

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
 Data File : VY020969.D
 Acq On : 29 Jan 2025 17:05
 Operator : SY/MD
 Sample : Q1220-01
 Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 TR-06-01292025

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\82Y012425S.M
 Title : SW846 8260

Signal : TIC: VY020969.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.299	256	259	268	rVB2	1200	2113	1.76%	0.169%
2	4.104	384	391	395	rBV4	1432	3574	2.98%	0.287%
3	6.878	836	846	855	rBV2	10384	31702	26.47%	2.542%
4	7.628	960	969	976	rBV3	19592	48133	40.19%	3.860%
5	7.707	976	982	994	rVB	37920	88842	74.17%	7.125%
6	8.061	1030	1040	1051	rVB2	19661	50645	42.28%	4.061%
7	8.616	1123	1131	1143	rBV	50901	104691	87.41%	8.396%
8	9.999	1354	1358	1363	rVB3	979	1612	1.35%	0.129%
9	10.103	1368	1375	1383	rBV	67009	119774	100.00%	9.605%
10	10.170	1383	1386	1393	rVB4	2716	4733	3.95%	0.380%
11	10.512	1436	1442	1448	rVB4	2694	5438	4.54%	0.436%
12	10.877	1496	1502	1506	rBV2	1043	2155	1.80%	0.173%
13	11.249	1558	1563	1569	rBV2	6493	10597	8.85%	0.850%
14	11.414	1584	1590	1601	rVB	68444	113618	94.86%	9.112%
15	11.627	1616	1625	1633	rBV7	1639	3944	3.29%	0.316%
16	11.956	1675	1679	1685	rVB4	1250	2334	1.95%	0.187%
17	12.414	1748	1754	1766	rVB2	51344	100185	83.65%	8.034%
18	12.597	1779	1784	1789	rBV	1574	2805	2.34%	0.225%
19	12.664	1789	1795	1798	rBV2	6539	10582	8.83%	0.849%
20	12.737	1803	1807	1812	rVB4	6864	10541	8.80%	0.845%
21	12.895	1827	1833	1838	rVB	4664	7007	5.85%	0.562%
22	12.968	1841	1845	1847	rBV3	1806	2385	1.99%	0.191%
23	13.048	1852	1858	1864	rVV	13251	20777	17.35%	1.666%
24	13.115	1864	1869	1871	rVB4	1214	1669	1.39%	0.134%
25	13.182	1871	1880	1882	rBV4	1781	3303	2.76%	0.265%
26	13.237	1882	1889	1894	rVB5	3538	6317	5.27%	0.507%
27	13.285	1894	1897	1901	rBV4	2029	2813	2.35%	0.226%
28	13.346	1901	1907	1917	rVV2	49560	89991	75.13%	7.217%
29	13.517	1931	1935	1936	rBV3	2762	3161	2.64%	0.253%
30	13.548	1936	1940	1944	rVV2	11810	20057	16.75%	1.608%
31	13.590	1944	1947	1956	rVV4	8942	18876	15.76%	1.514%
32	13.676	1956	1961	1965	rVV5	9059	15171	12.67%	1.217%
33	13.749	1969	1973	1978	rVV	5596	9300	7.76%	0.746%
34	13.810	1978	1983	1985	rVV2	6012	10062	8.40%	0.807%
35	13.840	1985	1988	1992	rVV2	7101	10965	9.15%	0.879%
36	13.895	1992	1997	2002	rVV2	11053	17625	14.72%	1.413%

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Peak Location: TOP

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
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37	13.950	2002	2006	2009	rVV6	2055	3572	2.98%	0.286%
38	14.005	2009	2015	2020	rVB	12004	19275	16.09%	1.546%
39	14.054	2020	2023	2024	rBV3	1570	1736	1.45%	0.139%
40	14.084	2025	2028	2031	rVV4	2090	3592	3.00%	0.288%
41	14.133	2031	2036	2040	rVV3	4390	9745	8.14%	0.782%
42	14.182	2040	2044	2047	rVV6	4513	8630	7.21%	0.692%
43	14.218	2047	2050	2053	rVV3	5807	9353	7.81%	0.750%
44	14.255	2053	2056	2061	rVV2	8731	12858	10.74%	1.031%
45	14.297	2061	2063	2067	rVV3	5054	7884	6.58%	0.632%
46	14.340	2067	2070	2077	rVV4	4393	6899	5.76%	0.553%
47	14.407	2077	2081	2083	rVV5	2069	3385	2.83%	0.271%
48	14.444	2084	2087	2090	rVV4	3547	5336	4.46%	0.428%
49	14.486	2090	2094	2099	rVV3	9636	16630	13.88%	1.334%
50	14.535	2099	2102	2107	rVB2	7344	10675	8.91%	0.856%
51	14.602	2107	2113	2119	rBV3	19009	40766	34.04%	3.269%
52	14.657	2119	2122	2130	rVV6	3572	6482	5.41%	0.520%
53	14.755	2132	2138	2144	rBV	15878	24465	20.43%	1.962%
54	14.822	2146	2149	2151	rVV3	1729	2242	1.87%	0.180%
55	14.864	2151	2156	2159	rVV5	7422	13745	11.48%	1.102%
56	14.895	2159	2161	2166	rVV4	4610	7041	5.88%	0.565%
57	14.968	2166	2173	2179	rVB5	6469	15035	12.55%	1.206%
58	15.047	2182	2186	2191	rVB5	1371	2551	2.13%	0.205%
59	15.133	2191	2200	2203	rBV7	1356	3654	3.05%	0.293%
60	15.212	2208	2213	2215	rBV4	5465	9896	8.26%	0.794%
61	15.279	2220	2224	2232	rVV5	4209	11374	9.50%	0.912%
62	15.346	2233	2235	2239	rVB3	1486	1420	1.19%	0.114%
63	15.431	2245	2249	2254	rVB5	2421	3696	3.09%	0.296%
64	15.596	2270	2276	2277	rBV5	1665	2633	2.20%	0.211%
65	15.626	2278	2281	2287	rVB4	6352	10571	8.83%	0.848%
66	15.791	2302	2308	2314	rVB8	1296	3196	2.67%	0.256%
67	15.931	2327	2331	2338	rVB5	2610	4509	3.76%	0.362%
68	16.120	2358	2362	2368	rVB5	2080	4180	3.49%	0.335%
69	16.187	2368	2373	2376	rBV4	1309	2433	2.03%	0.195%

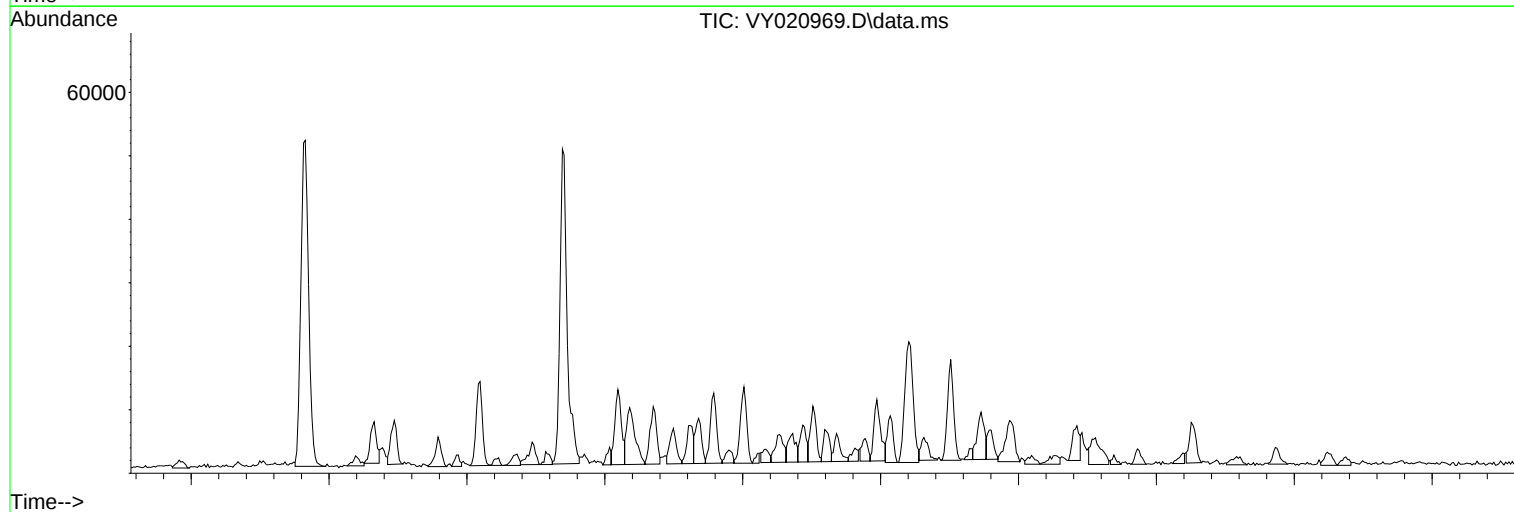
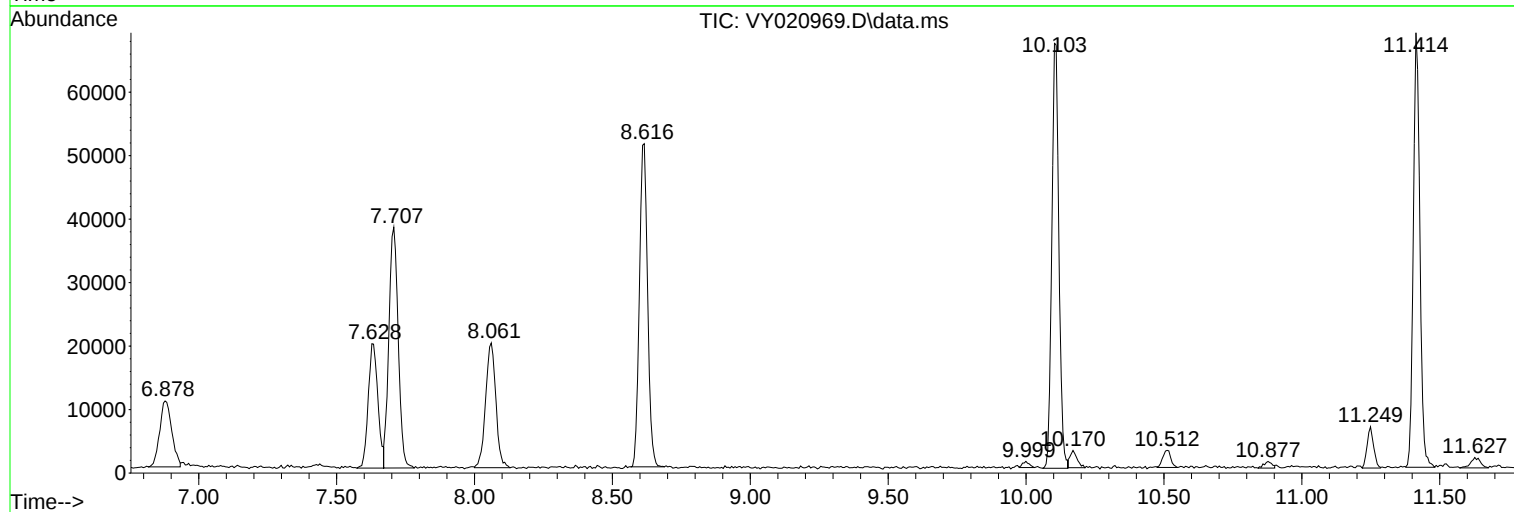
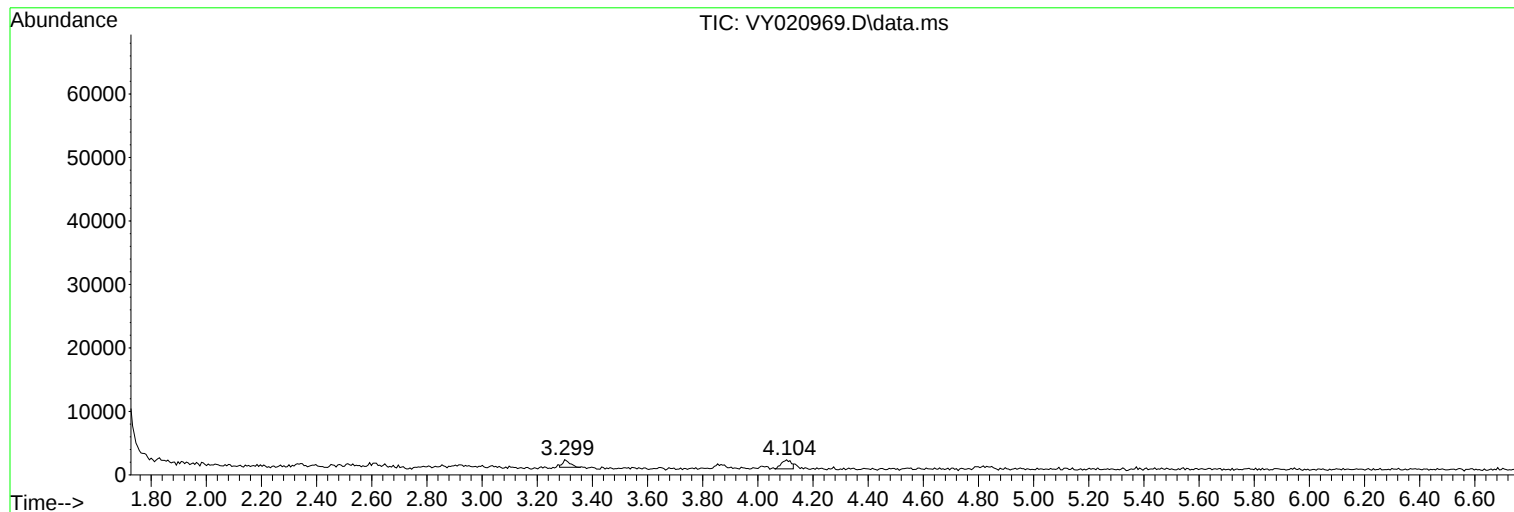
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Quant Title : SW846 8260

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



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ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

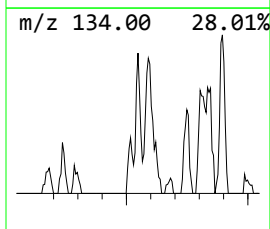
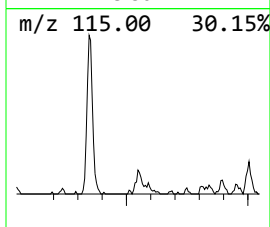
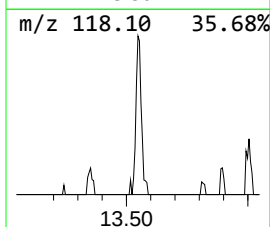
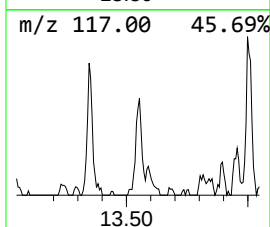
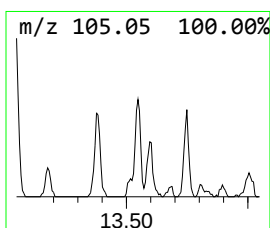
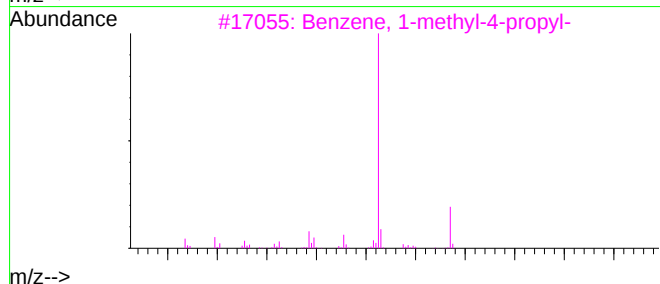
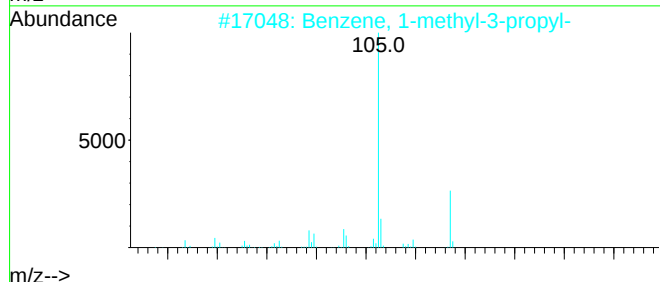
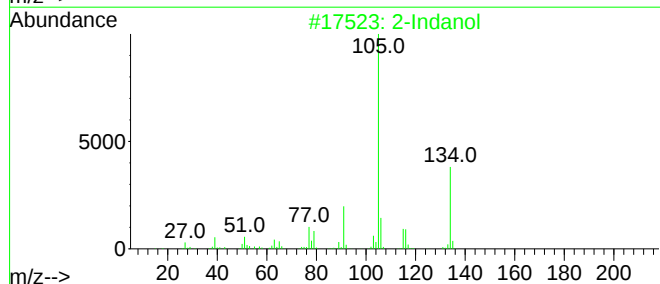
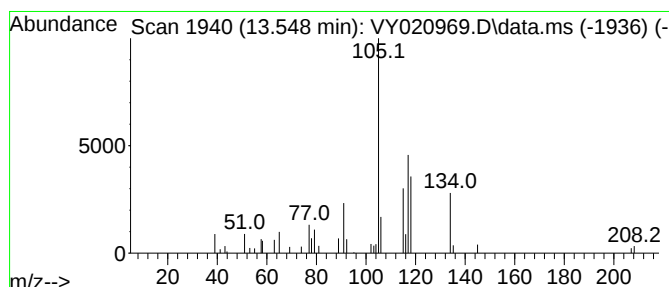
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown13.548 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.548	11.14 ug/l	20057	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Indanol	134	C9H10O	004254-29-9	41
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	30
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	27
4			2-Tolyloxirane	134	C9H10O	002783-26-8	25
5			Hydratropaldehyde	134	C9H10O	034713-70-7	22



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Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

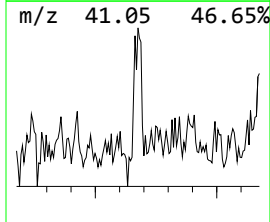
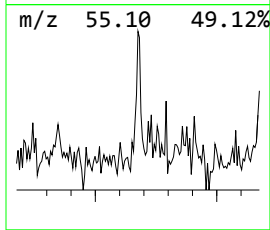
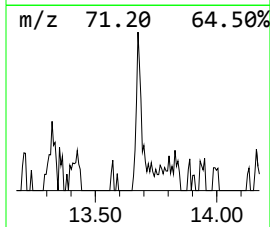
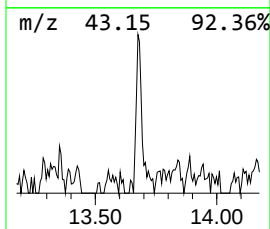
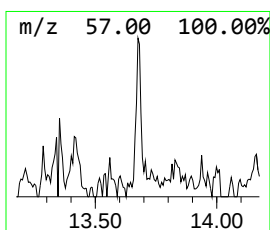
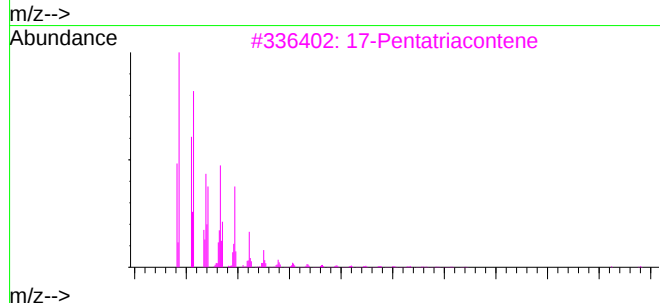
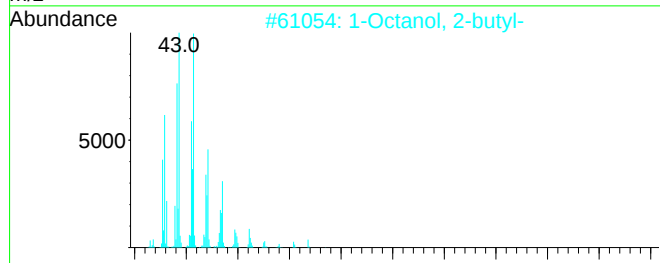
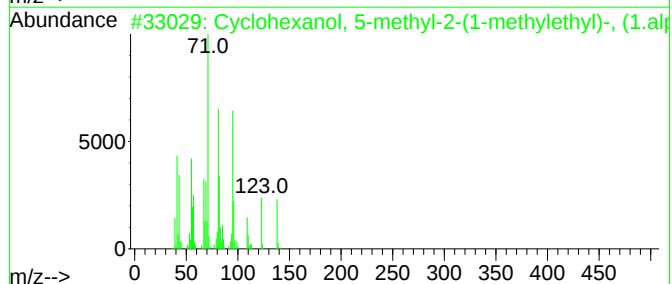
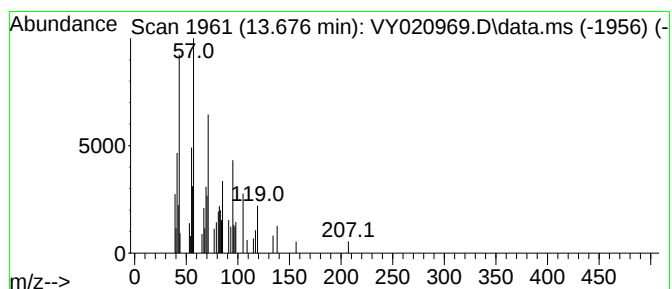
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown13.676 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.676	8.43 ug/l	15171	1,4-Dichlorobenzene-d4	13.352

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	000491-02-1	35
2	1-Octanol, 2-butyl-	186	C12H26O	003913-02-8	27
3	17-Pentatriacontene	491	C35H70	006971-40-0	27
4	Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	27
5	Carbonic acid, nonyl prop-1-en-2...	228	C13H24O3	1000382-53-9	27



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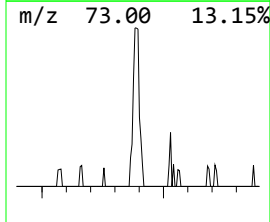
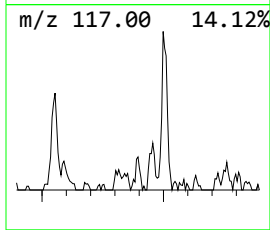
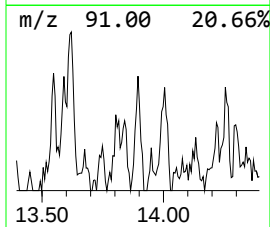
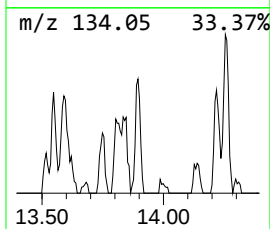
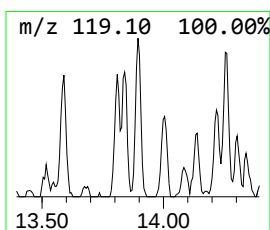
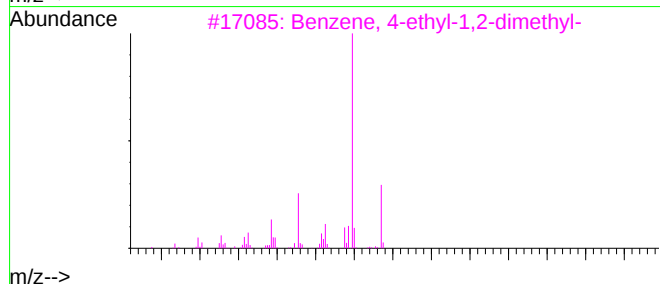
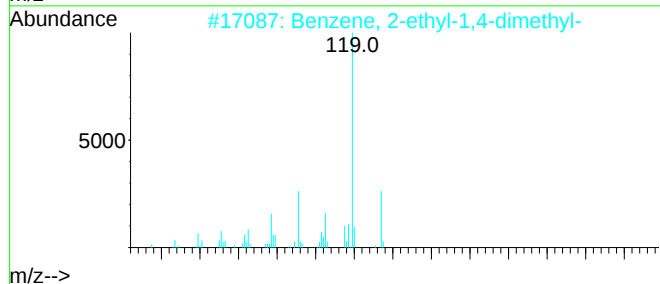
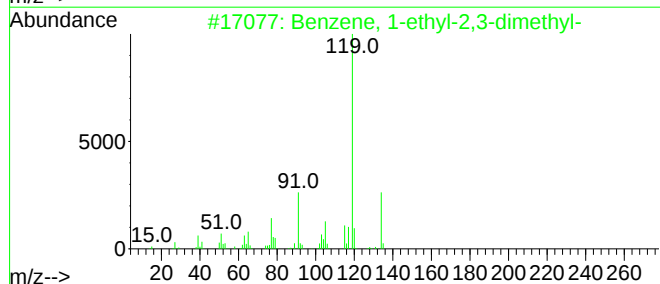
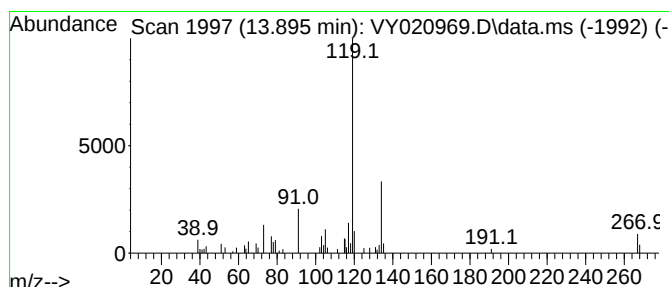
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.895	9.79 ug/l	17625	1,4-Dichlorobenzene-d4	13.352

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	90
4	1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	87
5	o-Cymene	134	C10H14	000527-84-4	81



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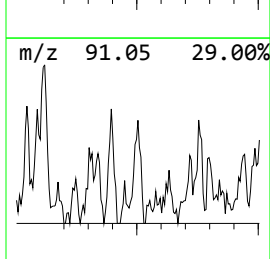
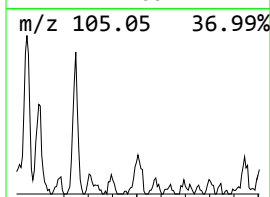
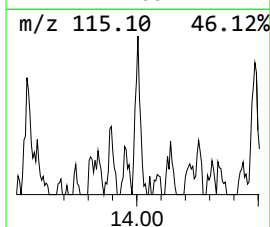
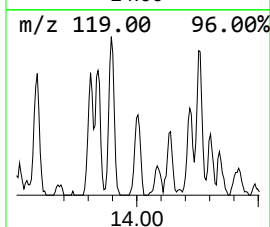
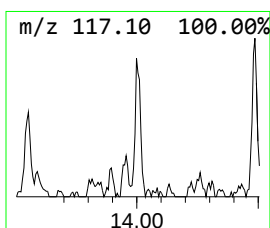
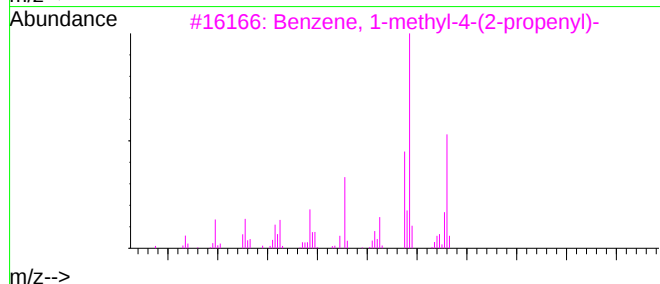
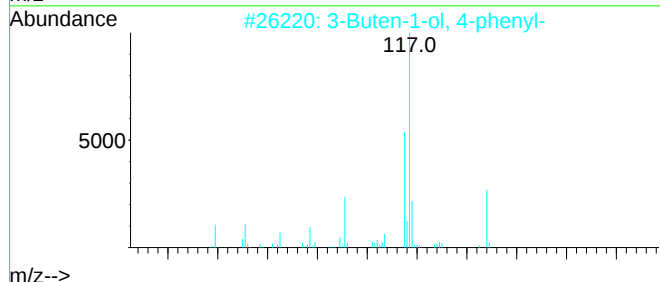
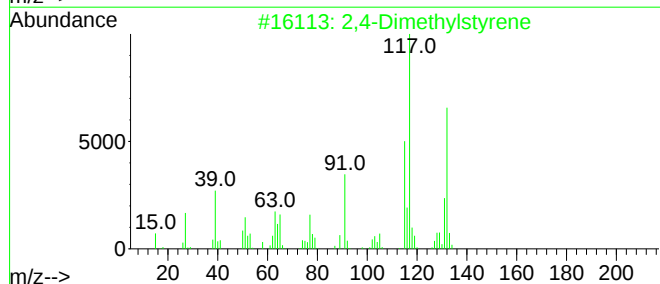
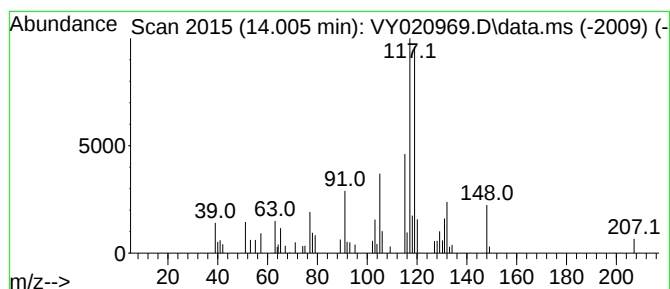
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Quant Title : SW846 8260

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

Peak Number 5 2,4-Dimethylstyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.005	10.71 ug/l	19275	1,4-Dichlorobenzene-d4	13.352

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dimethylstyrene	132	C10H12	002234-20-0	55
2	3-Buten-1-ol, 4-phenyl-	148	C10H12O	000937-58-6	55
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	55
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	53
5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
Data File : VY020969.D
Acq On : 29 Jan 2025 17:05
Operator : SY/MD
Sample : Q1220-01
Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 4 Sample Multiplier: 1

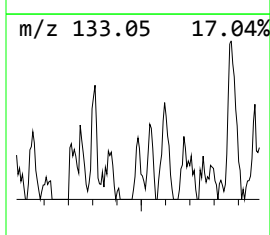
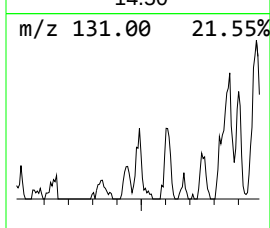
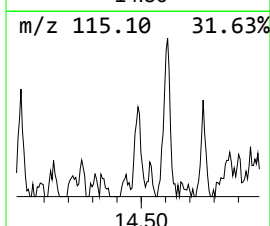
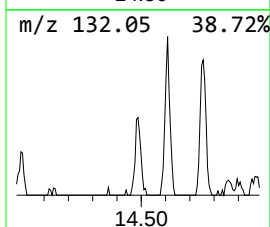
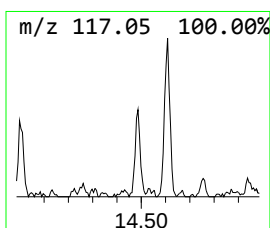
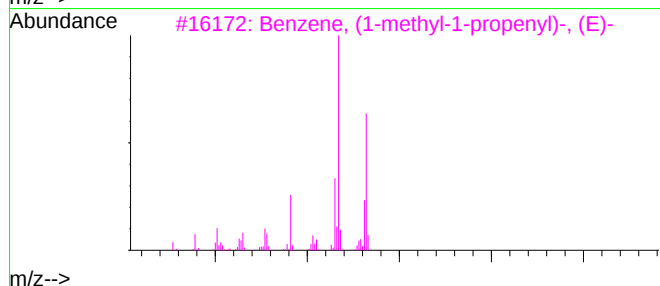
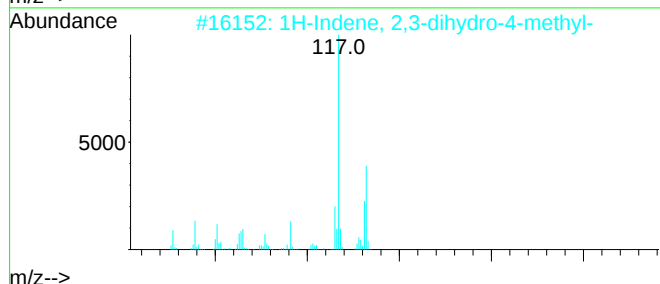
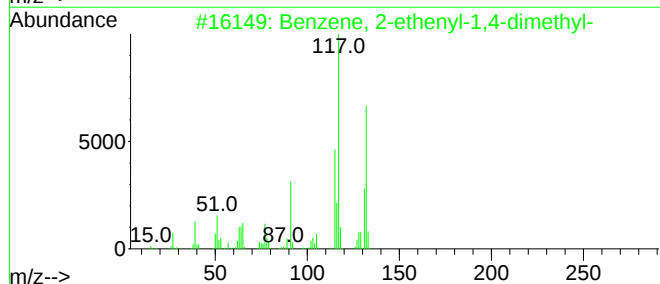
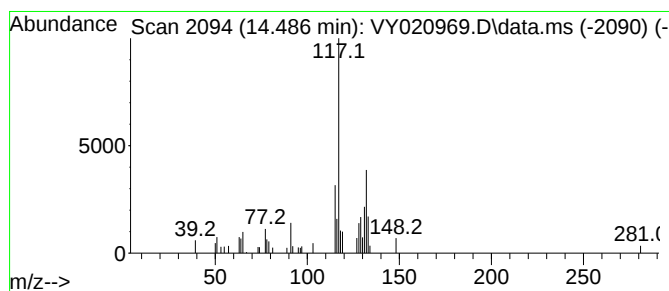
Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.486	9.24 ug/l	16630	1,4-Dichlorobenzene-d4	13.352	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	80
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	72
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	72
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
Data File : VY020969.D
Acq On : 29 Jan 2025 17:05
Operator : SY/MD
Sample : Q1220-01
Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 4 Sample Multiplier: 1

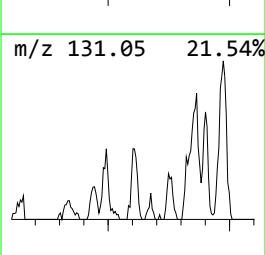
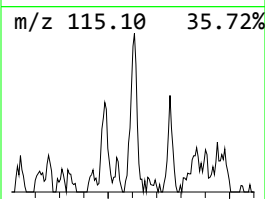
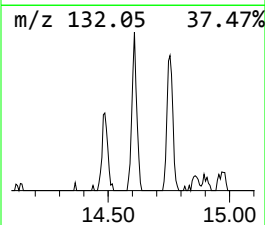
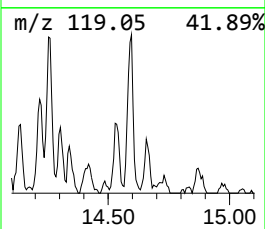
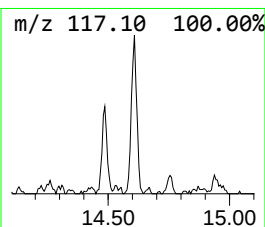
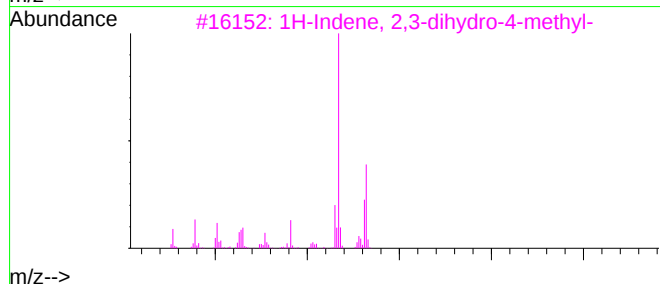
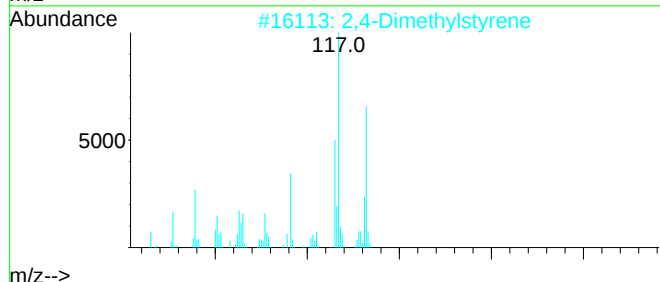
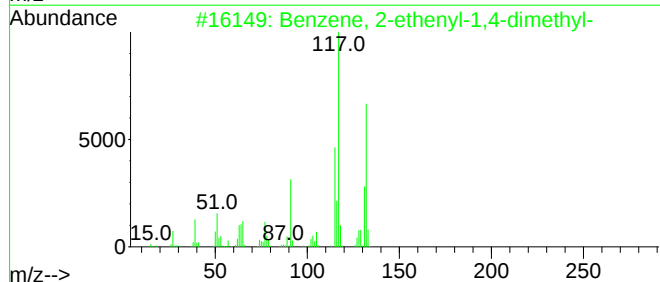
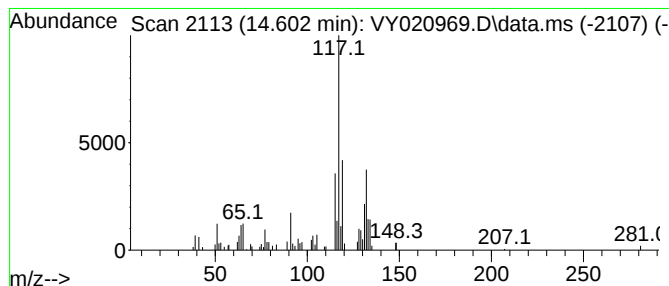
Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.602	22.65 ug/l	40766	1,4-Dichlorobenzene-d4	13.352	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
2	2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
Data File : VY020969.D
Acq On : 29 Jan 2025 17:05
Operator : SY/MD
Sample : Q1220-01
Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

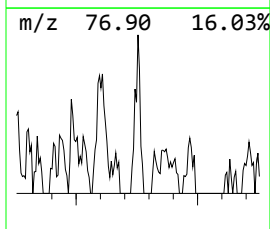
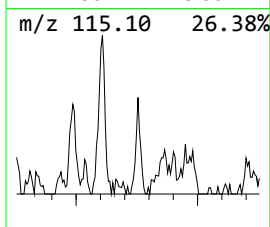
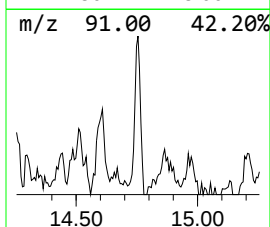
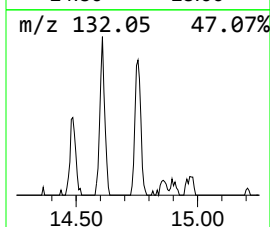
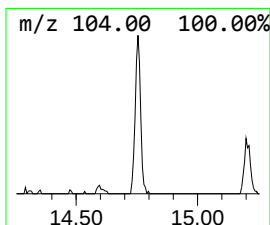
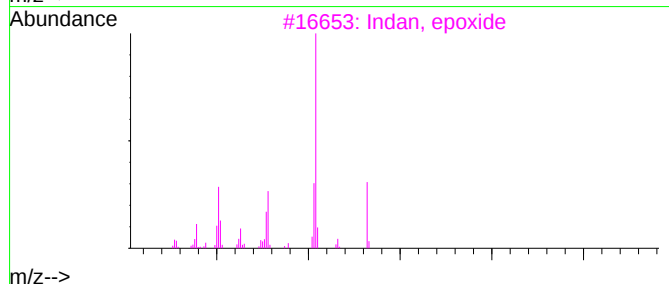
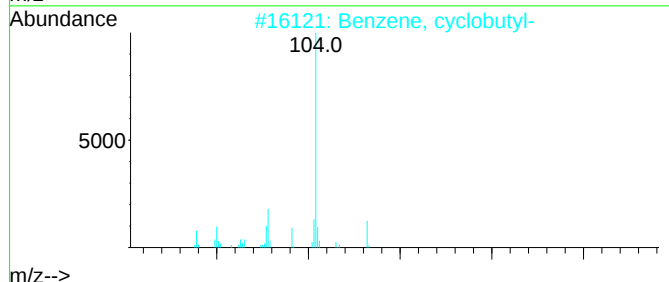
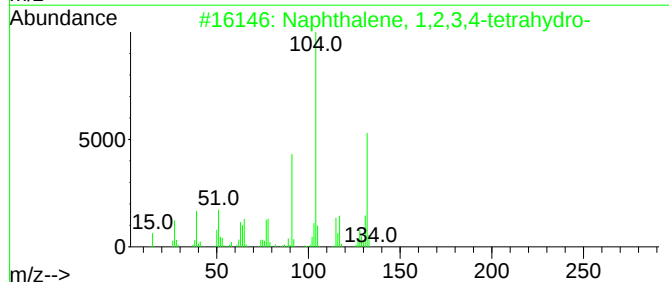
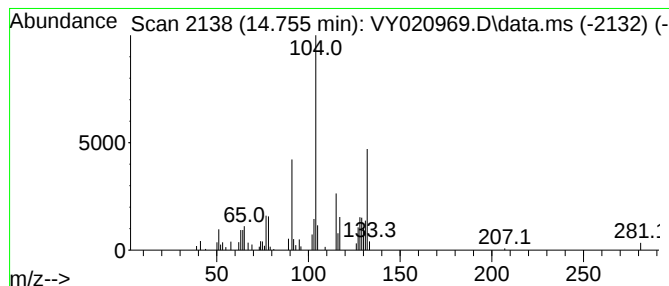
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.755	13.59 ug/l	24465	1,4-Dichlorobenzene-d4	13.352

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Benzene, cyclobutyl-	132	C10H12	004392-30-7	47
3	Indan, epoxide	132	C9H8O	000768-22-9	46
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	46
5	2-Methyl-4-phenylthiolane, 1,1-d...	210	C11H14O2S	058121-30-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
Data File : VY020969.D
Acq On : 29 Jan 2025 17:05
Operator : SY/MD
Sample : Q1220-01
Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

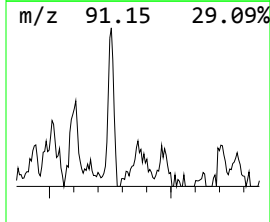
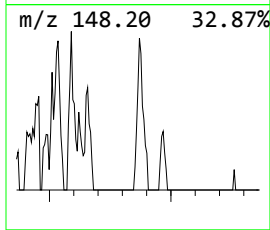
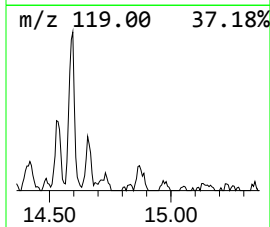
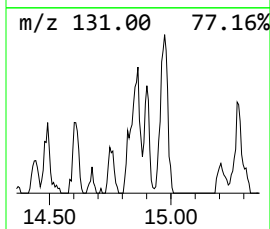
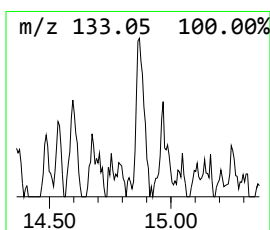
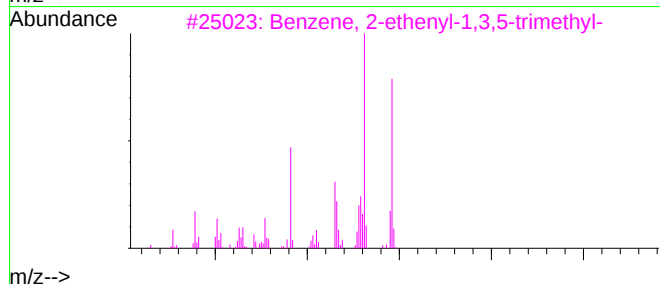
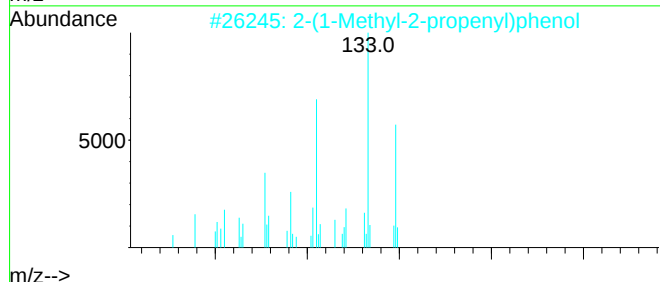
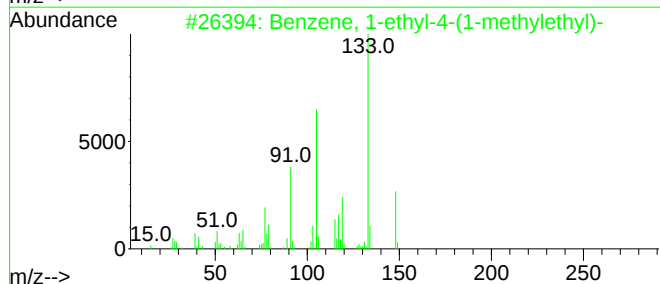
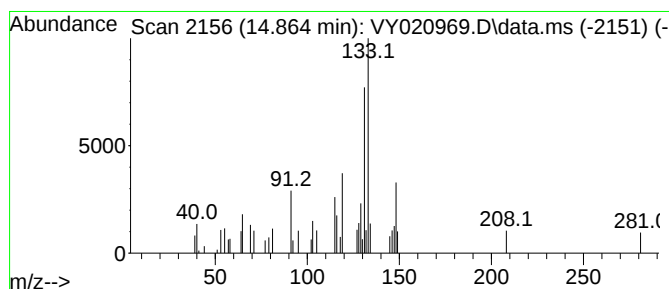
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown14.864 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.864	7.64 ug/l	13745	1,4-Dichlorobenzene-d4	13.352

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	38
2			2-(1-Methyl-2-propenyl)phenol	148	C10H12O	1000432-85-2	38
3			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	38
4			Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	35
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
Data File : VY020969.D
Acq On : 29 Jan 2025 17:05
Operator : SY/MD
Sample : Q1220-01
Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.M
Quant Title : SW846 8260

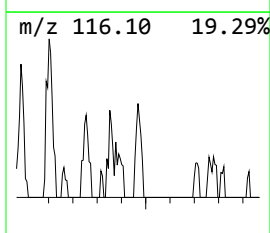
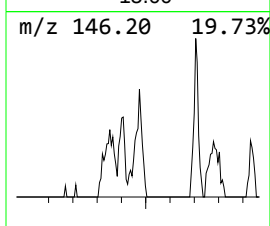
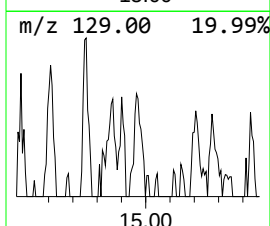
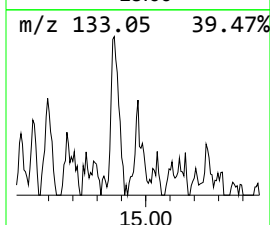
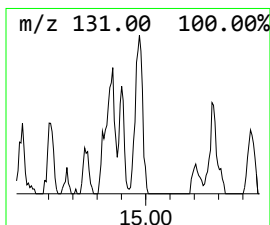
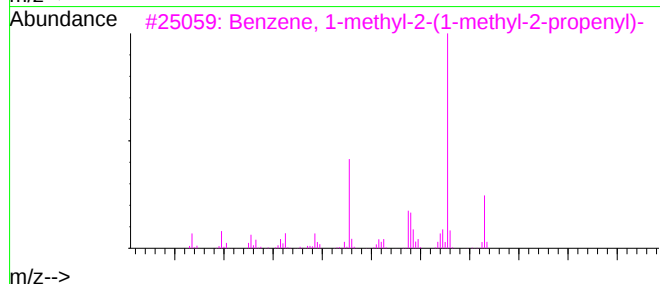
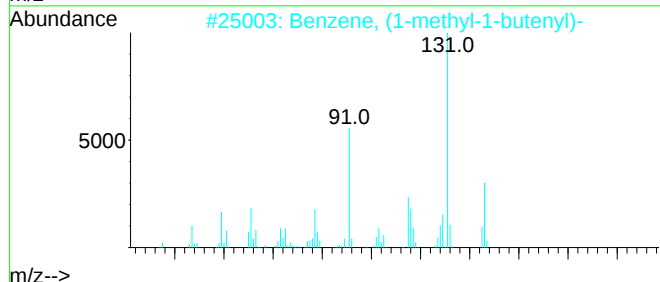
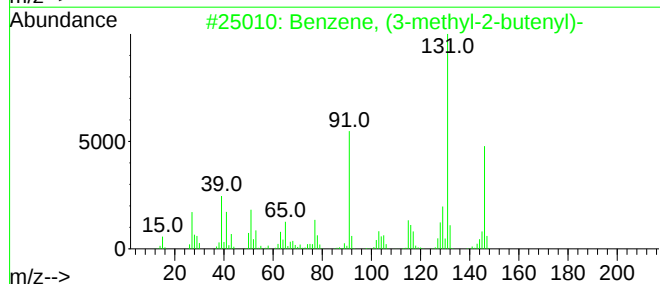
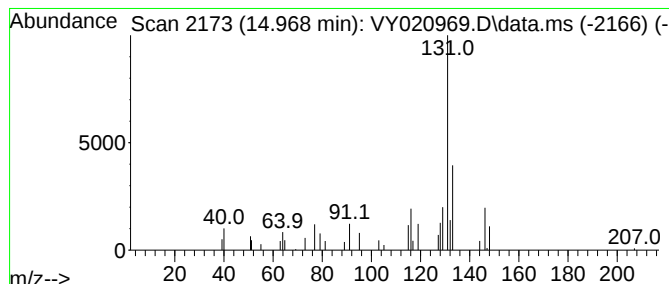
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, (3-methyl-2-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.968	8.35 ug/l	15035	1,4-Dichlorobenzene-d4	13.352

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	58
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	53
3	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	50
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	50
5	Benzene, (3-chloro-1-methyl-1-pr...	166	C10H11Cl	1000327-43-7	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY012925\
Data File : VY020969.D
Acq On : 29 Jan 2025 17:05
Operator : SY/MD
Sample : Q1220-01
Misc : 5.03g/5.0mL/MSVOA_Y/SOIL/B
ALS Vial : 4 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
TR-06-01292025

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y012425S.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
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unknown13.676	13.676	8.4	ug/l	15171	4	13.352	89991	50.0
Benzene, 1-ethy...	13.895	9.8	ug/l	17625	4	13.352	89991	50.0
2,4-Dimethylsty...	14.005	10.7	ug/l	19275	4	13.352	89991	50.0
Benzene, 2-ethe...	14.486	9.2	ug/l	16630	4	13.352	89991	50.0
1H-Indene, 2,3-...	14.602	22.6	ug/l	40766	4	13.352	89991	50.0
Naphthalene, 1,...	14.755	13.6	ug/l	24465	4	13.352	89991	50.0
unknown14.864	14.864	7.6	ug/l	13745	4	13.352	89991	50.0
Benzene, (3-met...	14.968	8.3	ug/l	15035	4	13.352	89991	50.0