

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
 Data File : VY009274.D
 Acq On : 22 Jun 2022 15:11
 Operator : KP/MD
 Sample : VIBLK
 Misc : 5.00/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 VIBLK

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\82Y062122S.M
 Title : SW846 8260

Signal : TIC: VY009274.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.899	169	177	185	rBV	13799	31860	1.32%	0.101%
2	3.247	225	234	242	rBV2	10463	24646	1.02%	0.078%
3	4.716	458	475	478	rBV6	26230	85560	3.55%	0.271%
4	5.228	549	559	573	rBV3	11572	40249	1.67%	0.127%
5	5.734	632	642	655	rBV4	14533	45293	1.88%	0.143%
6	6.783	805	814	825	rVB3	13877	44171	1.83%	0.140%
7	6.978	833	846	856	rBV2	11153	36883	1.53%	0.117%
8	7.709	956	966	972	rBV	200643	497403	20.64%	1.576%
9	7.783	972	978	987	rVV	340155	821152	34.07%	2.601%
10	7.862	987	991	1004	rVB3	24575	61245	2.54%	0.194%
11	7.996	1004	1013	1020	rBV2	28447	66092	2.74%	0.209%
12	8.136	1028	1036	1048	rBV	194723	432830	17.96%	1.371%
13	8.319	1051	1066	1076	rBV2	69921	204312	8.48%	0.647%
14	8.545	1093	1103	1115	rBV2	22566	51155	2.12%	0.162%
15	8.685	1115	1126	1139	rBV	485911	955814	39.66%	3.028%
16	9.185	1203	1208	1212	rVB2	22268	40189	1.67%	0.127%
17	9.240	1212	1217	1225	rVB3	17849	34625	1.44%	0.110%
18	9.612	1271	1278	1288	rVB	44547	87328	3.62%	0.277%
19	9.746	1288	1300	1307	rBV2	44634	106721	4.43%	0.338%
20	9.819	1307	1312	1316	rBV2	27066	49757	2.06%	0.158%
21	9.959	1325	1335	1346	rVB2	30086	70335	2.92%	0.223%
22	10.105	1354	1359	1364	rBV	16453	27167	1.13%	0.086%
23	10.178	1364	1371	1377	rBV	726679	1231674	51.10%	3.902%
24	10.239	1377	1381	1387	rVB	36320	61710	2.56%	0.195%
25	10.367	1394	1402	1408	rVB	39110	73373	3.04%	0.232%
26	10.947	1493	1497	1504	rVB	25975	48407	2.01%	0.153%
27	11.276	1542	1551	1562	rVB2	46952	103978	4.31%	0.329%
28	11.379	1562	1568	1578	rBV2	36308	64714	2.68%	0.205%
29	11.489	1578	1586	1595	rBV	720039	1188164	49.30%	3.764%
30	11.593	1595	1603	1610	rVB	175613	295897	12.28%	0.937%
31	11.697	1612	1620	1631	rBV	831571	1533170	63.61%	4.857%
32	12.026	1666	1674	1682	rVB	288408	476371	19.76%	1.509%
33	12.331	1718	1724	1727	rBV	33909	66748	2.77%	0.211%
34	12.477	1743	1748	1754	rVB	574648	918006	38.09%	2.908%
35	12.666	1774	1779	1784	rVB	166785	250755	10.40%	0.794%
36	12.733	1784	1790	1793	rBV	665255	1007383	41.80%	3.191%

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37	12.763	1793	1795	1799	rVV	318943	451257	18.72%	1.429%
38	12.812	1799	1803	1809	rVB	490444	739259	30.67%	2.342%
39	12.971	1823	1829	1837	rVB	252416	414333	17.19%	1.312%
40	13.117	1847	1853	1859	rVB	1595057	2410240	100.00%	7.635%
41	13.166	1859	1861	1867	rVB3	19105	28392	1.18%	0.090%
42	13.233	1867	1872	1874	rBV3	26120	48071	1.99%	0.152%
43	13.312	1882	1885	1889	rVB	46579	65050	2.70%	0.206%
44	13.361	1889	1893	1897	rBV2	36052	56929	2.36%	0.180%
45	13.422	1897	1903	1907	rVV	758575	1349374	55.99%	4.274%
46	13.452	1907	1908	1914	rVV	372868	450598	18.70%	1.427%
47	13.495	1914	1915	1923	rVB2	42733	56685	2.35%	0.180%
48	13.593	1925	1931	1932	rBV	145126	188048	7.80%	0.596%
49	13.623	1932	1936	1939	rVV2	513751	796308	33.04%	2.522%
50	13.666	1939	1943	1953	rVB3	539325	1022364	42.42%	3.239%
51	13.751	1953	1957	1962	rBV	81566	112646	4.67%	0.357%
52	13.824	1962	1969	1974	rVB	107528	170524	7.07%	0.540%
53	13.885	1974	1979	1981	rBV	273154	403375	16.74%	1.278%
54	13.916	1981	1984	1988	rVB	267681	371509	15.41%	1.177%
55	13.970	1988	1993	1998	rBV	514615	737812	30.61%	2.337%
56	14.031	1998	2003	2006	rVB	45391	64638	2.68%	0.205%
57	14.080	2006	2011	2016	rVB	195975	296395	12.30%	0.939%
58	14.153	2016	2023	2027	rVB5	25165	51825	2.15%	0.164%
59	14.214	2027	2033	2039	rBV	123929	254950	10.58%	0.808%
60	14.294	2040	2046	2049	rVV	340817	531140	22.04%	1.682%
61	14.336	2049	2053	2057	rVV	462120	677702	28.12%	2.147%
62	14.379	2057	2060	2064	rVV	149464	195183	8.10%	0.618%
63	14.422	2064	2067	2070	rVB	65568	77333	3.21%	0.245%
64	14.519	2079	2083	2086	rBV	102847	136541	5.67%	0.433%
65	14.562	2086	2090	2095	rVV	318391	537916	22.32%	1.704%
66	14.611	2095	2098	2102	rVB	195206	257763	10.69%	0.817%
67	14.684	2102	2110	2115	rBV2	449798	1024322	42.50%	3.245%
68	14.739	2115	2119	2125	rVB2	132854	229457	9.52%	0.727%
69	14.830	2131	2134	2138	rBV3	39215	58261	2.42%	0.185%
70	14.903	2142	2146	2148	rVV2	57131	86168	3.58%	0.273%
71	14.946	2148	2153	2164	rVV3	232063	621921	25.80%	1.970%
72	15.056	2164	2171	2180	rVV2	176723	388488	16.12%	1.231%
73	15.135	2181	2184	2189	rVB4	29182	42653	1.77%	0.135%
74	15.232	2189	2200	2207	rBV	548610	1036817	43.02%	3.284%
75	15.342	2213	2218	2223	rBV3	103322	163915	6.80%	0.519%
76	15.397	2223	2227	2230	rVV2	60313	103086	4.28%	0.327%

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77	15.434	2230	2233	2242	rVB3	117652	197709	8.20%	0.626%
78	15.525	2242	2248	2255	rBV	158492	313820	13.02%	0.994%
79	15.690	2265	2275	2284	rBV4	92297	321100	13.32%	1.017%
80	15.812	2290	2295	2301	rBV2	55561	99983	4.15%	0.317%
81	15.891	2301	2308	2324	rVB3	93588	269576	11.18%	0.854%
82	16.025	2324	2330	2335	rBV6	26266	57830	2.40%	0.183%
83	16.086	2336	2340	2346	rVB8	12313	24438	1.01%	0.077%
84	16.202	2353	2359	2360	rBV4	16725	27331	1.13%	0.087%
85	16.293	2367	2374	2381	rBV	792247	1532652	63.59%	4.855%
86	16.354	2381	2384	2391	rVB5	38145	70531	2.93%	0.223%
87	16.488	2398	2406	2417	rVB	380339	809440	33.58%	2.564%

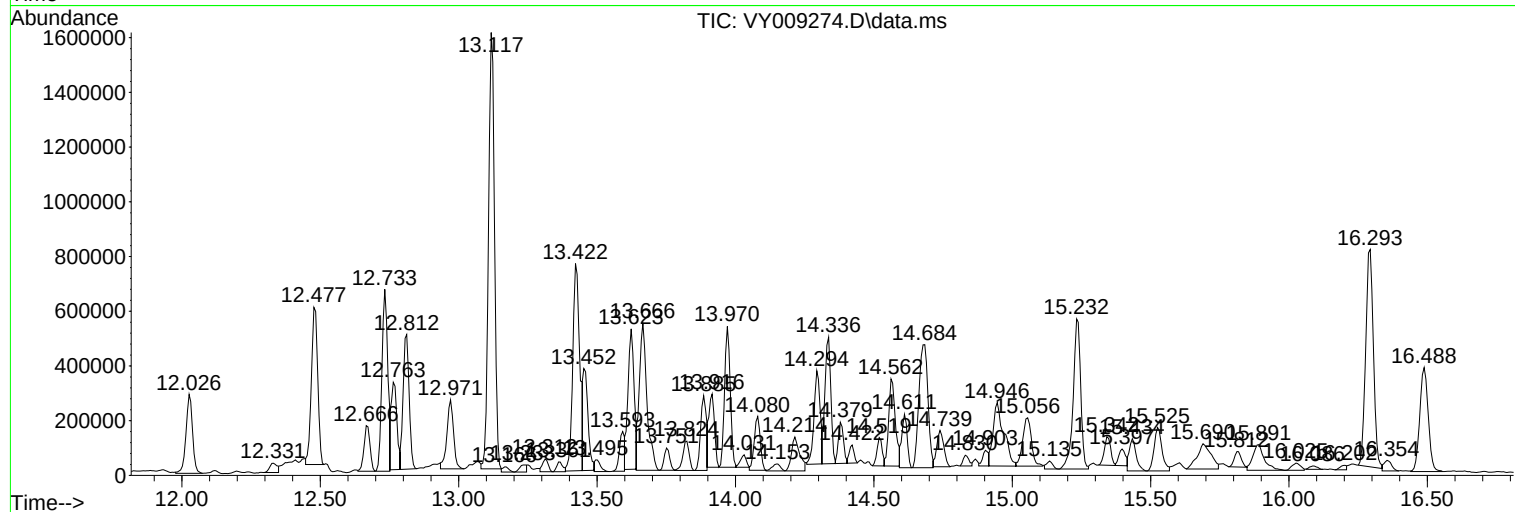
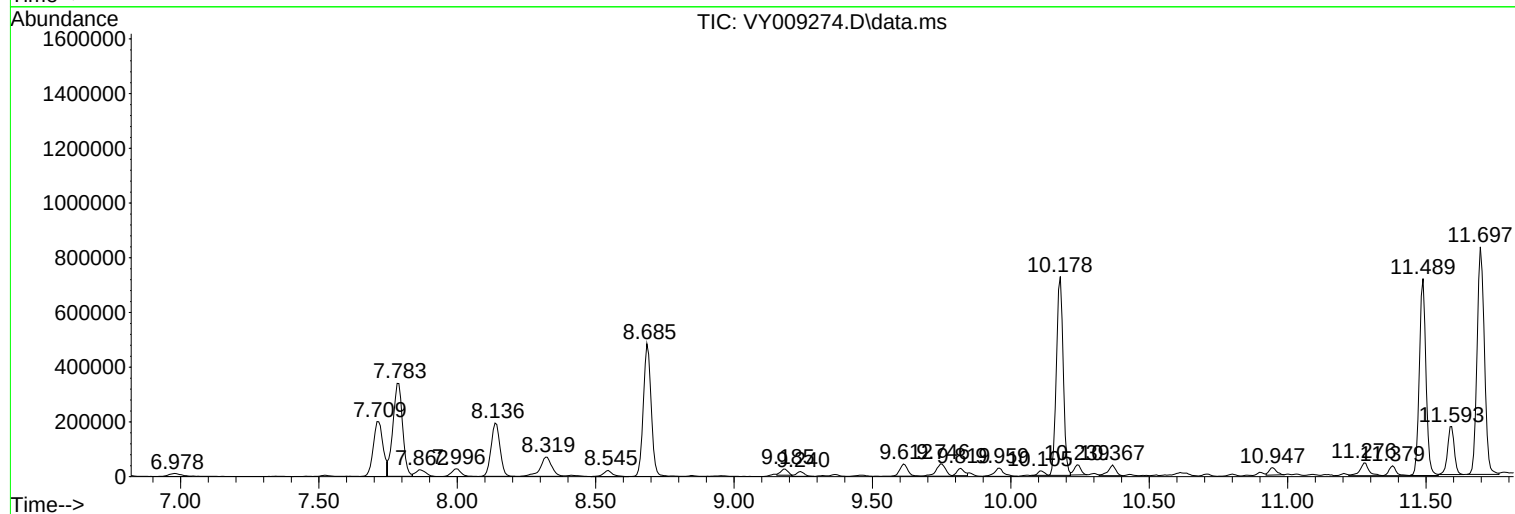
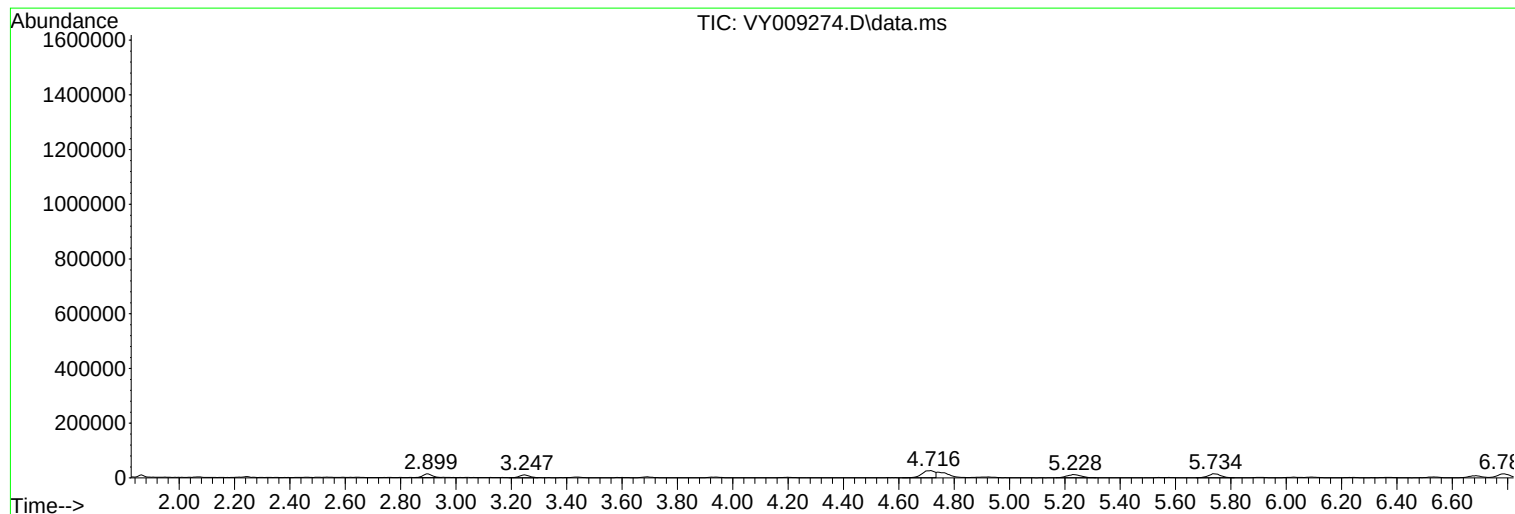
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Instrument :
MSVOA_Y
ClientSampleId :
VIBLK

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062122S.M
Quant Title : SW846 8260

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



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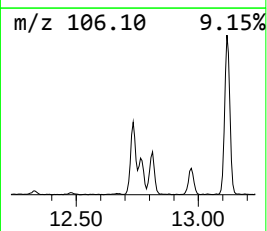
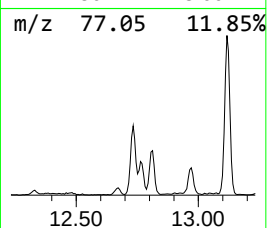
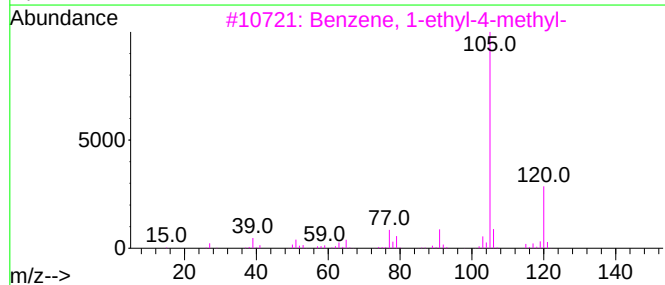
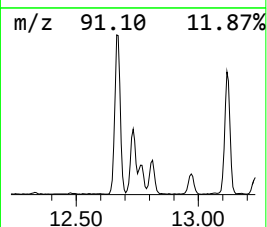
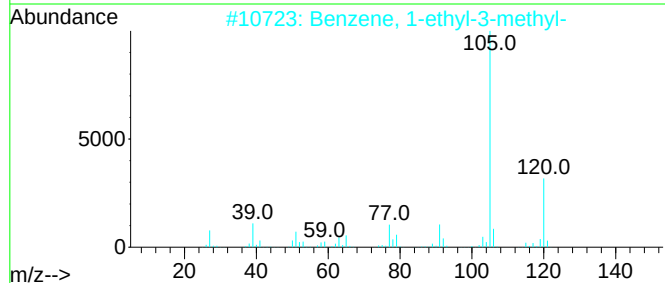
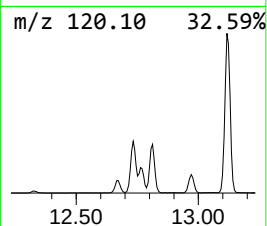
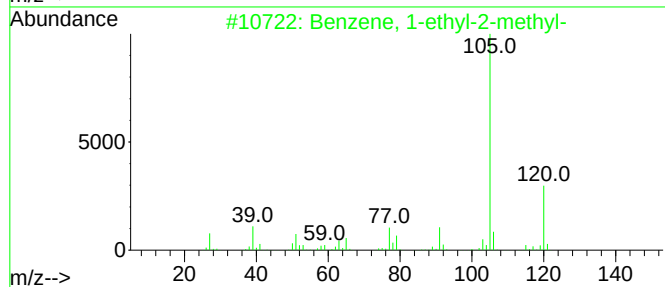
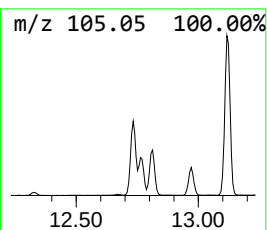
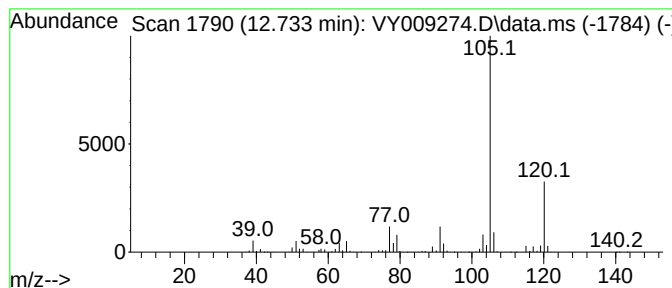
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062122S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	37.33 ug/l	1007380	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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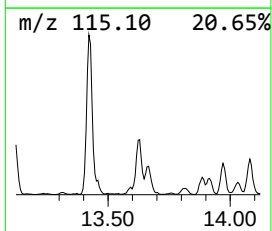
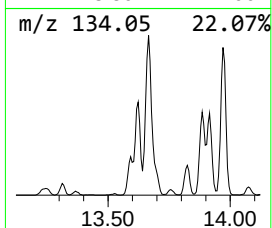
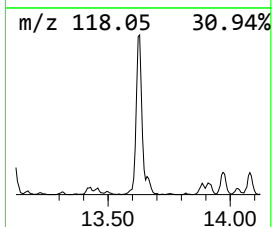
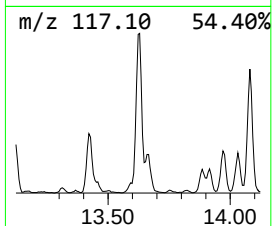
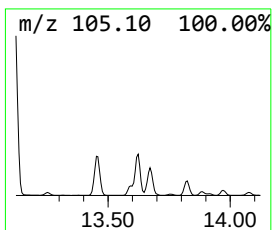
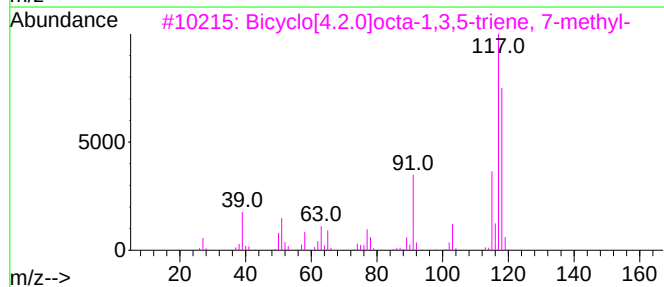
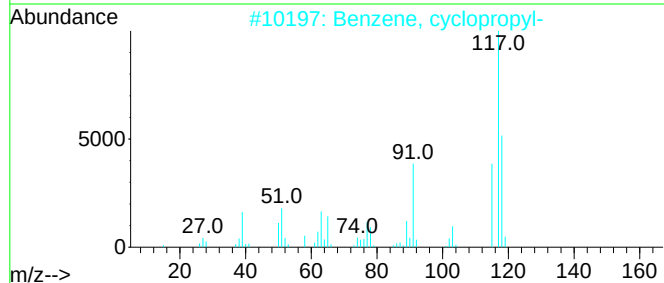
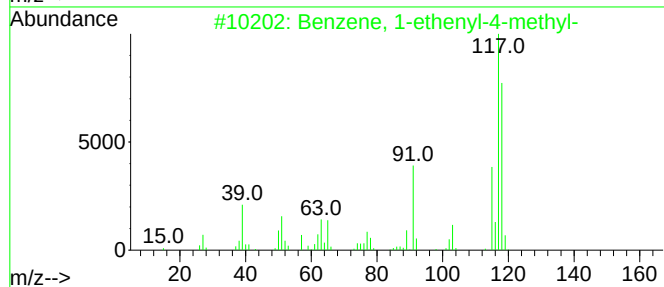
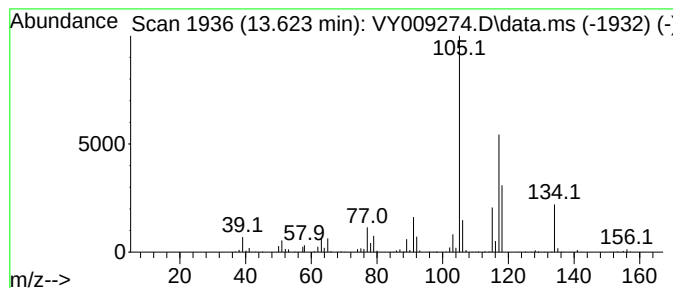
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062122S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethenyl-4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	29.51 ug/l	796308	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	50
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	46
3			Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	43
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	43
5			Indane	118	C9H10	000496-11-7	38



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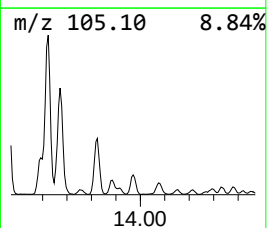
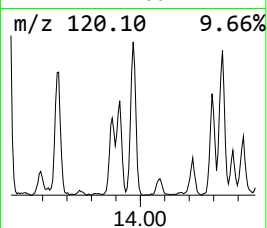
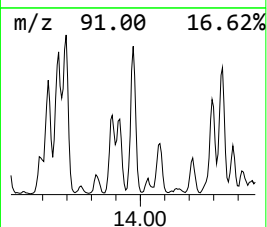
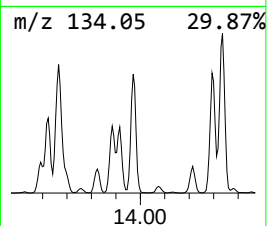
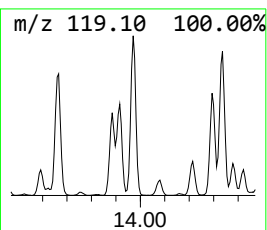
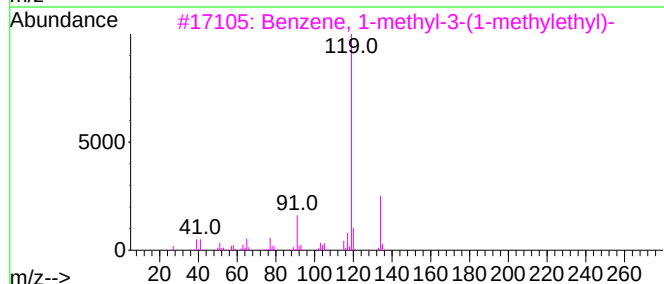
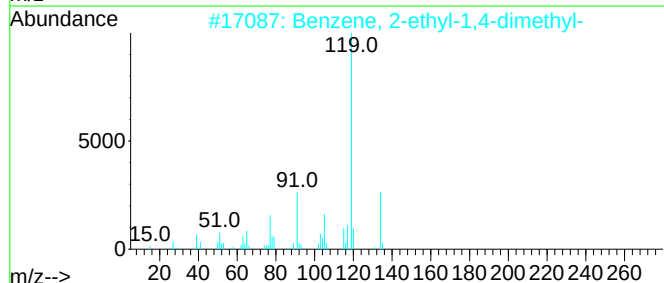
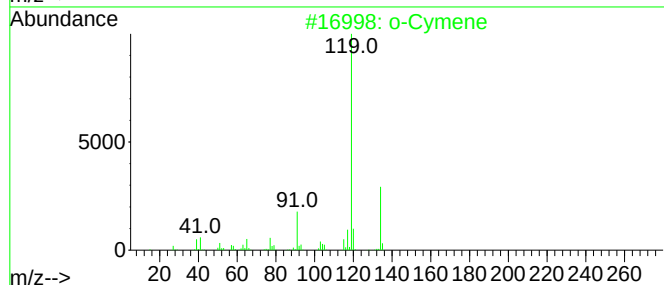
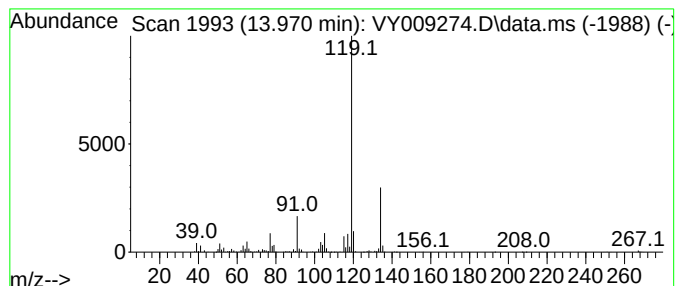
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062122S.M
Quant Title : SW846 8260

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

Peak Number 3 o-Cymene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.970	27.34 ug/l	737812	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 1-methyl-3-(1-methyleth...)	134	C10H14	000535-77-3	95
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



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ClientSampleId :
VIBLK

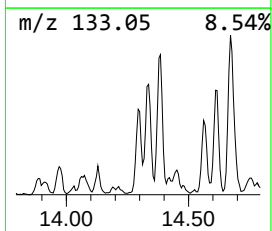
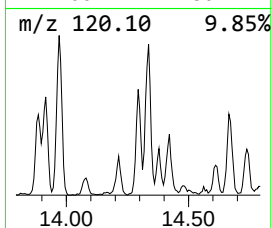
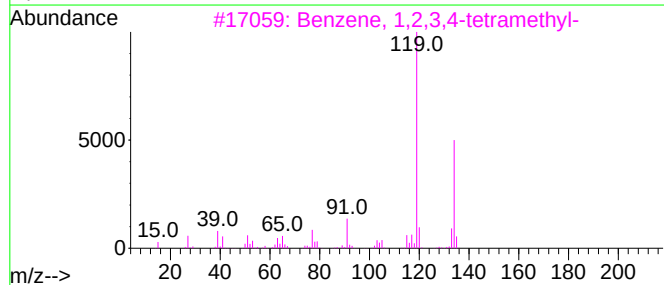
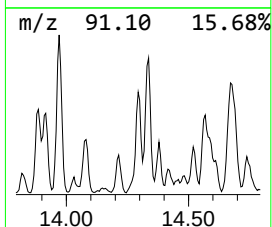
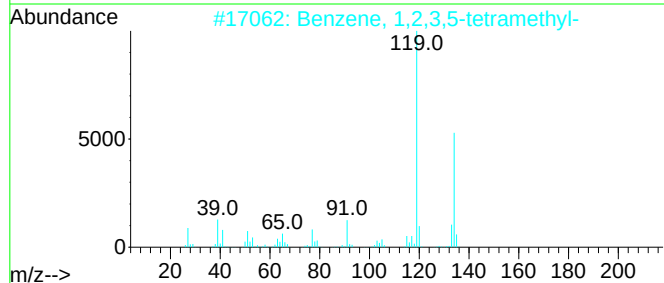
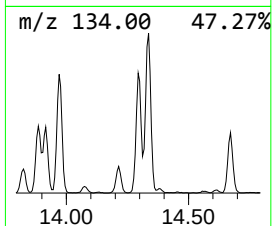
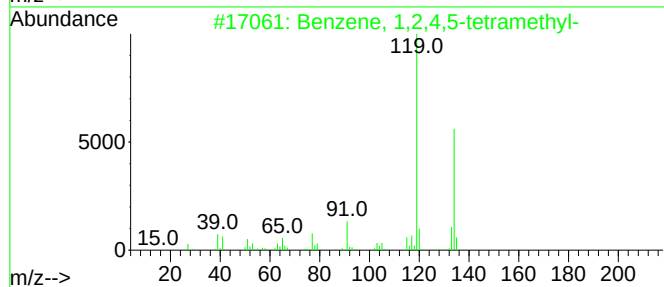
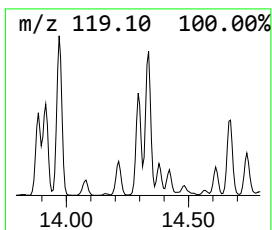
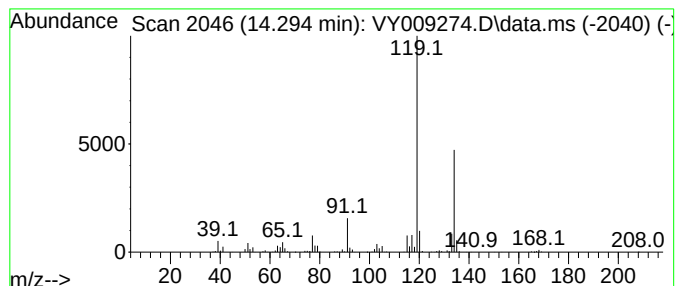
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.294	19.68 ug/l	531140	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VIBLK

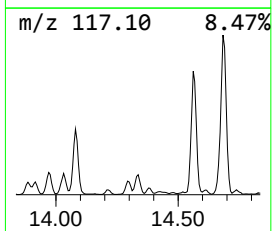
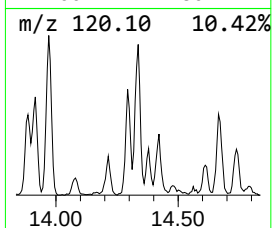
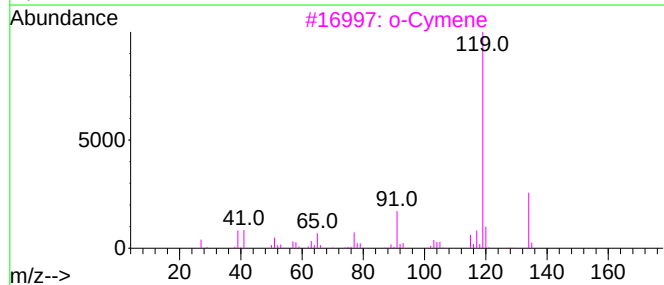
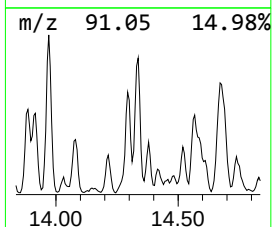
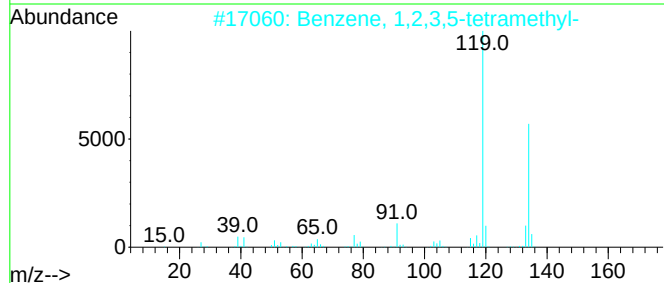
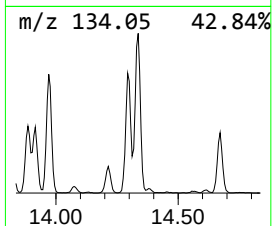
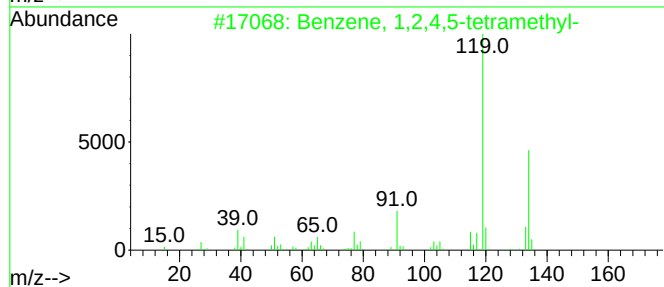
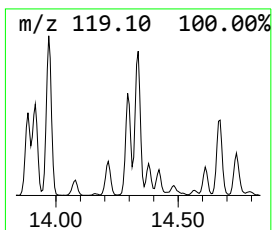
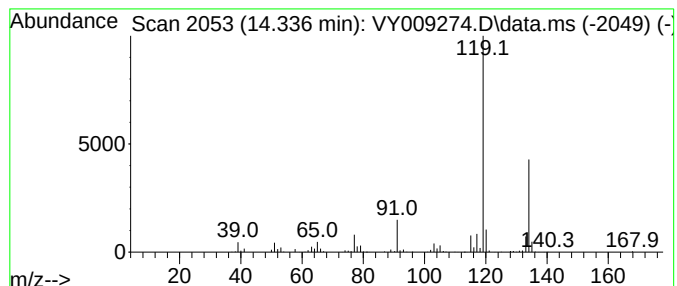
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.336	25.11 ug/l	677702	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VIBLK

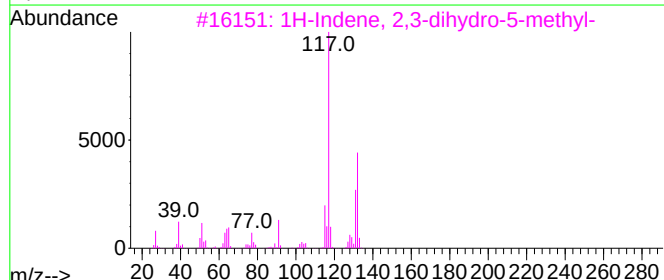
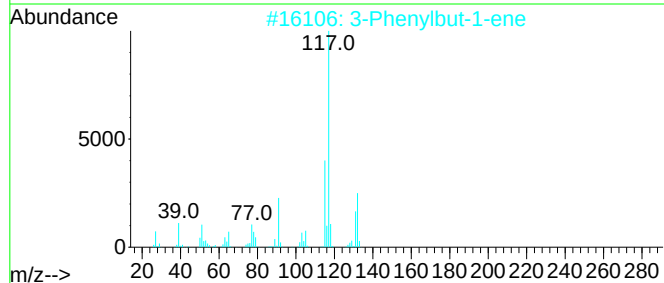
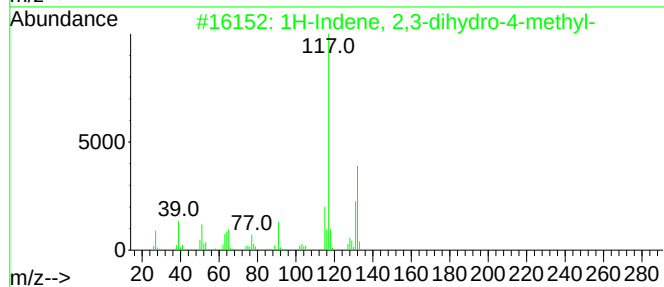
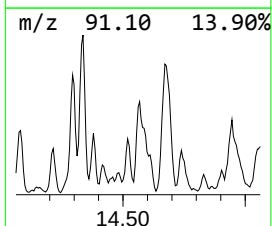
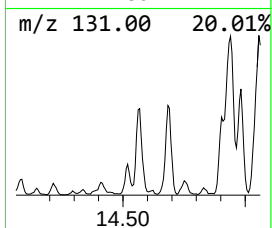
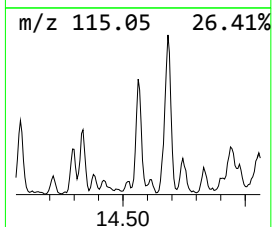
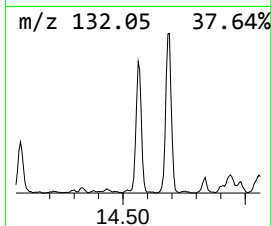
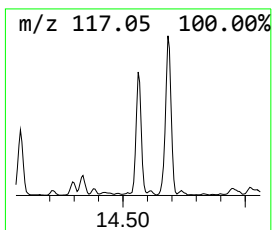
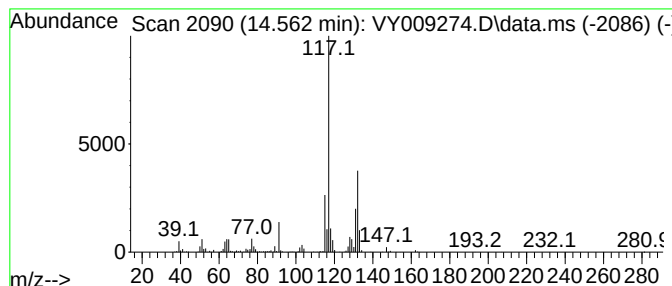
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.562	19.93 ug/l	537916	1,4-Dichlorobenzene-d4	13.422

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
4		Indan, 1-methyl-	132	C10H12	000767-58-8	81
5		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VIBLK

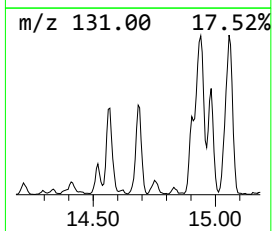
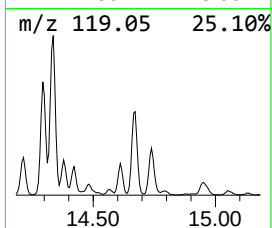
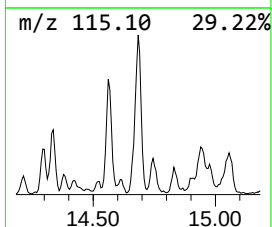
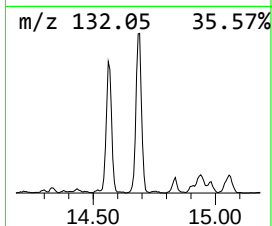
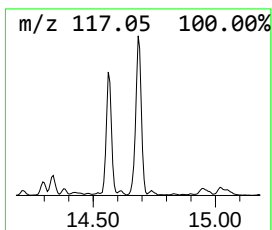
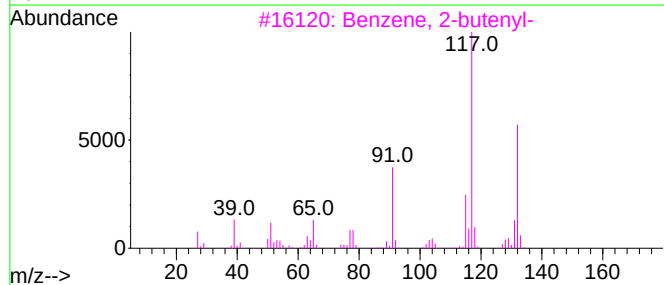
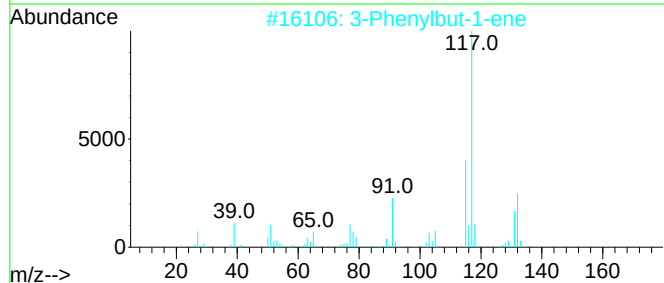
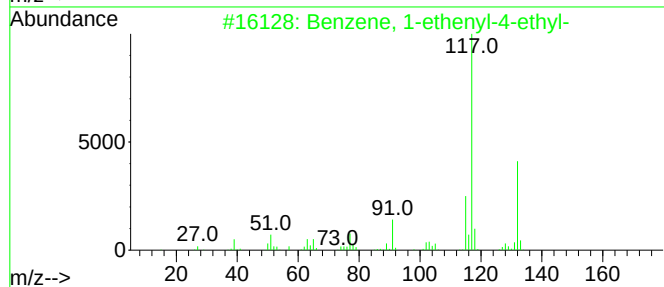
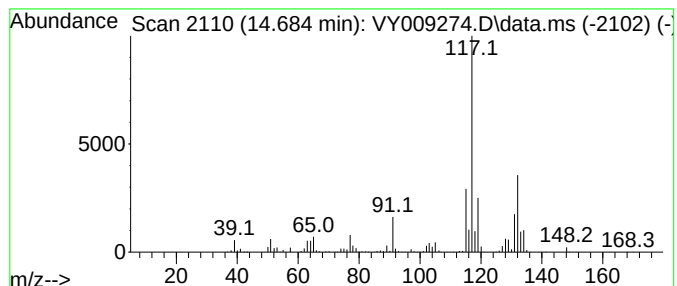
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	37.96 ug/l	1024320	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	81
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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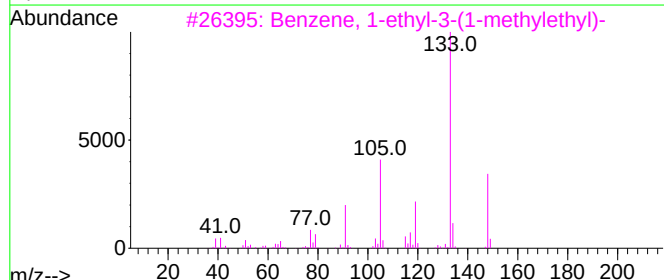
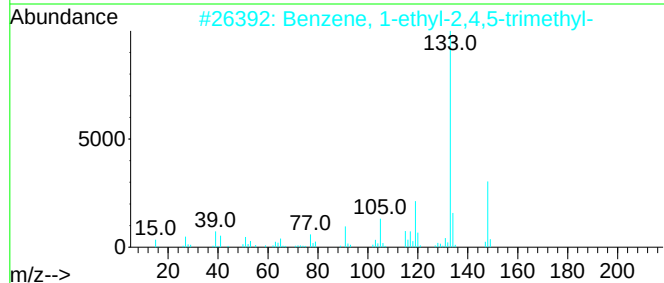
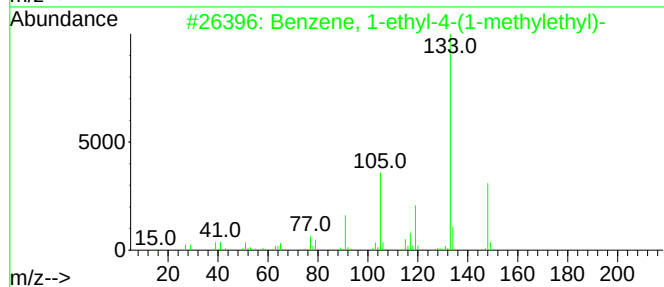
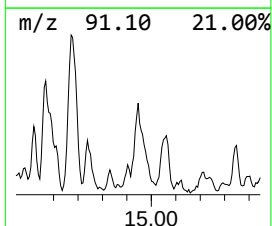
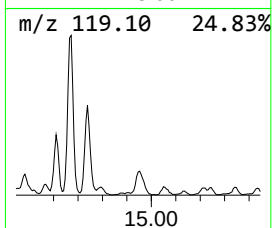
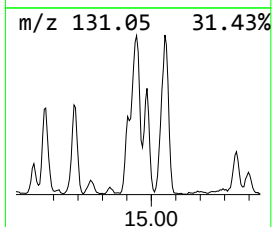
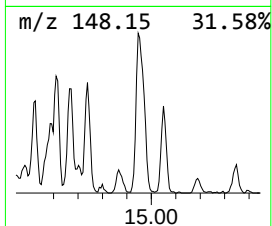
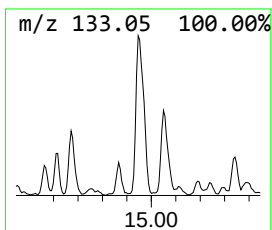
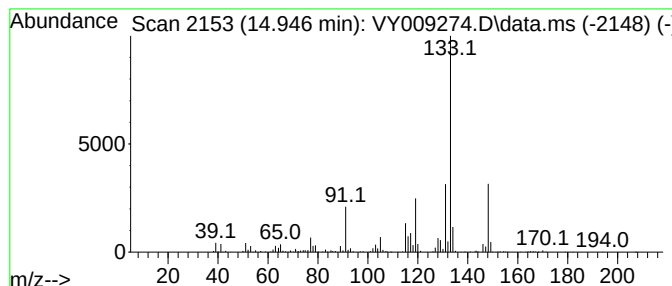
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.946	23.04 ug/l	621921	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	76
2			Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	74
3			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	68
4			4-tert-Butyltoluene	148	C11H16	000098-51-1	58
5			3,4-Dimethylcumene	148	C11H16	1000370-34-1	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VIBLK

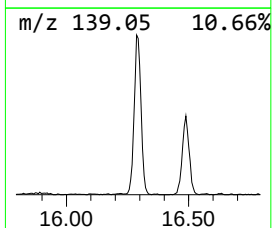
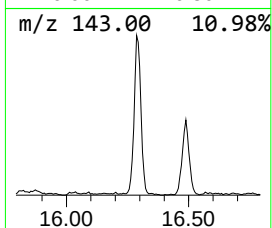
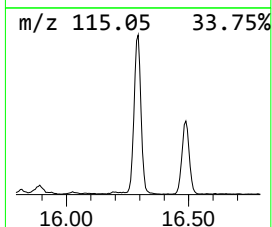
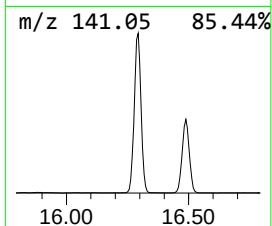
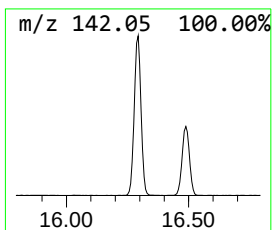
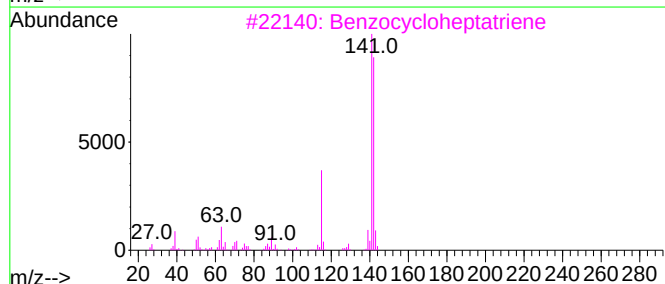
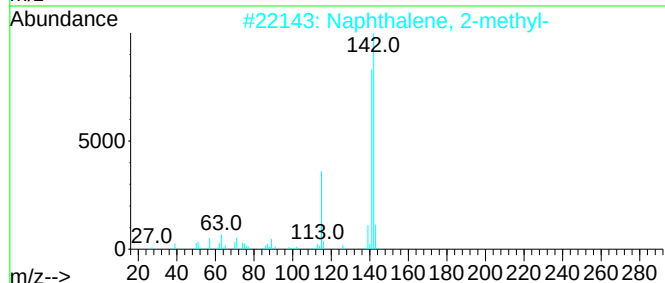
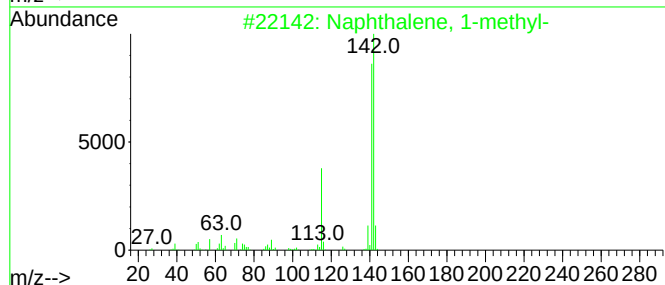
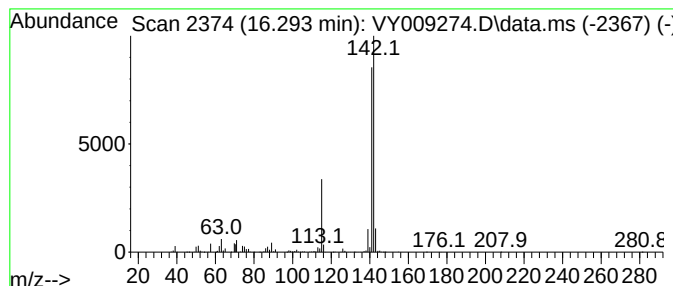
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.293	56.79 ug/l	1532650	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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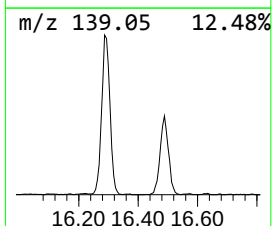
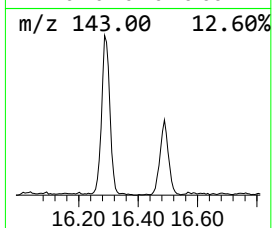
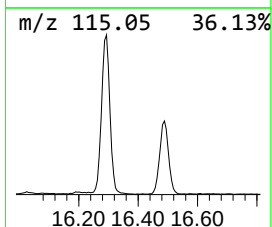
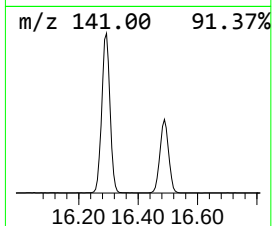
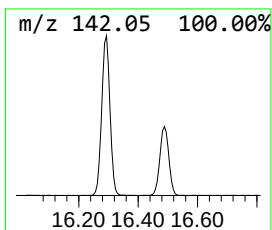
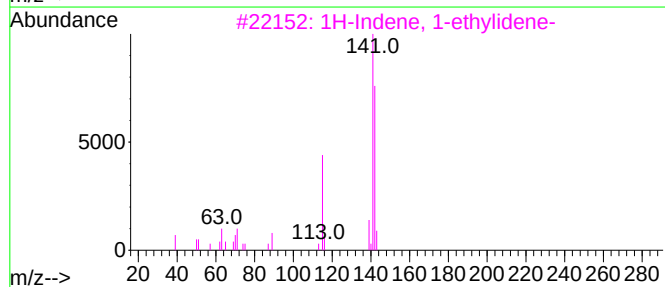
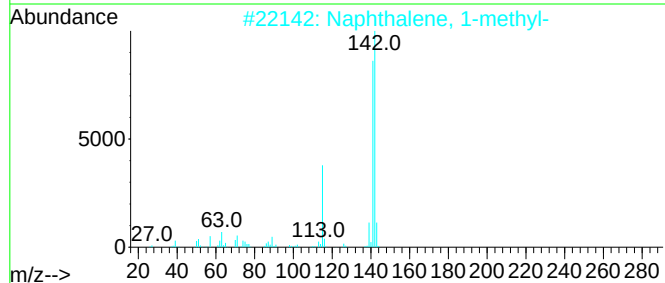
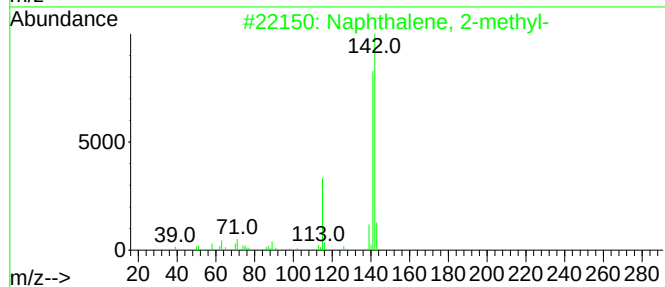
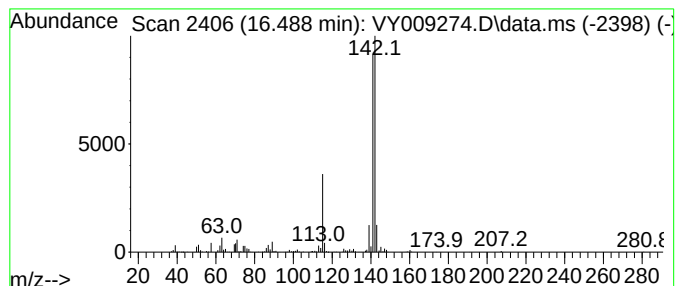
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 1-ethylidene- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.488	29.99 ug/l	809440	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		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	95
4		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5		Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY062222\
Data File : VY009274.D
Acq On : 22 Jun 2022 15:11
Operator : KP/MD
Sample : VIBLK
Misc : 5.00/5.0mL/MSVOA_Y/SOIL
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
VIBLK

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062122S.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
Benzene, 1-ethy...	12.733	37.3	ug/l	1007380	4	13.422	1349370	50.0
Benzene, 1-ethe...	13.623	29.5	ug/l	796308	4	13.422	1349370	50.0
o-Cymene	13.970	27.3	ug/l	737812	4	13.422	1349370	50.0
Benzene, 1,2,4,...	14.294	19.7	ug/l	531140	4	13.422	1349370	50.0
Benzene, 1,2,3,...	14.336	25.1	ug/l	677702	4	13.422	1349370	50.0
1H-Indene, 2,3-...	14.562	19.9	ug/l	537916	4	13.422	1349370	50.0
Benzene, 1-ethe...	14.684	38.0	ug/l	1024320	4	13.422	1349370	50.0
Benzene, 1-ethy...	14.946	23.0	ug/l	621921	4	13.422	1349370	50.0
Naphthalene, 1-...	16.293	56.8	ug/l	1532650	4	13.422	1349370	50.0
1H-Indene, 1-et...	16.488	30.0	ug/l	809440	4	13.422	1349370	50.0