

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
 Data File : VY014722.D
 Acq On : 20 Jul 2023 15:41
 Operator : SY/MD
 Sample : 03664-01
 Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 AUD-1021-1020-COMP

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\82Y071323S.M

Title : SW846 8260

Signal : TIC: VY014722.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.216	59	65	76	rBV	66910	112161	1.24%	0.214%
2	2.381	88	92	100	rVB	153432	249777	2.77%	0.475%
3	3.942	337	348	378	rVV	3273073	9026516	100.00%	17.183%
4	4.551	438	448	463	rVB3	74720	240724	2.67%	0.458%
5	4.710	463	474	483	rBV3	39676	150445	1.67%	0.286%
6	4.874	485	501	507	rVV	300824	1193504	13.22%	2.272%
7	4.948	508	513	531	rVB	226937	780020	8.64%	1.485%
8	5.740	628	643	661	rBV	701925	2223648	24.63%	4.233%
9	6.697	787	800	809	rBV2	153811	485820	5.38%	0.925%
10	6.978	830	846	859	rBV2	115593	342774	3.80%	0.653%
11	7.526	925	936	948	rVB2	37306	108572	1.20%	0.207%
12	7.715	955	967	969	rBV	187293	383968	4.25%	0.731%
13	7.770	970	976	986	rVV2	506308	1659030	18.38%	3.158%
14	7.868	986	992	1003	rVB	206034	520885	5.77%	0.992%
15	7.996	1003	1013	1028	rBV	527344	1244847	13.79%	2.370%
16	8.142	1028	1037	1047	rVB	189991	447065	4.95%	0.851%
17	8.270	1047	1058	1061	rBV3	82224	219065	2.43%	0.417%
18	8.325	1061	1067	1074	rVV	158880	430309	4.77%	0.819%
19	8.398	1074	1079	1093	rVB3	73644	224403	2.49%	0.427%
20	8.545	1093	1103	1111	rBV	409993	832406	9.22%	1.585%
21	8.685	1120	1126	1135	rVB	473639	921872	10.21%	1.755%
22	9.136	1193	1200	1203	rBV2	83939	167364	1.85%	0.319%
23	9.185	1203	1208	1213	rVV2	138420	285995	3.17%	0.544%
24	9.258	1213	1220	1233	rVV4	156562	476389	5.28%	0.907%
25	9.611	1271	1278	1289	rVB	203425	406106	4.50%	0.773%
26	9.746	1289	1300	1307	rBV	207198	458943	5.08%	0.874%
27	9.959	1325	1335	1346	rBV2	86628	222353	2.46%	0.423%
28	10.069	1346	1353	1357	rBV	113593	208743	2.31%	0.397%
29	10.178	1364	1371	1376	rBV	794756	1417957	15.71%	2.699%
30	10.239	1376	1381	1394	rVB	3650514	6377025	70.65%	12.139%
31	10.367	1394	1402	1408	rVB2	90601	176090	1.95%	0.335%
32	10.434	1408	1413	1416	rBV	62695	107876	1.20%	0.205%
33	10.483	1416	1421	1425	rBV4	42535	91971	1.02%	0.175%
34	10.611	1436	1442	1450	rVB5	69401	171964	1.91%	0.327%
35	10.709	1451	1458	1464	rVB3	43712	97090	1.08%	0.185%
36	10.928	1486	1494	1498	rBV	161795	300079	3.32%	0.571%

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

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37	11.032	1508	1511	1516	rVB	64567	91088	1.01%	0.173%
38	11.087	1516	1520	1526	rBV	82804	132401	1.47%	0.252%
39	11.203	1534	1539	1543	rBV3	57454	93415	1.03%	0.178%
40	11.282	1543	1552	1562	rVB3	206836	558806	6.19%	1.064%
41	11.379	1562	1568	1573	rBV	151014	255471	2.83%	0.486%
42	11.489	1579	1586	1592	rBV	680420	1125137	12.46%	2.142%
43	11.593	1595	1603	1606	rVV	235167	447248	4.95%	0.851%
44	11.623	1606	1608	1612	rVV	106706	148772	1.65%	0.283%
45	11.702	1612	1621	1631	rVV3	941843	2536648	28.10%	4.829%
46	11.776	1631	1633	1642	rVB2	88605	130736	1.45%	0.249%
47	12.026	1664	1674	1682	rVB3	468194	975728	10.81%	1.857%
48	12.129	1682	1691	1698	rBV3	152978	328842	3.64%	0.626%
49	12.202	1698	1703	1707	rVV2	147251	232775	2.58%	0.443%
50	12.245	1707	1710	1714	rVB3	66122	96270	1.07%	0.183%
51	12.324	1719	1723	1727	rBV	104688	158219	1.75%	0.301%
52	12.477	1742	1748	1753	rVB2	518808	919166	10.18%	1.750%
53	12.593	1759	1767	1771	rBV	227538	370298	4.10%	0.705%
54	12.666	1776	1779	1784	rVB	84873	128477	1.42%	0.245%
55	12.733	1784	1790	1793	rBV	129203	220555	2.44%	0.420%
56	12.763	1793	1795	1798	rVV	87255	118181	1.31%	0.225%
57	12.806	1798	1802	1815	rVB2	244595	506093	5.61%	0.963%
58	12.964	1823	1828	1832	rBV2	56883	95850	1.06%	0.182%
59	13.117	1848	1853	1858	rVV	187328	295286	3.27%	0.562%
60	13.275	1873	1879	1883	rBV4	76533	148470	1.64%	0.283%
61	13.367	1889	1894	1896	rBV3	89759	152497	1.69%	0.290%
62	13.422	1898	1903	1908	rVV	493896	743458	8.24%	1.415%
63	13.525	1915	1920	1926	rVB	1025189	1508315	16.71%	2.871%
64	13.623	1932	1936	1940	rVV	199142	333253	3.69%	0.634%
65	13.666	1940	1943	1952	rVB3	133853	309447	3.43%	0.589%
66	13.751	1952	1957	1961	rBV	199280	296162	3.28%	0.564%
67	13.824	1966	1969	1975	rVV	88624	148680	1.65%	0.283%
68	13.885	1975	1979	1981	rVV	66321	101632	1.13%	0.193%
69	13.909	1981	1983	1988	rVB3	83677	110666	1.23%	0.211%
70	13.970	1988	1993	1998	rBV	318643	462540	5.12%	0.880%
71	14.019	1998	2001	2006	rVB2	53005	100713	1.12%	0.192%
72	14.074	2006	2010	2016	rBV	174561	272765	3.02%	0.519%
73	14.141	2016	2021	2027	rBV7	35453	91382	1.01%	0.174%
74	14.245	2028	2038	2039	rBV5	69421	158029	1.75%	0.301%
75	14.293	2043	2046	2056	rVB2	175708	392134	4.34%	0.746%
76	14.379	2056	2060	2064	rBV2	159474	249810	2.77%	0.476%

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77	14.562	2086	2090	2095	rVV	208552	351114	3.89%	0.668%
78	14.611	2095	2098	2102	rVB	111010	147568	1.63%	0.281%
79	14.684	2103	2110	2115	rBV3	325080	688301	7.63%	1.310%
80	14.732	2115	2118	2124	rVB2	124713	209316	2.32%	0.398%
81	14.940	2148	2152	2156	rVV2	138246	227115	2.52%	0.432%
82	15.056	2164	2171	2176	rBV4	119956	250974	2.78%	0.478%
83	15.232	2191	2200	2206	rBV	318460	619136	6.86%	1.179%
84	15.342	2213	2218	2223	rBV3	100092	175746	1.95%	0.335%
85	15.427	2230	2232	2241	rVB4	57822	114255	1.27%	0.217%
86	15.525	2241	2248	2254	rBV	111872	246822	2.73%	0.470%
87	15.690	2266	2275	2278	rBV4	61101	168222	1.86%	0.320%
88	15.885	2302	2307	2313	rVB6	52696	108900	1.21%	0.207%
89	16.287	2367	2373	2384	rVB	316011	624666	6.92%	1.189%
90	16.482	2399	2405	2419	rVB	123515	290615	3.22%	0.553%

Sum of corrected areas: 52531921

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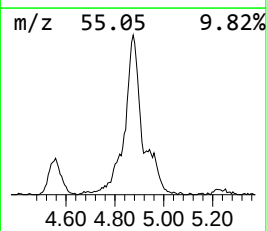
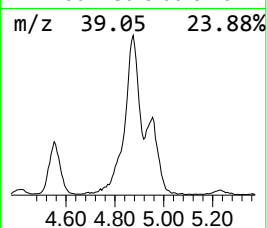
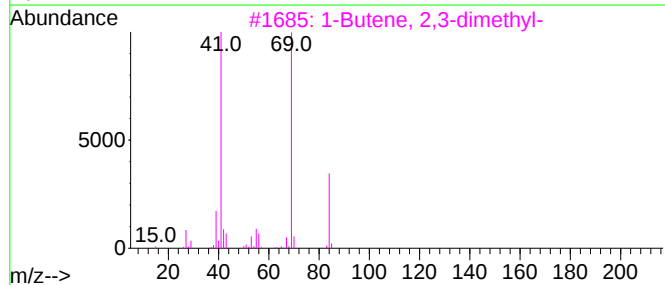
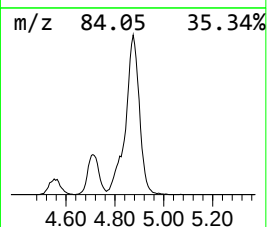
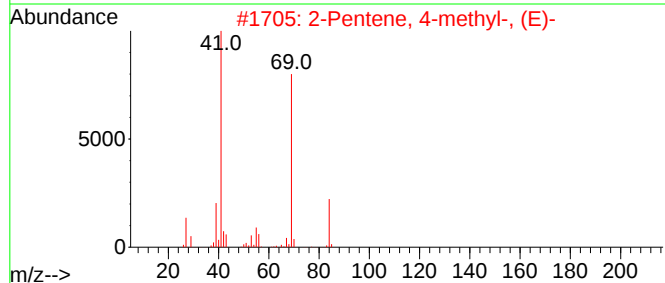
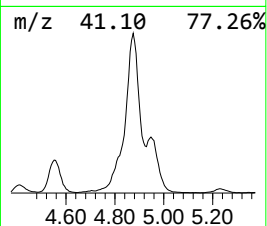
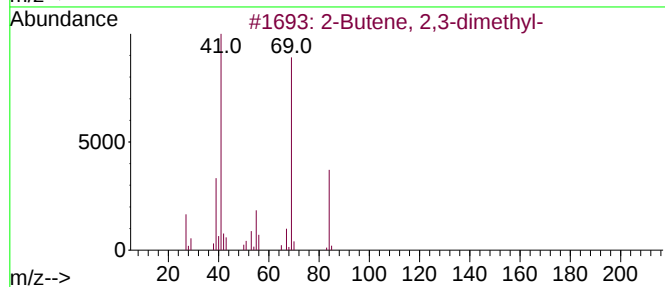
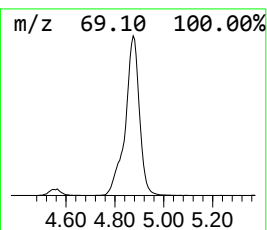
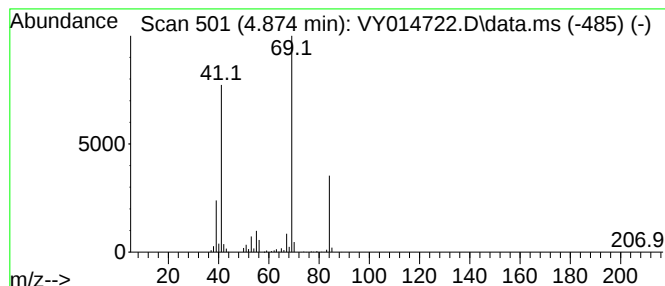
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071323S.M
Quant Title : SW846 8260

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

Peak Number 1 2-Butene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.874	35.97 ug/l	1193500	Pentafluorobenzene	7.789

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	91	
2	2-Pentene, 4-methyl-, (E)-	84	C6H12	000674-76-0	91	
3	1-Butene, 2,3-dimethyl-	84	C6H12	000563-78-0	90	
4	Cyclopropane, 1,1,2-trimethyl-	84	C6H12	004127-45-1	86	
5	2-Pentene, 2-methyl-	84	C6H12	000625-27-4	86	



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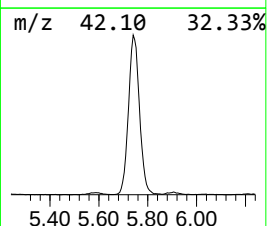
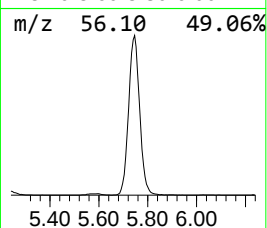
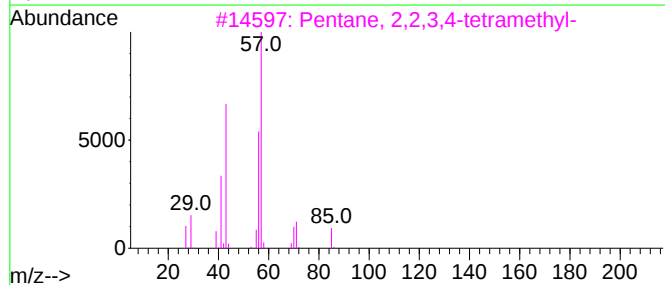
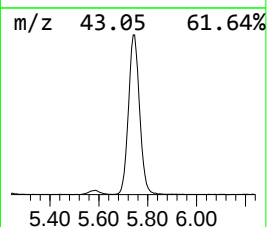
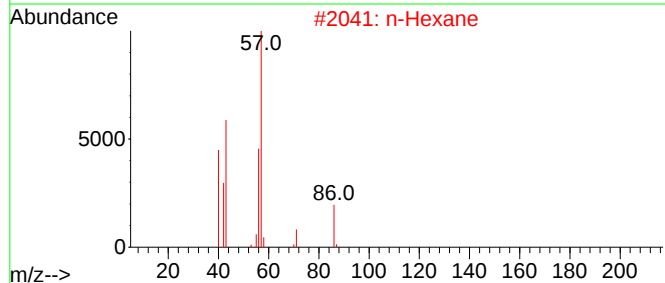
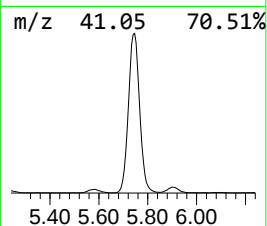
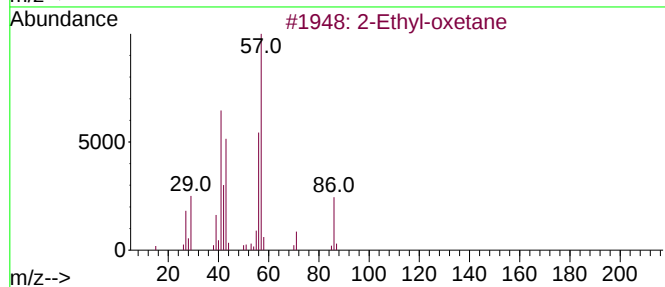
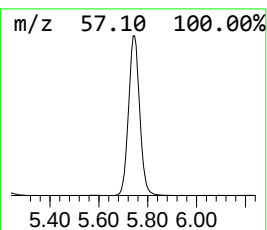
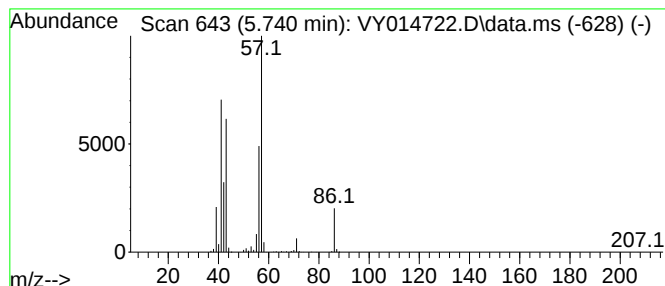
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Ethyl-oxetane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.740	67.02 ug/l	2223650	Pentafluorobenzene	7.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethyl-oxetane	86	C5H10O	1010386-40-2	90
2			n-Hexane	86	C6H14	000110-54-3	87
3			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	40
4			Hydroperoxide, hexyl	118	C6H14O2	004312-76-9	38
5			Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	38



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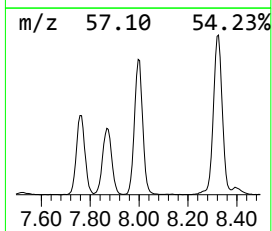
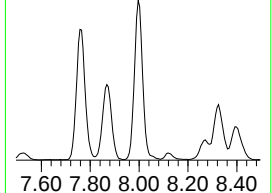
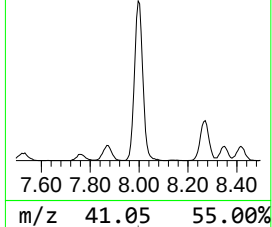
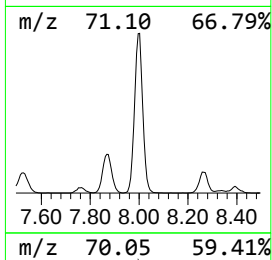
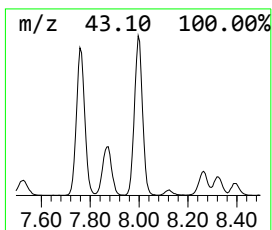
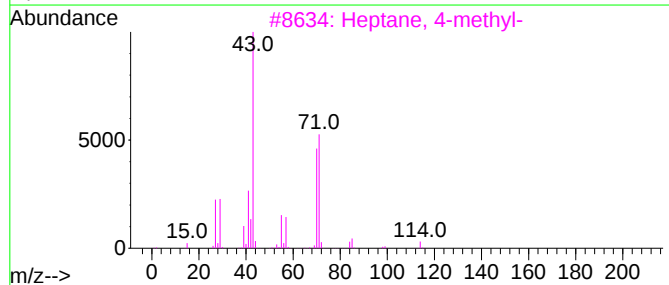
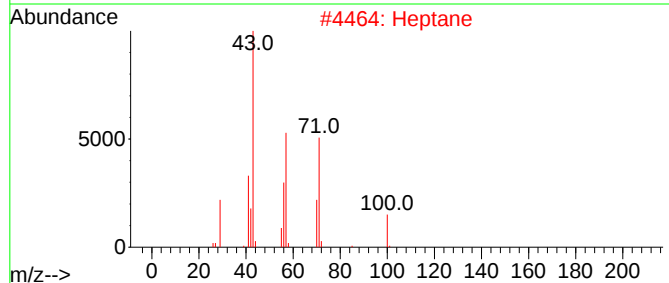
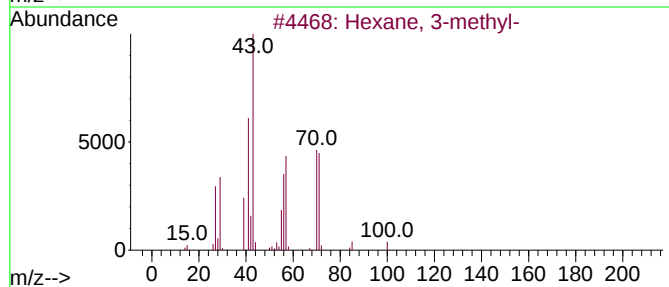
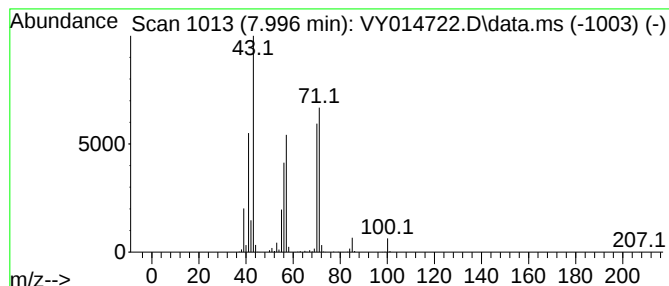
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 3 Hexane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.996	37.52 ug/l	1244850	Pentafluorobenzene	7.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	53



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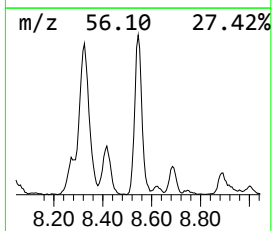
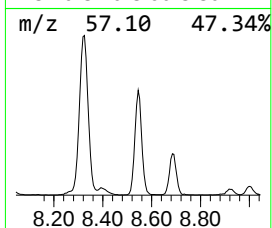
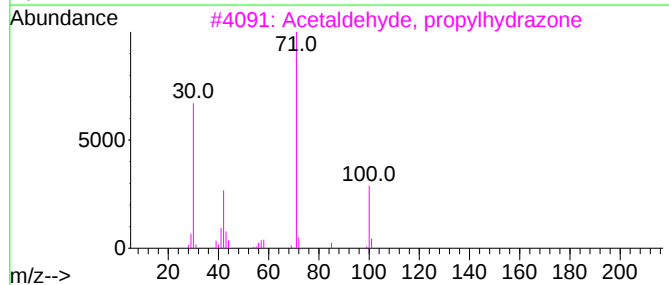
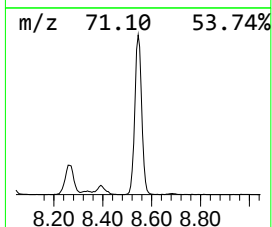
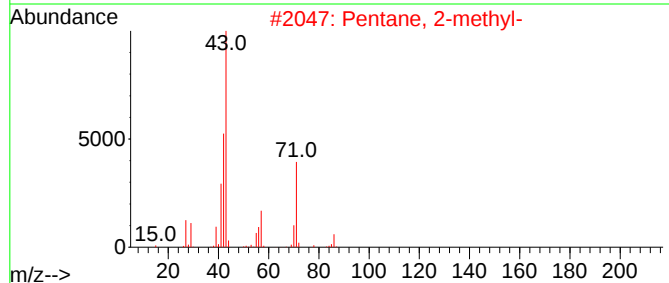
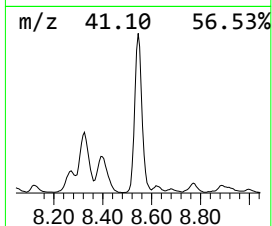
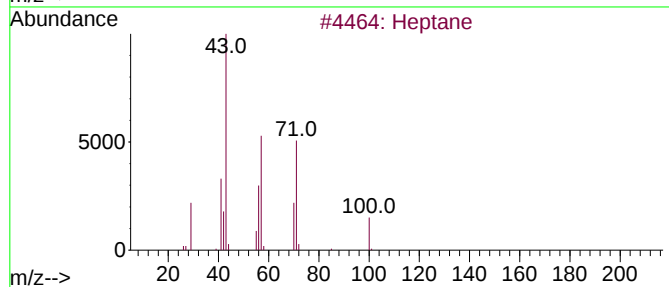
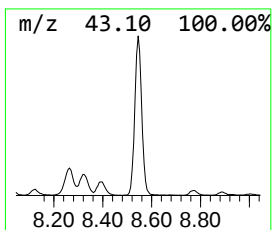
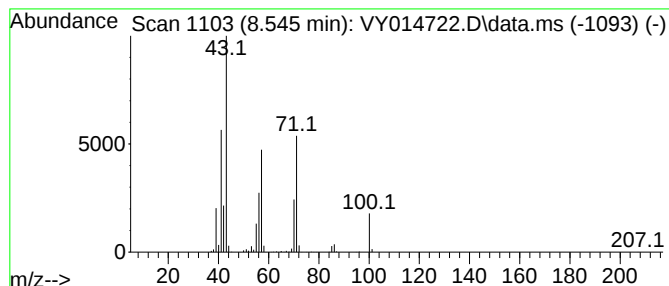
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Quant Title : SW846 8260

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

Peak Number 4 Heptane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.545	45.15 ug/l	832406	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	90
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	47
3			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	38
4			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	38
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
Acq On : 20 Jul 2023 15:41
Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
AUD-1021-1020-COMP

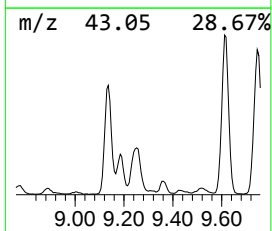
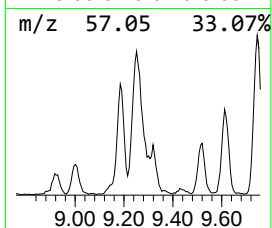
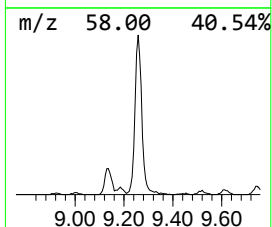
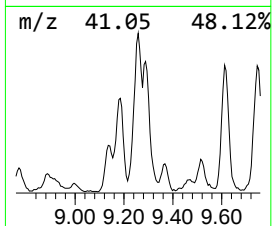
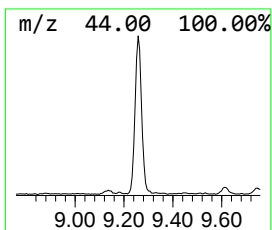
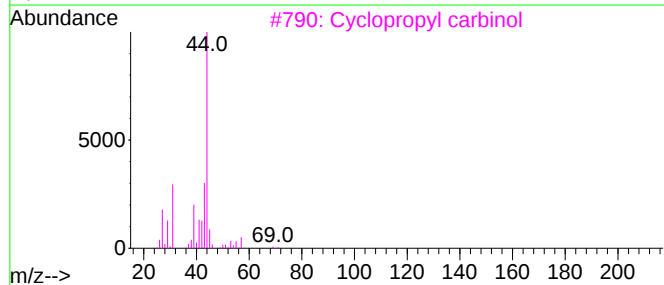
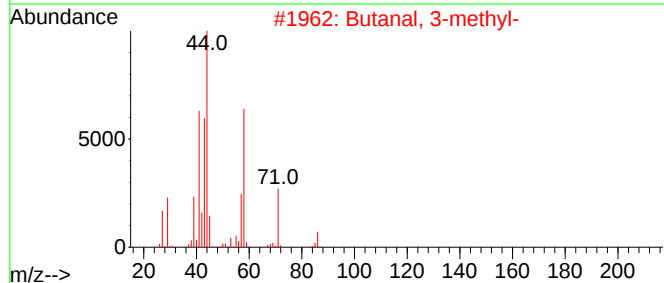
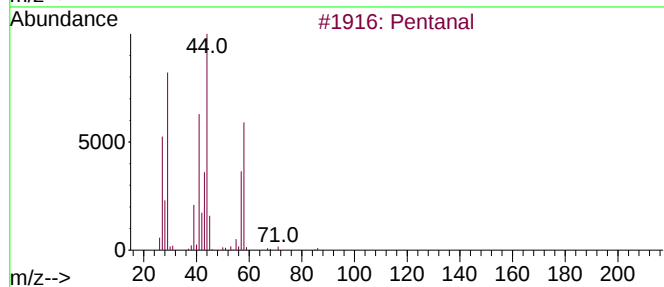
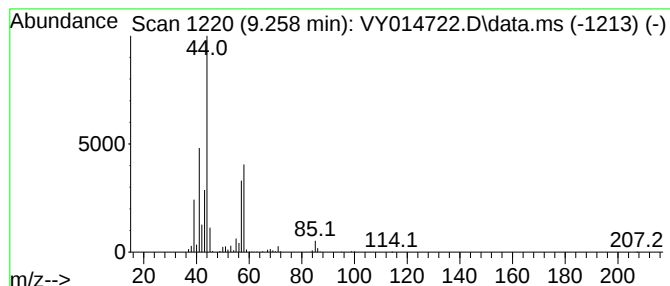
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Quant Title : SW846 8260

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

Peak Number 5 Pentanal Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.258	25.84 ug/l	476389	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanal	86	C5H10O	000110-62-3	78
2			Butanal, 3-methyl-	86	C5H10O	000590-86-3	64
3			Cyclopropyl carbinol	72	C4H8O	002516-33-8	33
4			Silacyclobutane	72	C3H8Si	1000434-34-1	33
5			Glutaraldehyde	100	C5H8O2	000111-30-8	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
Acq On : 20 Jul 2023 15:41
Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
AUD-1021-1020-COMP

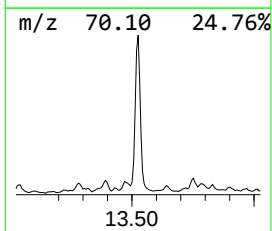
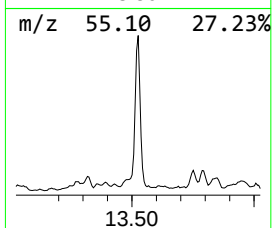
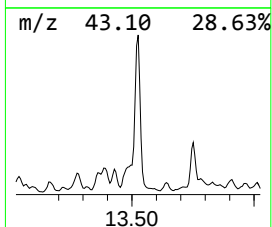
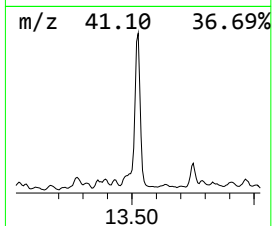
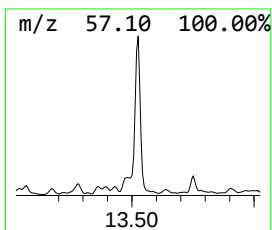
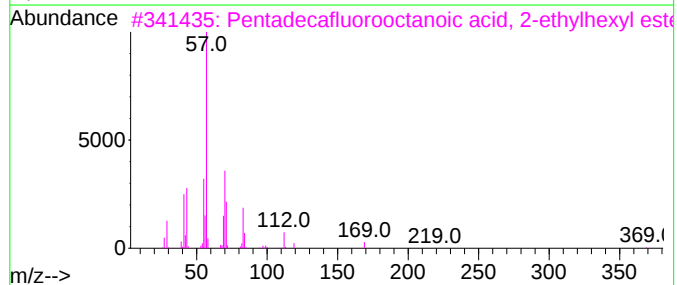
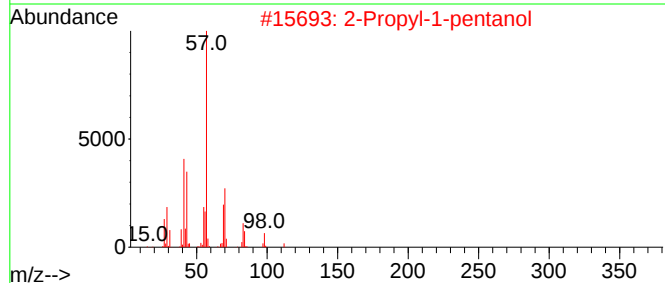
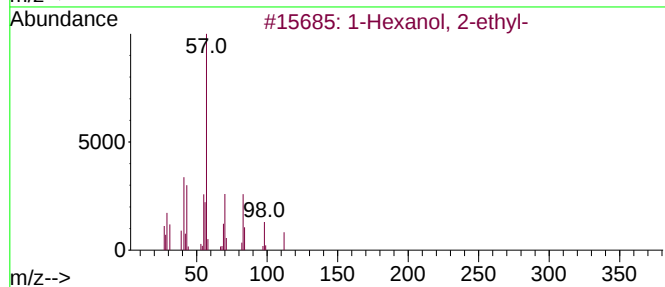
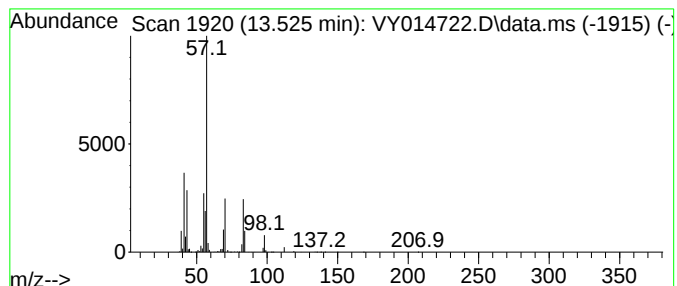
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Quant Title : SW846 8260

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

Peak Number 6 1-Hexanol, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.525	101.44 ug/l	1508320	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	74
2			2-Propyl-1-pentanol	130	C8H18O	058175-57-8	59
3			Pentadecafluorooctanoic acid, 2-...	526	C16H17F15O2	1000406-80-9	59
4			Carbonic acid, isobutyl 2-ethylh...	230	C13H26O3	1000383-13-3	50
5			Trichloroacetic acid, 2-ethylhex...	274	C10H17Cl3O2	016397-79-8	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
Acq On : 20 Jul 2023 15:41
Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
AUD-1021-1020-COMP

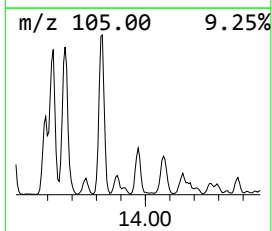
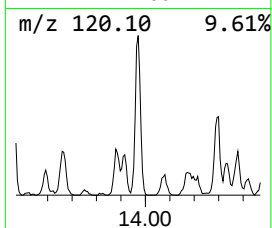
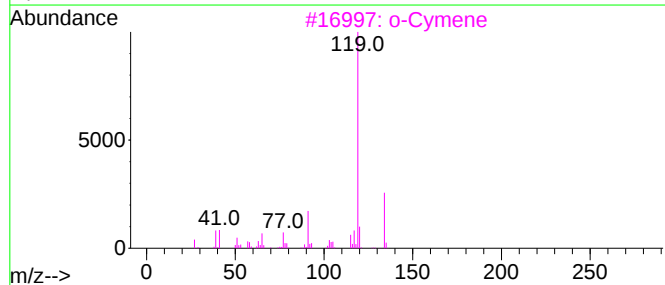
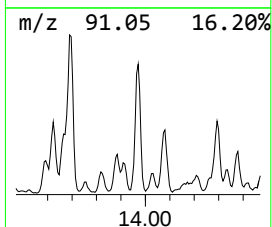
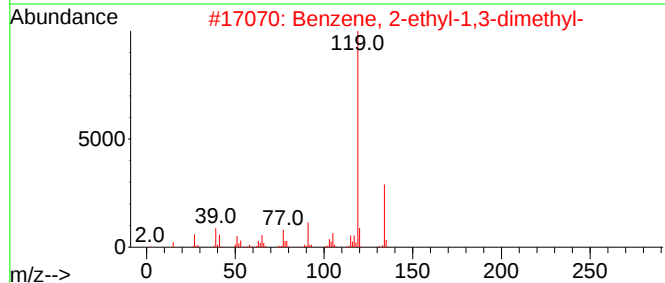
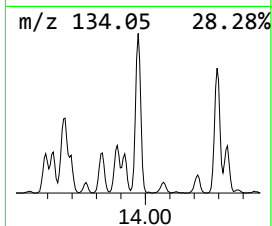
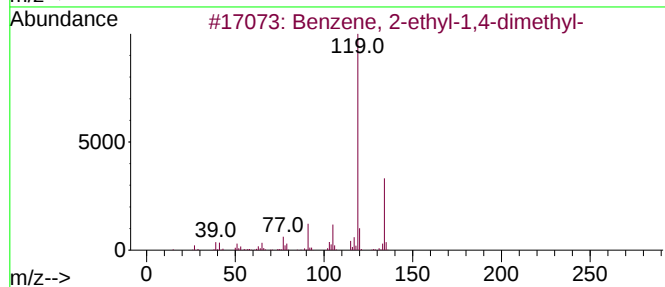
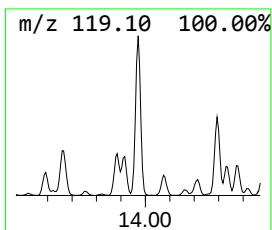
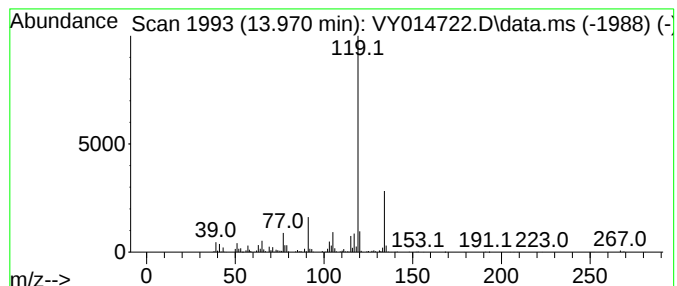
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.970	31.11 ug/l	462540	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
Acq On : 20 Jul 2023 15:41
Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
AUD-1021-1020-COMP

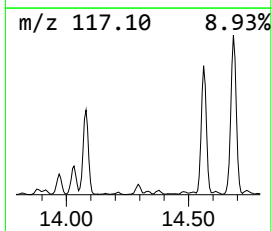
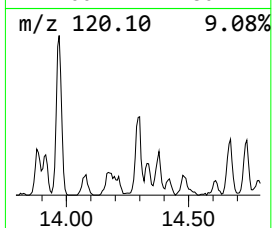
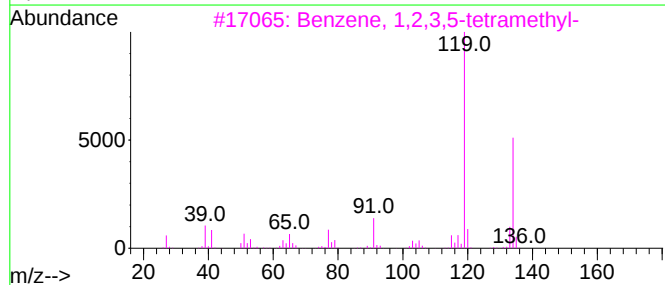
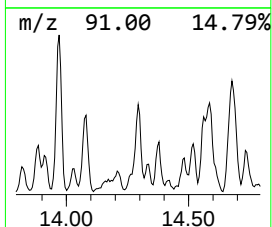
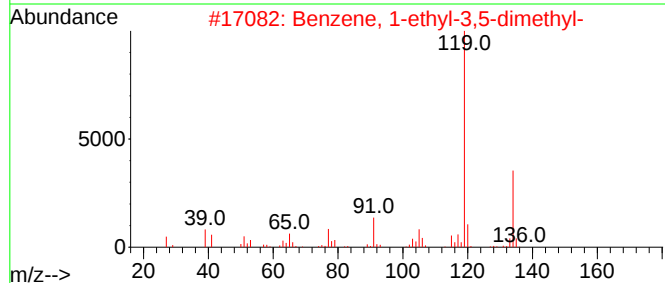
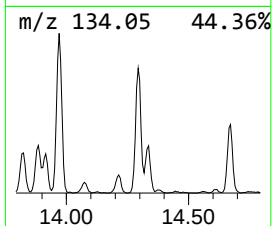
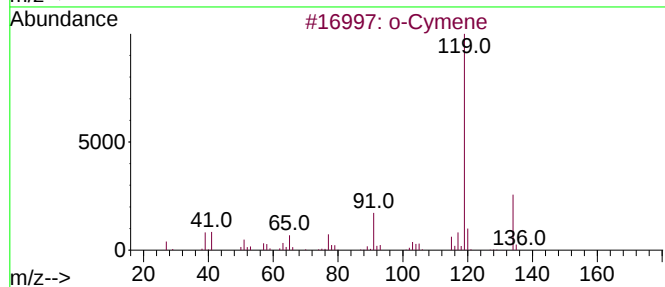
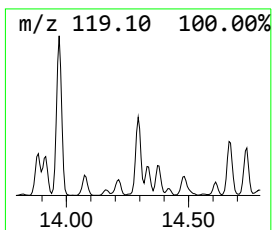
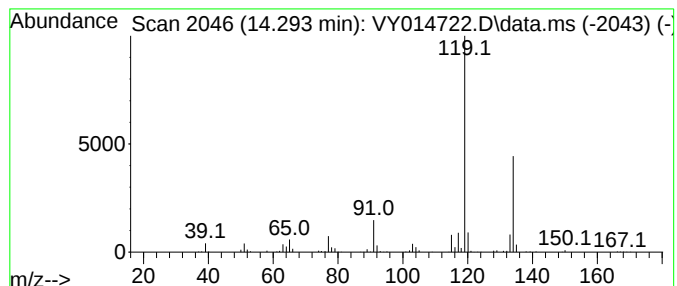
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Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.294	26.37 ug/l	392134	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
Acq On : 20 Jul 2023 15:41
Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
AUD-1021-1020-COMP

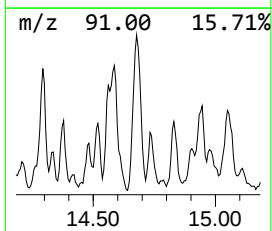
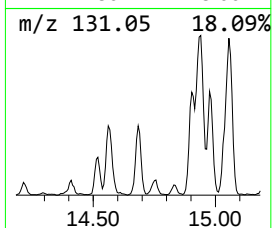
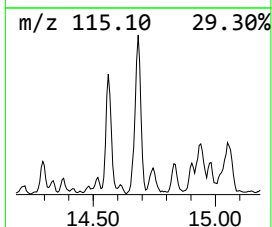
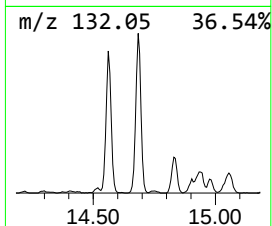
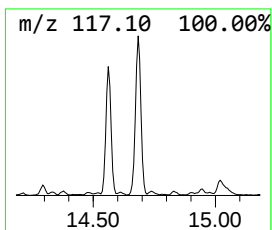
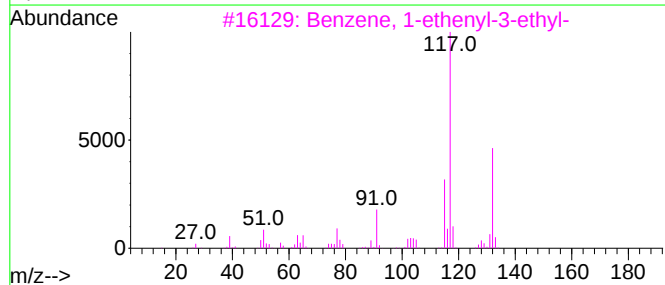
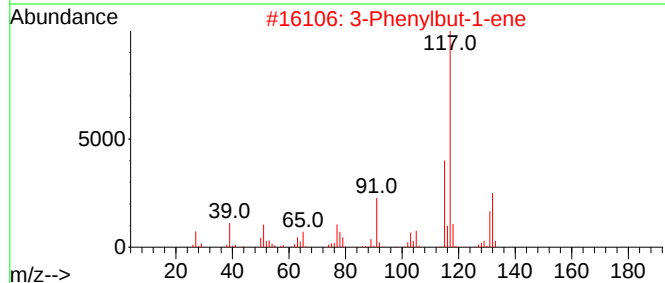
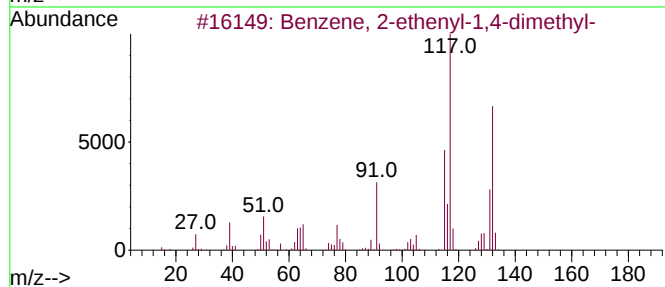
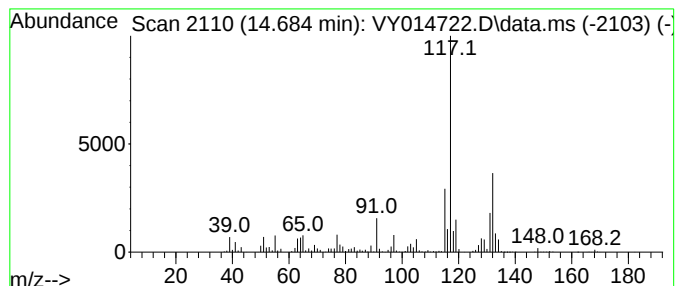
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	46.29 ug/l	688301	1,4-Dichlorobenzene-d4	13.422

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
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Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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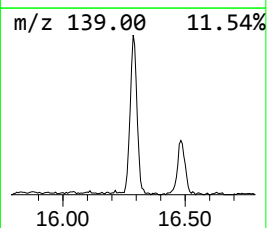
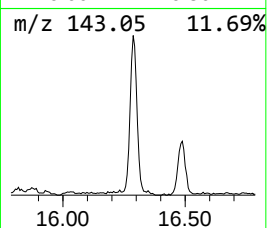
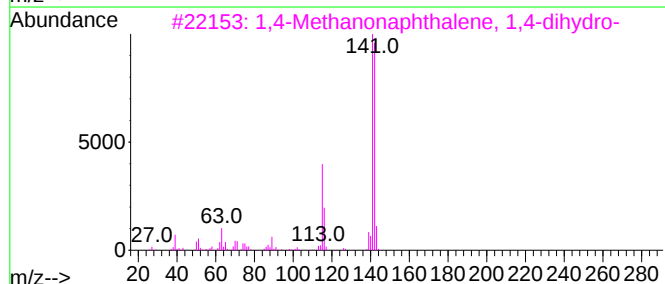
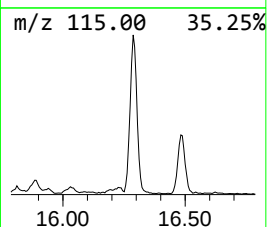
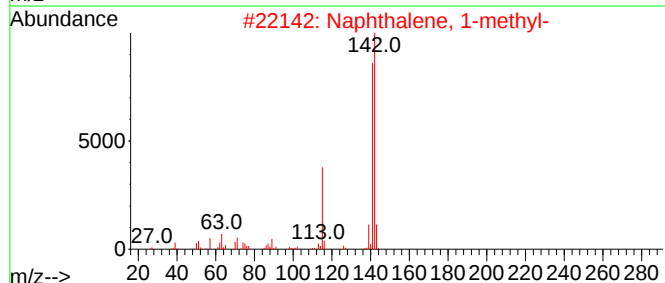
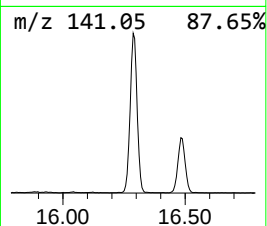
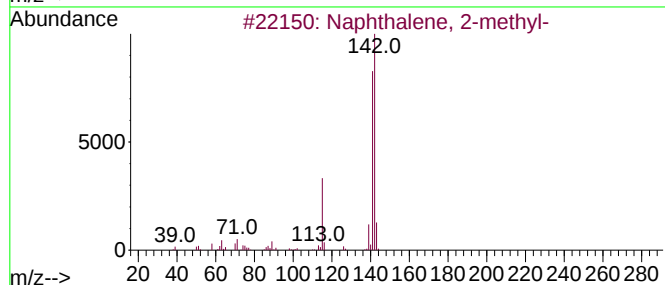
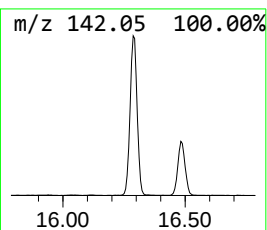
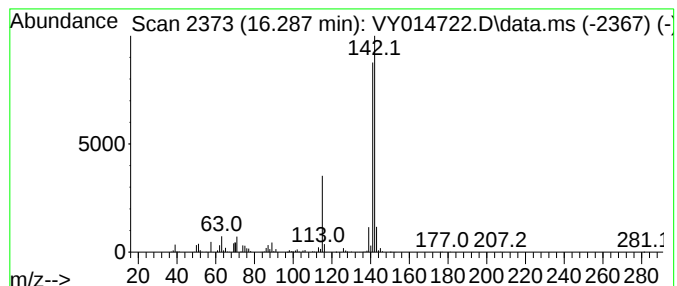
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.287	42.01 ug/l	624666	1,4-Dichlorobenzene-d4	13.422

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	94
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY072023\
Data File : VY014722.D
Acq On : 20 Jul 2023 15:41
Operator : SY/MD
Sample : 03664-01
Misc : 5.02g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
AUD-1021-1020-COMP

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y071323S.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
2-Butene, 2,3-d...	4.874	36.0	ug/l	1193500	1	7.789	1659030	50.0
2-Ethyl-oxetane	5.740	67.0	ug/l	2223650	1	7.789	1659030	50.0
Hexane, 3-methyl-	7.996	37.5	ug/l	1244850	1	7.789	1659030	50.0
Heptane	8.545	45.1	ug/l	832406	2	8.685	921872	50.0
Pentanal	9.258	25.8	ug/l	476389	2	8.685	921872	50.0
1-Hexanol, 2-et...	13.525	101.4	ug/l	1508320	4	13.422	743458	50.0
Benzene, 2-ethy...	13.970	31.1	ug/l	462540	4	13.422	743458	50.0
o-Cymene	14.294	26.4	ug/l	392134	4	13.422	743458	50.0
Benzene, 2-ethe...	14.684	46.3	ug/l	688301	4	13.422	743458	50.0
Naphthalene, 2-...	16.287	42.0	ug/l	624666	4	13.422	743458	50.0