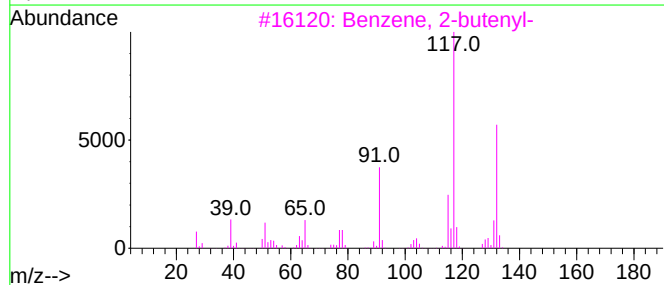
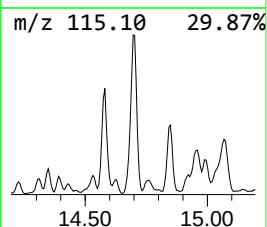
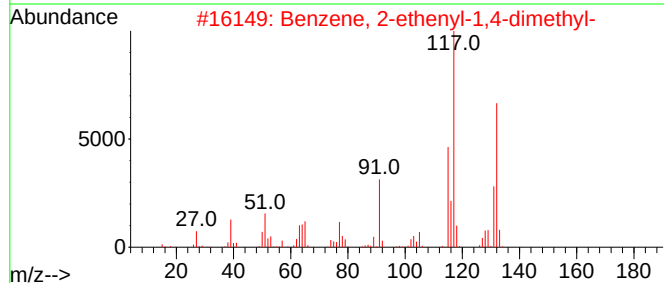
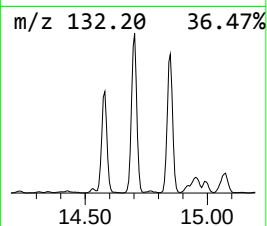
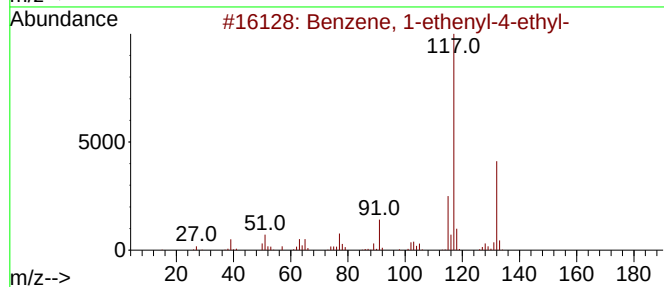
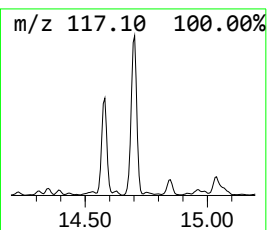
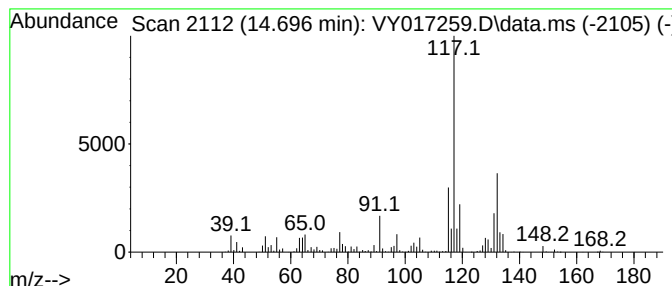
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.M
Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.696	161.17 ug/l	1593870	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
Data File : VY017259.D
Acq On : 06 Feb 2024 13:05
Operator : SY/MD
Sample : P1334-07
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

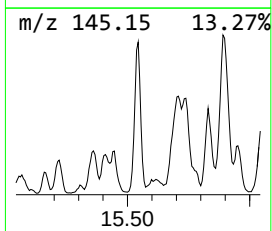
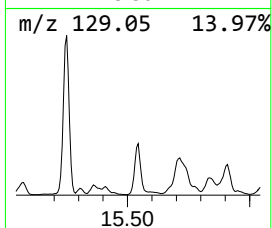
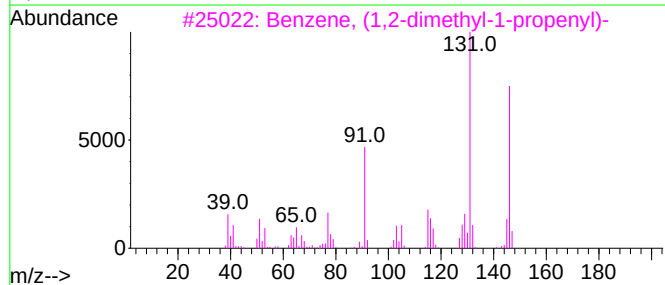
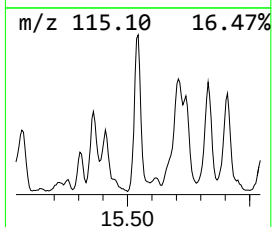
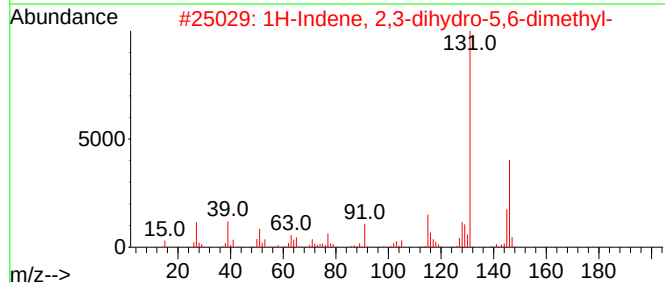
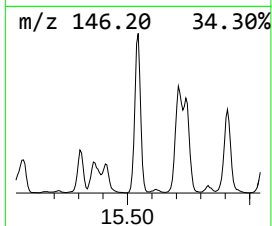
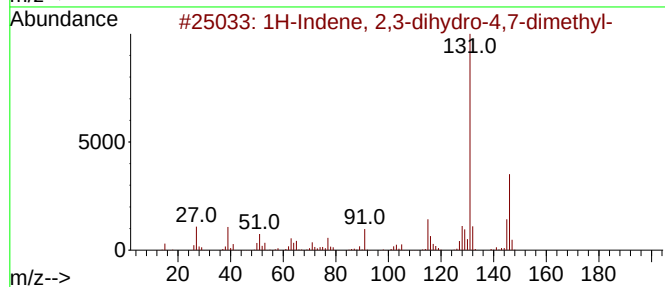
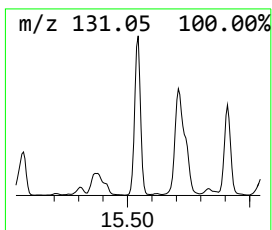
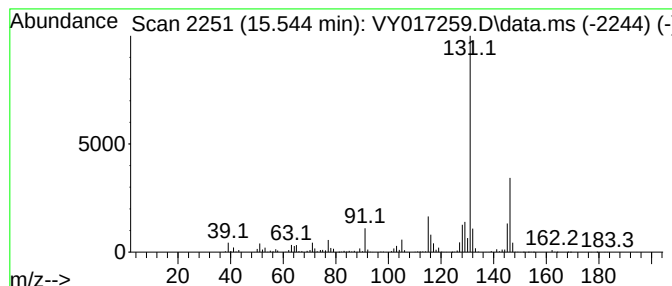
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.544	172.12 ug/l	1702170	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	97
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	94
3			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
Data File : VY017259.D
Acq On : 06 Feb 2024 13:05
Operator : SY/MD
Sample : P1334-07
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

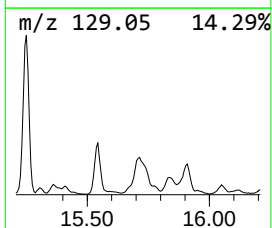
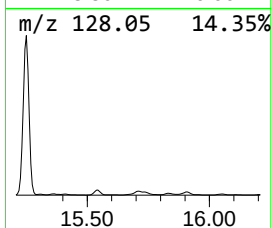
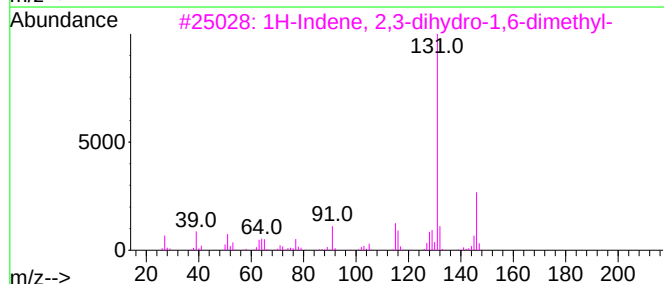
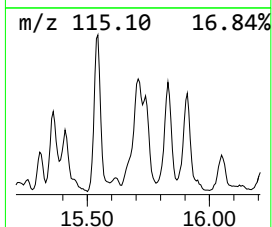
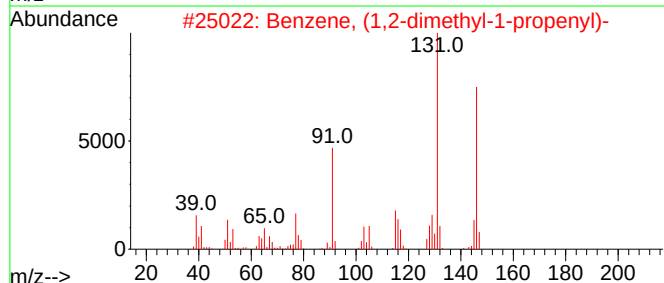
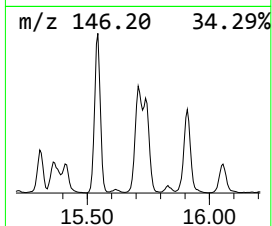
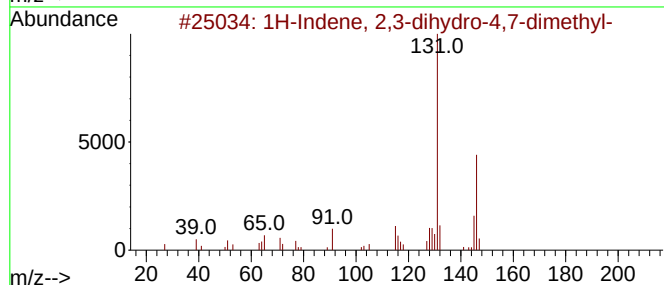
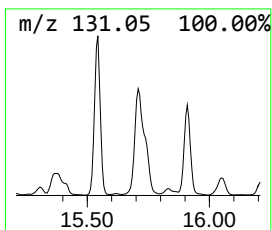
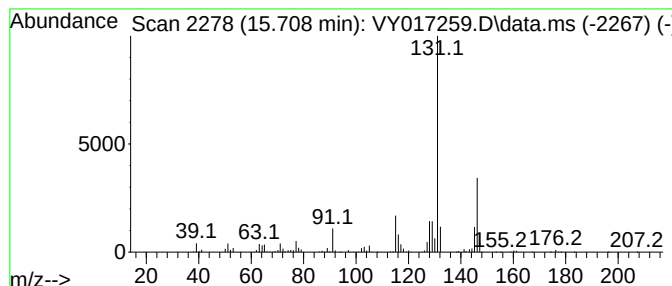
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.708	149.16 ug/l	1475090	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	96		
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94		
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94		
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91		
5	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	91		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
Data File : VY017259.D
Acq On : 06 Feb 2024 13:05
Operator : SY/MD
Sample : P1334-07
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

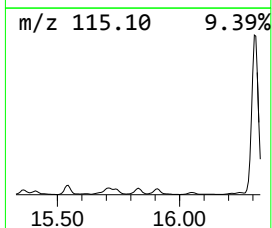
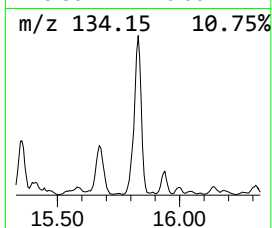
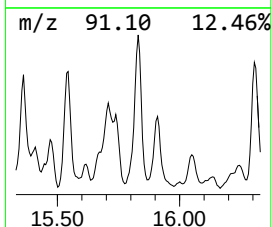
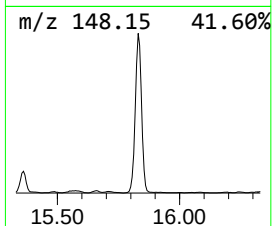
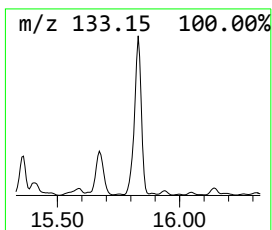
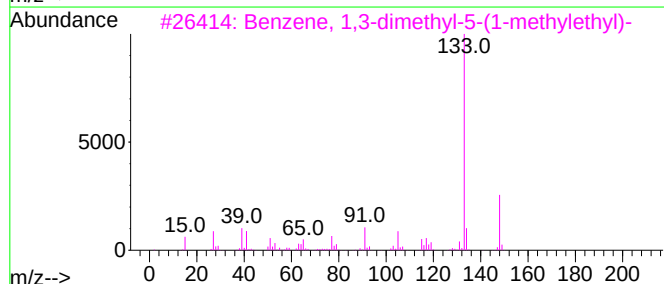
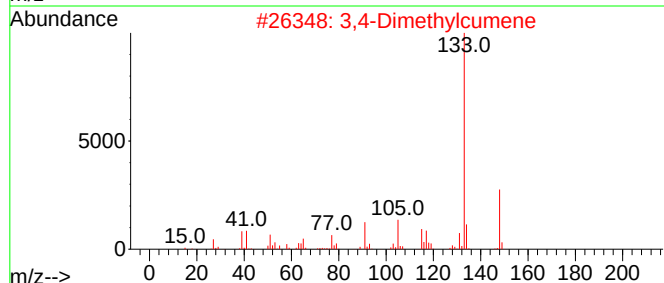
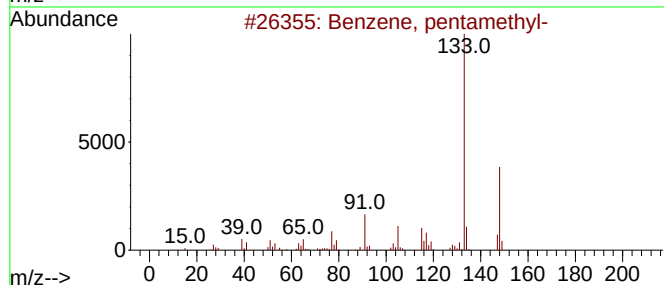
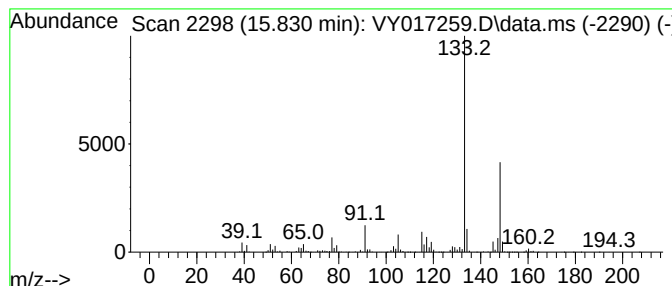
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, pentamethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.830	169.84 ug/l	1679600	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	95
2			3,4-Dimethylcumene	148	C11H16	1000370-34-1	90
3			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	87
4			Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	87
5			1,5,6,7-Tetramethylbicyclo[3.2.0...	148	C11H16	134329-46-7	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
Data File : VY017259.D
Acq On : 06 Feb 2024 13:05
Operator : SY/MD
Sample : P1334-07
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

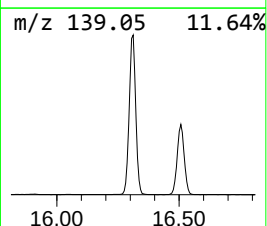
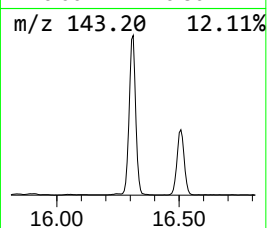
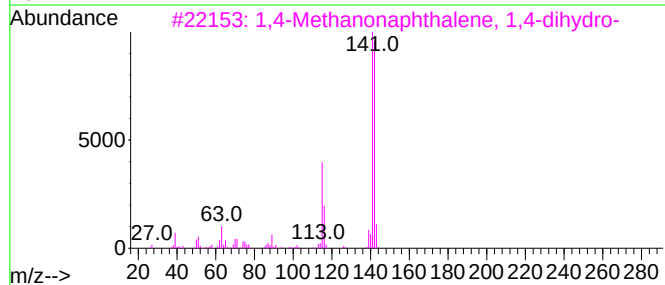
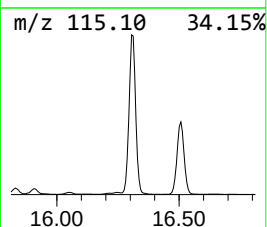
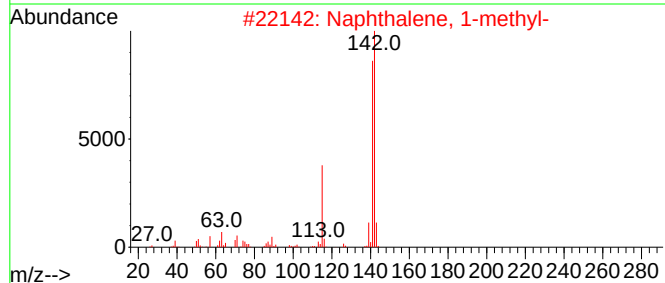
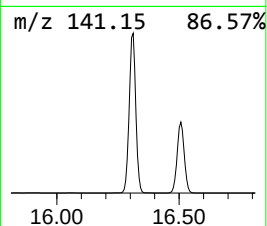
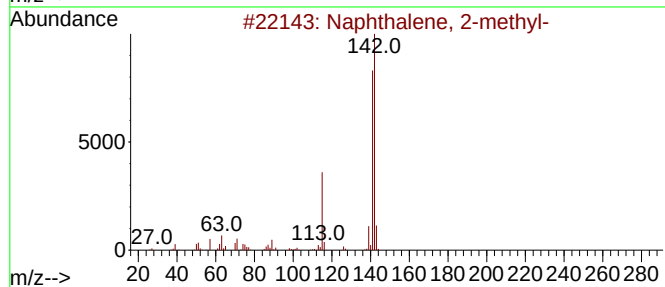
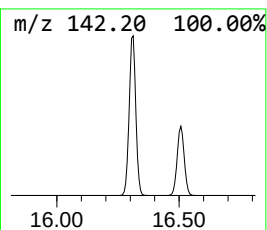
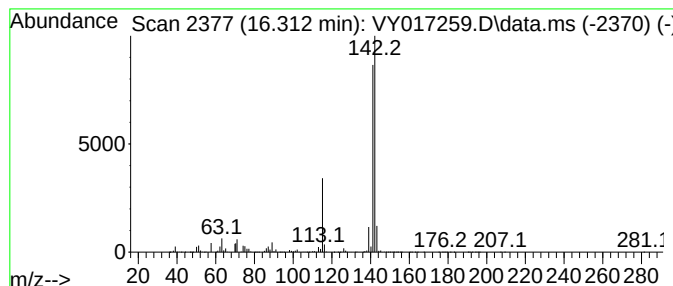
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.312	1362.45 ug/l	13473800	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
Data File : VY017259.D
Acq On : 06 Feb 2024 13:05
Operator : SY/MD
Sample : P1334-07
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

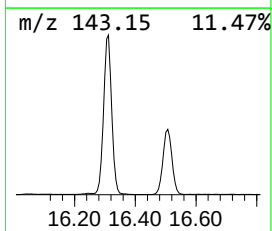
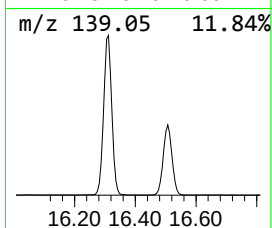
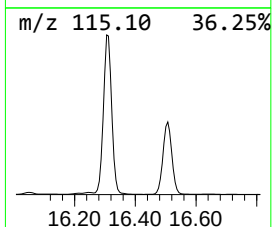
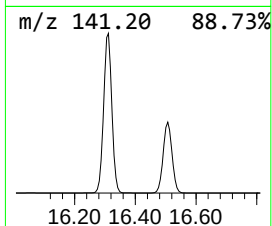
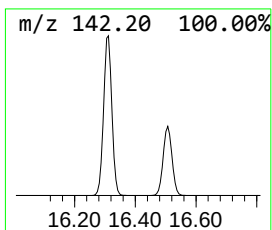
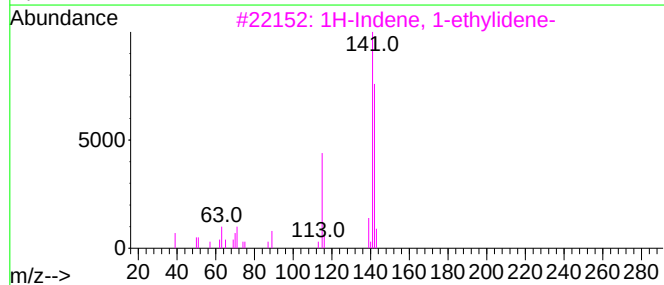
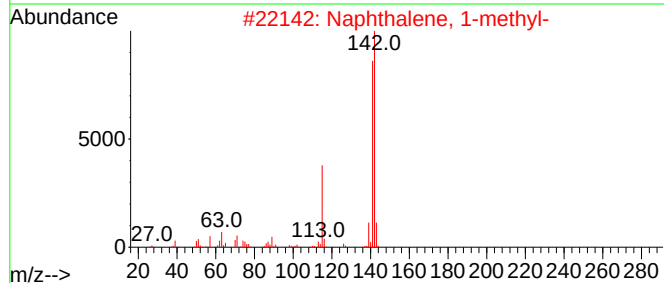
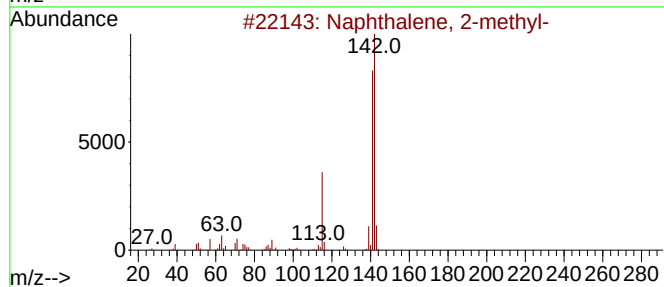
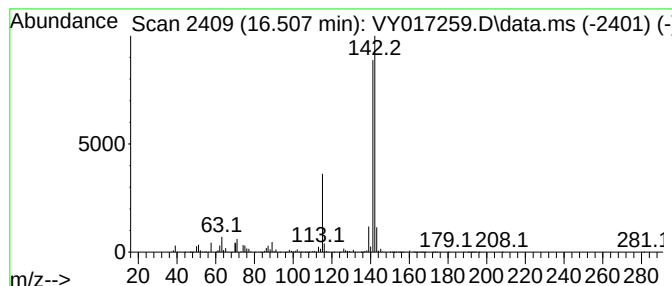
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 1-ethylidene- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.507	693.20 ug/l	6855300	1,4-Dichlorobenzene-d4	13.440

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	94
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY020624\
Data File : VY017259.D
Acq On : 06 Feb 2024 13:05
Operator : SY/MD
Sample : P1334-07
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
HR-04-020124

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y020124S.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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Benzene, 1-ethe...	13.635	123.4	ug/l	1220080	4	13.440	494471	50.0
Undecane	13.763	162.3	ug/l	1604740	4	13.440	494471	50.0
Benzene, 1-ethe...	14.696	161.2	ug/l	1593870	4	13.440	494471	50.0
1H-Indene, 2,3-...	15.544	172.1	ug/l	1702170	4	13.440	494471	50.0
Benzene, (1,2-d...	15.708	149.2	ug/l	1475090	4	13.440	494471	50.0
Benzene, pentam...	15.830	169.8	ug/l	1679600	4	13.440	494471	50.0
Naphthalene, 1-...	16.312	1362.5	ug/l	13473800	4	13.440	494471	50.0
1H-Indene, 1-et...	16.507	693.2	ug/l	6855300	4	13.440	494471	50.0