

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
 Data File : VY006784.D
 Acq On : 17 Nov 2021 13:33
 Operator : SY/MD
 Sample : M4693-03
 Misc : 5.49G/5ML/MSVOA_Y/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 UT-COMP-02

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\82Y111521S.M
 Title : SW846 8260

Signal : TIC: VY006784.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.869	4	8	18	rVB3	7183	13458	1.41%	0.153%
2	2.229	62	67	79	rBV4	4558	12740	1.34%	0.145%
3	2.534	110	117	125	rBV8	3600	11565	1.21%	0.131%
4	3.954	338	350	363	rBV2	39154	112446	11.80%	1.278%
5	4.204	381	391	402	rBV	4354	11768	1.23%	0.134%
6	4.729	465	477	490	rBV2	11259	37296	3.91%	0.424%
7	6.997	839	849	861	rBV	9467	27078	2.84%	0.308%
8	7.728	958	969	975	rBV	84543	204976	21.51%	2.329%
9	7.801	975	981	990	rVB	142818	320501	33.63%	3.642%
10	8.155	1028	1039	1050	rVB	101843	237291	24.90%	2.696%
11	8.697	1118	1128	1138	rBV	231213	449357	47.15%	5.106%
12	9.191	1196	1209	1215	rVV3	19532	50883	5.34%	0.578%
13	9.472	1244	1255	1260	rBV3	5958	15130	1.59%	0.172%
14	9.612	1272	1278	1287	rVB3	10468	22117	2.32%	0.251%
15	9.764	1292	1303	1307	rBV2	8962	19347	2.03%	0.220%
16	9.825	1307	1313	1317	rBV2	22692	41390	4.34%	0.470%
17	9.959	1325	1335	1348	rVB5	23135	74273	7.79%	0.844%
18	10.075	1348	1354	1364	rBV6	8052	22499	2.36%	0.256%
19	10.185	1364	1372	1388	rBV	449020	864897	90.75%	9.828%
20	10.368	1394	1402	1410	rVB3	31015	90131	9.46%	1.024%
21	10.526	1423	1428	1434	rVB2	61325	107538	11.28%	1.222%
22	10.624	1434	1444	1449	rBV2	33714	69634	7.31%	0.791%
23	10.715	1455	1459	1464	rVB4	10529	17090	1.79%	0.194%
24	10.807	1469	1474	1481	rVB	28659	43705	4.59%	0.497%
25	10.904	1481	1490	1497	rBV	13837	30034	3.15%	0.341%
26	10.978	1497	1502	1504	rBV3	10553	18164	1.91%	0.206%
27	11.008	1504	1507	1508	rVV	18088	23119	2.43%	0.263%
28	11.038	1508	1512	1517	rVV	64783	113243	11.88%	1.287%
29	11.093	1517	1521	1526	rVV	77943	133435	14.00%	1.516%
30	11.148	1526	1530	1534	rVV	14594	22005	2.31%	0.250%
31	11.209	1534	1540	1544	rVV2	21868	39062	4.10%	0.444%
32	11.295	1544	1554	1563	rVB4	43038	110626	11.61%	1.257%
33	11.386	1563	1569	1573	rBV	24390	39811	4.18%	0.452%
34	11.496	1580	1587	1596	rBV	328018	534878	56.12%	6.078%
35	11.593	1596	1603	1607	rBV3	32294	65835	6.91%	0.748%
36	11.630	1607	1609	1613	rVB2	14480	16252	1.71%	0.185%

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37	11.703	1613	1621	1625	rBV4	47401	125534	13.17%	1.427%
38	11.904	1648	1654	1663	rVB4	10421	22144	2.32%	0.252%
39	12.032	1663	1675	1683	rBV6	49032	166186	17.44%	1.888%
40	12.124	1687	1690	1693	rVB3	8784	11829	1.24%	0.134%
41	12.166	1693	1697	1699	rBV3	7956	11698	1.23%	0.133%
42	12.203	1699	1703	1707	rVV	25371	41822	4.39%	0.475%
43	12.246	1707	1710	1715	rVB5	14043	24658	2.59%	0.280%
44	12.337	1720	1725	1731	rVB2	110094	177202	18.59%	2.014%
45	12.489	1744	1750	1758	rBV2	540634	953038	100.00%	10.830%
46	12.587	1758	1766	1771	rVB3	19760	40165	4.21%	0.456%
47	12.648	1771	1776	1784	rVB3	20206	48025	5.04%	0.546%
48	12.739	1784	1791	1794	rBV	37270	65137	6.83%	0.740%
49	12.813	1798	1803	1813	rVV3	75657	152011	15.95%	1.727%
50	12.898	1813	1817	1822	rVB2	14818	26039	2.73%	0.296%
51	12.977	1822	1830	1836	rBV	27351	53033	5.56%	0.603%
52	13.044	1837	1841	1849	rVB5	14741	30009	3.15%	0.341%
53	13.123	1849	1854	1868	rVB2	78726	168225	17.65%	1.912%
54	13.251	1868	1875	1878	rBV6	13811	24349	2.55%	0.277%
55	13.319	1878	1886	1888	rVV4	21787	39703	4.17%	0.451%
56	13.367	1891	1894	1899	rVV	72332	111177	11.67%	1.263%
57	13.428	1899	1904	1912	rVV2	237222	418802	43.94%	4.759%
58	13.489	1912	1914	1919	rVB3	22875	31032	3.26%	0.353%
59	13.629	1933	1937	1940	rVV	58526	87505	9.18%	0.994%
60	13.666	1940	1943	1952	rVB2	33797	58365	6.12%	0.663%
61	13.751	1952	1957	1962	rBV3	33146	55246	5.80%	0.628%
62	13.825	1962	1969	1973	rVB2	13815	23679	2.48%	0.269%
63	13.886	1975	1979	1982	rBV	15760	25210	2.65%	0.286%
64	13.916	1982	1984	1988	rVV	18023	20858	2.19%	0.237%
65	13.971	1988	1993	1999	rVB2	48785	80374	8.43%	0.913%
66	14.081	2006	2011	2016	rBV3	31306	50797	5.33%	0.577%
67	14.215	2028	2033	2037	rBV2	13573	22451	2.36%	0.255%
68	14.264	2037	2041	2043	rBV4	7710	12535	1.32%	0.142%
69	14.312	2043	2049	2057	rVV2	54165	133678	14.03%	1.519%
70	14.385	2057	2061	2064	rVV2	13211	17792	1.87%	0.202%
71	14.422	2064	2067	2071	rVB4	9786	12049	1.26%	0.137%
72	14.520	2079	2083	2086	rVB3	6811	10400	1.09%	0.118%
73	14.568	2086	2091	2095	rVV	16703	26835	2.82%	0.305%
74	14.617	2095	2099	2103	rVB	15514	23297	2.44%	0.265%
75	14.684	2103	2110	2116	rBV3	40410	96296	10.10%	1.094%
76	14.739	2116	2119	2125	rVB6	11350	20318	2.13%	0.231%

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77	14.837	2130	2135	2139	rBV2	12257	17867	1.87%	0.203%
78	14.946	2148	2153	2158	rBV2	46859	76610	8.04%	0.871%
79	15.056	2166	2171	2177	rVB2	11498	25098	2.63%	0.285%
80	15.135	2181	2184	2190	rVB6	5525	10351	1.09%	0.118%
81	15.239	2191	2201	2207	rBV	507760	851380	89.33%	9.675%
82	15.336	2213	2217	2223	rBV2	13563	24367	2.56%	0.277%
83	15.434	2230	2233	2240	rVB7	7644	13566	1.42%	0.154%
84	15.532	2240	2249	2253	rBV4	7158	15596	1.64%	0.177%
85	15.690	2267	2275	2277	rVV6	5031	12578	1.32%	0.143%
86	15.727	2278	2281	2286	rVB4	8259	14247	1.49%	0.162%
87	16.038	2324	2332	2337	rBV9	6761	16555	1.74%	0.188%
88	16.227	2356	2363	2368	rBV8	5299	11822	1.24%	0.134%
89	16.294	2368	2374	2381	rVB	48920	95557	10.03%	1.086%
90	16.489	2399	2406	2414	rBV	34363	71394	7.49%	0.811%

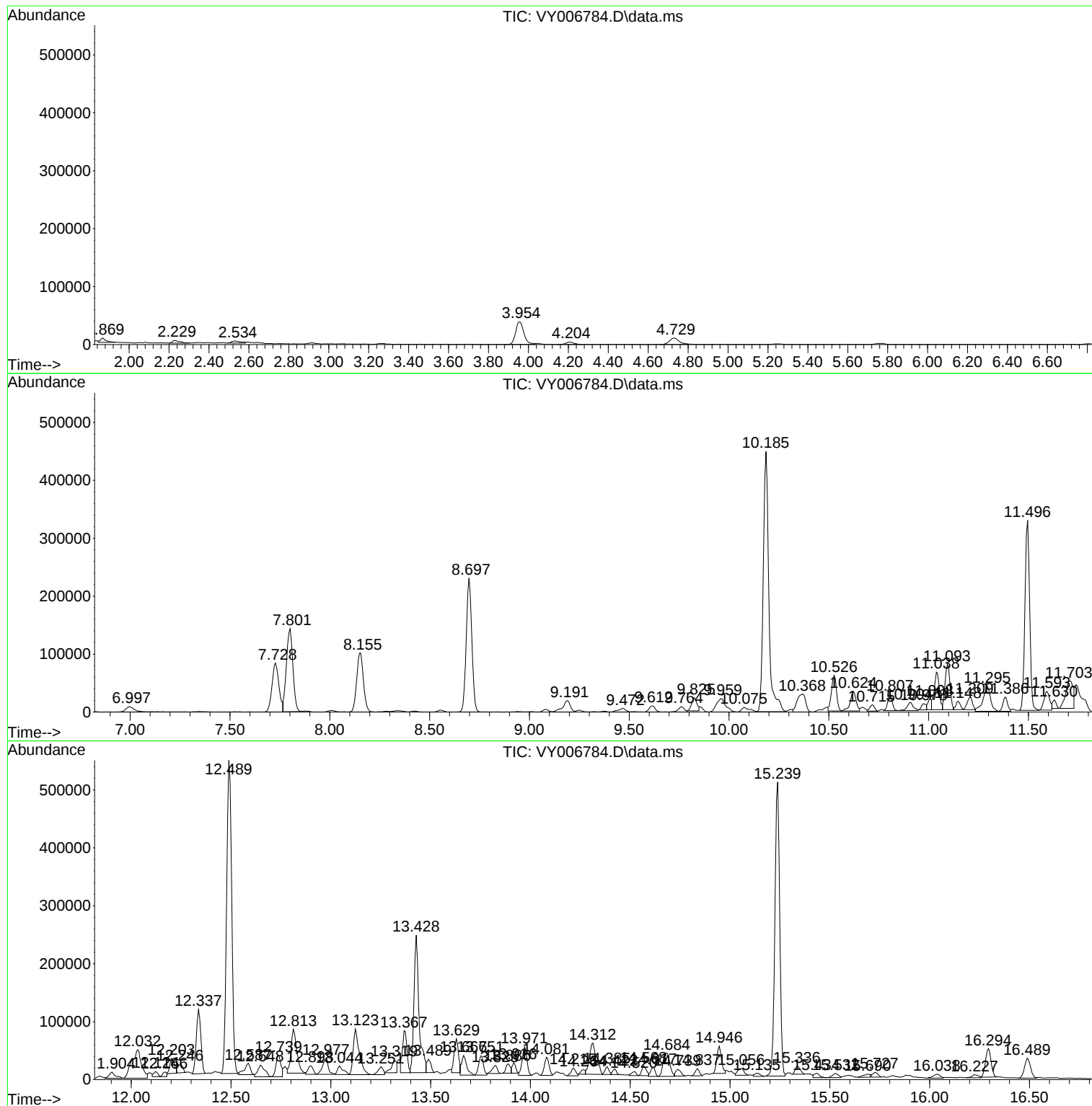
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y111521S.M
Quant Title : SW846 8260

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



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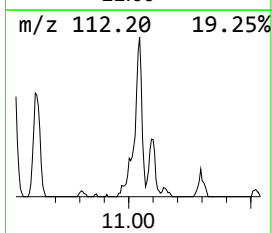
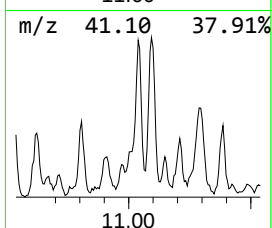
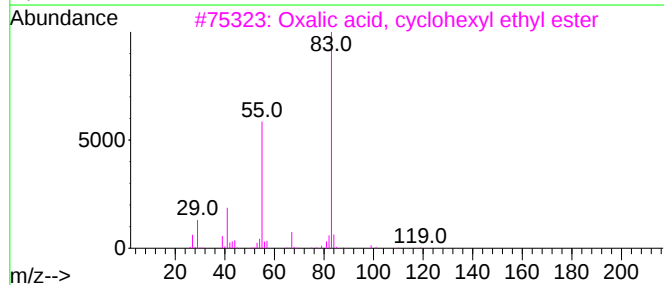
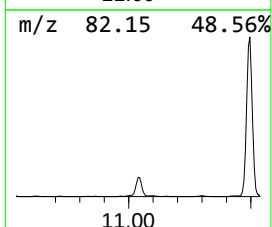
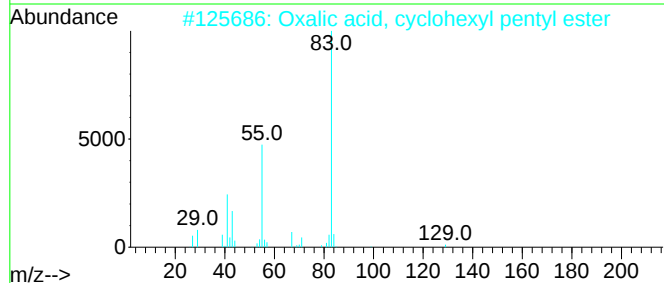
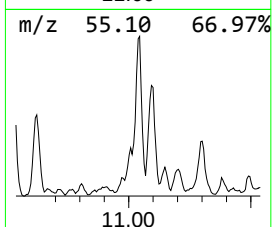
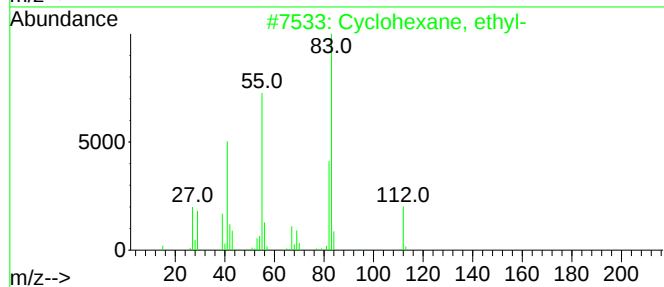
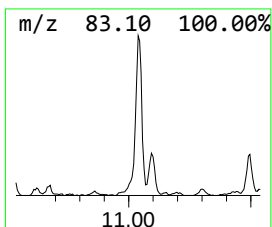
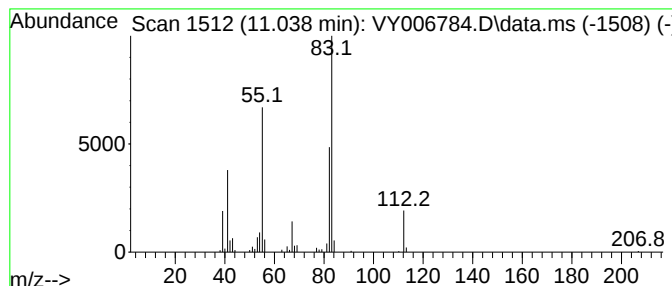
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y111521S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.039	10.59 ug/l	113243	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	90
2			Oxalic acid, cyclohexyl pentyl e...	242	C13H22O4	1000309-30-6	50
3			Oxalic acid, cyclohexyl ethyl ester	200	C10H16O4	1000309-30-2	50
4			Oxalic acid, cyclohexyl butyl ester	228	C12H20O4	1000309-30-5	50
5			Cyclohexanecarboxylic acid, ethe...	154	C9H14O2	004840-76-0	50



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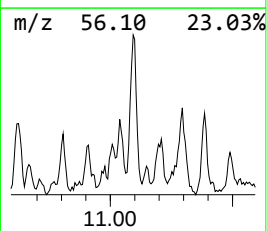
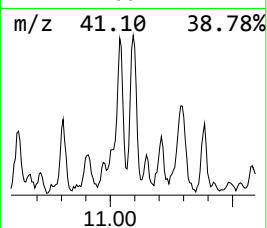
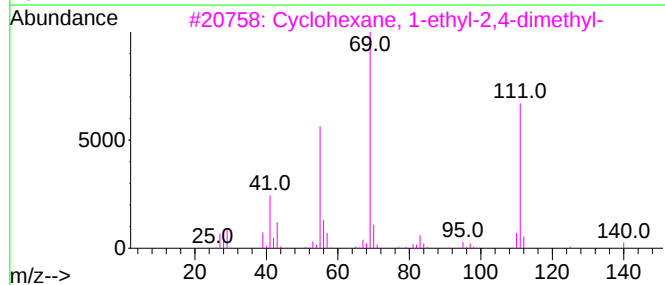
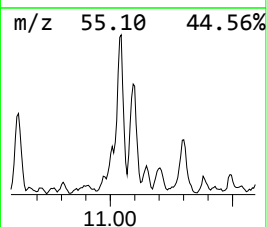
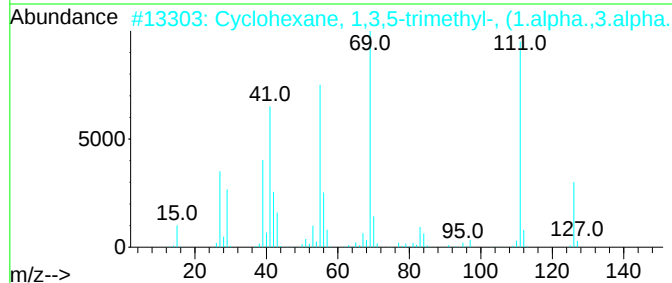
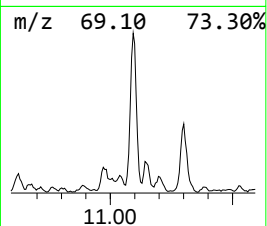
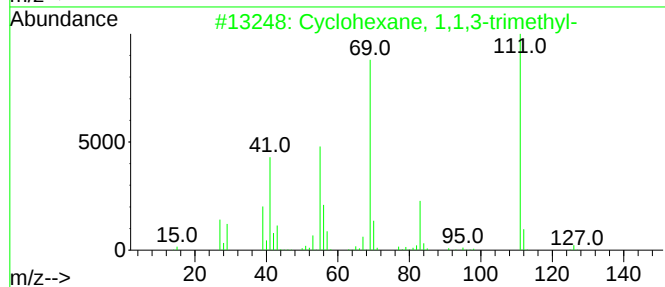
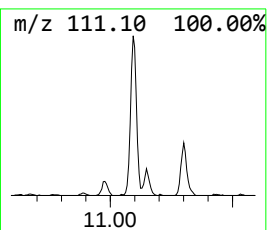
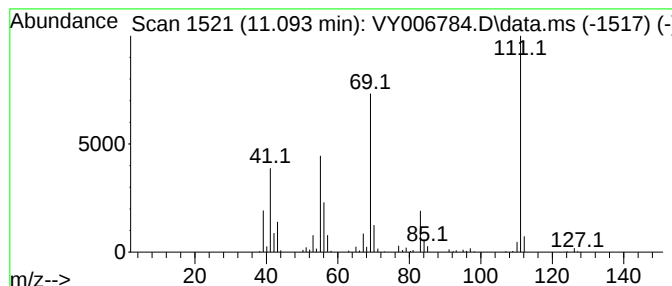
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y111521S.M
Quant Title : SW846 8260

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

Peak Number 2 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.093	12.47 ug/l	133435	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	95
2			Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-26-2	72
3			Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	72
4			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	64
5			Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-31-7	59



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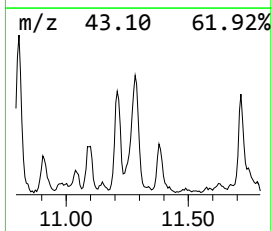
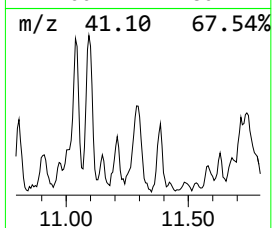
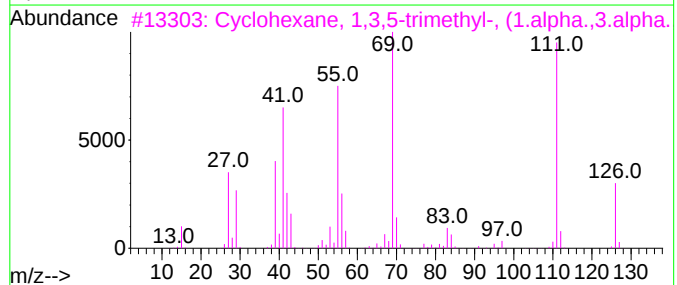
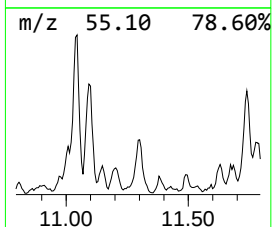
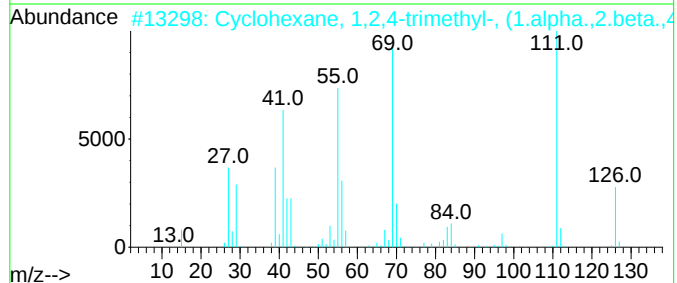
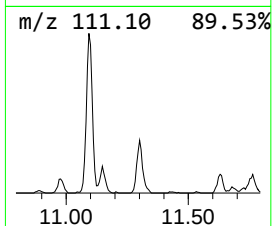
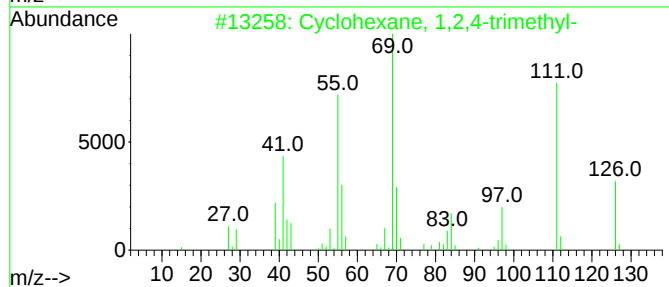
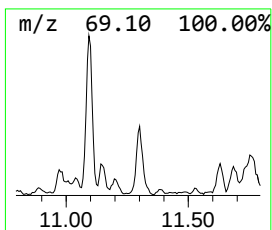
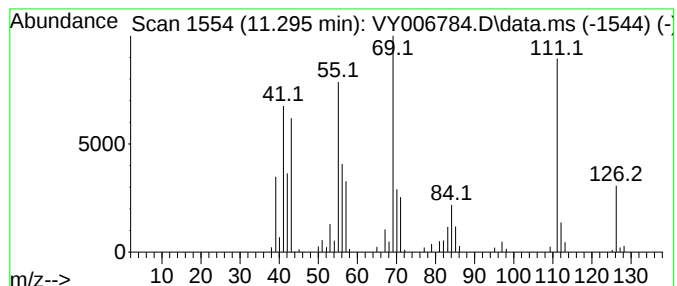
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y111521S.M
Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, 1,2,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.294	10.34 ug/l	110626	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	90
2			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	76
3			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	62
4			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	62
5			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	58



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ClientSampleId :
UT-COMP-02

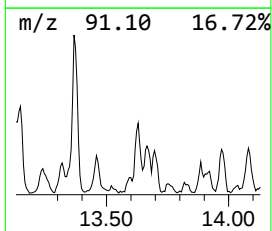
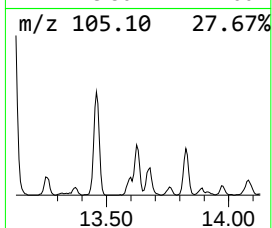
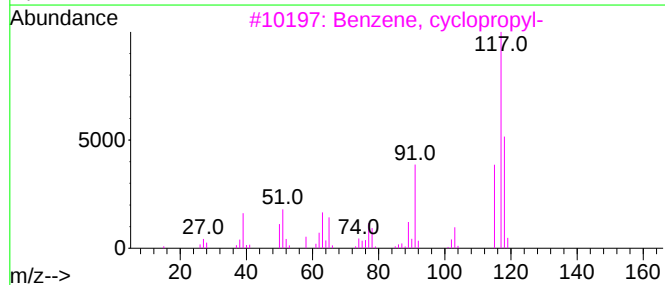
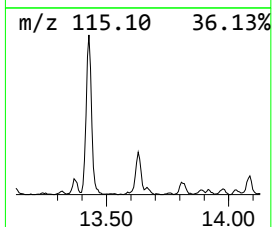
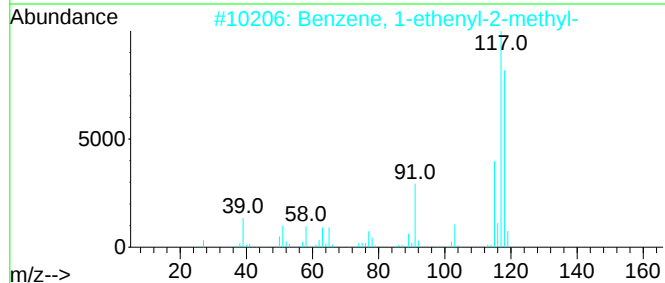
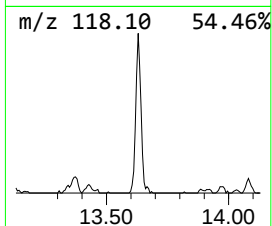
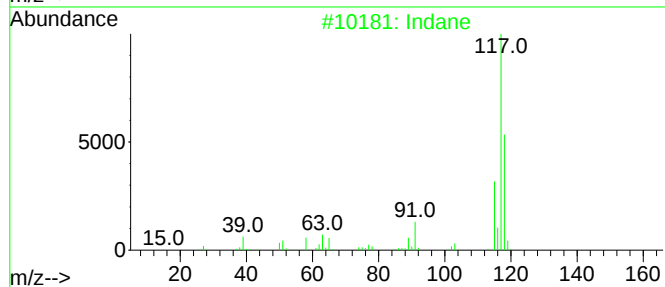
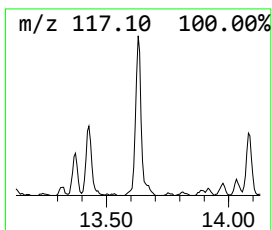
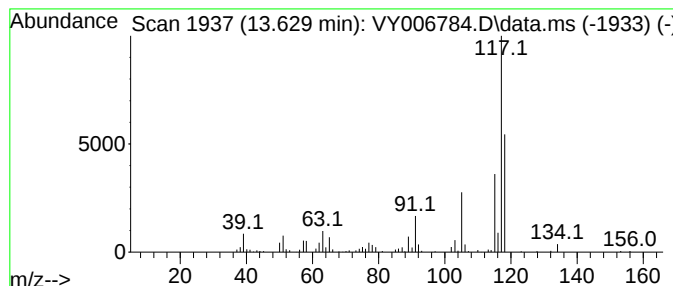
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Quant Title : SW846 8260

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

Peak Number 4 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.630	10.45 ug/l	87505	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	92
2			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	70
5			Benzene, 1-propenyl-	118	C9H10	000637-50-3	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
Data File : VY006784.D
Acq On : 17 Nov 2021 13:33
Operator : SY/MD
Sample : M4693-03
Misc : 5.49G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
UT-COMP-02

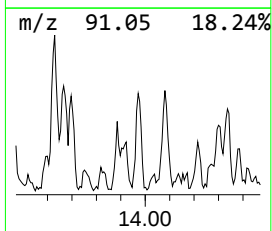
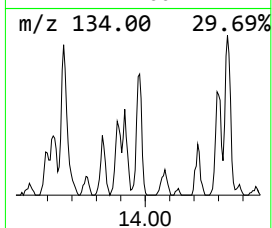
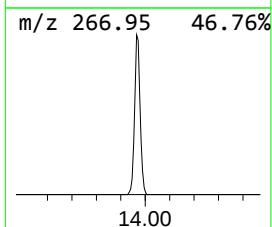
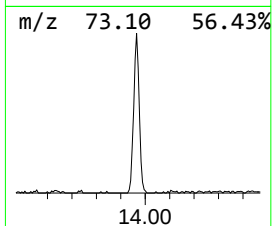
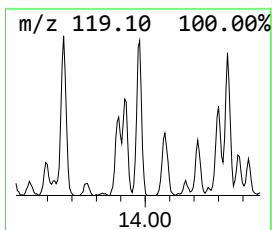
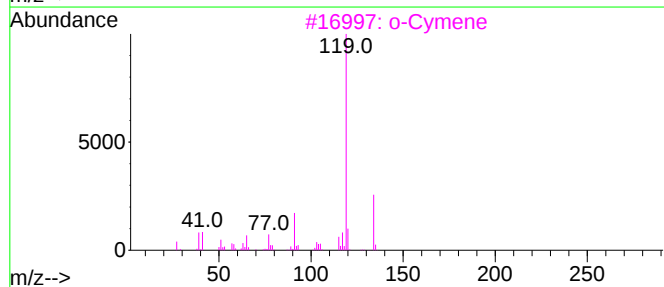
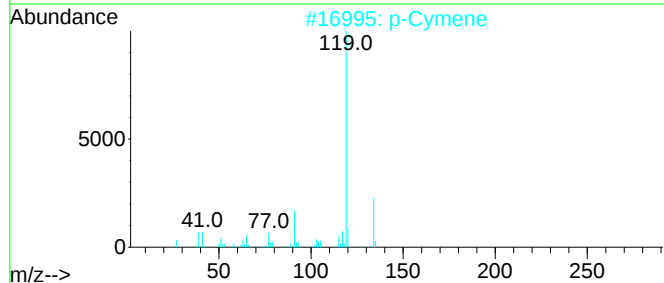
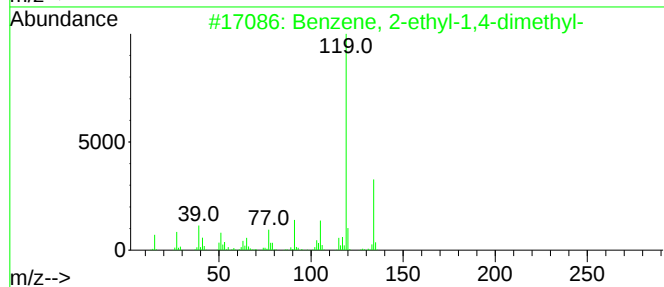
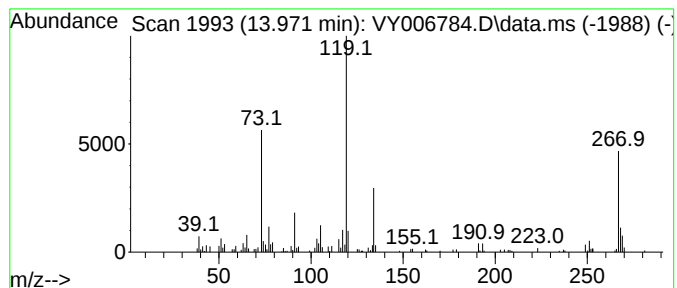
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	9.60 ug/l	80374	1,4-Dichlorobenzene-d4	13.428

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	92
3			o-Cymene	134	C10H14	000527-84-4	92
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
Data File : VY006784.D
Acq On : 17 Nov 2021 13:33
Operator : SY/MD
Sample : M4693-03
Misc : 5.49G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
UT-COMP-02

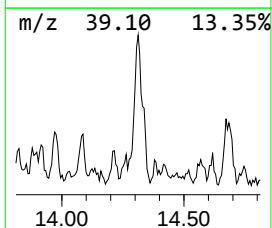
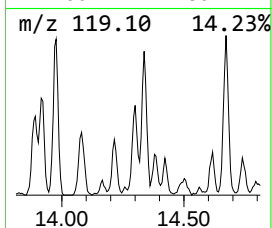
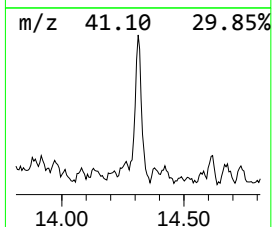
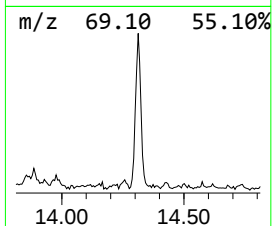
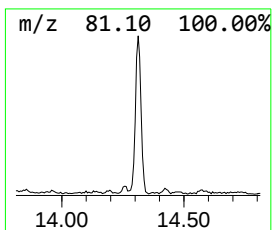
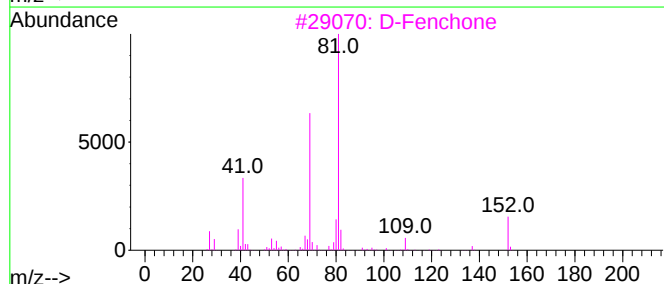
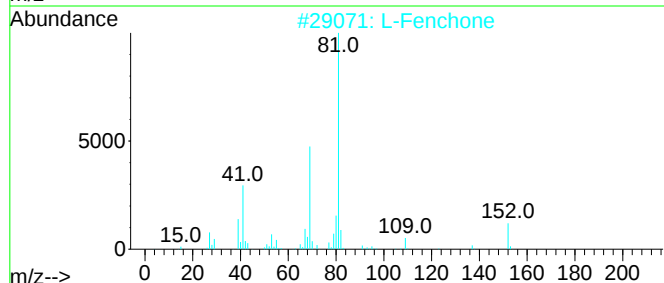
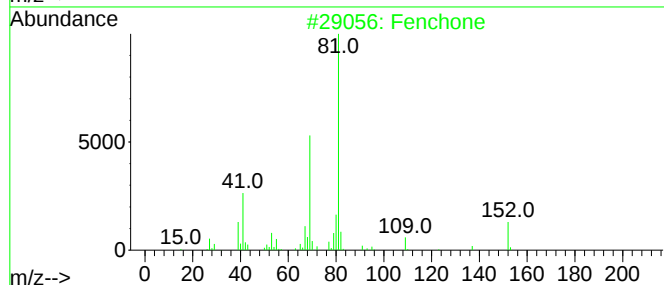
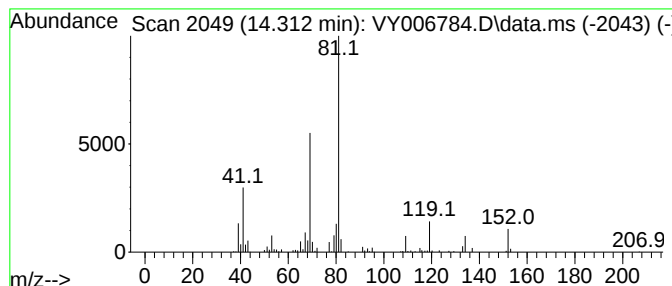
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Quant Title : SW846 8260

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

Peak Number 6 Fenchone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.312	15.96 ug/l	133678	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fenchone	152	C10H16O	001195-79-5	94
2			L-Fenchone	152	C10H16O	007787-20-4	93
3			D-Fenchone	152	C10H16O	004695-62-9	42
4			2,4-Dimethyl 1,4-pentadiene	96	C7H12	004161-65-3	38
5			Cyclohexene,3-(2-propenyl)-	122	C9H14	015232-95-8	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
Data File : VY006784.D
Acq On : 17 Nov 2021 13:33
Operator : SY/MD
Sample : M4693-03
Misc : 5.49G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
UT-COMP-02

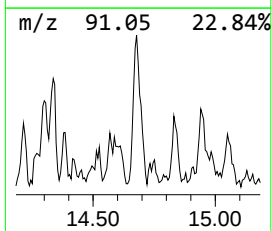
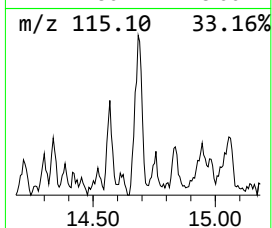
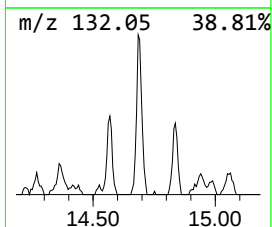
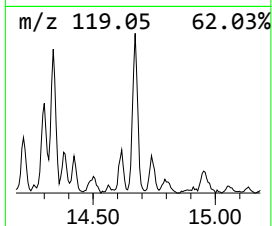
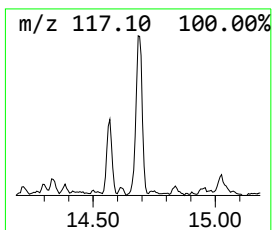
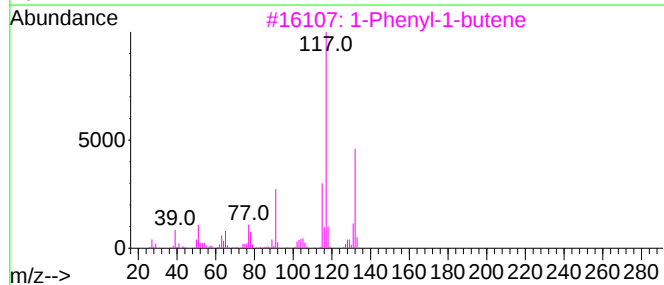
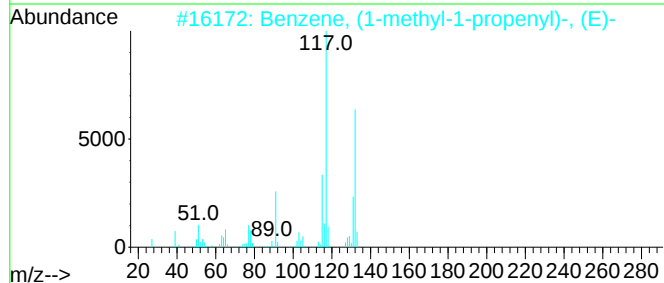
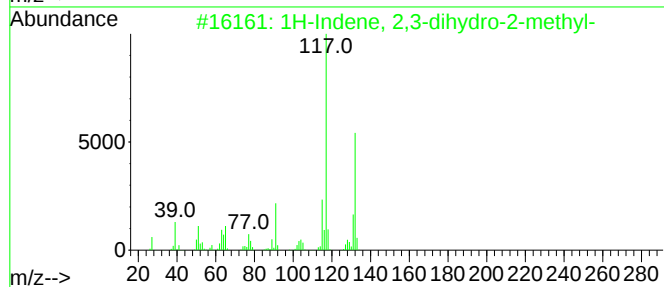
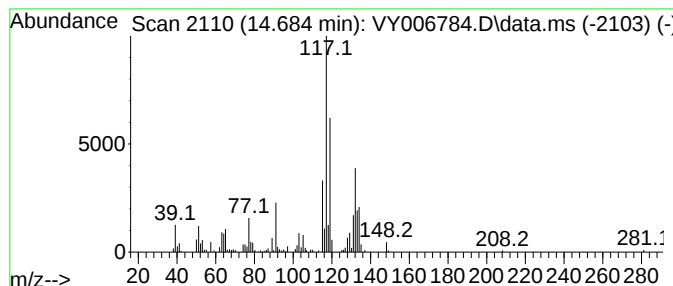
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	11.50 ug/l	96296	1,4-Dichlorobenzene-d4	13.428

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	70
2		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	70
4		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	70
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
Data File : VY006784.D
Acq On : 17 Nov 2021 13:33
Operator : SY/MD
Sample : M4693-03
Misc : 5.49G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
UT-COMP-02

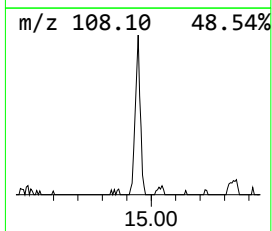
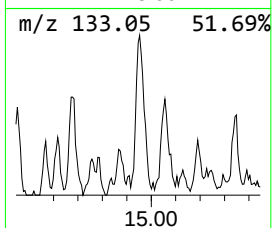
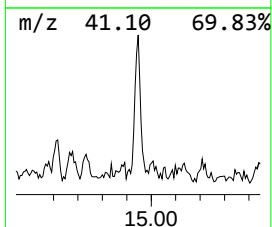
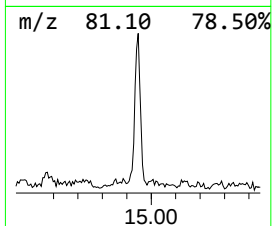
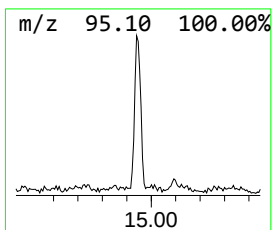
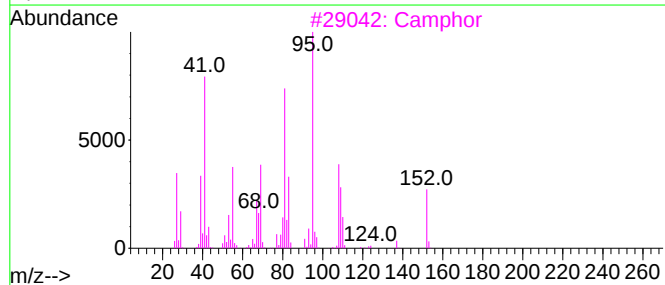
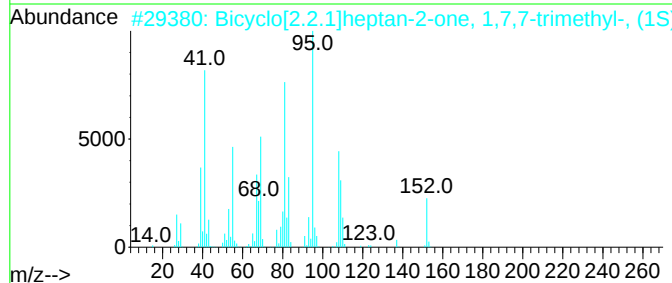
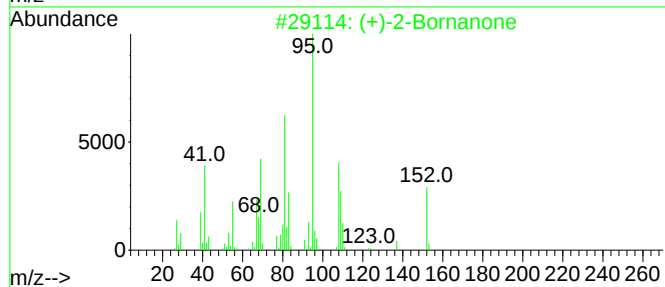
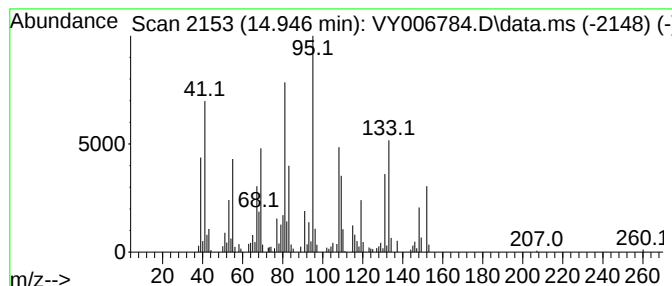
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Quant Title : SW846 8260

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

Peak Number 8 (+)-2-Bornanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.946	9.15 ug/l	76610	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(+)-2-Bornanone	152	C10H16O	000464-49-3	93
2			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	92
3			Camphor	152	C10H16O	000076-22-2	86
4			2-Butanone, 4-(5-methyl-2-furanyl)-	152	C9H12O2	013679-56-6	18
5			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	18



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
Data File : VY006784.D
Acq On : 17 Nov 2021 13:33
Operator : SY/MD
Sample : M4693-03
Misc : 5.49G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
UT-COMP-02

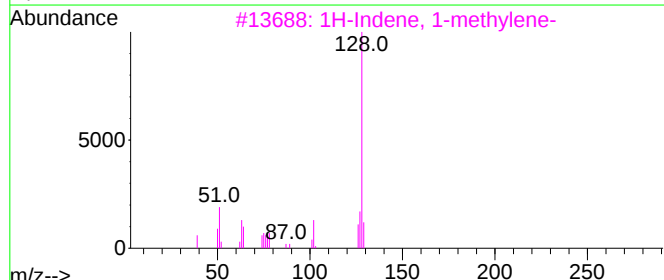
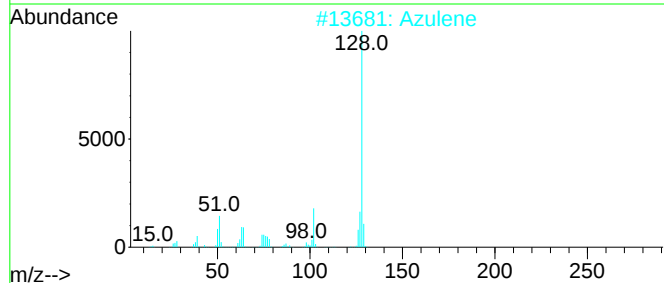
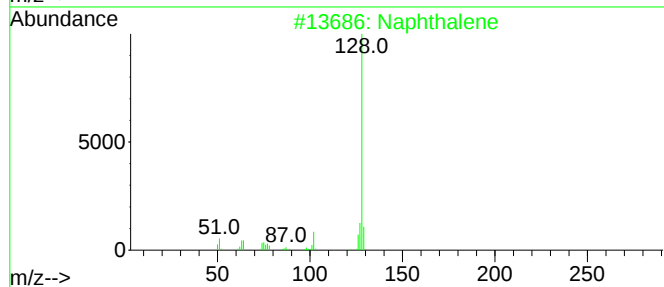
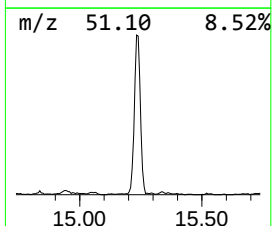
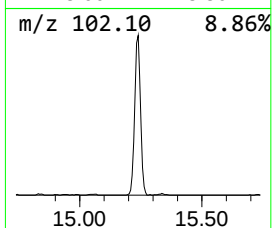
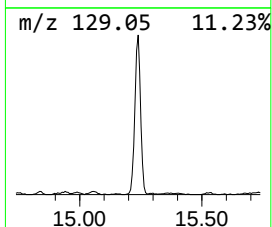
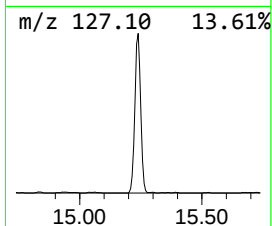
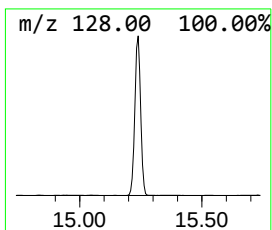
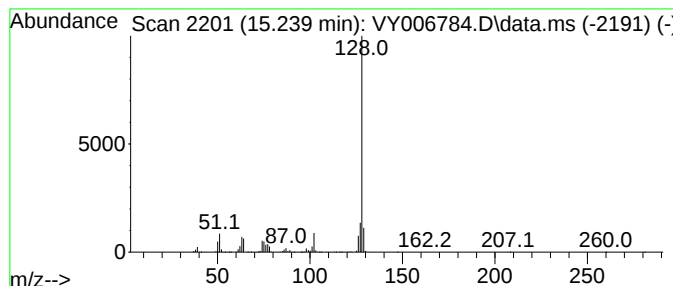
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Quant Title : SW846 8260

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

Peak Number 9 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.239	101.65 ug/l	851380	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	91
3			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	87
4			1,4-Benzenedicarbonitrile	128	C8H4N2	000623-26-7	59
5			4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY111721\
Data File : VY006784.D
Acq On : 17 Nov 2021 13:33
Operator : SY/MD
Sample : M4693-03
Misc : 5.49G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
UT-COMP-02

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y111521S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Cyclohexane, 1,...	11.093	12.5	ug/l	133435	3	11.496	534878	50.0
Cyclohexane, 1,...	11.294	10.3	ug/l	110626	3	11.496	534878	50.0
Indane	13.630	10.4	ug/l	87505	4	13.428	418802	50.0
Benzene, 2-ethy...	13.971	9.6	ug/l	80374	4	13.428	418802	50.0
Fenchone	14.312	16.0	ug/l	133678	4	13.428	418802	50.0
1H-Indene, 2,3-...	14.684	11.5	ug/l	96296	4	13.428	418802	50.0
(+)-2-Bornanone	14.946	9.2	ug/l	76610	4	13.428	418802	50.0
Azulene	15.239	101.7	ug/l	851380	4	13.428	418802	50.0