

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
 Data File : VY019764.D
 Acq On : 02 Oct 2024 15:00
 Operator : SY/MD
 Sample : P4232-03
 Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TP-R-VOC

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\82Y090924S.M

Title : SW846 8260

Signal : TIC: VY019764.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.634	959	970	976	rBV	235559	575041	2.76%	0.428%
2	7.707	976	982	991	rVB	372495	855690	4.10%	0.637%
3	8.061	1029	1040	1053	rBV	208449	473143	2.27%	0.352%
4	8.250	1053	1071	1081	rBV	353191	877220	4.20%	0.653%
5	8.616	1121	1131	1144	rBV	644193	1282809	6.15%	0.956%
6	9.116	1199	1213	1217	rBV	159388	334817	1.60%	0.249%
7	9.170	1217	1222	1229	rVB	174047	330718	1.58%	0.246%
8	9.542	1275	1283	1294	rBV	701886	1311438	6.28%	0.977%
9	9.676	1294	1305	1312	rBV2	761858	1683675	8.07%	1.254%
10	9.750	1312	1317	1320	rBV	253415	430701	2.06%	0.321%
11	9.890	1329	1340	1351	rBV	362908	906777	4.35%	0.675%
12	10.042	1357	1365	1370	rBV	401715	721955	3.46%	0.538%
13	10.103	1370	1375	1388	rVB	1068574	1876347	8.99%	1.398%
14	10.298	1399	1407	1413	rVB	349026	635290	3.04%	0.473%
15	10.560	1440	1450	1458	rVB5	133896	442238	2.12%	0.329%
16	10.731	1468	1478	1482	rBV	124814	234984	1.13%	0.175%
17	10.829	1489	1494	1501	rBV	192307	357660	1.71%	0.266%
18	11.207	1547	1556	1567	rVB4	569810	1426275	6.83%	1.062%
19	11.304	1567	1572	1581	rBV	485015	882244	4.23%	0.657%
20	11.414	1584	1590	1596	rBV	1037146	1708749	8.19%	1.273%
21	11.518	1604	1607	1615	rVB4	166143	302833	1.45%	0.226%
22	11.603	1615	1621	1623	rVV3	142248	290934	1.39%	0.217%
23	11.640	1623	1627	1635	rVV6	281674	826487	3.96%	0.616%
24	11.926	1668	1674	1681	rBV5	149787	390480	1.87%	0.291%
25	12.249	1721	1727	1733	rBV	2041675	3060584	14.67%	2.280%
26	12.402	1743	1752	1757	rBV2	659773	1202829	5.76%	0.896%
27	12.591	1772	1783	1790	rBV	9030713	13522987	64.80%	10.073%
28	12.725	1800	1805	1812	rVB4	133859	239017	1.15%	0.178%
29	12.889	1824	1832	1836	rBV	175189	295264	1.41%	0.220%
30	13.170	1870	1878	1886	rBV2	1425163	3621028	17.35%	2.697%
31	13.340	1901	1906	1909	rBV	651712	1037791	4.97%	0.773%
32	13.371	1909	1911	1917	rVB	570664	750697	3.60%	0.559%
33	13.444	1918	1923	1927	rVB	228735	345932	1.66%	0.258%
34	13.511	1927	1934	1937	rBV	5850133	8949114	42.89%	6.666%
35	13.542	1937	1939	1943	rVB	2488812	3031835	14.53%	2.258%
36	13.590	1943	1947	1949	rBV	2264231	3381150	16.20%	2.518%

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
 Title : SW846 8260

37	13.670	1956	1960	1966	rVB	952963	1446551	6.93%	1.077%
38	13.737	1966	1971	1977	rVB	391020	579514	2.78%	0.432%
39	13.816	1977	1984	1989	rBV2	531828	923709	4.43%	0.688%
40	13.889	1989	1996	2001	rBV	14971285	20867355	100.00%	15.543%
41	13.944	2001	2005	2009	rVV	954111	1466203	7.03%	1.092%
42	13.993	2009	2013	2019	rVB	3930667	6185026	29.64%	4.607%
43	14.066	2019	2025	2030	rBV3	309797	517942	2.48%	0.386%
44	14.127	2030	2035	2040	rVB3	343679	537918	2.58%	0.401%
45	14.212	2040	2049	2054	rBV	6924690	10583341	50.72%	7.883%
46	14.291	2058	2062	2066	rBV	874399	1202606	5.76%	0.896%
47	14.365	2071	2074	2077	rVV	523248	765082	3.67%	0.570%
48	14.432	2081	2085	2088	rVV2	364662	595923	2.86%	0.444%
49	14.480	2088	2093	2098	rVV	2612849	4218268	20.21%	3.142%
50	14.529	2098	2101	2105	rVV	538375	706561	3.39%	0.526%
51	14.596	2105	2112	2117	rVV3	5307640	10516103	50.39%	7.833%
52	14.651	2117	2121	2127	rVB	805959	1336463	6.40%	0.995%
53	14.743	2133	2136	2141	rVB	258001	347255	1.66%	0.259%
54	14.816	2141	2148	2149	rVV3	344248	498367	2.39%	0.371%
55	14.852	2149	2154	2158	rVV2	1380907	2753160	13.19%	2.051%
56	14.889	2158	2160	2165	rVV	600909	882657	4.23%	0.657%
57	14.962	2165	2172	2180	rVB3	1114921	2322355	11.13%	1.730%
58	15.139	2196	2201	2206	rVV	502009	851687	4.08%	0.634%
59	15.194	2206	2210	2214	rVV	180664	291582	1.40%	0.217%
60	15.249	2214	2219	2223	rVV2	357635	695039	3.33%	0.518%
61	15.297	2223	2227	2240	rVB	298304	604198	2.90%	0.450%
62	15.425	2241	2248	2254	rBV	343757	595052	2.85%	0.443%
63	15.584	2264	2274	2277	rBV3	232029	534401	2.56%	0.398%
64	15.620	2277	2280	2289	rVV3	251225	473321	2.27%	0.353%
65	15.706	2289	2294	2300	rVV	129282	232801	1.12%	0.173%
66	15.779	2300	2306	2320	rVB2	163254	436732	2.09%	0.325%
67	15.925	2320	2330	2341	rBV2	184028	382165	1.83%	0.285%
68	16.175	2365	2371	2386	rVB	274673	573910	2.75%	0.427%
69	16.364	2394	2402	2410	rBV	353492	734958	3.52%	0.547%

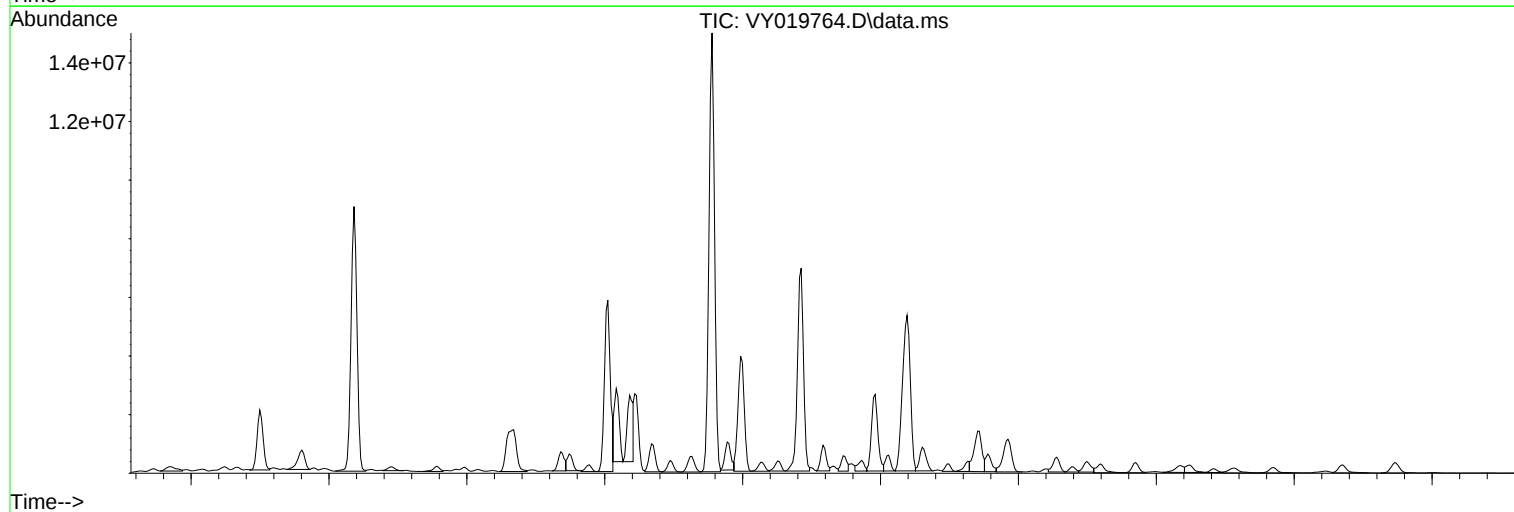
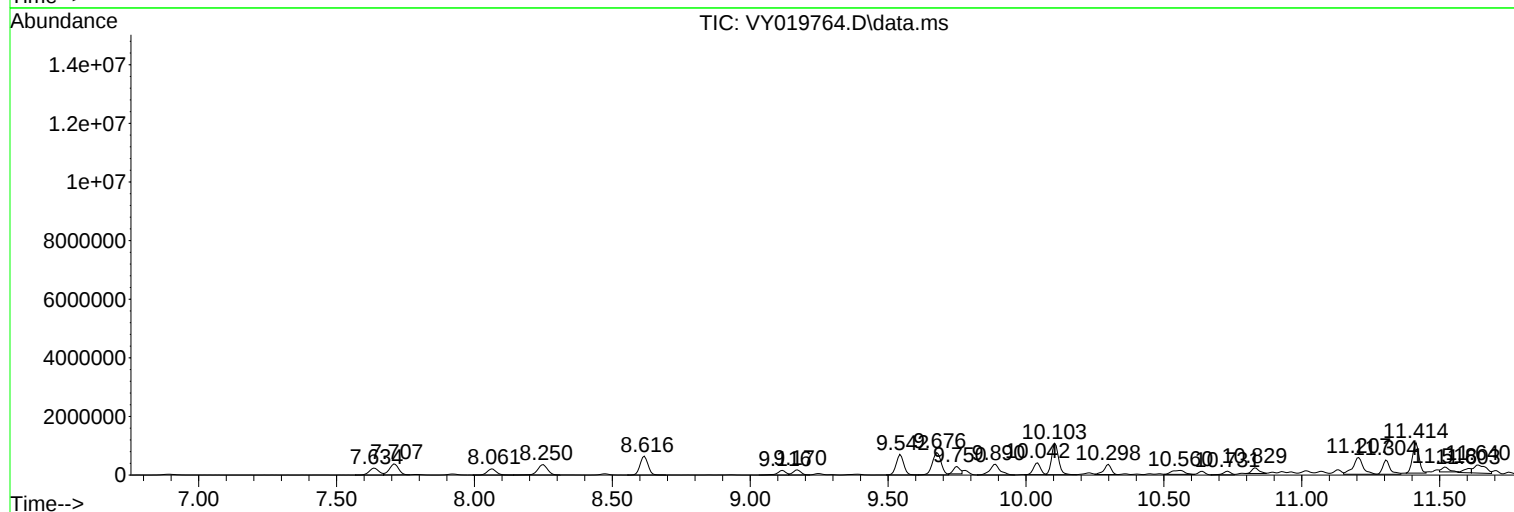
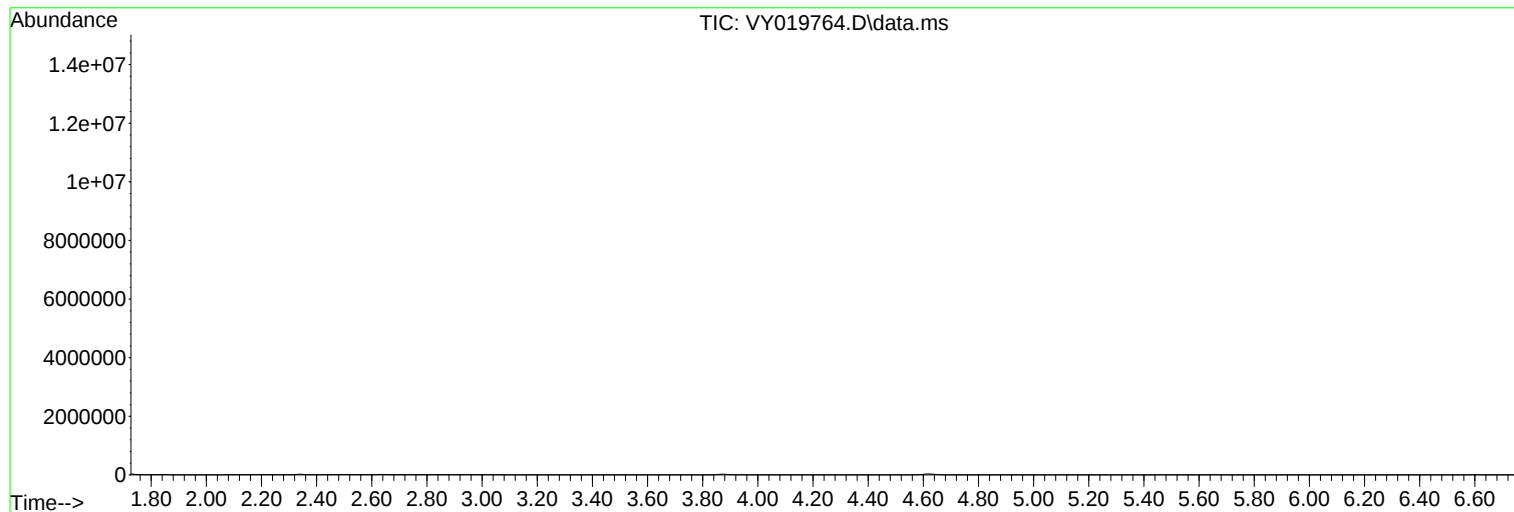
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ClientSampleId :
TP-R-VOC

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.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 :
TP-R-VOC

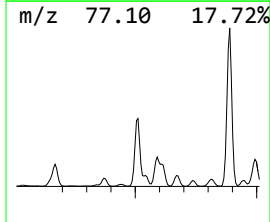
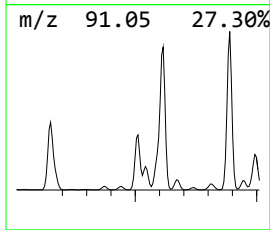
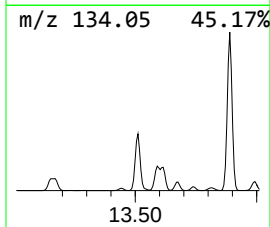
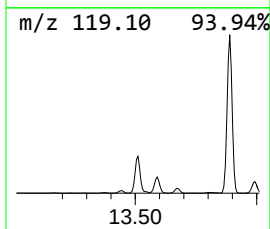
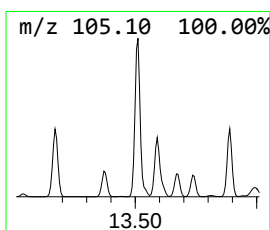
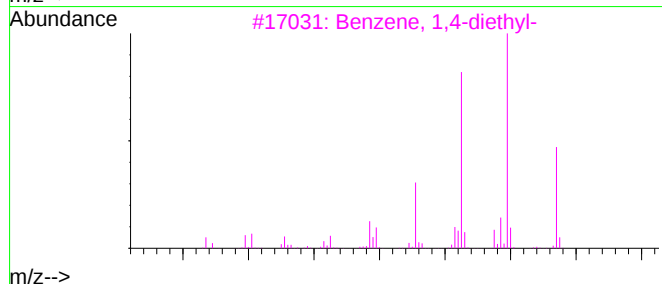
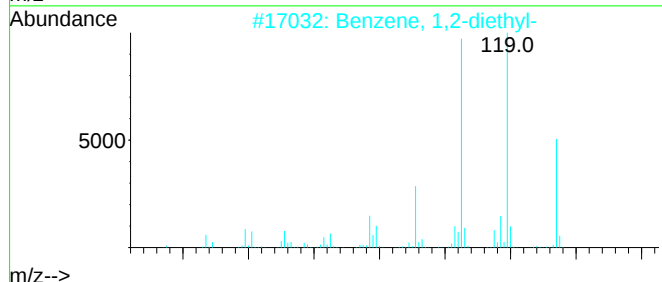
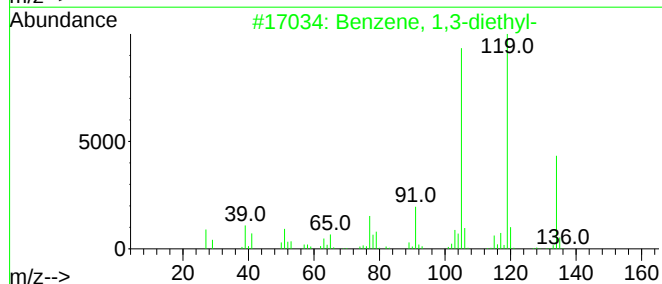
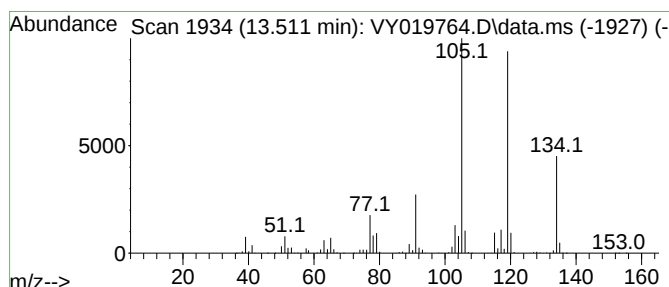
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,3-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.511	431.16 ug/l	8949110	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



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Instrument :
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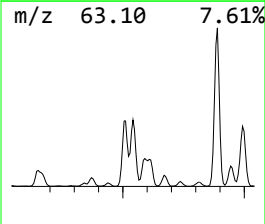
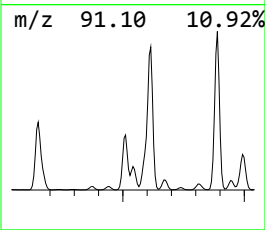
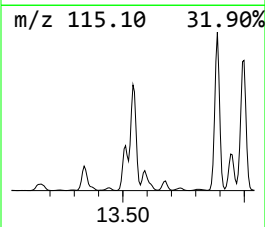
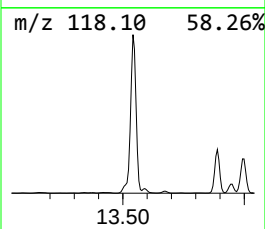
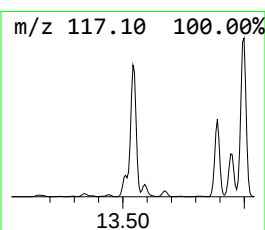
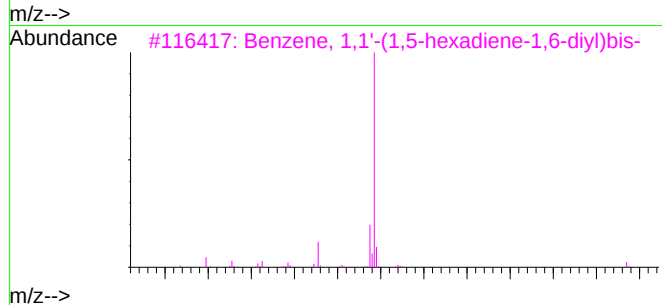
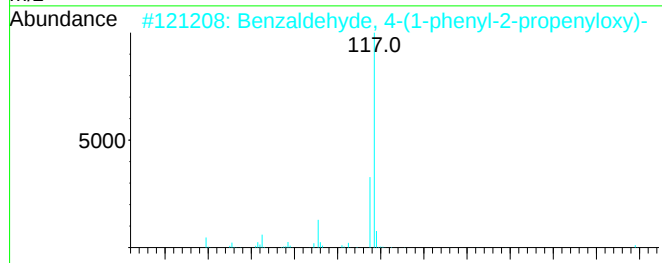
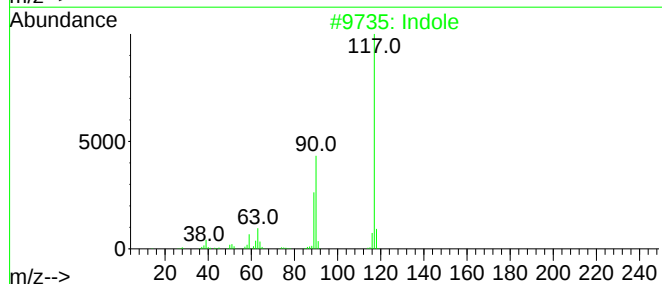
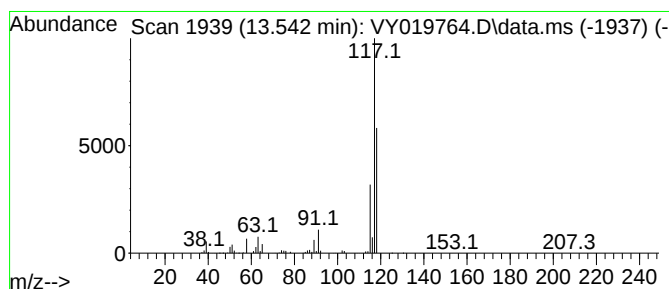
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown13.542 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.542	146.07 ug/l	3031840	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indole	117	C8H7N	000120-72-9	46
2	Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	42
3	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	40
4	Benzyl nitrile	117	C8H7N	000140-29-4	37
5	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	35



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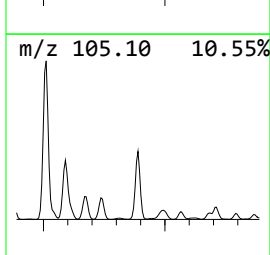
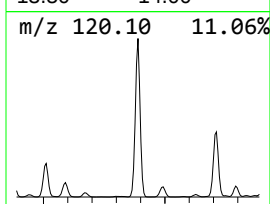
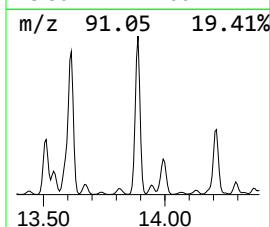
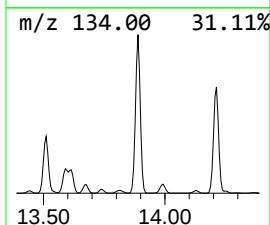
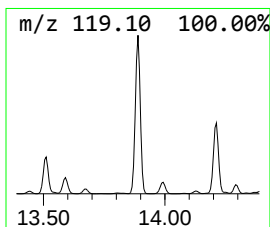
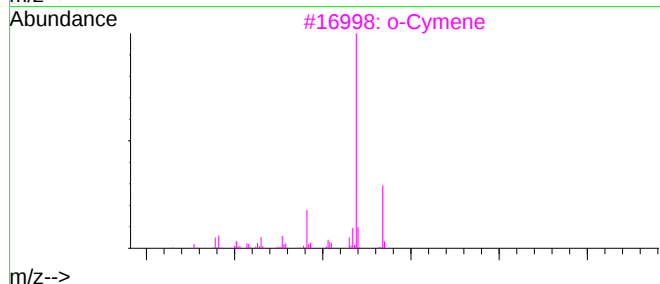
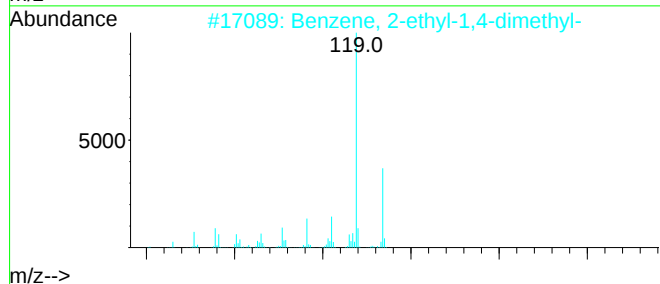
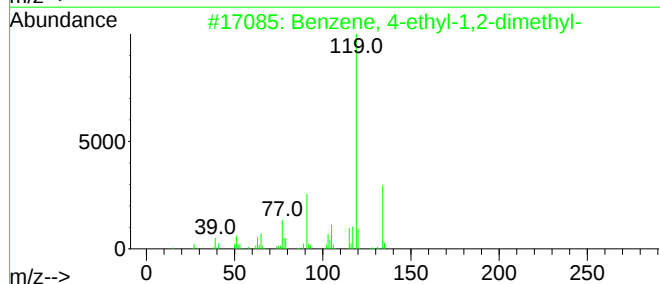
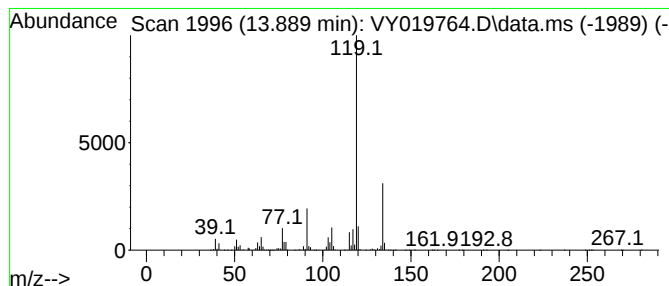
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.889	1005.37 ug/l	20867400	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	p-Cymene	134	C10H14	000099-87-6	95



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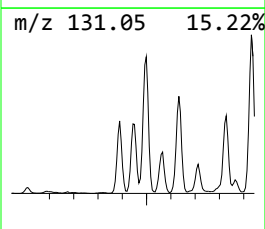
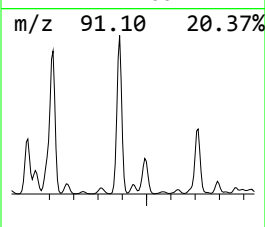
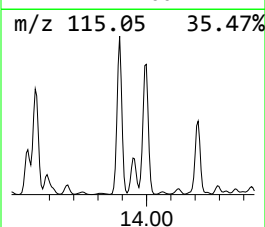
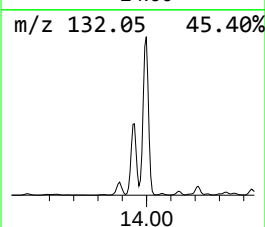
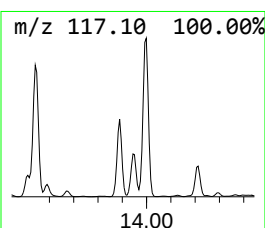
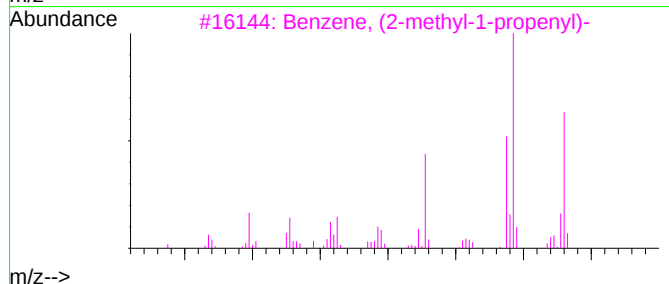
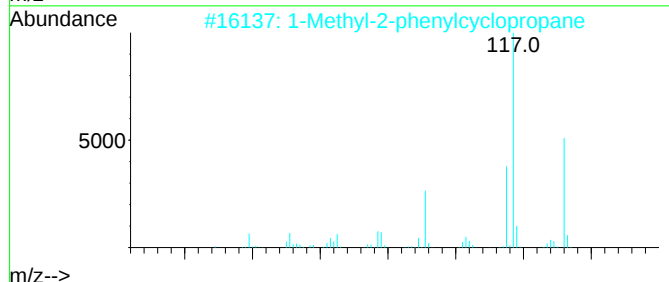
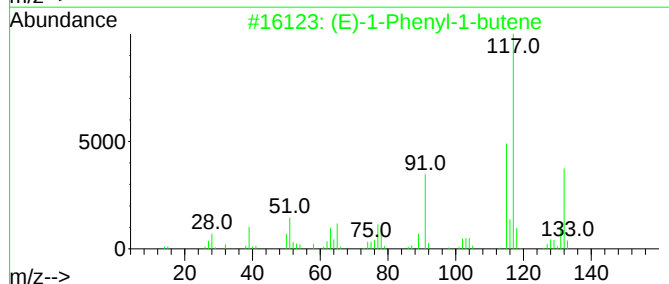
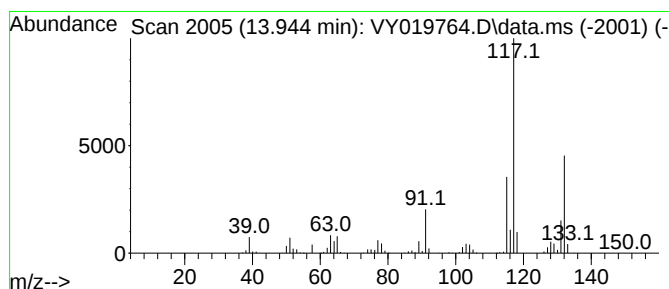
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 4 (E)-1-Phenyl-1-butene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.944	70.64 ug/l	1466200	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	95
2	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	94
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	94
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	94
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/soil/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
Quant Title : SW846 8260

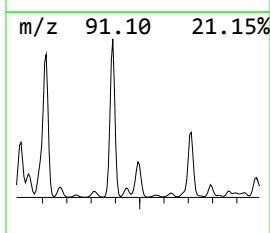
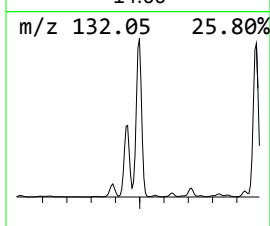
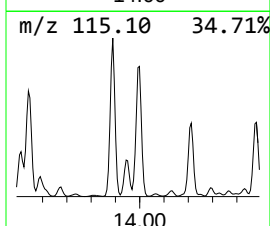
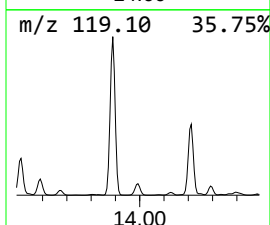
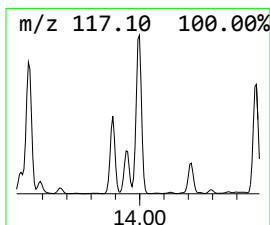
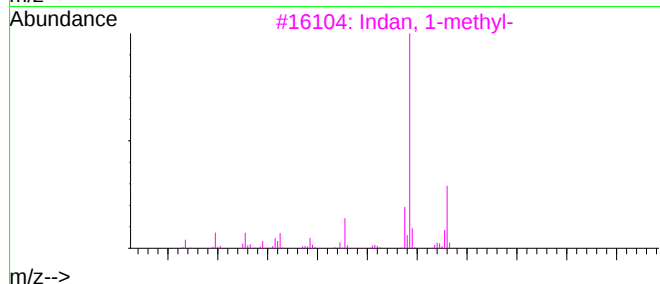
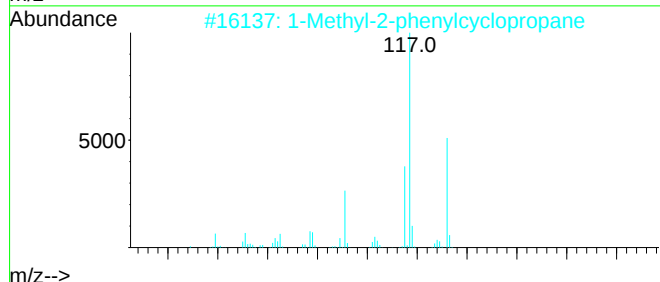
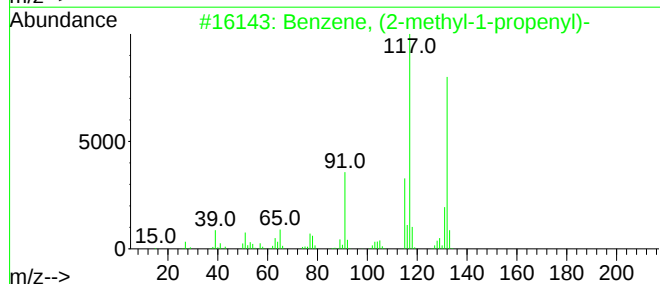
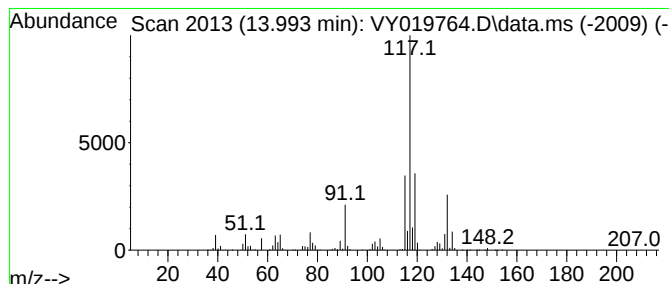
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, (2-methyl-1-propen... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.993	297.99 ug/l	6185030	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
2	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	76
3	Indan, 1-methyl-	132	C10H12	000767-58-8	76
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
Quant Title : SW846 8260

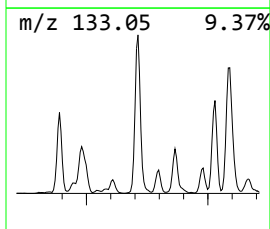
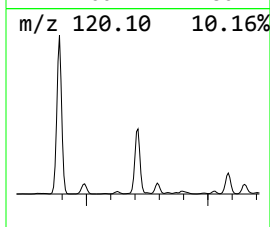
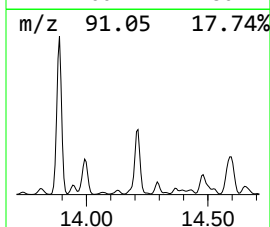
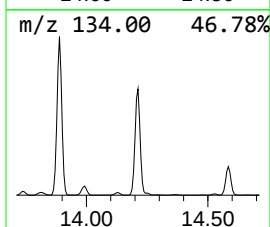
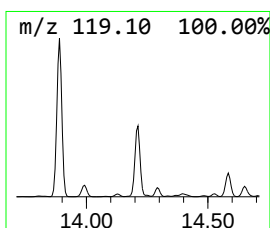
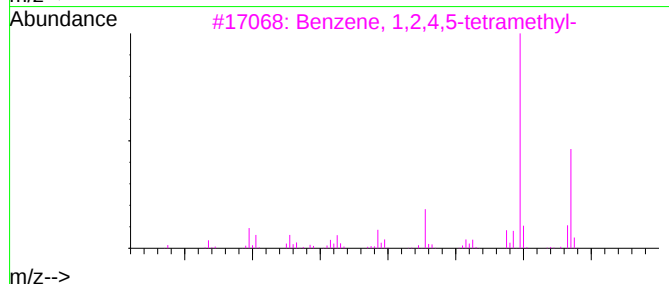
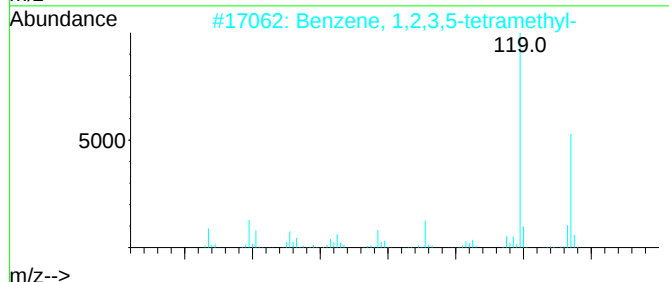
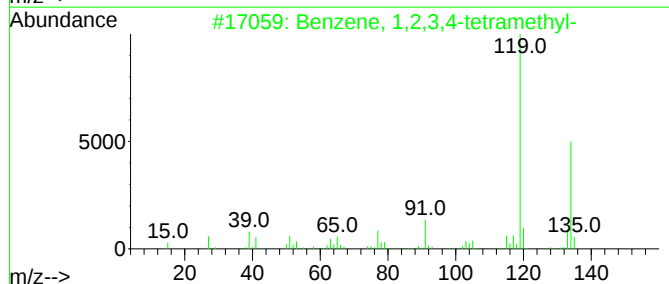
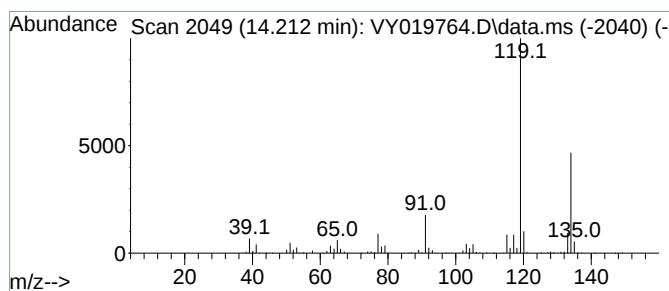
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.212	509.90 ug/l	10583300	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	o-Cymene	134	C10H14	000527-84-4	95
5	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.M
Quant Title : SW846 8260

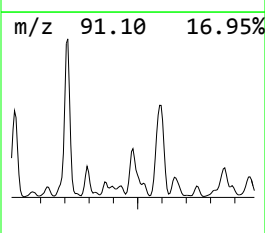
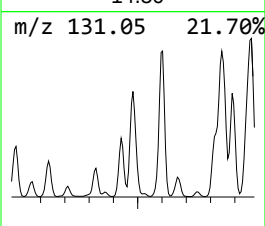
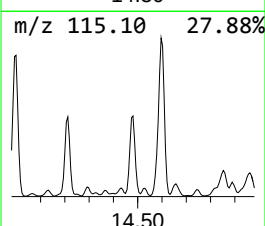
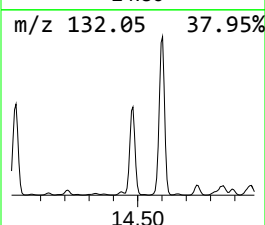
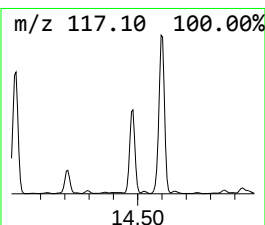
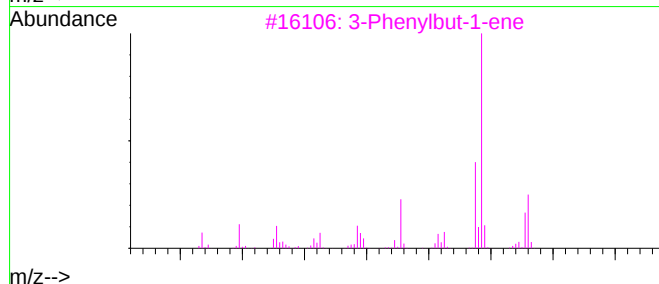
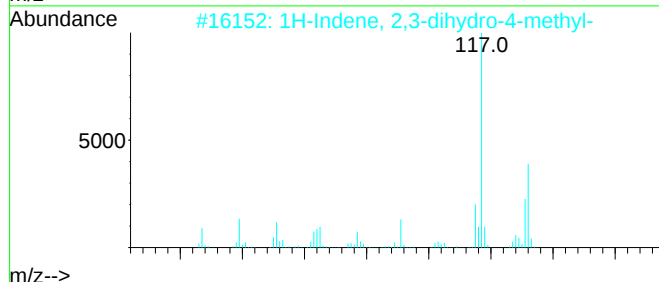
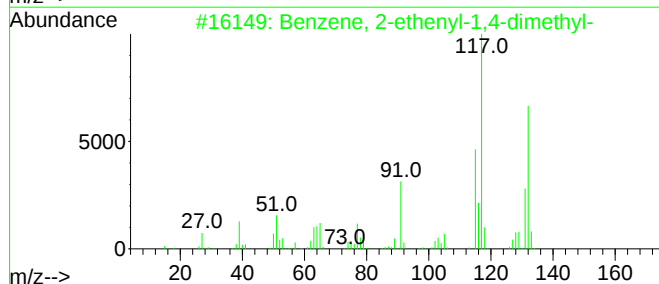
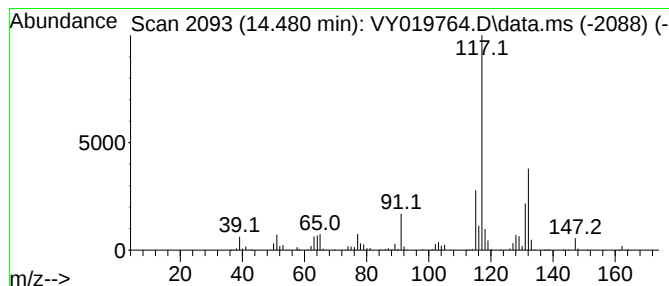
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.480	203.23 ug/l	4218270	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

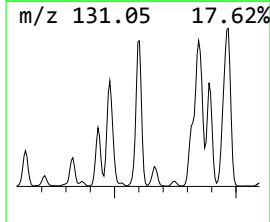
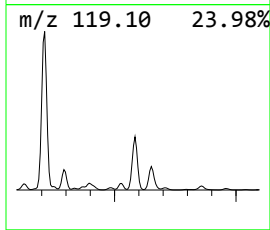
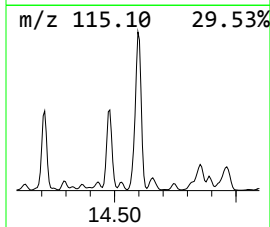
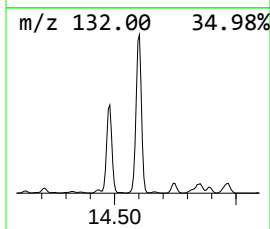
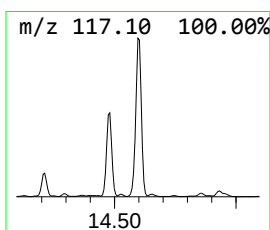
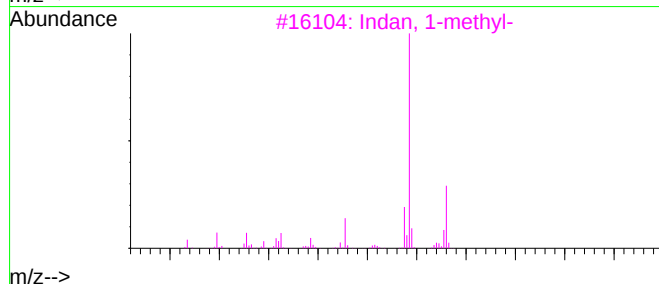
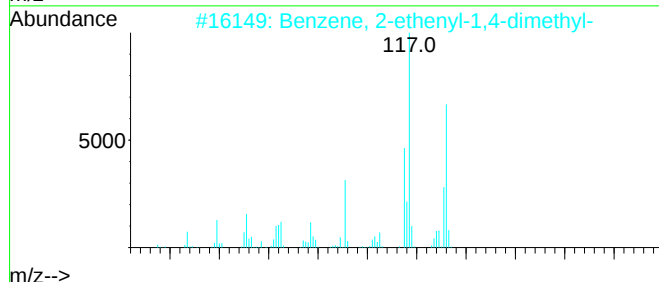
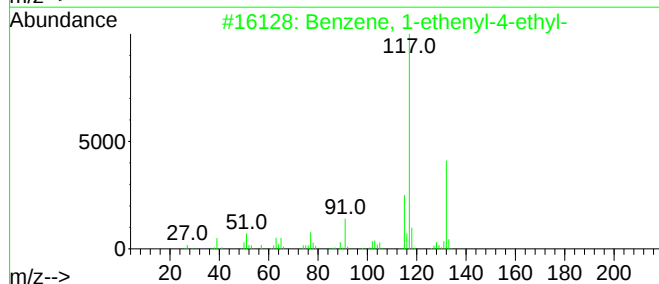
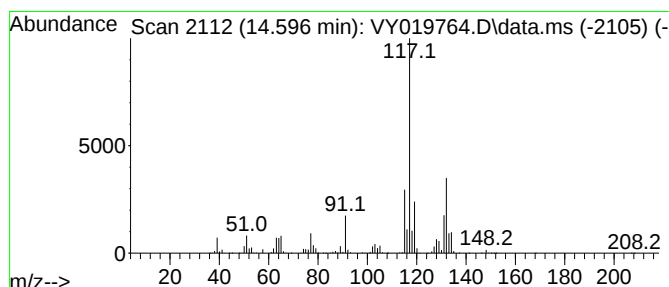
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.596	506.66 ug/l	10516100	1,4-Dichlorobenzene-d4	13.340

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	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

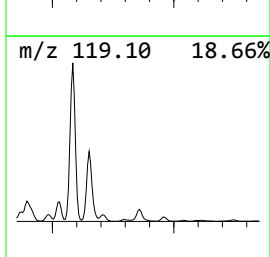
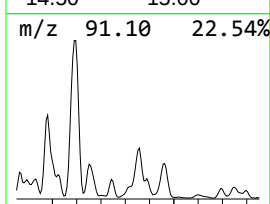
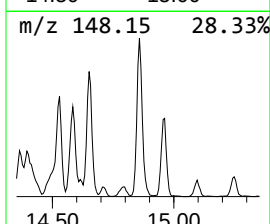
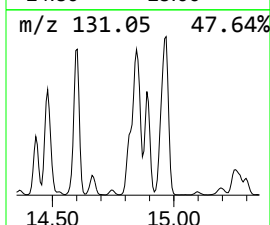
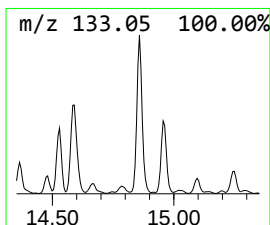
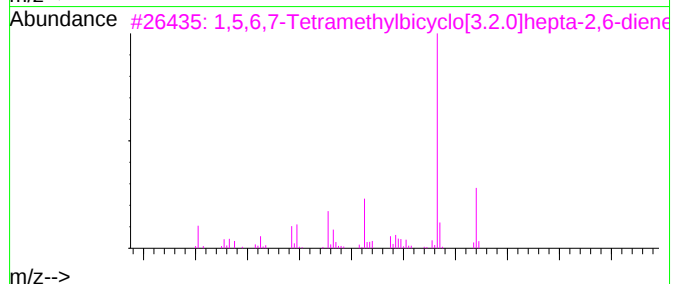
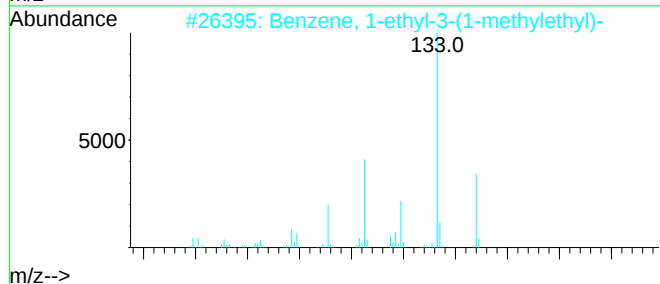
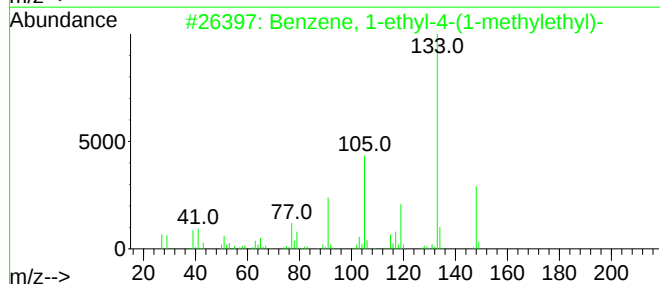
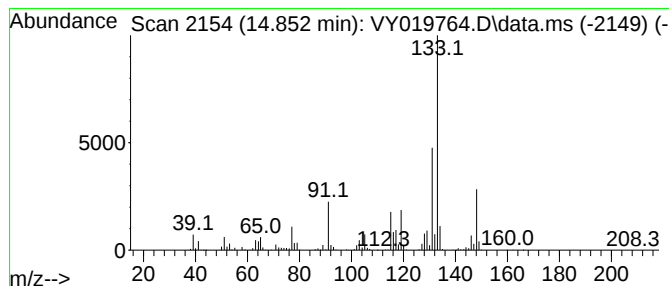
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.852	132.65 ug/l	2753160	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	70
2	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	58
3	1,5,6,7-Tetramethylbicyclo[3.2.0]hepta-2,6-diene	148	C11H16	134329-46-7	58
4	Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	58
5	4-tert-Butyltoluene	148	C11H16	000098-51-1	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

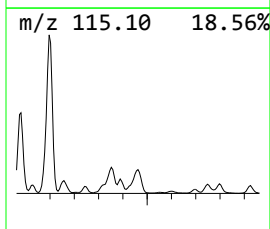
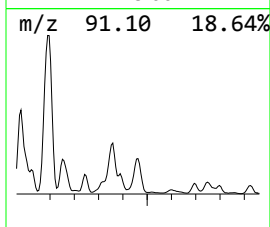
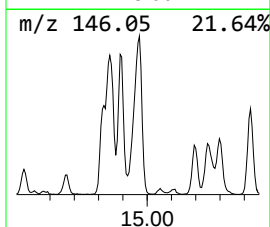
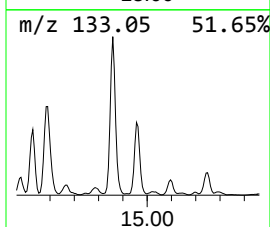
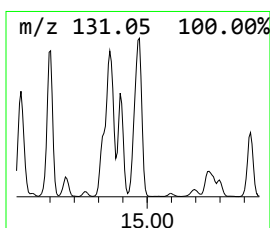
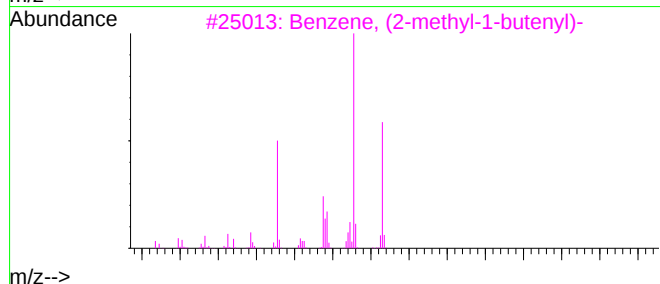
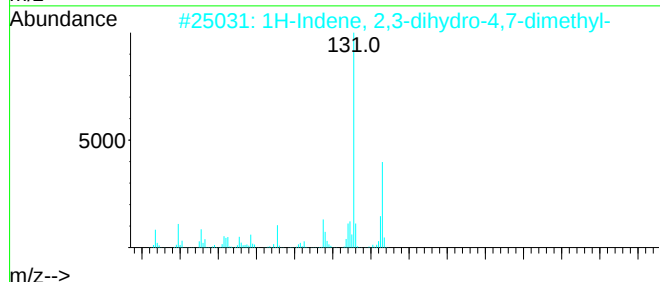
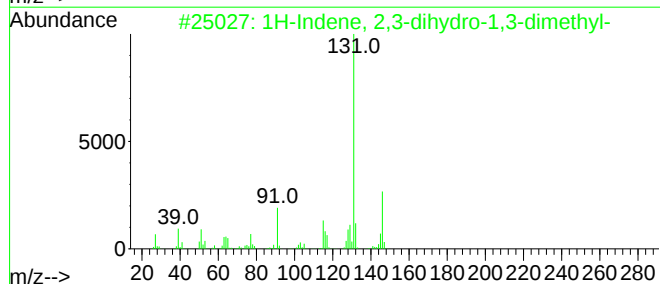
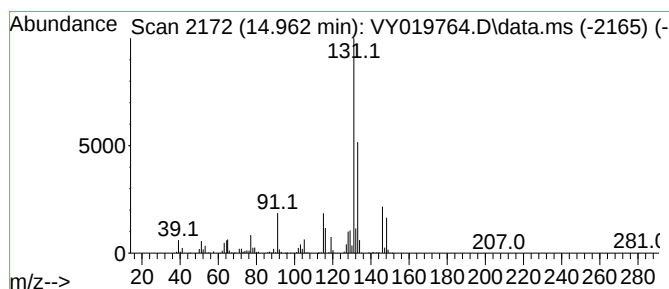
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.962	111.89 ug/l	2322360	1,4-Dichlorobenzene-d4	13.340

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	70
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
3	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	70
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY100224\
Data File : VY019764.D
Acq On : 02 Oct 2024 15:00
Operator : SY/MD
Sample : P4232-03
Misc : 6.12g/5.0mL/MSVOA_Y/SOIL/A
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TP-R-VOC

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y090924S.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, 1,3-di...	13.511	431.2	ug/l	8949110	4	13.340	1037790	50.0
unknown13.542	13.542	146.1	ug/l	3031840	4	13.340	1037790	50.0
Benzene, 4-ethy...	13.889	1005.4	ug/l	20867400	4	13.340	1037790	50.0
(E)-1-Phenyl-1-...	13.944	70.6	ug/l	1466200	4	13.340	1037790	50.0
Benzene, (2-met...	13.993	298.0	ug/l	6185030	4	13.340	1037790	50.0
Benzene, 1,2,3,...	14.212	509.9	ug/l	10583300	4	13.340	1037790	50.0
Benzene, 2-ethe...	14.480	203.2	ug/l	4218270	4	13.340	1037790	50.0
Benzene, 1-ethe...	14.596	506.7	ug/l	10516100	4	13.340	1037790	50.0
Benzene, 1-ethy...	14.852	132.7	ug/l	2753160	4	13.340	1037790	50.0
1H-Indene, 2,3-...	14.962	111.9	ug/l	2322360	4	13.340	1037790	50.0