

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
 Data File : VX032223.D
 Acq On : 19 Oct 2022 21:14
 Operator : JC/MD
 Sample : N5170-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TRE-1661

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\82X101922W.M
 Title : SW846 8260

Signal : TIC: VX032223.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.294	30	34	37	rBV	4457	7632	1.11%	0.103%
2	2.386	205	213	225	rVB	9182	17904	2.61%	0.241%
3	5.385	694	705	721	rBV	83858	242532	35.39%	3.265%
4	5.556	721	733	746	rVV	139034	394390	57.55%	5.310%
5	5.958	787	799	812	rBV	94220	247656	36.14%	3.334%
6	6.763	921	931	946	rBV	210787	500190	72.99%	6.734%
7	8.647	1234	1240	1249	rBV	436027	685243	100.00%	9.226%
8	10.055	1465	1471	1485	rVB	428731	569096	83.05%	7.662%
9	11.079	1634	1639	1653	rBV	364917	469322	68.49%	6.319%
10	11.231	1660	1664	1671	rVB2	5375	8949	1.31%	0.120%
11	11.402	1681	1692	1695	rBV	6382	13225	1.93%	0.178%
12	11.670	1732	1736	1741	rBV2	5397	7526	1.10%	0.101%
13	12.024	1788	1794	1801	rBV	412975	518510	75.67%	6.981%
14	12.085	1801	1804	1808	rBV	9927	12556	1.83%	0.169%
15	12.268	1828	1834	1838	rVB3	11060	17223	2.51%	0.232%
16	12.317	1838	1842	1849	rVB3	29061	47046	6.87%	0.633%
17	12.408	1853	1857	1863	rBV5	6474	12084	1.76%	0.163%
18	12.463	1863	1866	1873	rVB	8979	11915	1.74%	0.160%
19	12.615	1885	1891	1894	rBV	45400	56122	8.19%	0.756%
20	12.725	1899	1909	1914	rBV4	19443	44471	6.49%	0.599%
21	12.798	1916	1921	1925	rBV2	15761	24377	3.56%	0.328%
22	12.847	1925	1929	1934	rVB3	21573	33172	4.84%	0.447%
23	12.920	1934	1941	1945	rBV	29411	42582	6.21%	0.573%
24	12.963	1945	1948	1955	rVB2	13057	17811	2.60%	0.240%
25	13.042	1955	1961	1964	rBV3	22042	50531	7.37%	0.680%
26	13.067	1964	1965	1968	rVB	10223	8657	1.26%	0.117%
27	13.140	1974	1977	1978	rBV2	15895	16550	2.42%	0.223%
28	13.213	1984	1989	1992	rBV3	19869	25125	3.67%	0.338%
29	13.262	1992	1997	1999	rBV4	24631	39676	5.79%	0.534%
30	13.292	1999	2002	2007	rVB3	33427	51623	7.53%	0.695%
31	13.371	2012	2015	2020	rVV	27092	38493	5.62%	0.518%
32	13.426	2020	2024	2033	rVB3	20556	42984	6.27%	0.579%
33	13.505	2033	2037	2040	rBV4	15288	24471	3.57%	0.329%
34	13.548	2040	2044	2046	rVV	24949	33135	4.84%	0.446%
35	13.579	2046	2049	2055	rVV3	76540	143534	20.95%	1.933%
36	13.640	2056	2059	2060	rVV	38524	44108	6.44%	0.594%

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Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

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Stop Thrs : 0

Peak Location: TOP

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
 Title : SW846 8260

37	13.664	2060	2063	2069	rVV2	60596	104633	15.27%	1.409%
38	13.719	2069	2072	2074	rVV	9321	12423	1.81%	0.167%
39	13.749	2075	2077	2081	rVB2	12061	13979	2.04%	0.188%
40	13.798	2081	2085	2088	rBV3	15901	22162	3.23%	0.298%

41	13.853	2088	2094	2099	rVB3	66605	104170	15.20%	1.403%
42	13.926	2099	2106	2111	rBV3	102424	159383	23.26%	2.146%
43	13.969	2111	2113	2117	rVV2	19043	25202	3.68%	0.339%
44	14.005	2117	2119	2122	rVV	15312	18862	2.75%	0.254%
45	14.036	2122	2124	2127	rVB2	18987	18203	2.66%	0.245%

46	14.072	2127	2130	2134	rBV3	30703	40475	5.91%	0.545%
47	14.164	2142	2145	2148	rVB2	13526	14068	2.05%	0.189%
48	14.213	2148	2153	2163	rBV2	120669	250880	36.61%	3.378%
49	14.286	2163	2165	2167	rVV	11376	10568	1.54%	0.142%
50	14.322	2167	2171	2175	rVV	48573	59850	8.73%	0.806%

51	14.371	2175	2179	2183	rVV	106158	130994	19.12%	1.764%
52	14.420	2183	2187	2192	rVB2	28874	45685	6.67%	0.615%
53	14.481	2192	2197	2202	rBV2	130296	190561	27.81%	2.566%
54	14.524	2202	2204	2208	rVV3	10653	10542	1.54%	0.142%
55	14.615	2215	2219	2225	rBV3	131526	217656	31.76%	2.930%

56	14.670	2225	2228	2233	rVB	103267	139546	20.36%	1.879%
57	14.731	2233	2238	2240	rBV	32034	46319	6.76%	0.624%
58	14.755	2240	2242	2245	rVB3	22504	21088	3.08%	0.284%
59	14.816	2245	2252	2255	rBV2	72721	123605	18.04%	1.664%
60	14.847	2255	2257	2261	rVB2	58072	65158	9.51%	0.877%

61	14.902	2261	2266	2273	rBV2	67437	129271	18.86%	1.740%
62	15.054	2284	2291	2295	rBV2	30269	54858	8.01%	0.739%
63	15.109	2296	2300	2302	rVV4	5193	7143	1.04%	0.096%
64	15.182	2307	2312	2320	rVV	125001	258273	37.69%	3.477%
65	15.255	2320	2324	2331	rVV7	23179	59362	8.66%	0.799%

66	15.328	2331	2336	2339	rVB5	10343	16478	2.40%	0.222%
67	15.377	2339	2344	2352	rBV4	29577	53231	7.77%	0.717%
68	15.481	2357	2361	2365	rVV2	52562	90430	13.20%	1.218%
69	15.529	2365	2369	2373	rVV2	35453	65498	9.56%	0.882%
70	15.572	2373	2376	2378	rVV3	15746	23324	3.40%	0.314%

71	15.609	2378	2382	2387	rVV3	35421	66858	9.76%	0.900%
72	15.651	2387	2389	2397	rVB6	20378	35234	5.14%	0.474%
73	15.725	2397	2401	2403	rBV5	5634	9040	1.32%	0.122%
74	15.755	2403	2406	2409	rBV3	6846	9914	1.45%	0.133%
75	15.847	2417	2421	2425	rBV5	15829	27763	4.05%	0.374%

76	15.889	2425	2428	2434	rVB4	12427	22834	3.33%	0.307%
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ClientSampleId :
TRE-1661

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

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Peak separation: 5

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Title : SW846 8260

77	16.023	2447	2450	2456	rVB3	15740	25021	3.65%	0.337%
78	16.090	2456	2461	2464	rBV2	15323	27164	3.96%	0.366%
79	16.298	2491	2495	2499	rVB6	7446	12754	1.86%	0.172%
80	16.377	2500	2508	2512	rBV4	11946	28652	4.18%	0.386%
81	16.517	2528	2531	2534	rBV2	6021	10052	1.47%	0.135%
82	16.596	2541	2544	2548	rVB3	7653	12414	1.81%	0.167%
83	16.657	2549	2554	2560	rVB4	8139	15331	2.24%	0.206%

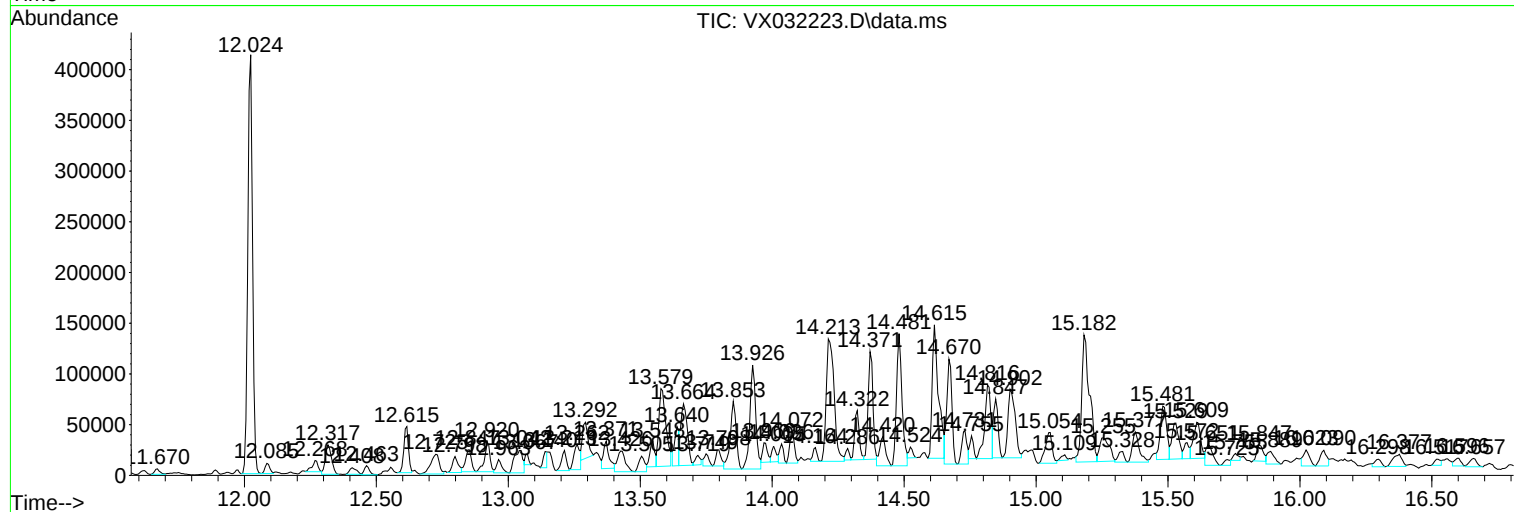
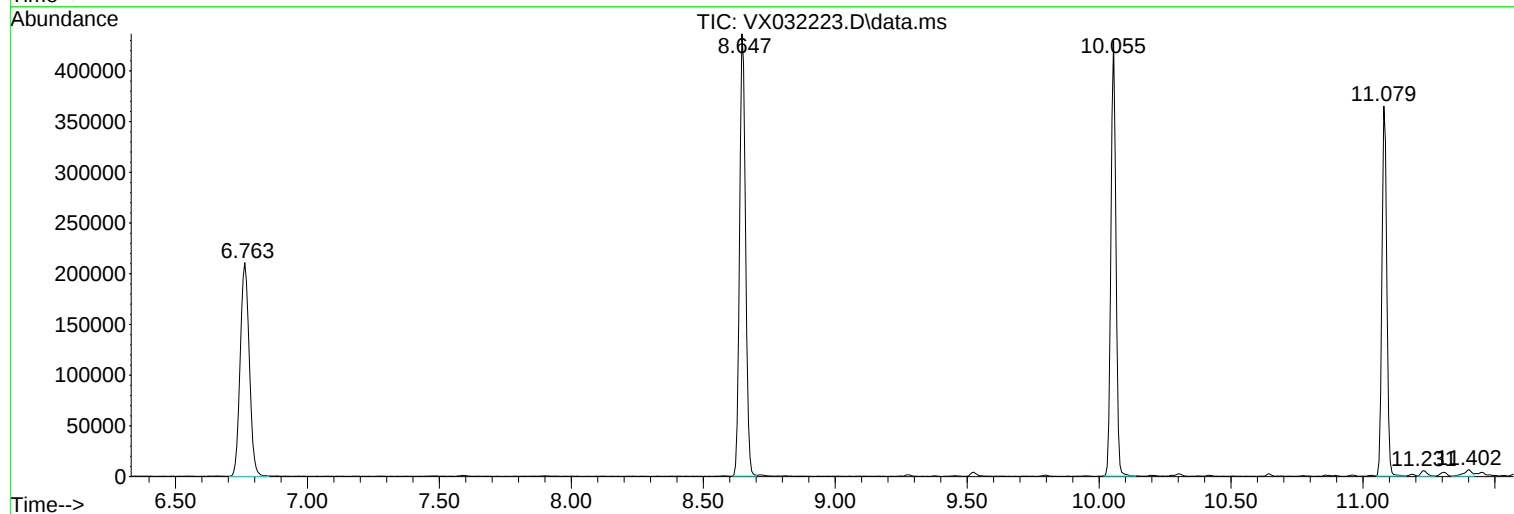
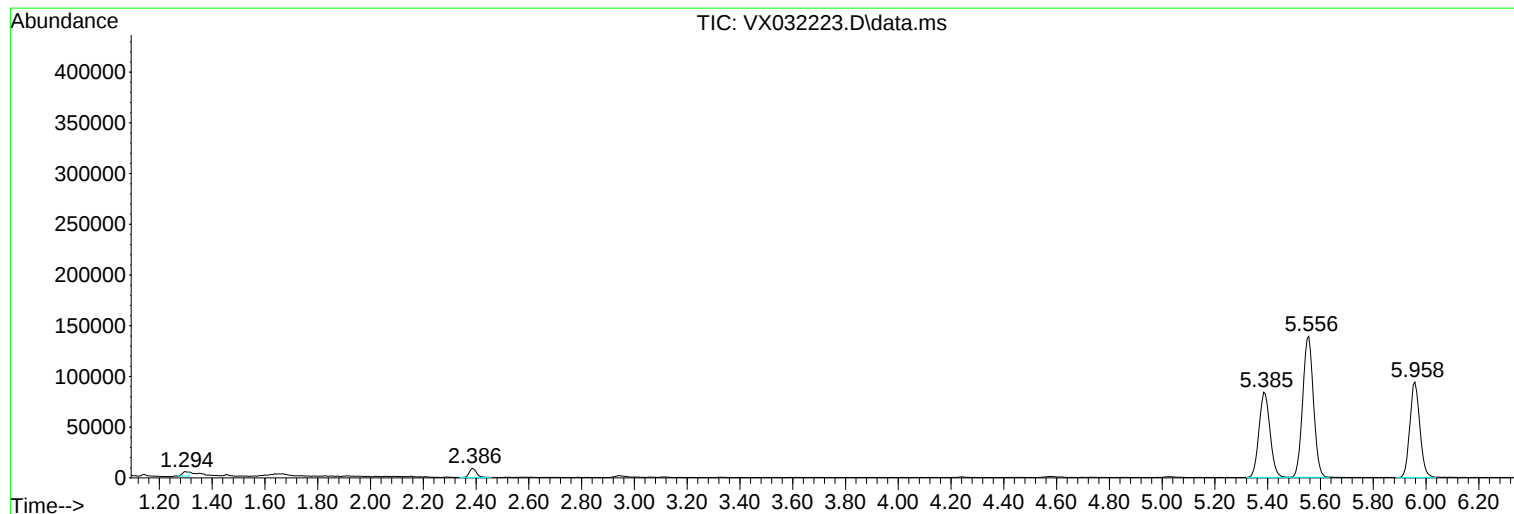
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Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
Quant Title : SW846 8260

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



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ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

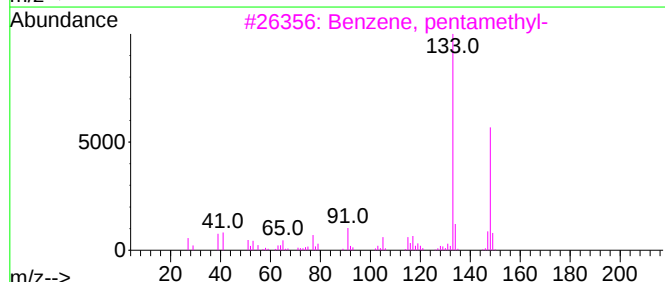
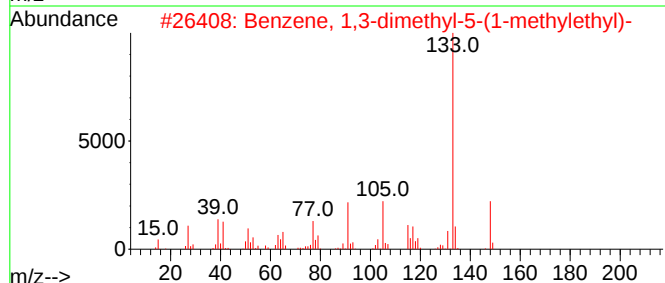
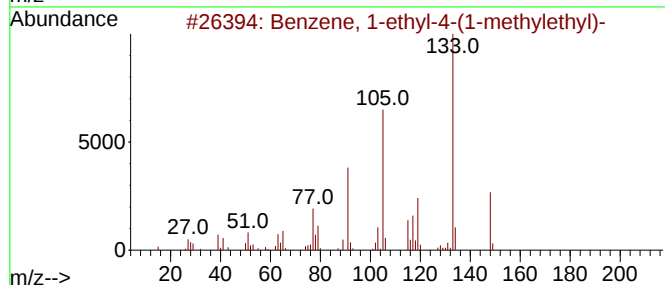
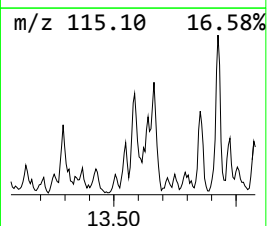
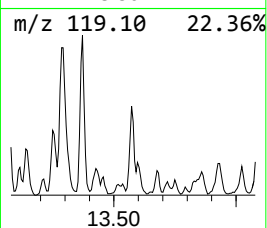
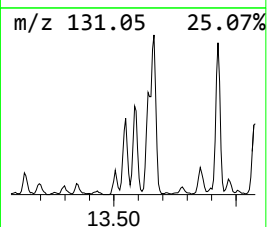
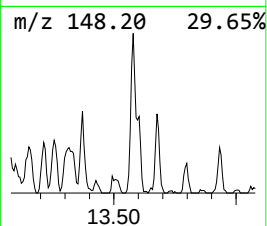
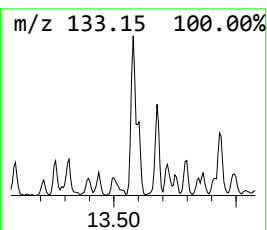
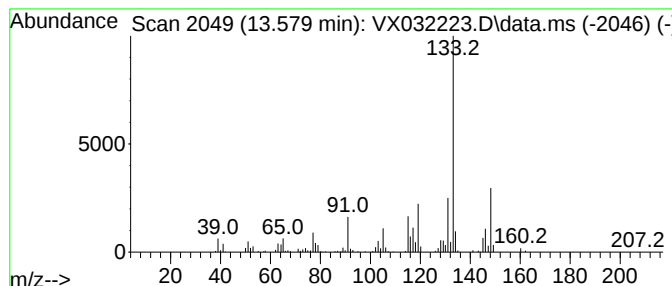
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-4-(1-methyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.579	13.84 ug/l	143534	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	76
2			Benzene, 1,3-dimethyl-5-(1-methyl-...	148	C11H16	004706-90-5	62
3			Benzene, pentamethyl-	148	C11H16	000700-12-9	58
4			3,4-Dimethylcumene	148	C11H16	1000370-34-1	58
5			1,5,6,7-Tetramethylbicyclo[3.2.0...	148	C11H16	134329-46-7	52



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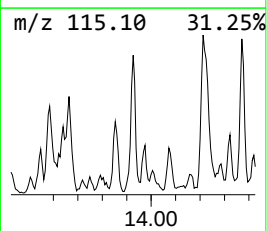
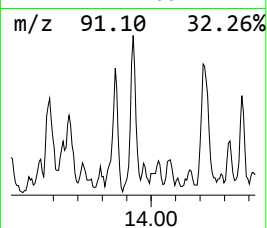
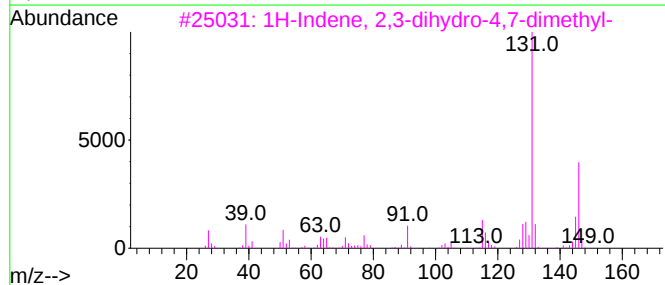
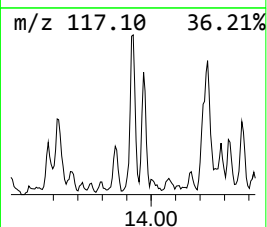
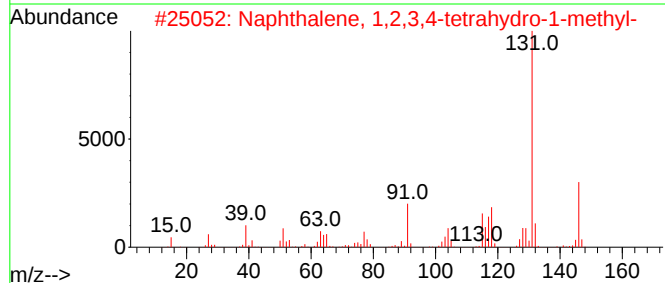
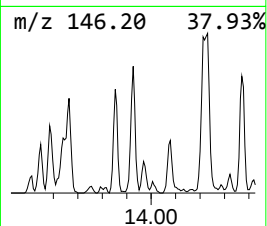
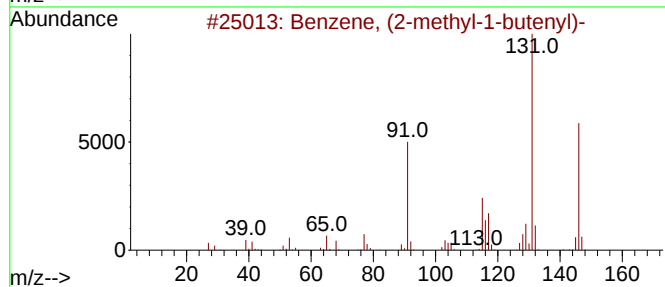
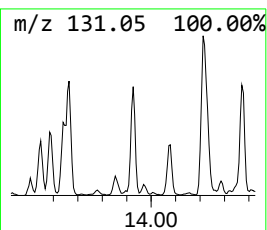
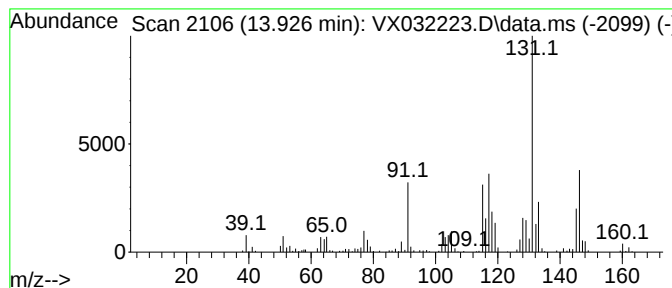
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	15.37 ug/l	159383	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			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	92
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	92
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	89
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83



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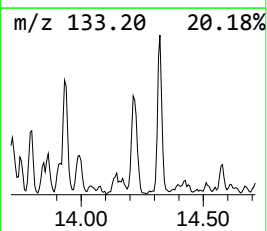
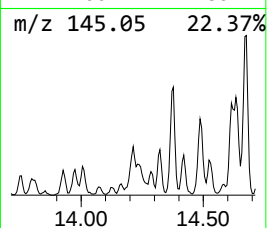
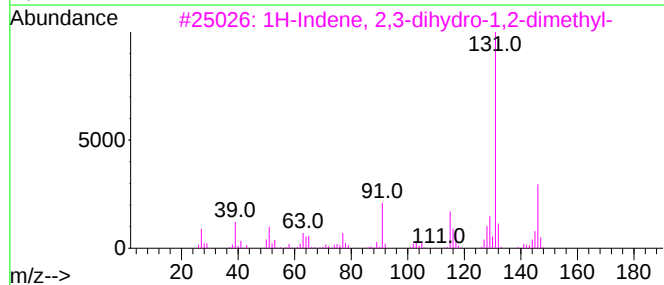
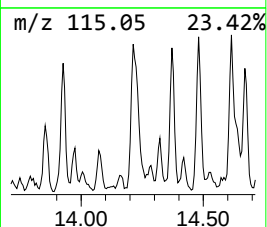
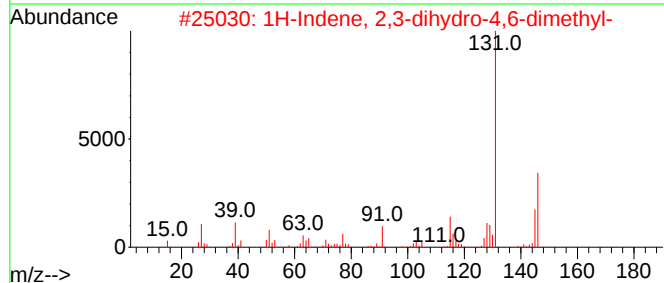
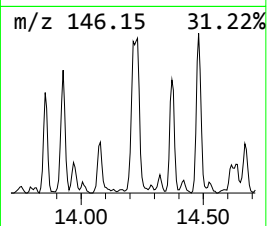
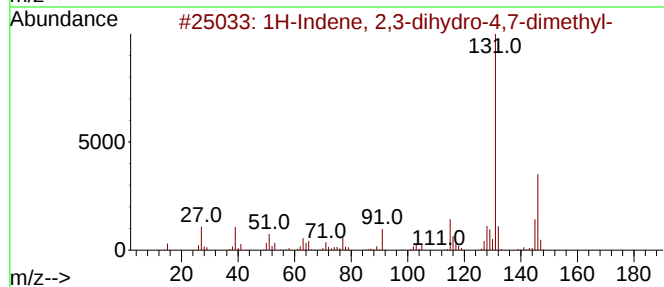
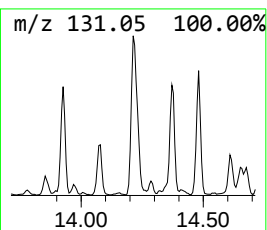
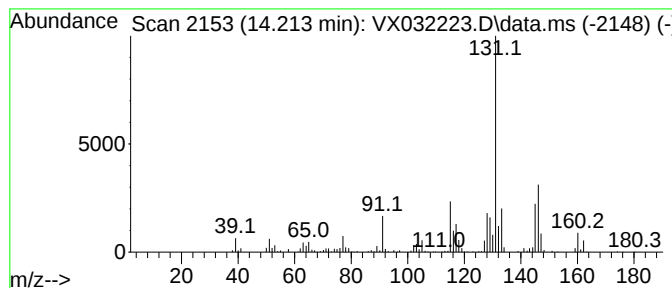
Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.213	24.19 ug/l	250880	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...		146	C11H14	006682-71-9	76
2	1H-Indene, 2,3-dihydro-4,6-dimet...		146	C11H14	001685-82-1	76
3	1H-Indene, 2,3-dihydro-1,2-dimet...		146	C11H14	017057-82-8	74
4	Benzene, (2-methyl-1-butenyl)-		146	C11H14	056253-64-6	74
5	Benzene, (1-methyl-1-butenyl)-		146	C11H14	053172-84-2	74



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MSVOA_X
ClientSampleId :
TRE-1661

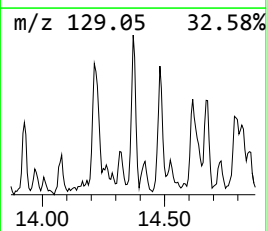
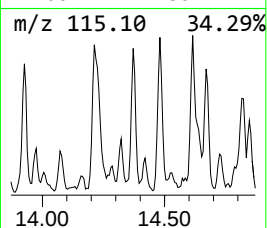
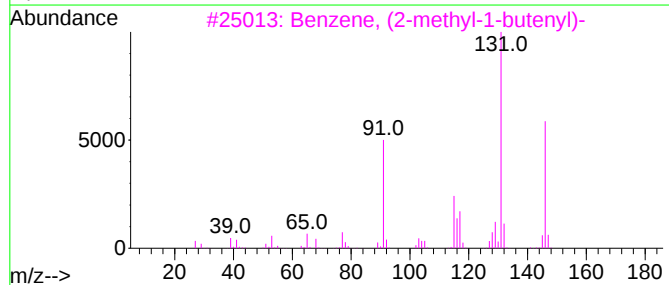
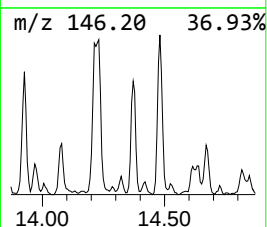
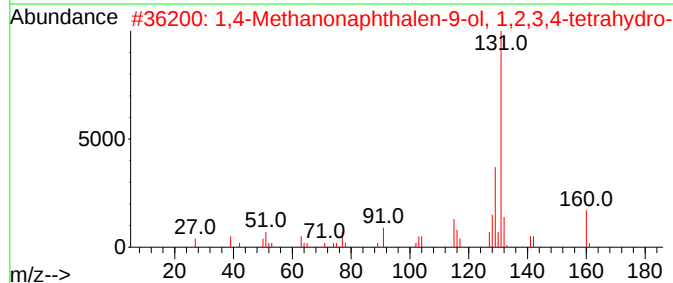
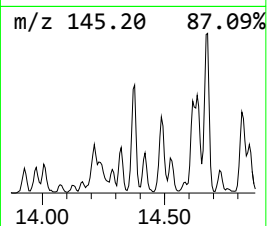
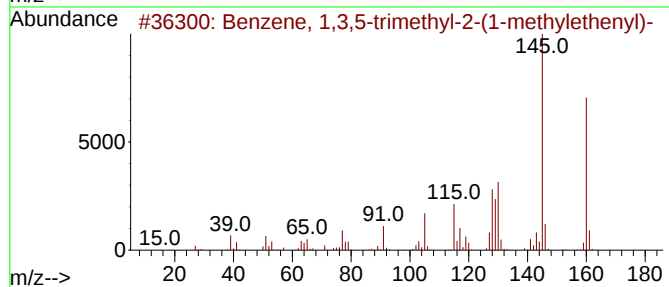
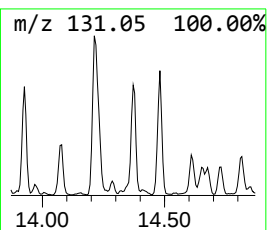
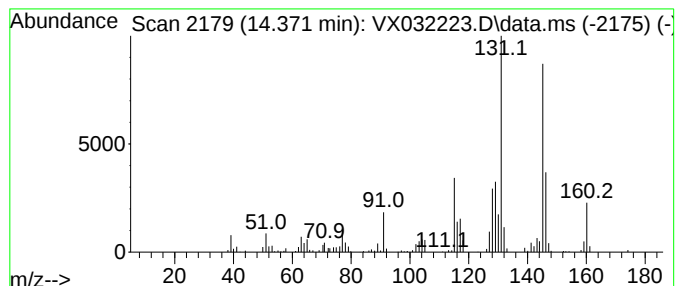
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.371	12.63 ug/l	130994	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	78
2			1,4-Methanonaphthalen-9-ol, 1,2,...	160	C11H12O	055255-94-2	46
3			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	45
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	43
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

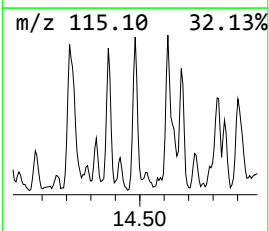
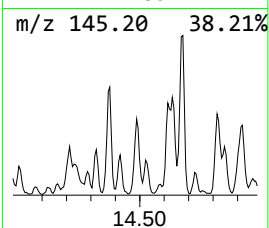
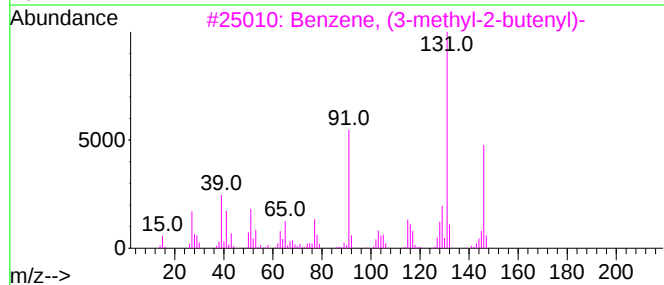
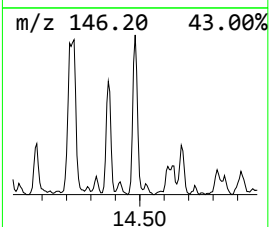
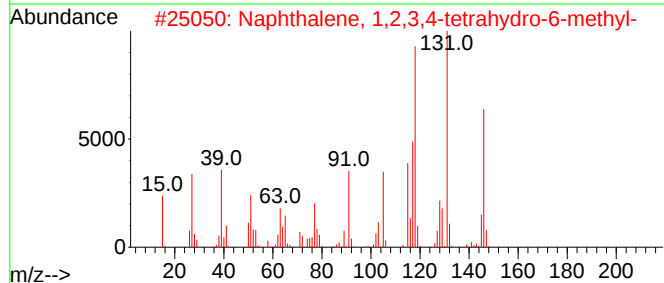
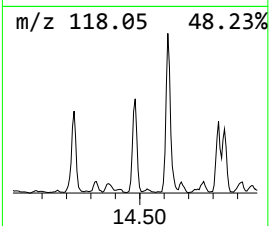
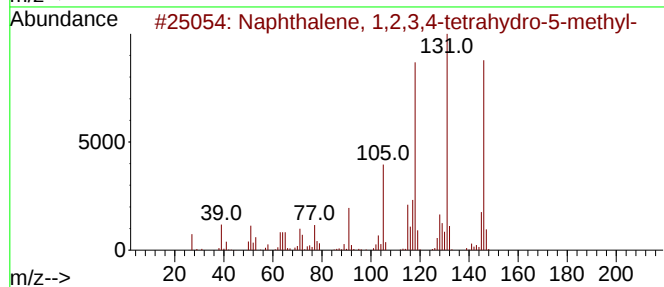
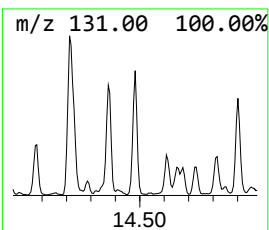
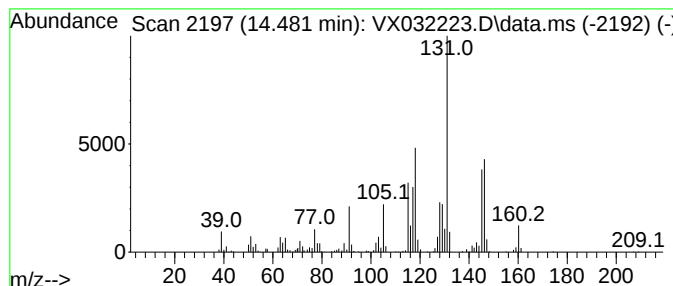
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, 1,2,3,4-tetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	18.38 ug/l	190561	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	68
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	55
3			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	53
4			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	53
5			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

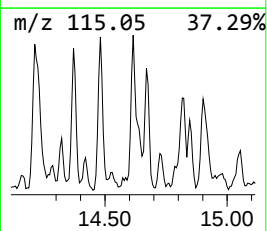
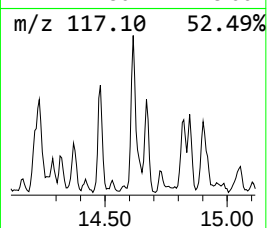
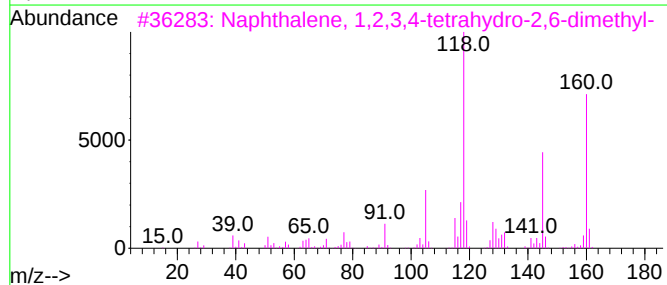
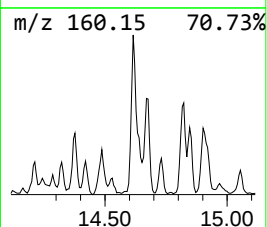
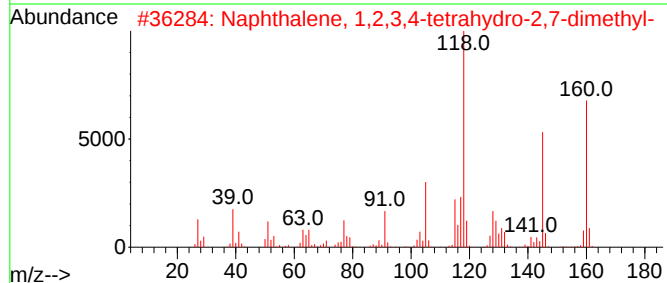
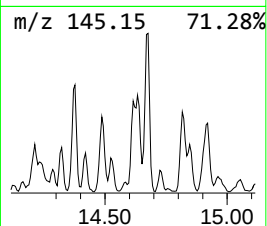
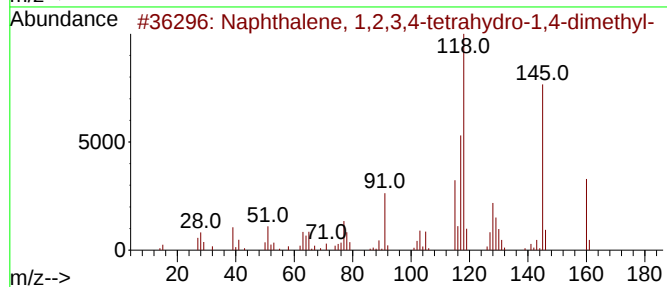
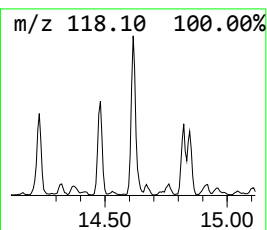
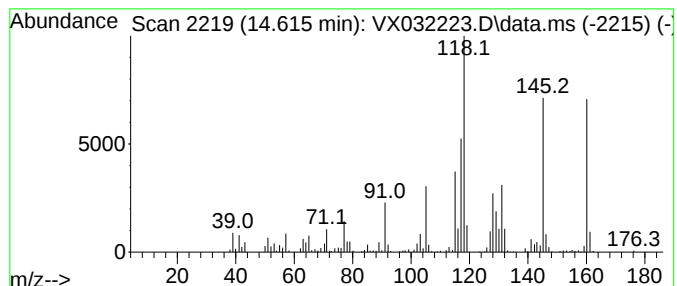
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.615	20.99 ug/l	217656	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	76
4			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	64
5			1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

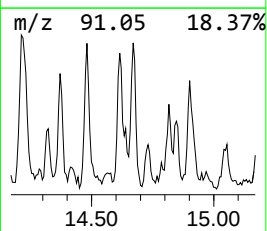
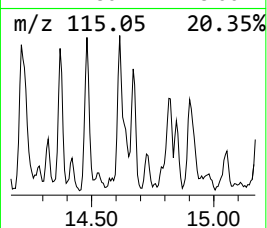
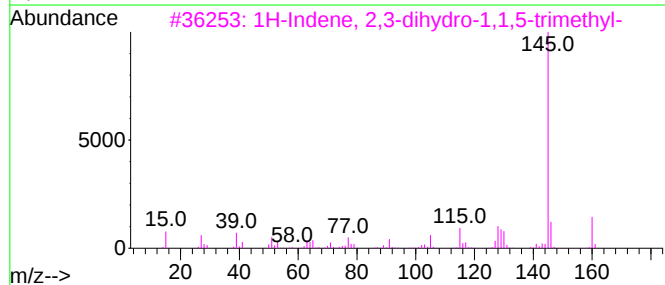
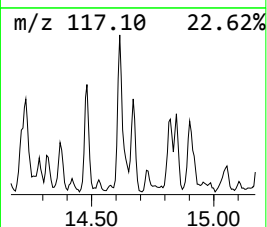
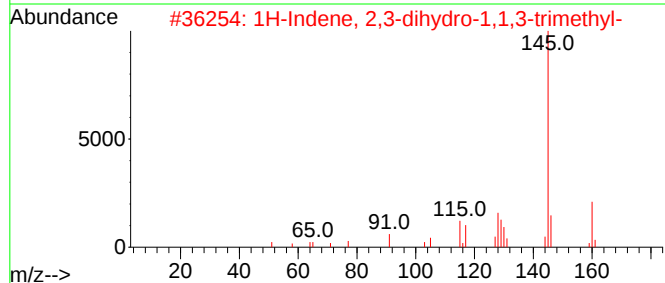
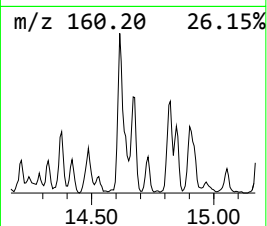
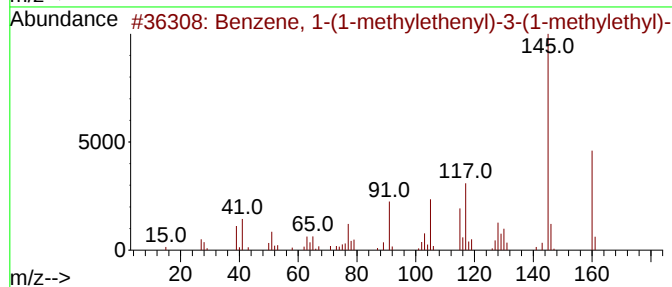
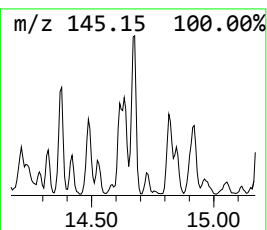
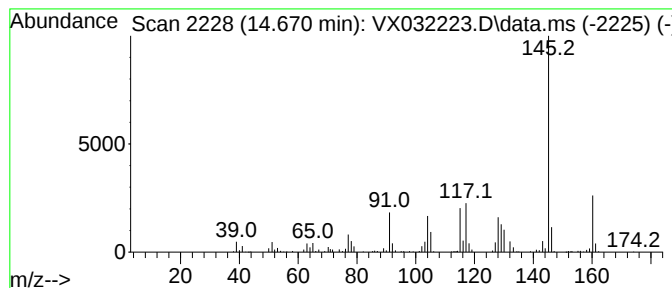
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-(1-methylethenyl)... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.670	13.46 ug/l	139546	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	90
2			1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	90
3			1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	87
4			Benzene, (2,2-dimethyl-1-methyle...	160	C12H16	005676-29-9	87
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

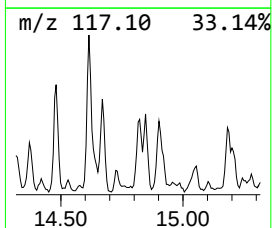
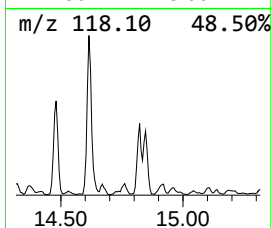
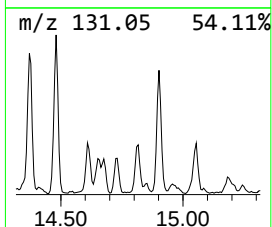
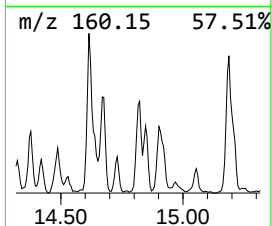
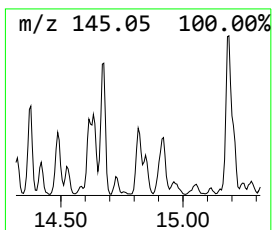
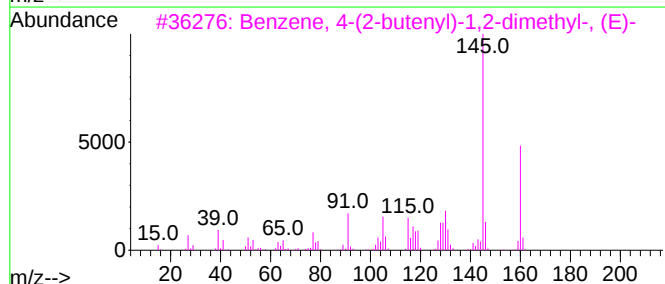
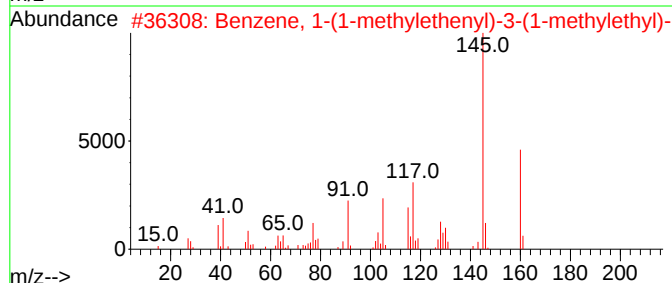
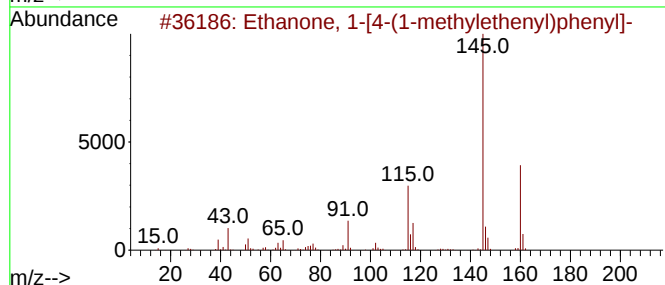
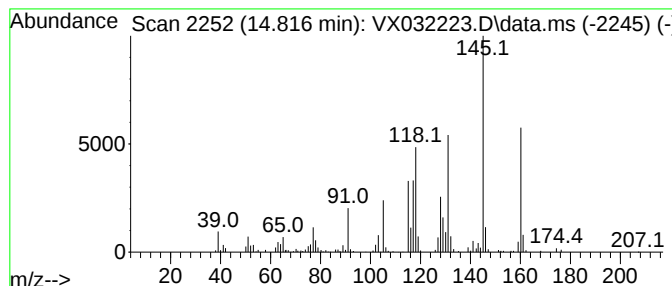
Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
Quant Title : SW846 8260

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

Peak Number 8 Ethanone, 1-[4-(1-methyleth... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.816	11.92 ug/l	123605	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Ethanone, 1-[4-(1-methylethenyl)...		160	C11H12O	005359-04-6	87
2	Benzene, 1-(1-methylethenyl)-3-(...		160	C12H16	001129-29-9	64
3	Benzene, 4-(2-butenyl)-1,2-dimet...		160	C12H16	054340-86-2	64
4	Naphthalene, 1,2,3,4-tetrahydro-...		160	C12H16	004175-54-6	62
5	1H-Indene, 2,3-dihydro-1,1,5-tri...		160	C12H16	040650-41-7	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

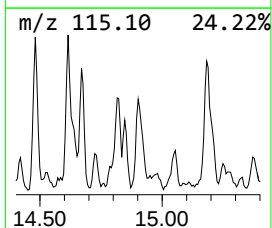
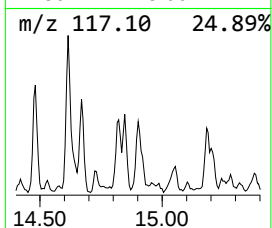
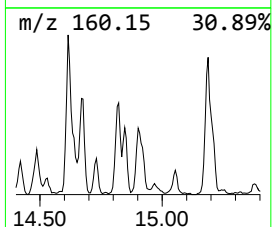
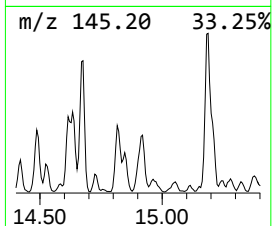
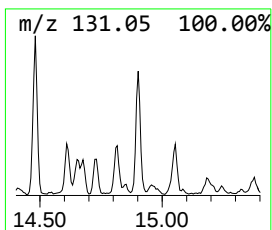
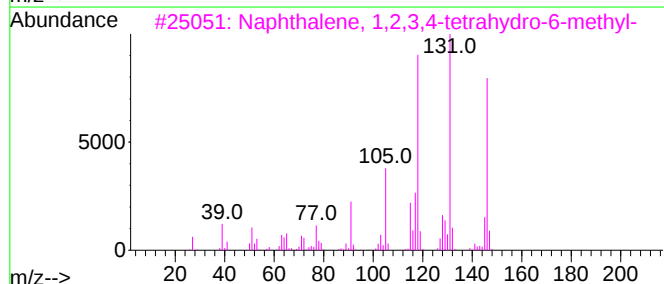
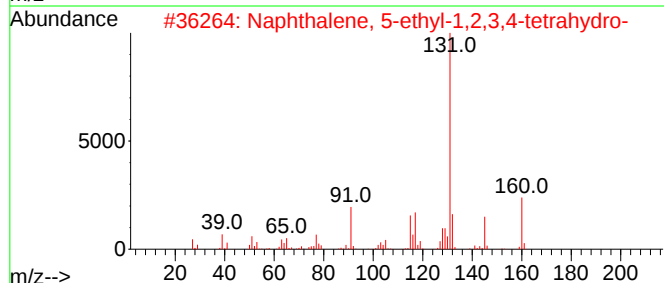
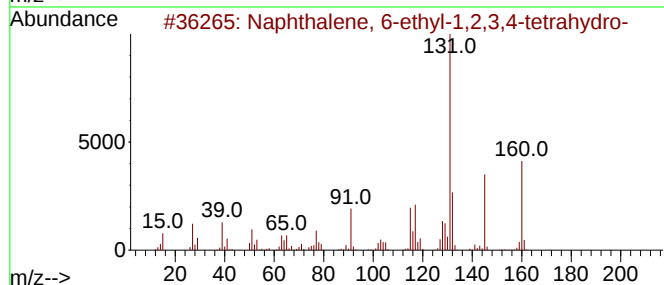
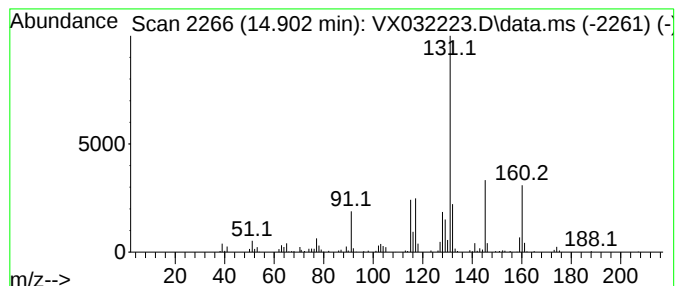
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 6-ethyl-1,2,3,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.902	12.47 ug/l	129271	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	94
2			Naphthalene, 5-ethyl-1,2,3,4-tet...	160	C12H16	042775-75-7	91
3			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	59
4			1H-Indene, 2,3-dihydro-4-propyl-	160	C12H16	092013-16-6	58
5			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

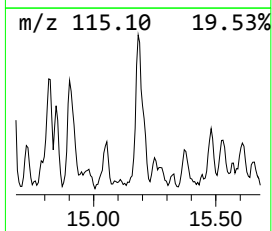
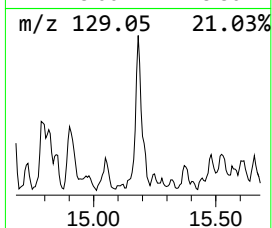
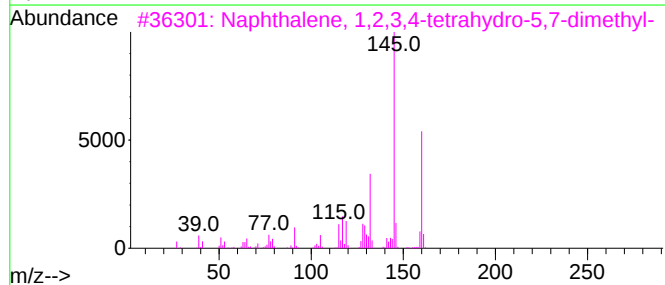
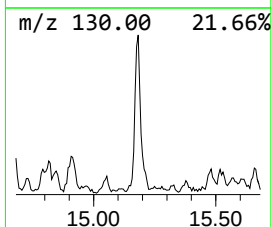
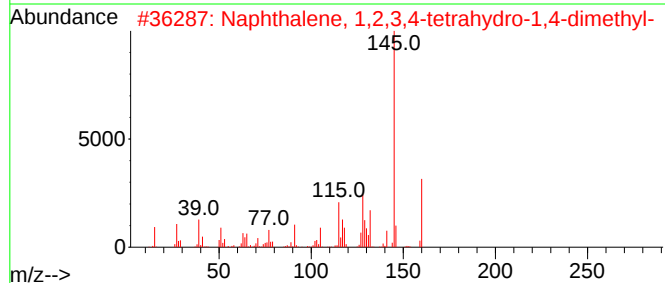
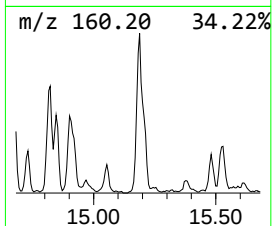
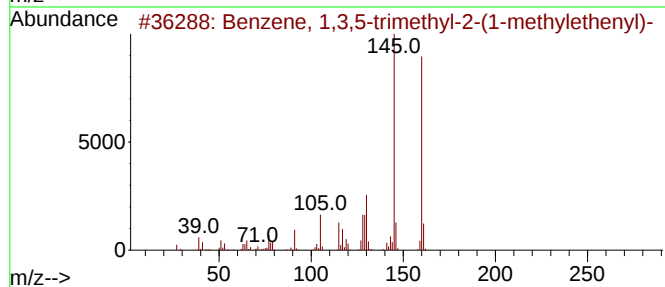
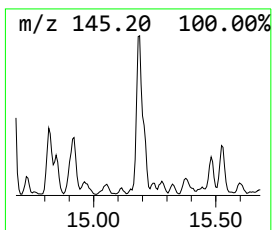
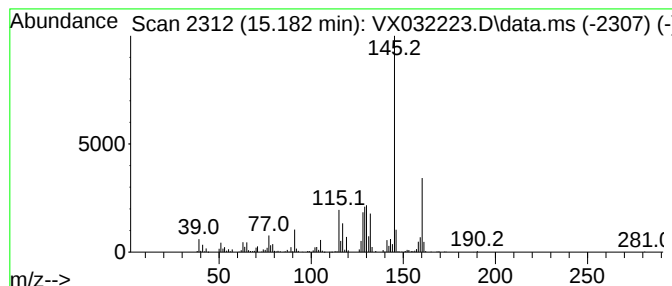
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.182	24.91 ug/l	258273	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	91
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	83
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	83
5			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
Data File : VX032223.D
Acq On : 19 Oct 2022 21:14
Operator : JC/MD
Sample : N5170-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TRE-1661

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.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...	13.579	13.8	ug/l	143534	4	12.024	518510	50.0
Benzene, (2-met...	13.926	15.4	ug/l	159383	4	12.024	518510	50.0
1H-Indene, 2,3-...	14.213	24.2	ug/l	250880	4	12.024	518510	50.0
Benzene, 1,3,5-...	14.371	12.6	ug/l	130994	4	12.024	518510	50.0
Naphthalene, 1,...	14.481	18.4	ug/l	190561	4	12.024	518510	50.0
Naphthalene, 1,...	14.615	21.0	ug/l	217656	4	12.024	518510	50.0
Benzene, 1-(1-m...	14.670	13.5	ug/l	139546	4	12.024	518510	50.0
Ethanone, 1-[4-...	14.816	11.9	ug/l	123605	4	12.024	518510	50.0
Naphthalene, 6-...	14.902	12.5	ug/l	129271	4	12.024	518510	50.0
Naphthalene, 1,...	15.182	24.9	ug/l	258273	4	12.024	518510	50.0