

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
 Data File : VY020033.D
 Acq On : 28 Oct 2024 11:56
 Operator : SY/MD
 Sample : P4578-03
 Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 WB-305-TOP

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\82Y100924S.M
 Title : SW846 8260

Signal : TIC: VY020033.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.080	58	59	63	rBV4	2426	3689	1.17%	0.090%
2	2.476	118	124	126	rBV4	2502	3344	1.06%	0.081%
3	2.586	135	142	143	rBV4	2079	4719	1.50%	0.115%
4	2.647	150	152	155	rVB4	3399	3152	1.00%	0.077%
5	4.110	388	392	400	rVB4	1403	3393	1.08%	0.083%
6	4.397	437	439	449	rVB3	3557	7744	2.46%	0.189%
7	4.628	471	477	481	rBV6	2221	4513	1.43%	0.110%
8	7.640	960	971	976	rBV2	19704	46816	14.88%	1.140%
9	7.713	977	983	995	rVB2	40678	96029	30.53%	2.339%
10	8.073	1032	1042	1051	rBV2	17535	40460	12.86%	0.986%
11	8.622	1125	1132	1143	rBV2	59337	119919	38.12%	2.921%
12	9.122	1210	1214	1218	rVB4	2928	5152	1.64%	0.125%
13	10.109	1369	1376	1383	rBV	84969	154882	49.23%	3.773%
14	10.176	1383	1387	1396	rVB	25834	45967	14.61%	1.120%
15	10.451	1428	1432	1437	rVB6	2382	4241	1.35%	0.103%
16	10.554	1443	1449	1453	rVB8	1768	4169	1.33%	0.102%
17	10.731	1474	1478	1484	rBV5	4217	6370	2.02%	0.155%
18	10.841	1493	1496	1500	rVB5	2875	4189	1.33%	0.102%
19	10.969	1513	1517	1521	rVB5	3371	5360	1.70%	0.131%
20	11.024	1521	1526	1531	rVB4	6646	10948	3.48%	0.267%
21	11.079	1531	1535	1539	rVB4	2381	3297	1.05%	0.080%
22	11.133	1539	1544	1549	rBV5	4210	8140	2.59%	0.198%
23	11.219	1551	1558	1564	rVB8	5719	14856	4.72%	0.362%
24	11.310	1567	1573	1577	rBV3	5237	8371	2.66%	0.204%
25	11.420	1584	1591	1601	rBV	68108	115897	36.84%	2.823%
26	11.639	1617	1627	1630	rBV7	9496	25693	8.17%	0.626%
27	11.713	1636	1639	1645	rVB6	3966	4936	1.57%	0.120%
28	11.932	1669	1675	1676	rBV5	7255	9984	3.17%	0.243%
29	12.054	1691	1695	1703	rVB7	4830	6675	2.12%	0.163%
30	12.133	1703	1708	1710	rBV4	4966	7585	2.41%	0.185%
31	12.170	1710	1714	1721	rVV6	9047	15750	5.01%	0.384%
32	12.255	1725	1728	1733	rVV4	6556	9846	3.13%	0.240%
33	12.310	1733	1737	1741	rVV7	3398	6941	2.21%	0.169%
34	12.347	1741	1743	1748	rVB3	2986	3756	1.19%	0.091%
35	12.414	1748	1754	1764	rBV3	53185	97177	30.89%	2.367%
36	12.505	1764	1769	1771	rBV6	2401	3188	1.01%	0.078%

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37	12.572	1777	1780	1788	rVB5	5643	12332	3.92%	0.300%
38	12.664	1788	1795	1796	rBV7	6356	10561	3.36%	0.257%
39	12.694	1796	1800	1802	rVV5	5030	7383	2.35%	0.180%
40	12.737	1802	1807	1813	rVB7	15378	29757	9.46%	0.725%
41	12.901	1827	1834	1838	rBV2	11905	22428	7.13%	0.546%
42	12.968	1842	1845	1848	rVV5	5688	7576	2.41%	0.185%
43	13.048	1852	1858	1864	rBV	38498	57828	18.38%	1.409%
44	13.170	1871	1878	1883	rBV	4940	10825	3.44%	0.264%
45	13.243	1885	1890	1894	rVV6	10678	19461	6.19%	0.474%
46	13.292	1894	1898	1902	rVV6	7251	13867	4.41%	0.338%
47	13.353	1902	1908	1910	rVV2	37946	67478	21.45%	1.644%
48	13.383	1910	1913	1919	rVB2	43581	62775	19.95%	1.529%
49	13.548	1931	1940	1944	rBV2	70262	136438	43.37%	3.324%
50	13.590	1944	1947	1956	rVB5	78411	149479	47.52%	3.641%
51	13.676	1956	1961	1965	rBV5	19614	32904	10.46%	0.802%
52	13.749	1969	1973	1978	rVV2	24634	37272	11.85%	0.908%
53	13.810	1978	1983	1985	rVV	68174	99767	31.71%	2.430%
54	13.840	1985	1988	1992	rVV	83117	117987	37.51%	2.874%
55	13.895	1992	1997	2003	rVB2	208244	314584	100.00%	7.663%
56	13.950	2003	2006	2009	rVB5	4409	5331	1.69%	0.130%
57	14.005	2009	2015	2020	rBV3	70244	112183	35.66%	2.733%
58	14.078	2020	2027	2031	rBV7	12585	33495	10.65%	0.816%
59	14.139	2031	2037	2041	rBV	52165	88858	28.25%	2.164%
60	14.218	2041	2050	2053	rVV	128018	232976	74.06%	5.675%
61	14.255	2053	2056	2061	rVV	187833	281603	89.52%	6.860%
62	14.304	2061	2064	2068	rVV2	64928	94943	30.18%	2.313%
63	14.340	2068	2070	2077	rVV2	37152	55549	17.66%	1.353%
64	14.401	2078	2080	2083	rVB3	8267	9275	2.95%	0.226%
65	14.444	2083	2087	2090	rBV3	20228	27058	8.60%	0.659%
66	14.486	2090	2094	2098	rVV	89523	135130	42.96%	3.292%
67	14.535	2098	2102	2106	rVB	91636	123528	39.27%	3.009%
68	14.602	2106	2113	2118	rBV3	144366	305033	96.96%	7.430%
69	14.657	2118	2122	2128	rVB4	47176	84601	26.89%	2.061%
70	14.724	2128	2133	2135	rBV6	7114	12805	4.07%	0.312%
71	14.864	2151	2156	2163	rBV3	19217	36633	11.64%	0.892%
72	14.925	2164	2166	2167	rVV2	6434	4665	1.48%	0.114%
73	14.968	2167	2173	2177	rVV6	14446	31077	9.88%	0.757%
74	15.011	2177	2180	2184	rVV5	9951	18681	5.94%	0.455%
75	15.053	2184	2187	2193	rVB5	21101	37779	12.01%	0.920%
76	15.121	2193	2198	2206	rBV3	32630	78428	24.93%	1.910%

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77	15.346	2231	2235	2242	rVB	62033	91934	29.22%	2.239%
78	15.425	2244	2248	2249	rBV4	4581	6016	1.91%	0.147%
79	15.791	2304	2308	2309	rBV4	2840	3654	1.16%	0.089%
80	15.864	2317	2320	2325	rVB7	3065	4818	1.53%	0.117%
81	15.919	2325	2329	2331	rBV5	4733	7894	2.51%	0.192%
82	15.980	2337	2339	2346	rVB6	5794	9587	3.05%	0.234%
83	16.053	2348	2351	2357	rVB7	6801	12892	4.10%	0.314%
84	16.254	2379	2384	2390	rBV	12702	21258	6.76%	0.518%
85	16.364	2399	2402	2405	rBV4	3727	6563	2.09%	0.160%
86	16.492	2416	2423	2427	rBV9	1180	3152	1.00%	0.077%

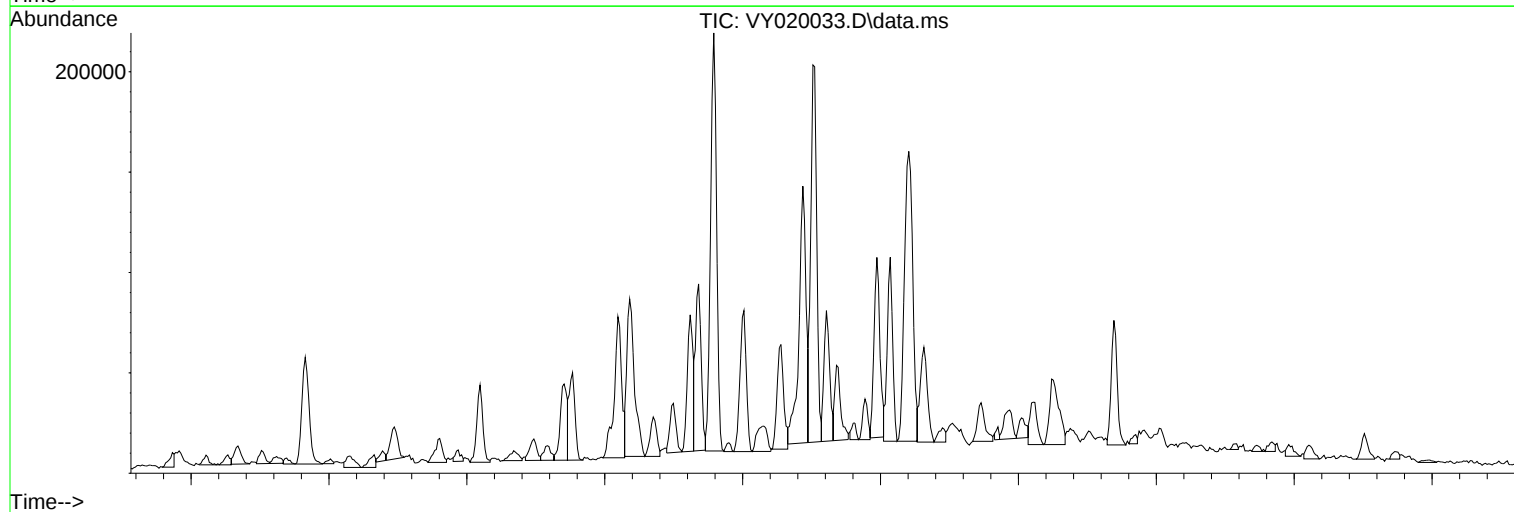
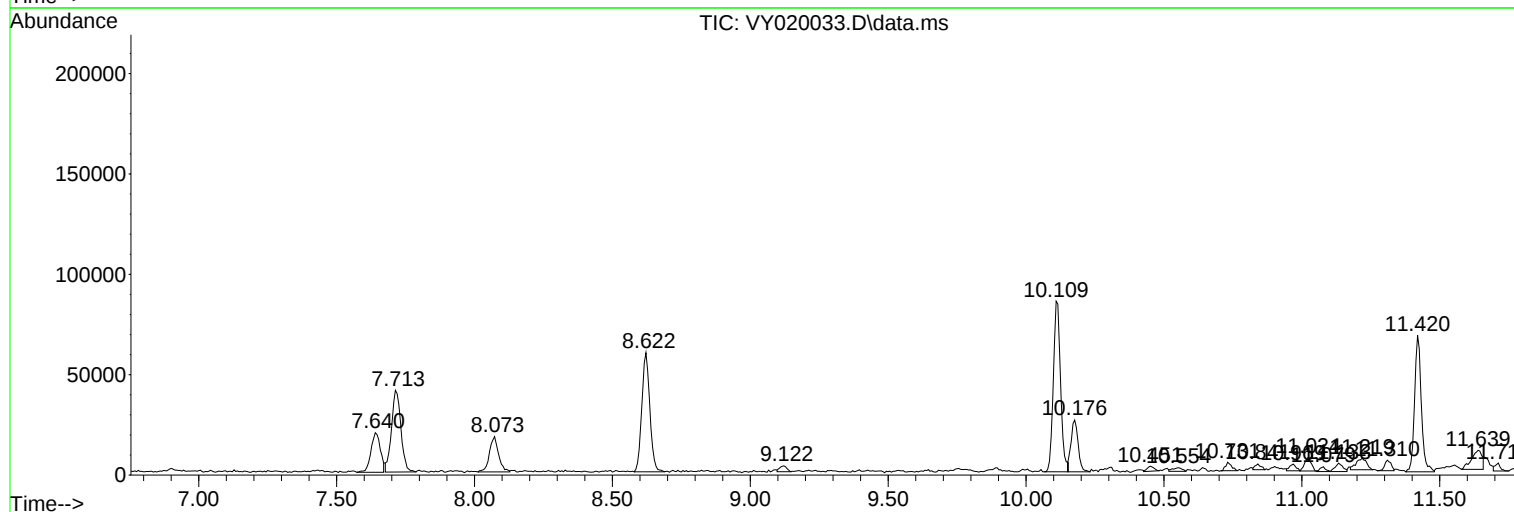
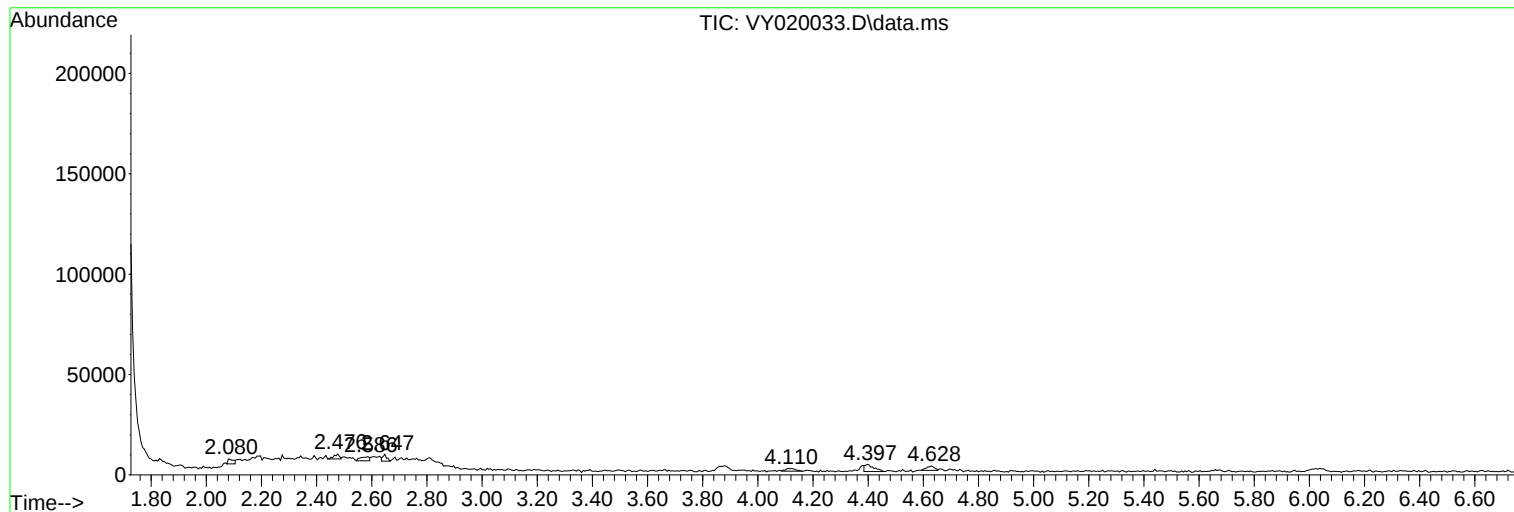
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

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



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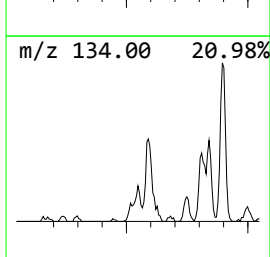
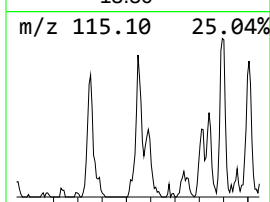
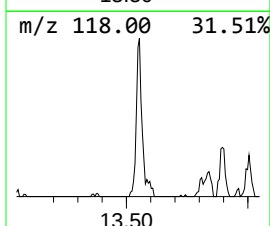
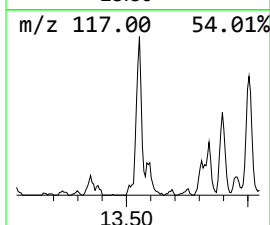
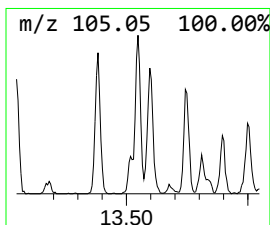
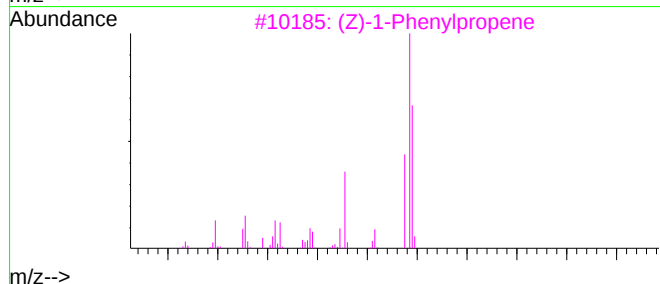
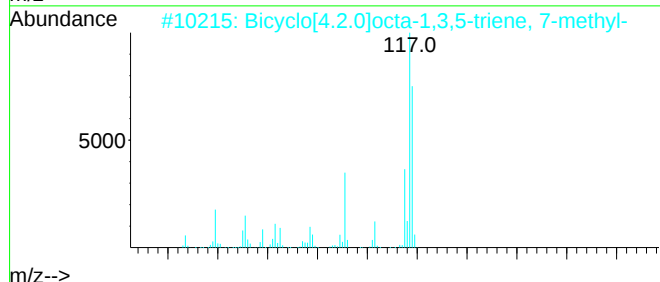
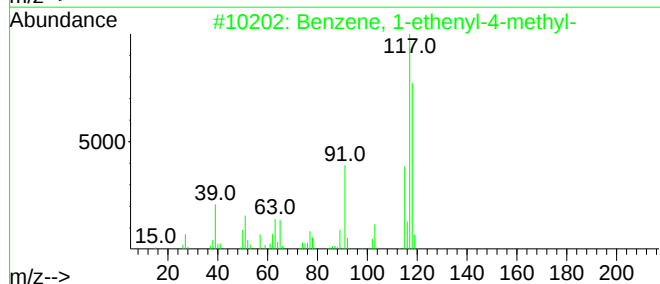
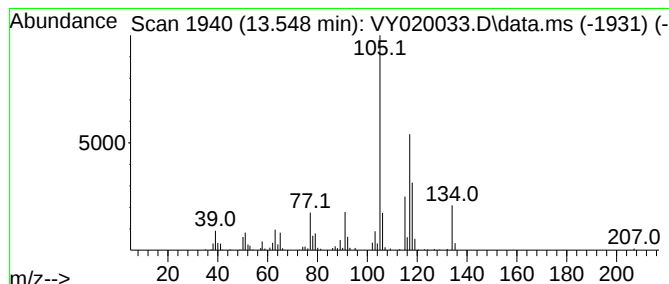
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, 1-ethenyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.548	101.10 ug/l	136438	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	86
2	Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	50
3	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	50
4	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	50
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	45



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

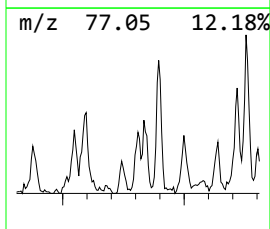
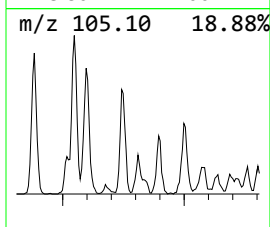
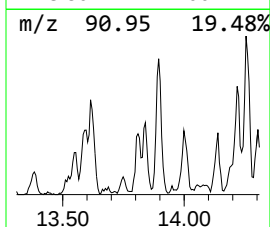
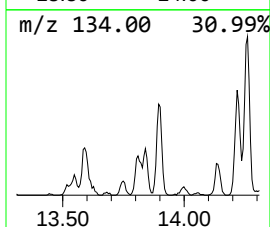
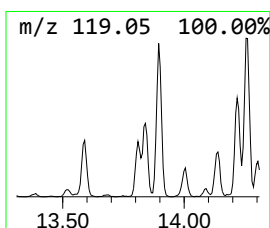
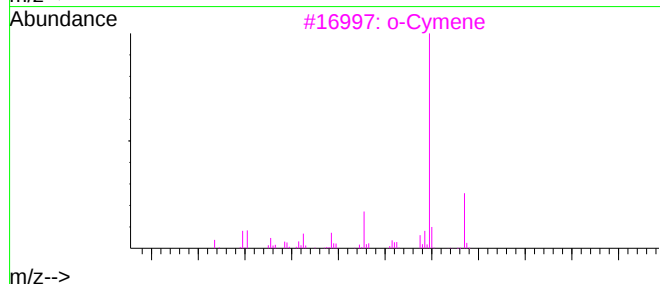
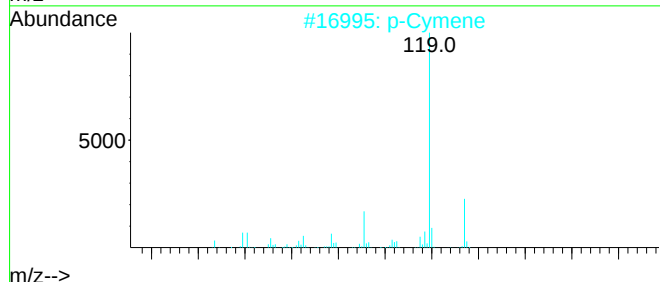
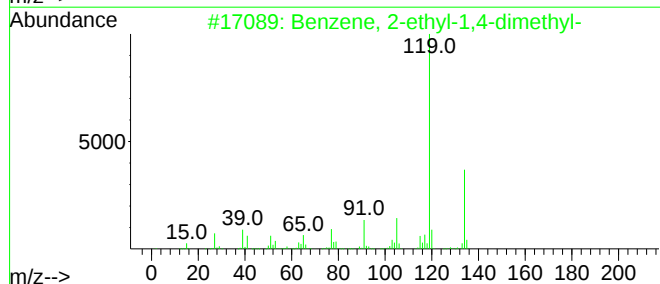
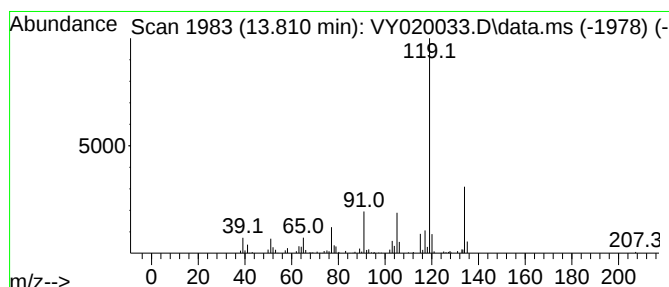
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.810	73.93 ug/l	99767	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	p-Cymene	134	C10H14	000099-87-6	94
3	o-Cymene	134	C10H14	000527-84-4	93
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



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Quant Title : SW846 8260

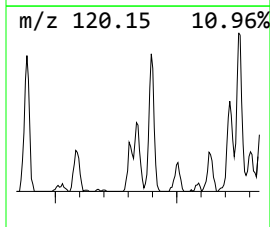
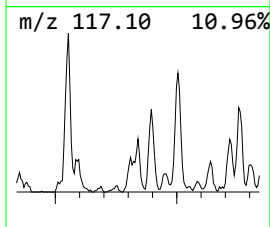
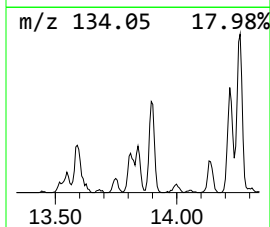
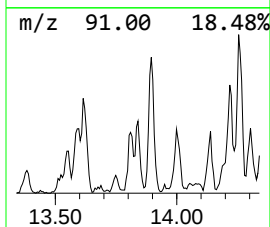
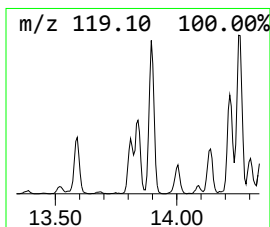
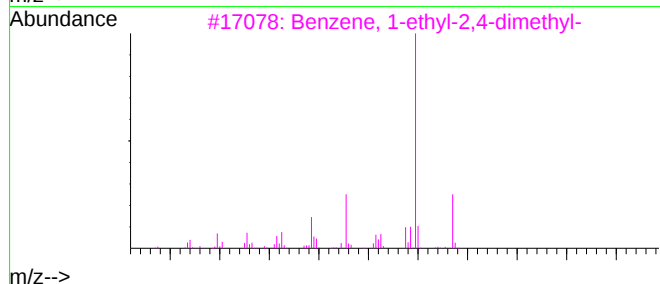
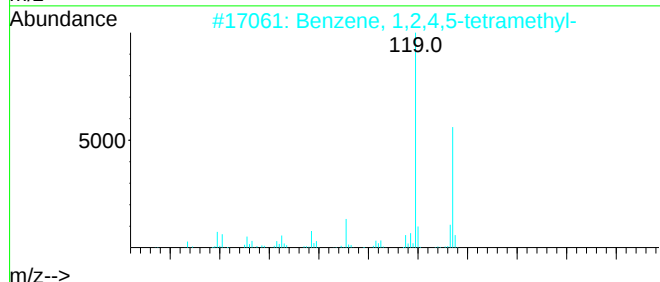
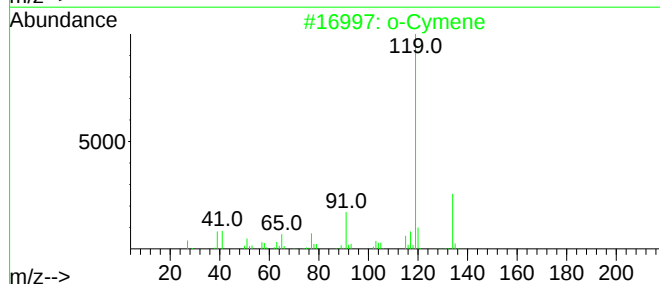
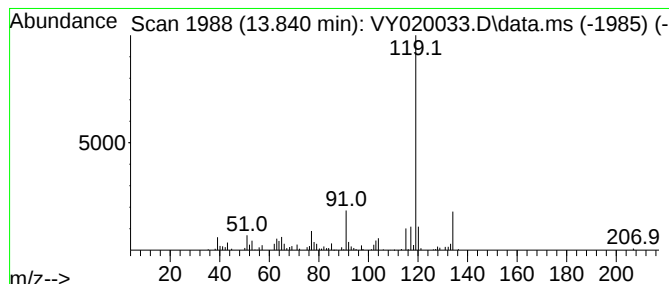
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.840	87.43 ug/l	117987	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	91
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91



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Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

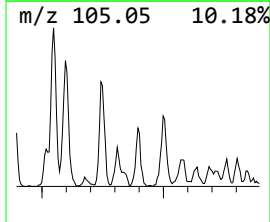
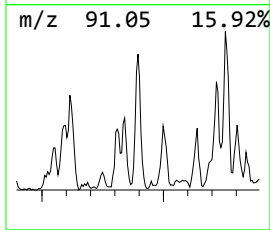
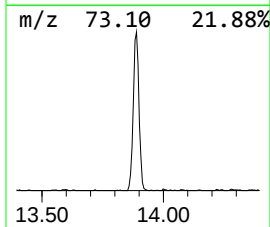
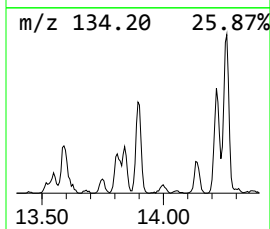
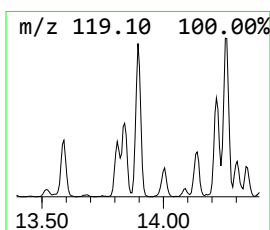
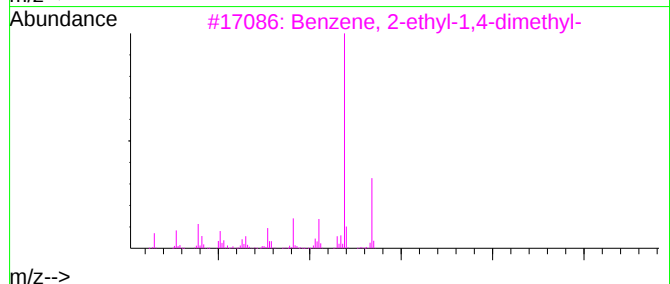
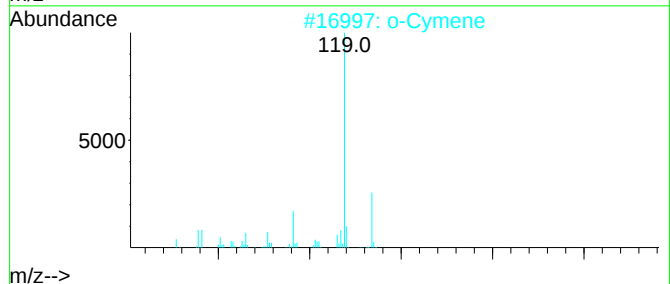
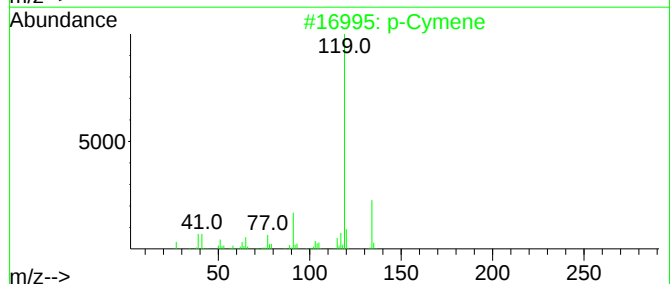
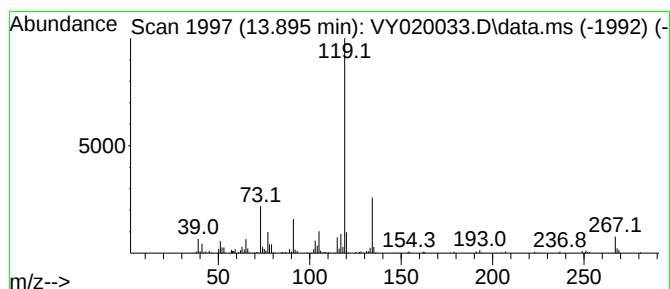
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.895	233.10 ug/l	314584	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
Acq On : 28 Oct 2024 11:56
Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

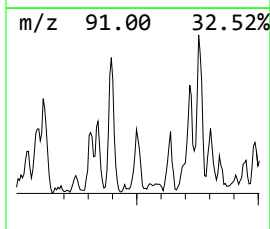
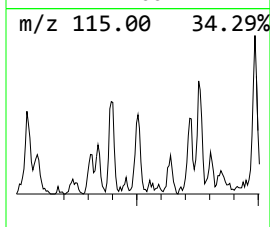
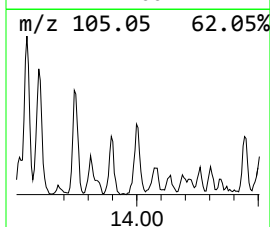
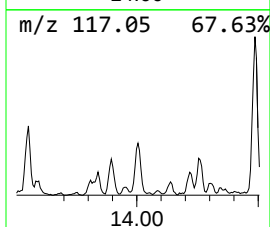
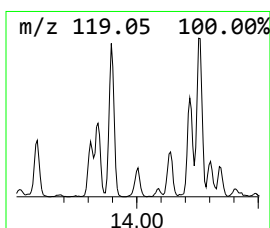
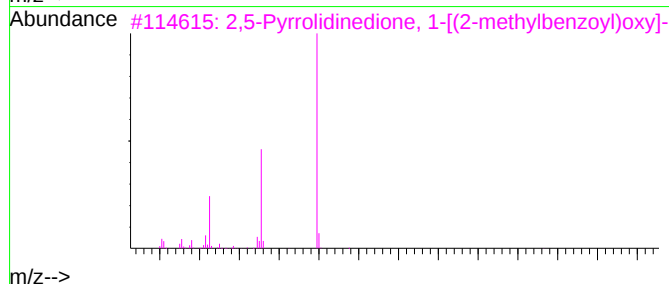
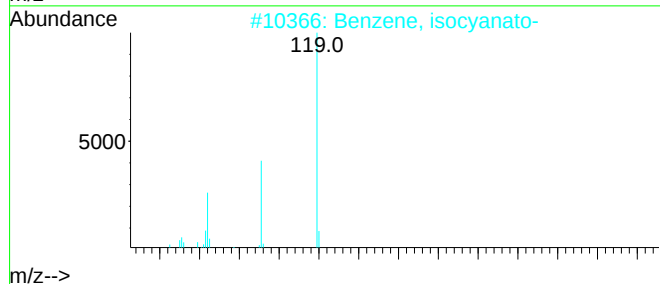
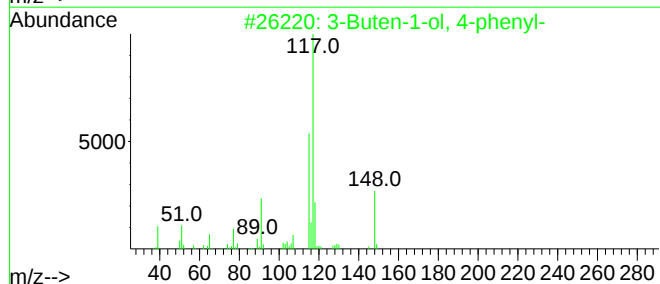
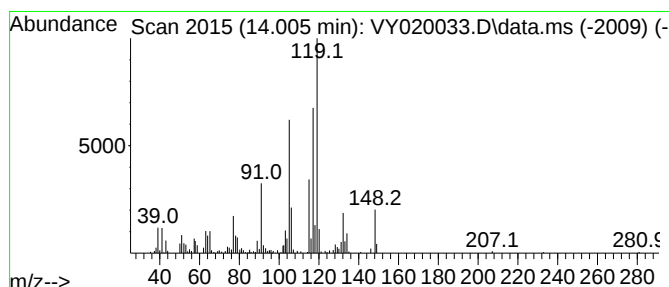
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Quant Title : SW846 8260

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

Peak Number 5 unknown14.005 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.005	83.13 ug/l	112183	1,4-Dichlorobenzene-d4	13.359

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Buten-1-ol, 4-phenyl-	148	C10H12O	000937-58-6	42
2			Benzene, isocyanato-	119	C7H5NO	000103-71-9	38
3			2,5-Pyrrolidinedione, 1-[(2-meth...	233	C12H11NO4	110920-14-4	38
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	35
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
Acq On : 28 Oct 2024 11:56
Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

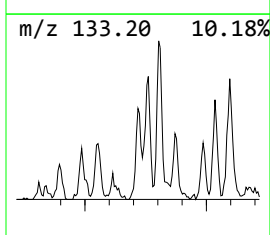
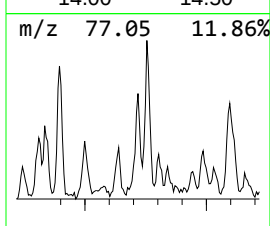
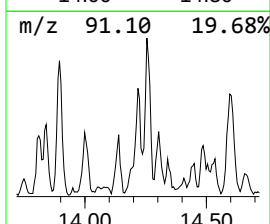
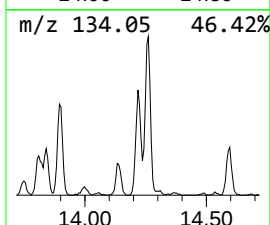
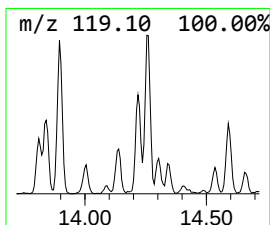
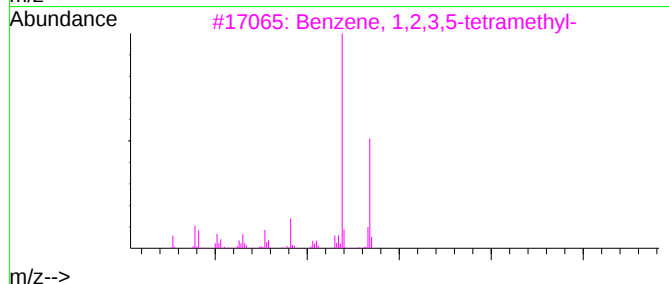
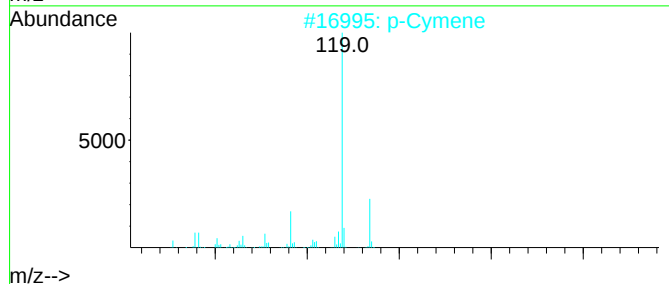
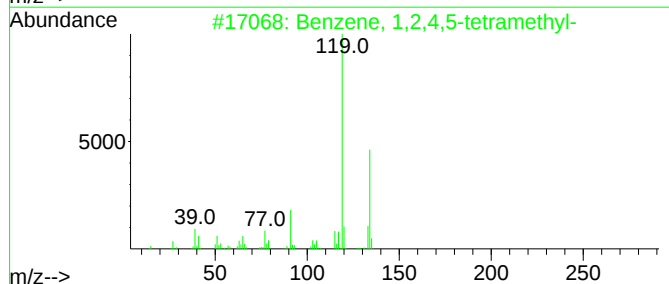
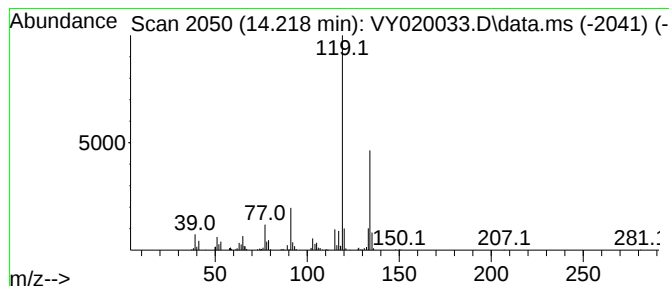
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.218	172.63 ug/l	232976	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	p-Cymene	134	C10H14	000099-87-6	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5	o-Cymene	134	C10H14	000527-84-4	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
Acq On : 28 Oct 2024 11:56
Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

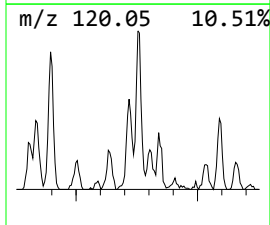
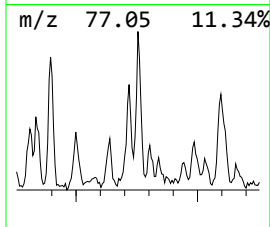
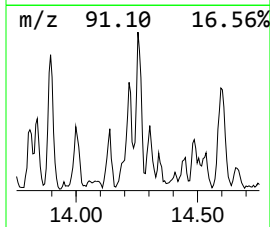
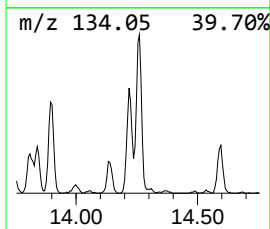
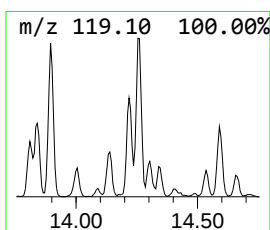
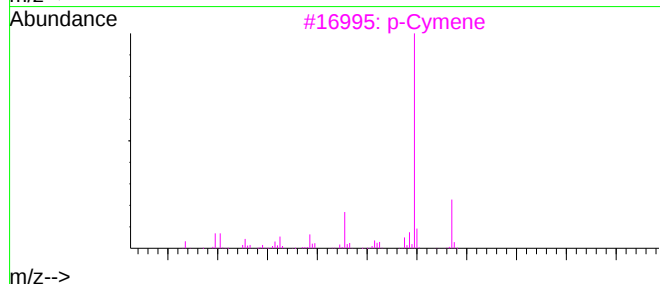
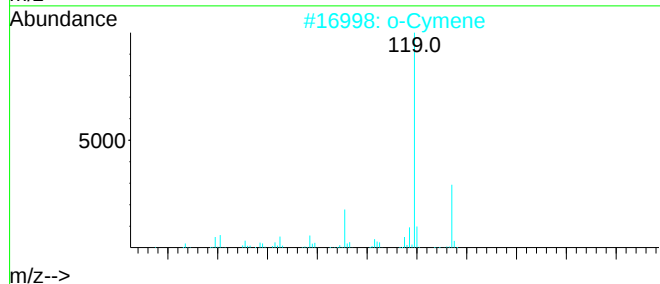
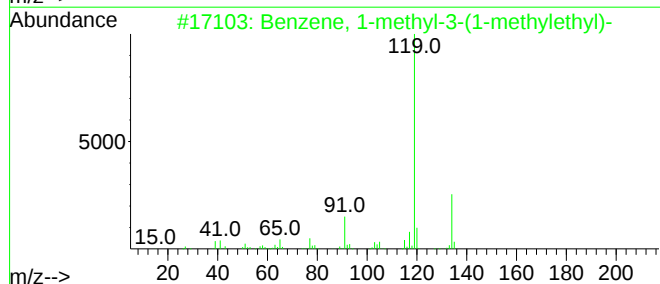
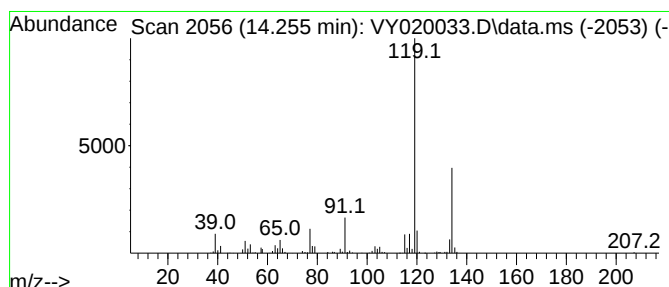
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1-methyl-3-(1-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.255	208.66 ug/l	281603	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
2	o-Cymene	134	C10H14	000527-84-4	91
3	p-Cymene	134	C10H14	000099-87-6	91
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
Acq On : 28 Oct 2024 11:56
Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.M
Quant Title : SW846 8260

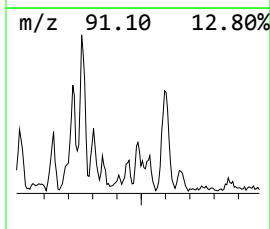
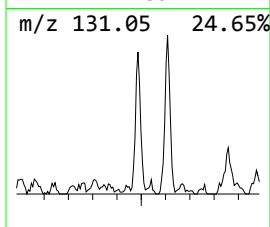
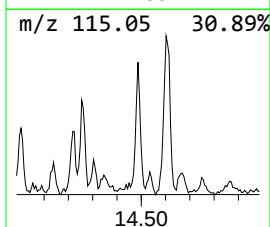
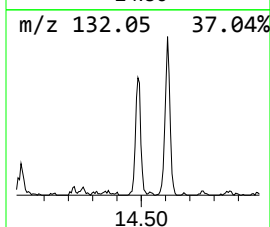
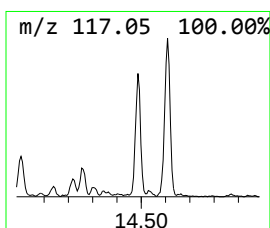
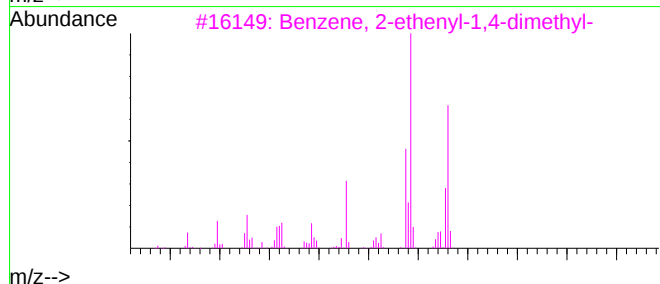
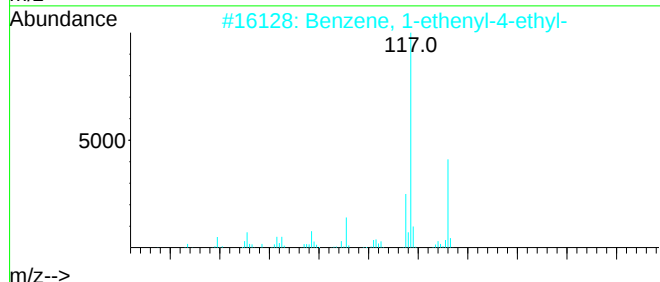
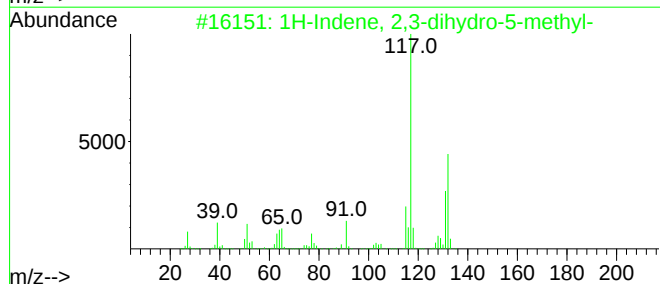
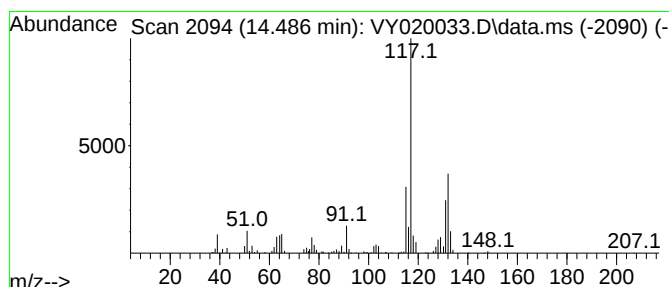
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.486	100.13 ug/l	135130	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
Acq On : 28 Oct 2024 11:56
Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

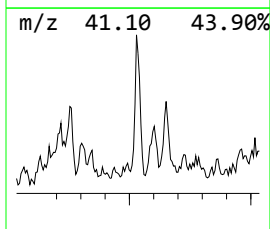
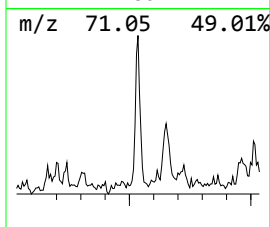
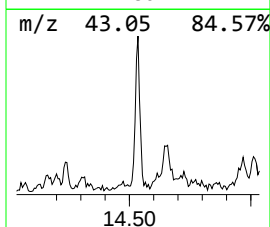
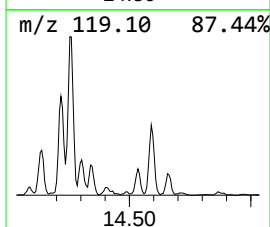
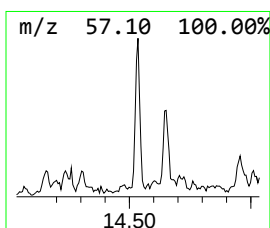
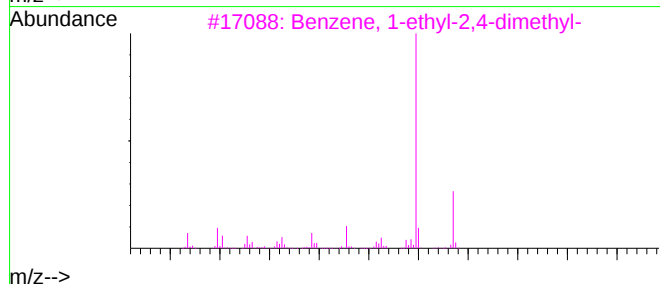
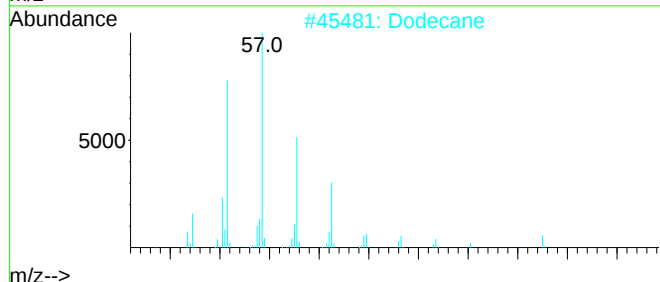
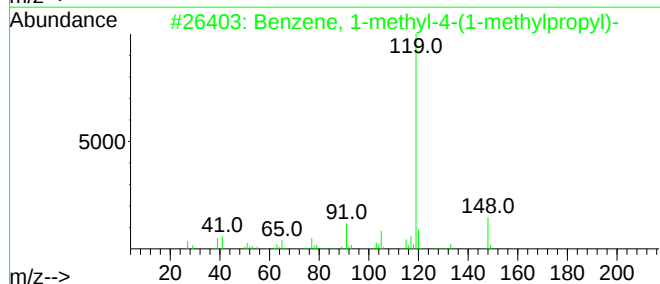
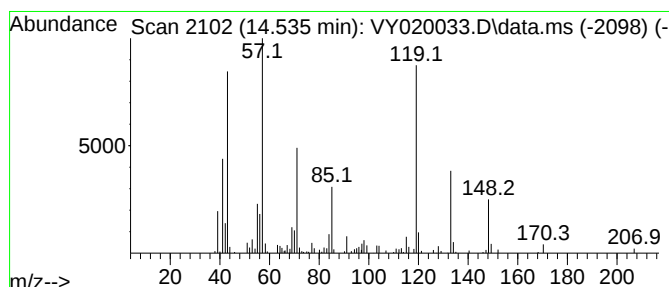
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Quant Title : SW846 8260

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

Peak Number 9 unknown14.535 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.535	91.53 ug/l	123528	1,4-Dichlorobenzene-d4	13.359

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	38
2	Dodecane	170	C12H26	000112-40-3	30
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	30
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	30
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
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Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

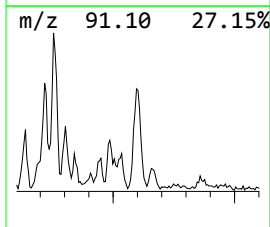
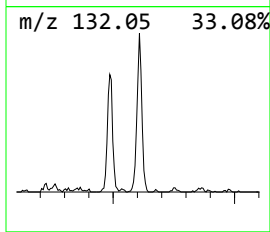
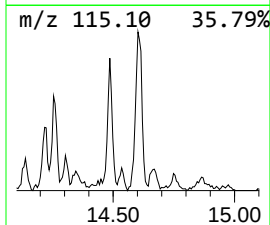
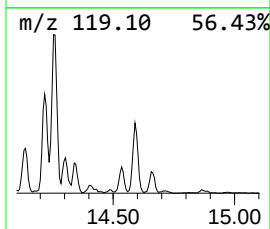
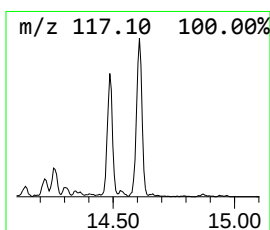
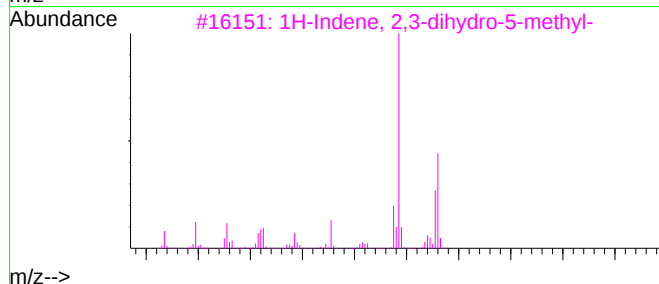
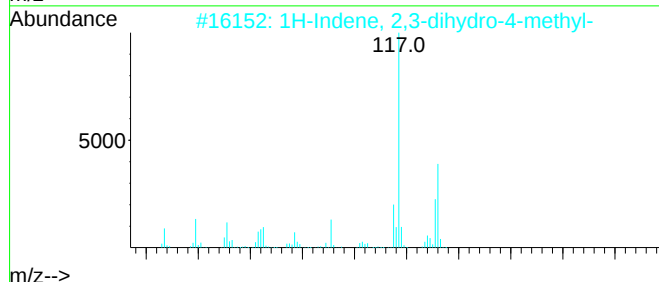
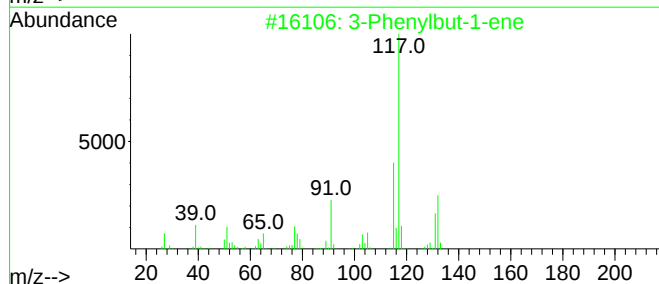
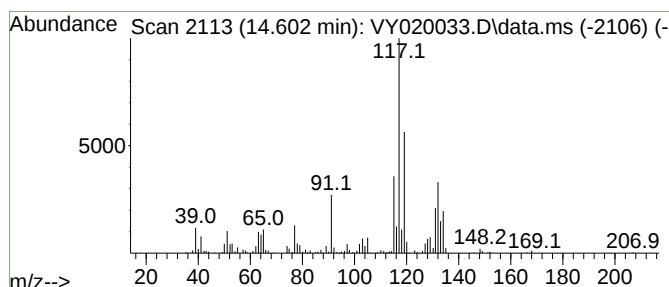
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Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 3-Phenylbut-1-ene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.602	226.02 ug/l	305033	1,4-Dichlorobenzene-d4		13.359	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene		132	C10H12	000934-10-1	92
2	1H-Indene, 2,3-dihydro-4-methyl-		132	C10H12	000824-22-6	70
3	1H-Indene, 2,3-dihydro-5-methyl-		132	C10H12	000874-35-1	64
4	Benzene, (1-methyl-1-propenyl)-,...		132	C10H12	000767-99-7	64
5	Benzene, 2-butenyl-		132	C10H12	001560-06-1	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102824\
Data File : VY020033.D
Acq On : 28 Oct 2024 11:56
Operator : SY/MD
Sample : P4578-03
Misc : 5.99g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
WB-305-TOP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y100924S.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
Benzene, 1-ethe...	13.548	101.1	ug/l	136438	4	13.359	67478	50.0
Benzene, 2-ethy...	13.810	73.9	ug/l	99767	4	13.359	67478	50.0
o-Cymene	13.840	87.4	ug/l	117987	4	13.359	67478	50.0
Benzene, 2-ethy...	13.895	233.1	ug/l	314584	4	13.359	67478	50.0
unknown14.005	14.005	83.1	ug/l	112183	4	13.359	67478	50.0
Benzene, 1,2,4,...	14.218	172.6	ug/l	232976	4	13.359	67478	50.0
Benzene, 1-meth...	14.255	208.7	ug/l	281603	4	13.359	67478	50.0
1H-Indene, 2,3-...	14.486	100.1	ug/l	135130	4	13.359	67478	50.0
unknown14.535	14.535	91.5	ug/l	123528	4	13.359	67478	50.0
3-Phenylbut-1-ene	14.602	226.0	ug/l	305033	4	13.359	67478	50.0