

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
 Data File : VX032561.D
 Acq On : 03 Nov 2022 17:24
 Operator : JC/MD
 Sample : N5454-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1674

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_X\Method\82X102622W.M
 Title : SW846 8260

Signal : TIC: VX032561.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.398	208	215	227	rVB	8377	17119	1.68%	0.110%
2	5.385	693	705	720	rBV2	87490	256521	25.22%	1.641%
3	5.550	720	732	749	rVB	148526	417257	41.02%	2.670%
4	5.958	789	799	815	rBV	97955	259845	25.55%	1.663%
5	6.757	920	930	949	rBV	227864	544542	53.54%	3.485%
6	8.647	1233	1240	1248	rBV	471355	737915	72.55%	4.722%
7	9.006	1290	1299	1301	rBV4	15414	30860	3.03%	0.197%
8	9.519	1378	1383	1402	rVB	31715	60739	5.97%	0.389%
9	10.055	1465	1471	1483	rBV	489725	643898	63.31%	4.120%
10	10.195	1488	1494	1504	rBV	18849	25669	2.52%	0.164%
11	10.299	1504	1511	1518	rBV	63318	90237	8.87%	0.577%
12	10.640	1562	1567	1577	rVB	75230	98162	9.65%	0.628%
13	10.963	1615	1620	1626	rVB	13084	15644	1.54%	0.100%
14	11.079	1634	1639	1648	rBV	419696	536923	52.79%	3.436%
15	11.305	1670	1676	1680	rBV	29085	36696	3.61%	0.235%
16	11.378	1683	1688	1696	rVV	99687	192512	18.93%	1.232%
17	11.451	1696	1700	1709	rVB	78562	112540	11.06%	0.720%
18	11.616	1721	1727	1733	rVB	94592	126916	12.48%	0.812%
19	11.750	1745	1749	1759	rVB	140126	185107	18.20%	1.185%
20	11.890	1765	1772	1776	rVB	21547	31436	3.09%	0.201%
21	11.969	1782	1785	1789	rVV	34249	40610	3.99%	0.260%
22	12.024	1789	1794	1800	rVB	487709	626387	61.59%	4.008%
23	12.085	1800	1804	1810	rBV	211483	266069	26.16%	1.703%
24	12.177	1816	1819	1823	rVB2	8591	11363	1.12%	0.073%
25	12.244	1823	1830	1832	rBV2	142572	185167	18.21%	1.185%
26	12.268	1832	1834	1838	rVB	90102	94015	9.24%	0.602%
27	12.317	1838	1842	1849	rVB4	161475	249289	24.51%	1.595%
28	12.420	1849	1859	1863	rBV2	50641	93487	9.19%	0.598%
29	12.463	1863	1866	1872	rVB	82145	99131	9.75%	0.634%
30	12.530	1872	1877	1879	rBV	75538	97735	9.61%	0.625%
31	12.555	1879	1881	1885	rVB	80590	76189	7.49%	0.488%
32	12.616	1886	1891	1894	rBV	161504	187427	18.43%	1.199%
33	12.646	1894	1896	1900	rVB	42808	43244	4.25%	0.277%
34	12.701	1900	1905	1908	rBV	109988	176812	17.38%	1.131%
35	12.798	1917	1921	1923	rBV	30071	40557	3.99%	0.260%
36	12.847	1926	1929	1933	rVB	78986	93010	9.14%	0.595%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
 Data File : VX032561.D
 Acq On : 03 Nov 2022 17:24
 Operator : JC/MD
 Sample : N5454-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1674

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_X\Method\82X102622W.M
 Title : SW846 8260

37	12.920	1933	1941	1944	rBV	137263	190142	18.69%	1.217%
38	12.963	1944	1948	1954	rVB	197855	258452	25.41%	1.654%
39	13.042	1954	1961	1964	rBV4	58888	119023	11.70%	0.762%
40	13.073	1964	1966	1968	rVV	25798	23882	2.35%	0.153%
41	13.097	1968	1970	1974	rVB2	8490	12139	1.19%	0.078%
42	13.170	1974	1982	1986	rBV4	156414	256276	25.20%	1.640%
43	13.213	1986	1989	1992	rVB2	47878	53165	5.23%	0.340%
44	13.292	1992	2002	2007	rBV2	618927	1017099	100.00%	6.508%
45	13.335	2007	2009	2012	rVV3	25896	33568	3.30%	0.215%
46	13.372	2012	2015	2019	rVV2	35465	51394	5.05%	0.329%
47	13.420	2019	2023	2032	rVB	402131	567805	55.83%	3.633%
48	13.506	2032	2037	2040	rBV2	61953	94721	9.31%	0.606%
49	13.542	2040	2043	2046	rBV	84661	106167	10.44%	0.679%
50	13.585	2046	2050	2055	rVV3	147629	252583	24.83%	1.616%
51	13.640	2056	2059	2060	rVV	51630	61404	6.04%	0.393%
52	13.664	2060	2063	2068	rVB2	131390	193773	19.05%	1.240%
53	13.798	2083	2085	2088	rVB3	31161	30235	2.97%	0.193%
54	13.853	2088	2094	2099	rVB2	218648	329538	32.40%	2.109%
55	13.926	2099	2106	2110	rBV2	238503	355015	34.90%	2.272%
56	13.969	2110	2113	2117	rBV	68666	71212	7.00%	0.456%
57	14.006	2117	2119	2121	rBV	15748	17354	1.71%	0.111%
58	14.036	2121	2124	2127	rVB	112693	115585	11.36%	0.740%
59	14.073	2127	2130	2134	rBV2	154626	178296	17.53%	1.141%
60	14.176	2142	2147	2149	rBV2	26207	46531	4.57%	0.298%
61	14.231	2149	2156	2160	rVV	465882	773035	76.00%	4.947%
62	14.323	2167	2171	2175	rBV	93394	123031	12.10%	0.787%
63	14.371	2175	2179	2183	rVV	189343	231662	22.78%	1.482%
64	14.414	2183	2186	2192	rVB4	41767	74171	7.29%	0.475%
65	14.481	2192	2197	2206	rBV2	429271	626511	61.60%	4.009%
66	14.615	2215	2219	2221	rBV	264655	339165	33.35%	2.170%
67	14.670	2225	2228	2233	rVB	149513	192963	18.97%	1.235%
68	14.725	2233	2237	2239	rBV2	59302	77992	7.67%	0.499%
69	14.755	2239	2242	2245	rBV	132026	121438	11.94%	0.777%
70	14.822	2249	2253	2255	rBV2	59875	75967	7.47%	0.486%
71	14.847	2255	2257	2262	rVB3	87674	88154	8.67%	0.564%
72	14.902	2262	2266	2274	rVB2	90363	147407	14.49%	0.943%
73	14.987	2277	2280	2284	rVB2	40534	56487	5.55%	0.361%
74	15.048	2284	2290	2294	rBV2	45229	75024	7.38%	0.480%
75	15.103	2295	2299	2302	rBV5	23883	36745	3.61%	0.235%
76	15.139	2302	2305	2308	rBV3	17426	25247	2.48%	0.162%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
 Data File : VX032561.D
 Acq On : 03 Nov 2022 17:24
 Operator : JC/MD
 Sample : N5454-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1674

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_X\Method\82X102622W.M
 Title : SW846 8260

77	15.182	2308	2312	2315	rBV	189376	288027	28.32%	1.843%
78	15.249	2320	2323	2325	rBV4	20808	28003	2.75%	0.179%
79	15.310	2331	2333	2339	rVB5	18604	32617	3.21%	0.209%
80	15.377	2340	2344	2352	rVB	58773	98963	9.73%	0.633%
81	15.487	2356	2362	2366	rBV3	181187	298460	29.34%	1.910%
82	15.566	2373	2375	2378	rBV3	13402	19983	1.96%	0.128%
83	15.615	2378	2383	2386	rVV2	91058	160080	15.74%	1.024%
84	15.652	2386	2389	2397	rVB2	60688	106068	10.43%	0.679%
85	15.841	2417	2420	2424	rBV3	20783	32871	3.23%	0.210%
86	15.993	2440	2445	2448	rBV5	16991	32547	3.20%	0.208%
87	16.023	2448	2450	2455	rVB3	26590	37821	3.72%	0.242%
88	16.091	2456	2461	2469	rBV4	34642	85593	8.42%	0.548%
89	16.377	2502	2508	2513	rVB2	59554	104905	10.31%	0.671%

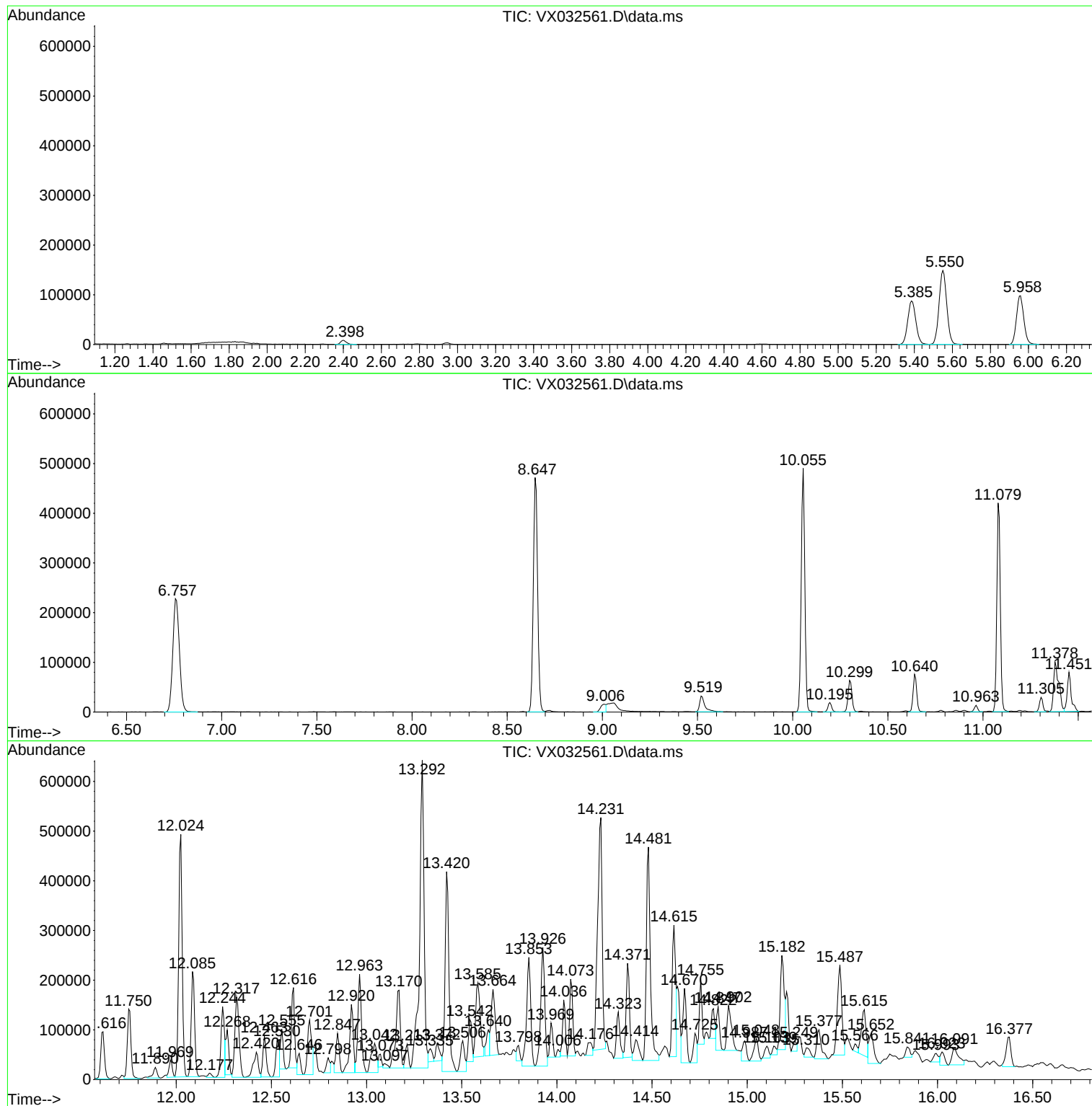
Sum of corrected areas: 15627326

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

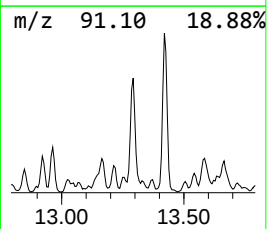
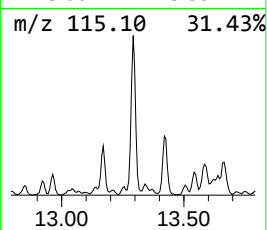
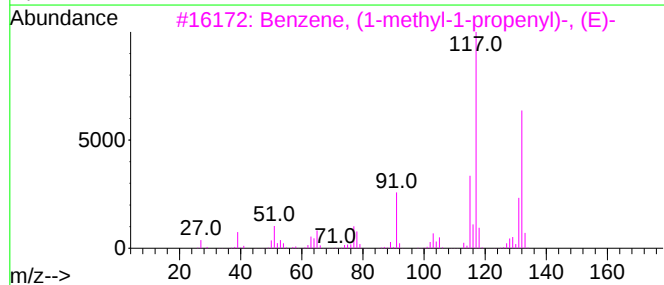
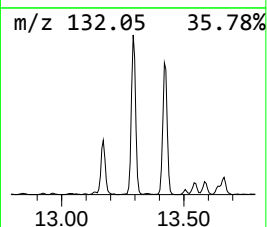
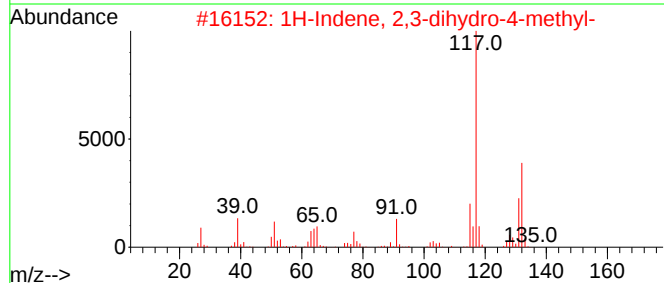
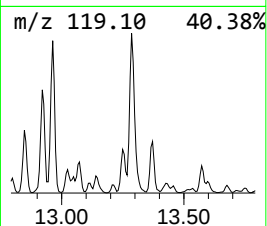
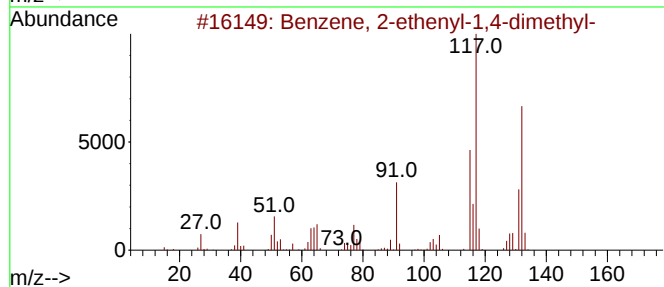
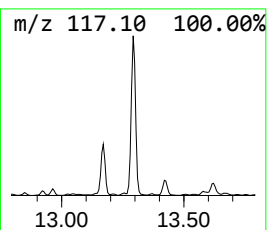
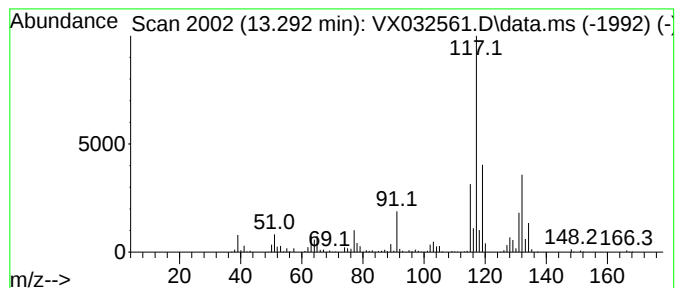
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	81.19 ug/l	1017100	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	92
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	89
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
5			1-Phenyl-1-butene	132	C10H12	000824-90-8	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

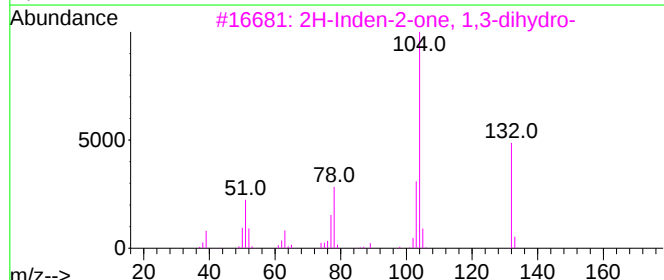
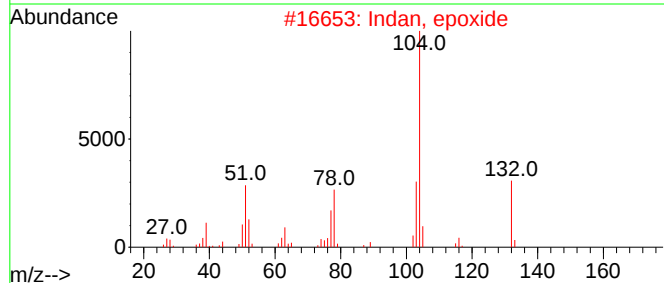
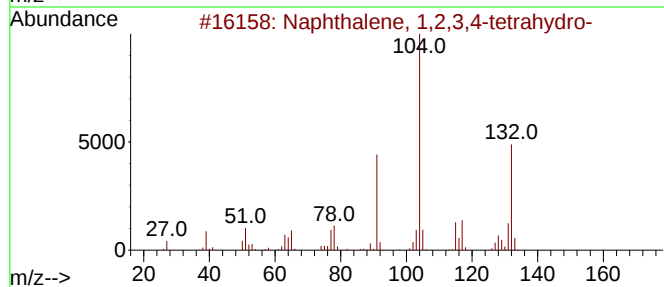
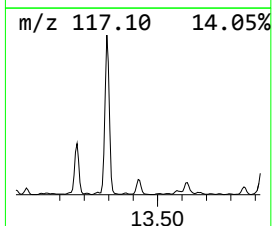
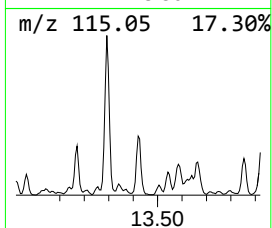
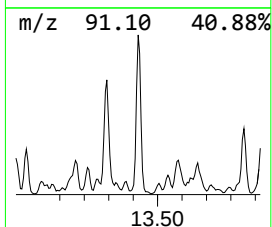
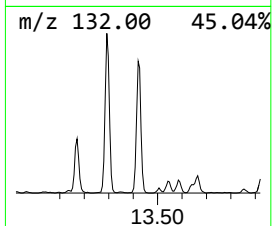
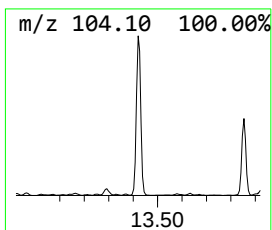
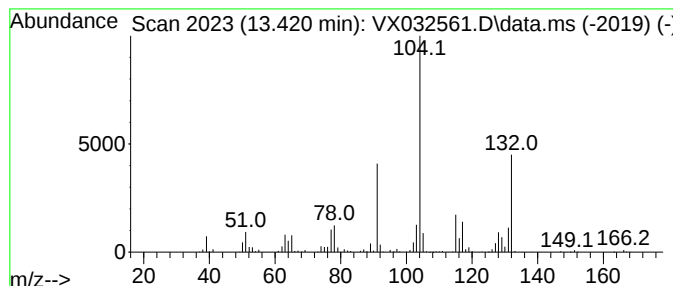
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	45.32 ug/l	567805	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	96
2			Indan, epoxide	132	C9H8O	000768-22-9	58
3			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
4			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	43
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-92-1	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

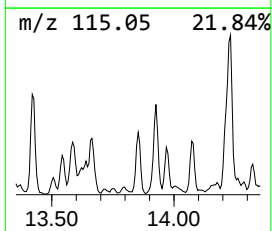
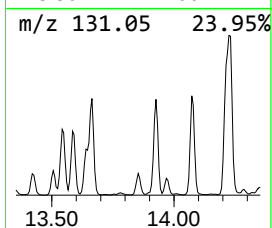
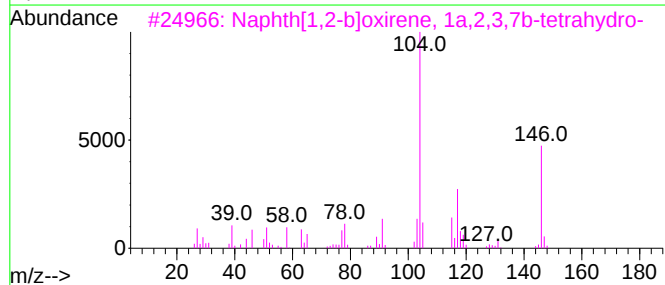
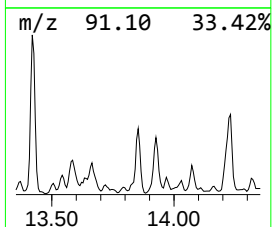
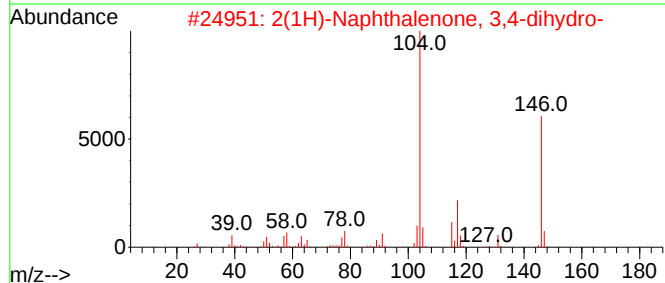
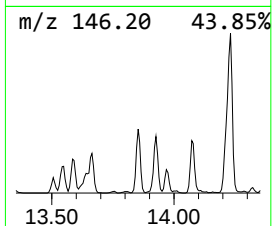
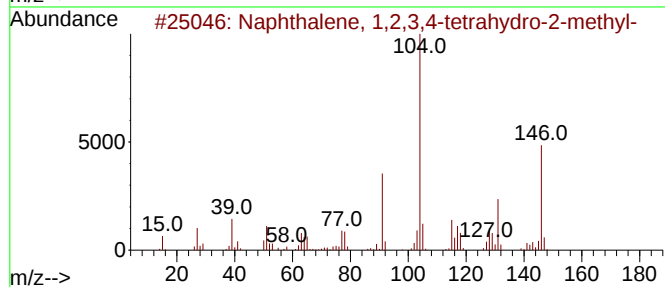
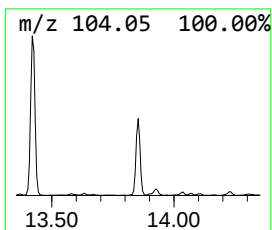
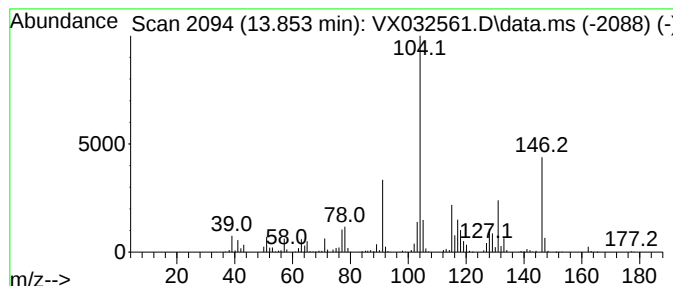
Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 4 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.853	26.30 ug/l	329538	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	003877-19-8	93
2	2(1H)-Naphthalenone, 3,4-dihydro-		146	C10H10O	000530-93-8	64
3	Naphth[1,2-b]oxirene, 1a,2,3,7b-...		146	C10H10O	002461-34-9	58
4	2-Methylbenzyl cyanide		131	C9H9N	022364-68-7	38
5	2-Azetidinone, 4-phenyl-		147	C9H9NO	005661-55-2	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

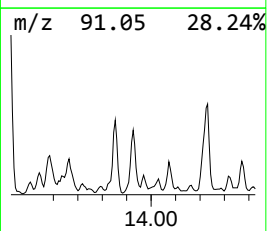
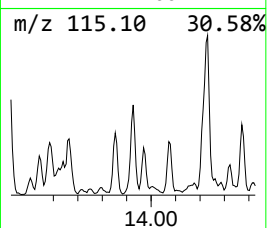
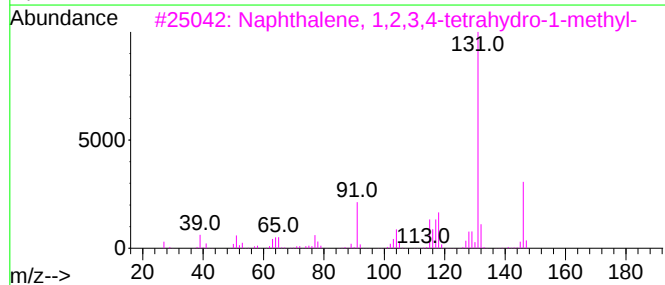
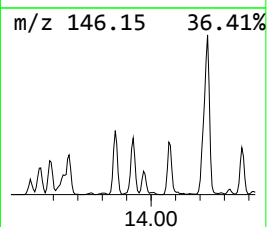
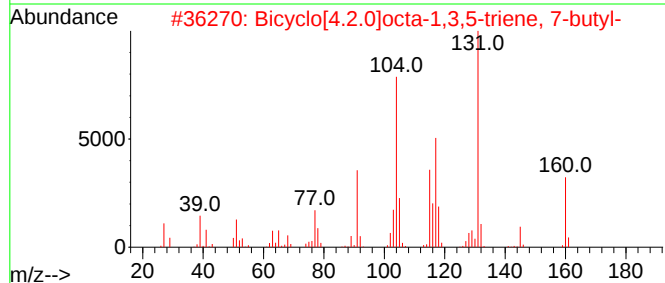
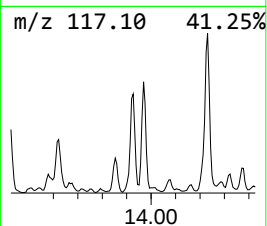
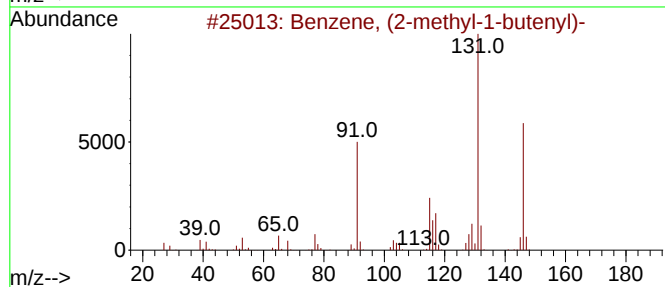
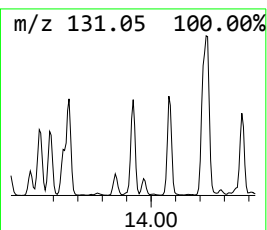
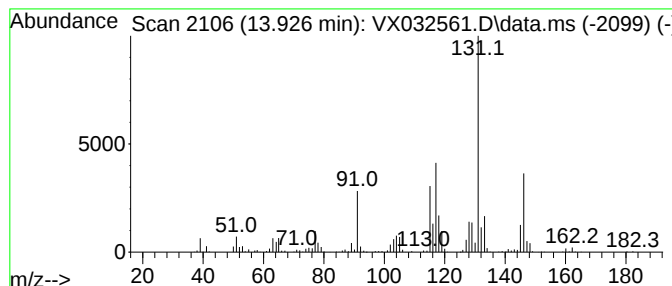
Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.926	28.34 ug/l	355015	1,4-Dichlorobenzene-d4		12.024	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-		146	C11H14	056253-64-6	93
2	Bicyclo[4.2.0]octa-1,3,5-triene,...		160	C12H16	078920-29-3	93
3	Naphthalene, 1,2,3,4-tetrahydro-...		146	C11H14	001559-81-5	92
4	Benzene, (3-methyl-2-butenyl)-		146	C11H14	004489-84-3	83
5	1H-Inden-1-one, 2,3-dihydro-2-me...		146	C10H10O	017496-14-9	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

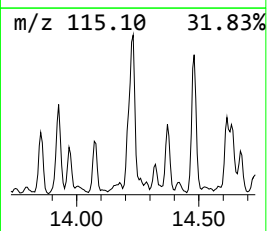
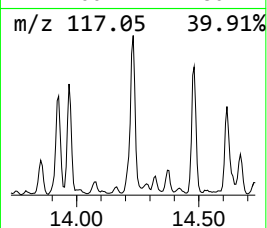
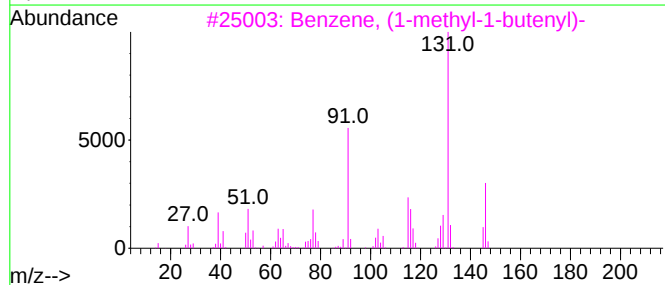
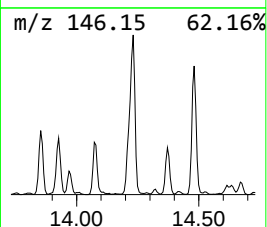
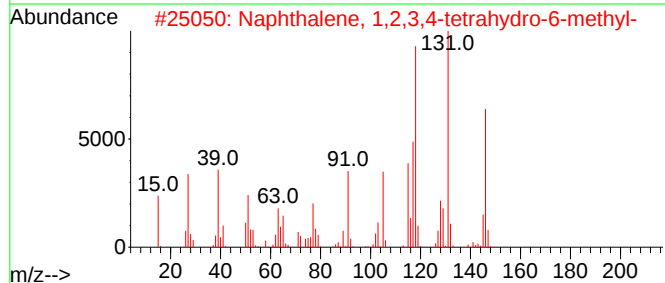
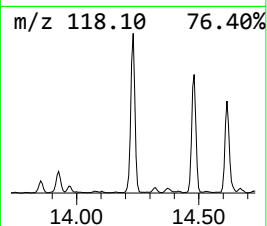
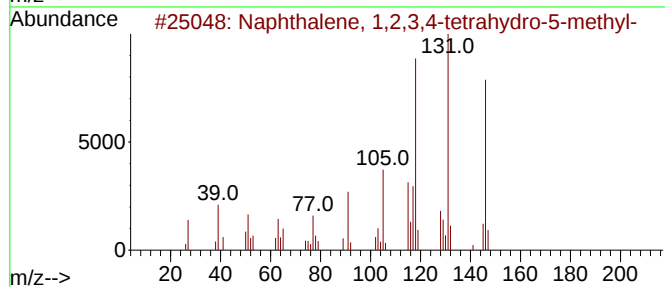
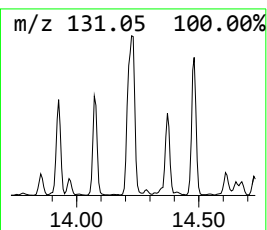
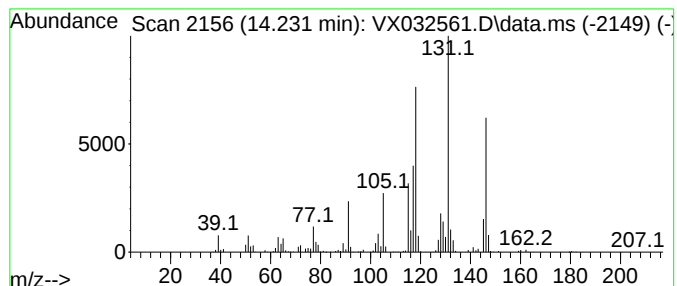
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.231	61.71 ug/l	773035	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	98
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

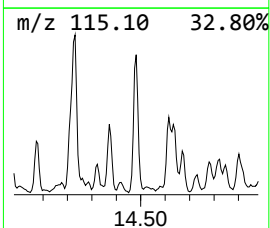
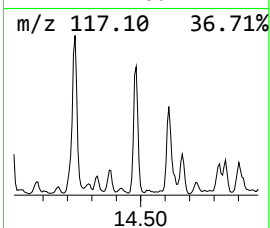
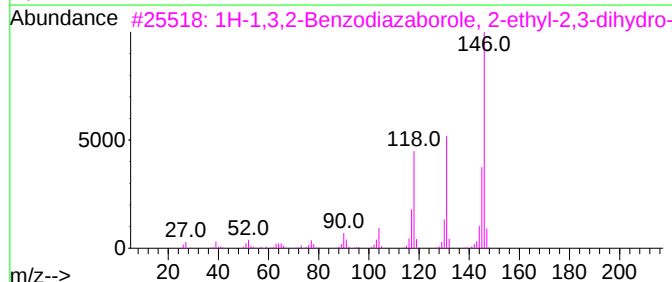
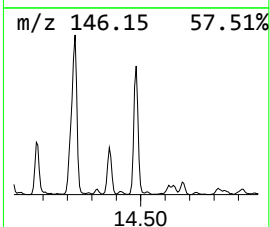
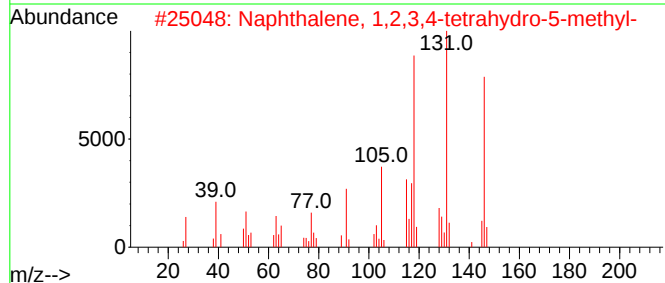
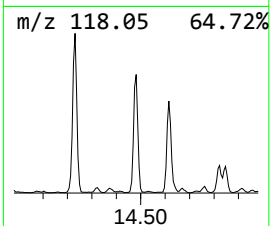
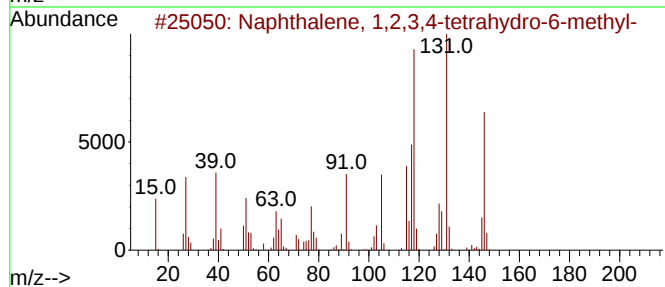
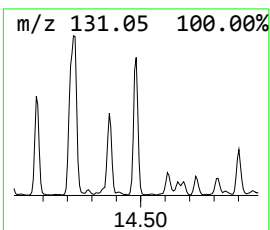
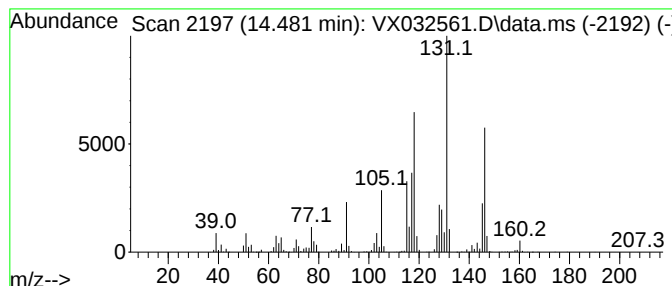
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	50.01 ug/l	626511	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3			1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	81
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5			2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

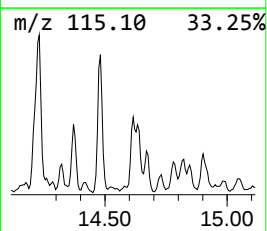
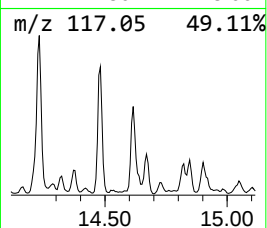
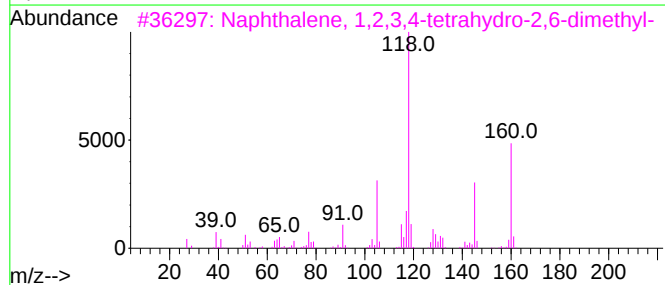
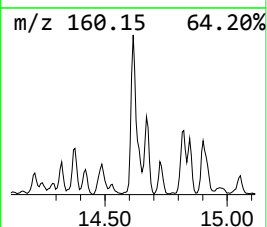
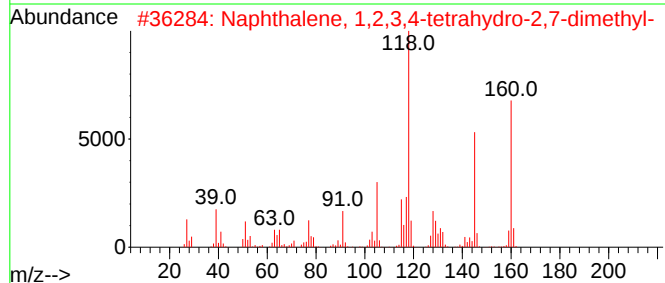
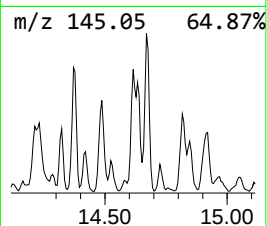
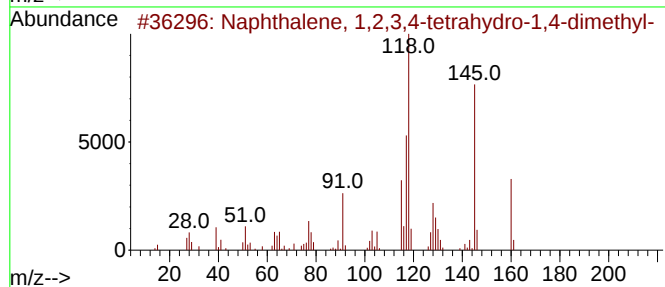
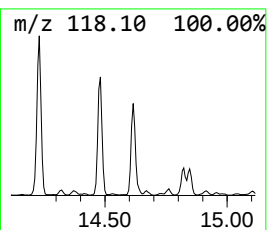
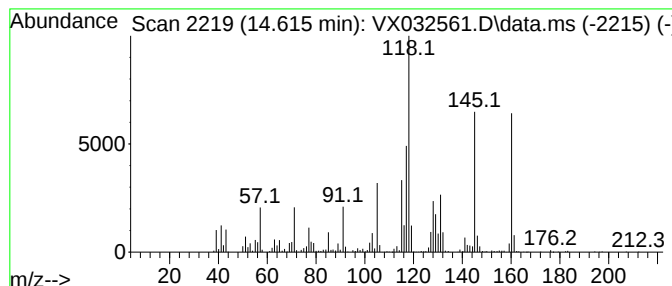
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.615	27.07 ug/l	339165	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	89
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	58
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	58
5			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

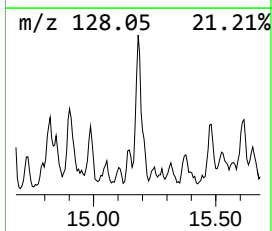
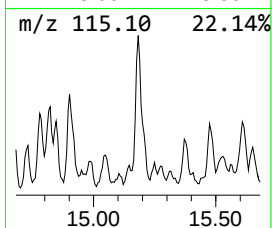
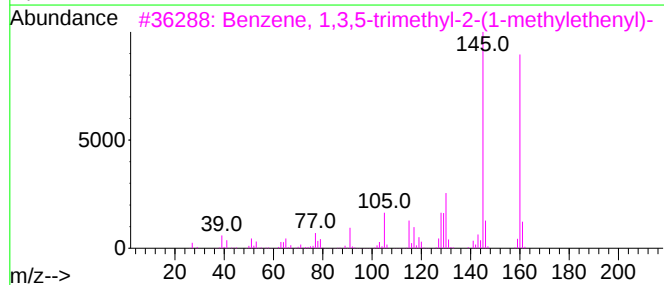
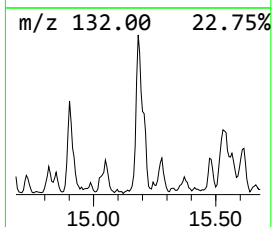
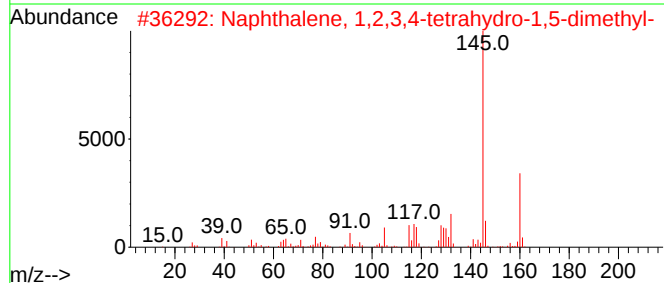
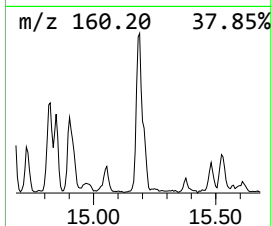
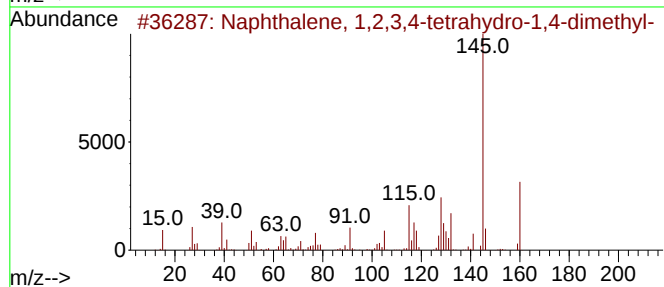
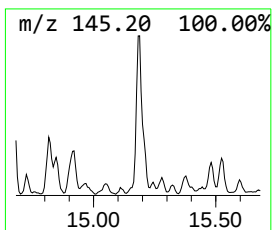
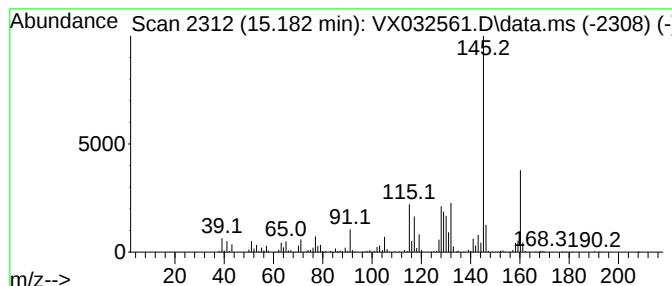
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.182	22.99 ug/l	288027	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	93
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	90
3			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	89
4			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	89
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

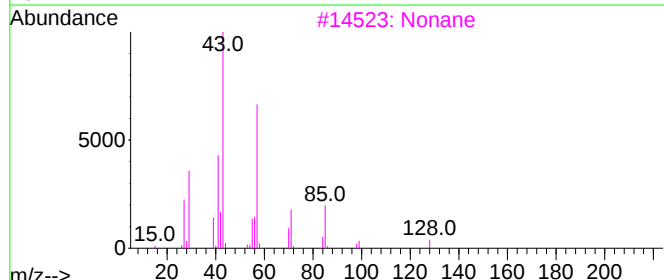
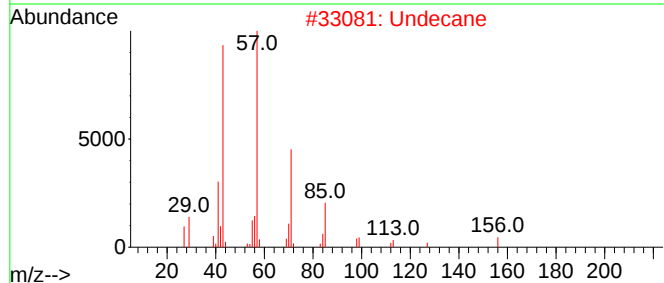
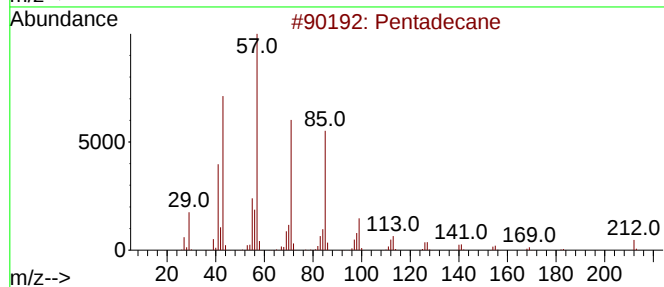
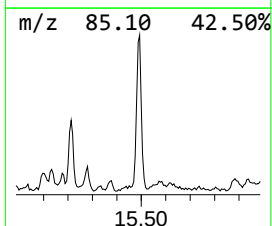
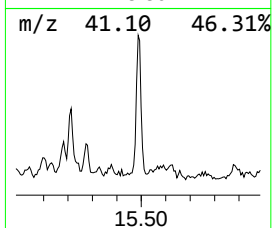
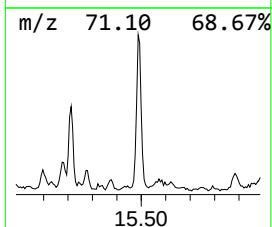
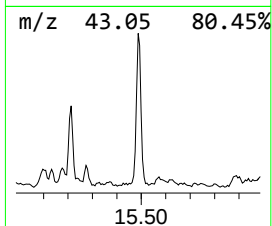
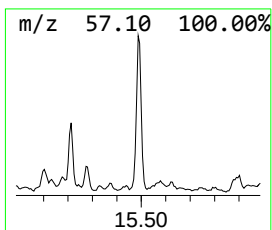
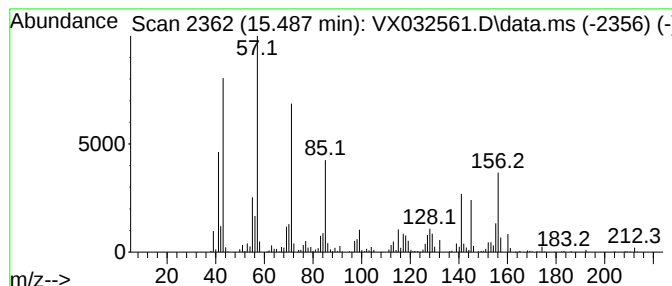
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.M
Quant Title : SW846 8260

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

Peak Number 10 Pentadecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.487	23.82 ug/l	298460	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane	212	C15H32	000629-62-9	95
2			Undecane	156	C11H24	001120-21-4	52
3			Nonane	128	C9H20	000111-84-2	50
4			Nonadecane	268	C19H40	000629-92-5	49
5			Hexadecane	226	C16H34	000544-76-3	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX110322\
Data File : VX032561.D
Acq On : 03 Nov 2022 17:24
Operator : JC/MD
Sample : N5454-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1674

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X102622W.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, 2-ethe...	13.292	81.2	ug/l	1017100	4	12.024	626387	50.0
Naphthalene, 1,...	13.420	45.3	ug/l	567805	4	12.024	626387	50.0
Naphthalene, 1,...	13.853	26.3	ug/l	329538	4	12.024	626387	50.0
Benzene, (2-met...	13.926	28.3	ug/l	355015	4	12.024	626387	50.0
Naphthalene, 1,...	14.231	61.7	ug/l	773035	4	12.024	626387	50.0
Naphthalene, 1,...	14.481	50.0	ug/l	626511	4	12.024	626387	50.0
Naphthalene, 1,...	14.615	27.1	ug/l	339165	4	12.024	626387	50.0
Naphthalene, 1,...	15.182	23.0	ug/l	288027	4	12.024	626387	50.0
Pentadecane	15.487	23.8	ug/l	298460	4	12.024	626387	50.0