LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060324\
Data File : VX041707.D
Acq On : 03 Jun 2024 16:17
Operator : JC/MD
Sample : P2660-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

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\82X052924W.M
Title : SW846 8260

Signal : TIC: VX041707.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.240	22	25	32	rBV2	17233	28277	4.19%	0.382%
2	1.605	69	85	89	rBV2	15642	85788	12.70%	1.160%
3	1.715	99	103	111	rVB	21299	31797	4.71%	0.430%
4	2.325	197	203	210	rVB	16069	29968	4.44%	0.405%
5	2.770	267	276	286	rBV	34602	77931	11.54%	1.054%
6	3.087	320	328	336	rBV	4568	10398	1.54%	0.141%
7	4.203	502	511	519	rBV2	5571	15837	2.34%	0.214%
8	5.123	651	662	669	rBV4	1946	6775	1.00%	0.092%
9	5.385	694	705	722	rBV2	80632	242806	35.94%	3.283%
10	5.550	722	732	756	rVB	136891	432445	64.01%	5.847%
11	5.836	772	779	790	rVB6	2353	6901	1.02%	0.093%
12	5.958	790	799	815	rBV	83101	225125	33.32%	3.044%
13	6.245	826	846	855	rVV2	7714	28151	4.17%	0.381%
14	6.355	855	864	873	rVB4	2562	7923	1.17%	0.107%
15	6.763	922	931	954	rBV	204204	497724	73.67%	6.729%
16	7.354	1019	1028	1037	rVB6	4985	12609	1.87%	0.170%
17	7.982	1123	1131	1140	rVB	9161	18037	2.67%	0.244%
18	8.104	1142	1151	1161	rBV3	18596	39876	5.90%	0.539%
19	8.647	1234	1240	1255	rVB	431307	675580	100.00%	9.134%
20	8.964	1287	1292	1299	rVB3	10647	21355	3.16%	0.289%
21	9.049	1299	1306	1311	rBV2	10023	17681	2.62%	0.239%
22	9.098	1311	1314	1319	rVB2	4543	7159	1.06%	0.097%
23	9.159	1319	1324	1328	rBV2	5736	8715	1.29%	0.118%
24	10.055	1465	1471	1481	rBV	457326	624922	92.50%	8.449%
25	10.586	1551	1558	1565	rBV7	4737	11326	1.68%	0.153%
26	10.781	1581	1590	1593	rBV6	3828	7730	1.14%	0.105%
27	10.964	1616	1620	1625	rBV	8715	12449	1.84%	0.168%
28	11.079	1634	1639	1645	rBV	389814	514594	76.17%	6.957%
29	11.305	1672	1676	1684	rVB2	7841	13603	2.01%	0.184%
30	11.616	1722	1727	1734	rVB5	5990	10686	1.58%	0.144%
31	11.713	1734	1743	1745	rBV6	4815	8845	1.31%	0.120%
32	11.750	1745	1749	1755	rVB4	4438	8908	1.32%	0.120%
33	11.866	1762	1768	1770	rBV	16052	21694	3.21%	0.293%
34	11.890	1770	1772	1779	rVB	9539	11853	1.75%	0.160%
35	12.024	1788	1794	1802	rBV	468190	609930	90.28%	8.246%
36	12.177	1812	1819	1824	rVB	22325	31779	4.70%	0.430%

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MSVOA_X
ClientSampleId :
MW-10

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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Title : SW846 8260

37	12.250	1826	1831	1835	rVV	52966	68810	10.19%	0.930%
38	12.311	1837	1841	1852	rVV	30231	42688	6.32%	0.577%
39	12.402	1852	1856	1863	rVV3	67371	93642	13.86%	1.266%
40	12.469	1863	1867	1873	rVB	57552	76699	11.35%	1.037%
41	12.597	1882	1888	1893	rBV5	3606	8058	1.19%	0.109%
42	12.658	1893	1898	1900	rBV5	8868	12061	1.79%	0.163%
43	12.701	1900	1905	1912	rVV	324970	493310	73.02%	6.670%
44	12.756	1912	1914	1918	rVV	17256	21258	3.15%	0.287%
45	12.799	1918	1921	1924	rVV2	14140	21576	3.19%	0.292%
46	12.841	1924	1928	1933	rVB	43419	58049	8.59%	0.785%
47	12.896	1933	1937	1939	rBV	29138	37200	5.51%	0.503%
48	12.920	1939	1941	1947	rVV	35552	43885	6.50%	0.593%
49	13.030	1954	1959	1965	rBV2	64883	94521	13.99%	1.278%
50	13.140	1968	1977	1981	rBV3	101715	164442	24.34%	2.223%
51	13.195	1981	1986	1990	rVV	57378	78248	11.58%	1.058%
52	13.244	1990	1994	2000	rVB4	25821	47530	7.04%	0.643%
53	13.298	2000	2003	2007	rVB2	9278	10953	1.62%	0.148%
54	13.347	2007	2011	2015	rBV	38096	49105	7.27%	0.664%
55	13.390	2015	2018	2021	rBV	27067	34795	5.15%	0.470%
56	13.426	2021	2024	2029	rVB2	39900	57297	8.48%	0.775%
57	13.512	2034	2038	2041	rBV	20210	30866	4.57%	0.417%
58	13.548	2041	2044	2047	rVB	13899	14612	2.16%	0.198%
59	13.579	2047	2049	2051	rBV2	9274	7611	1.13%	0.103%
60	13.640	2051	2059	2063	rBV2	173698	268124	39.69%	3.625%
61	13.750	2073	2077	2081	rBV	32996	41478	6.14%	0.561%
62	13.798	2081	2085	2090	rVB3	28354	53159	7.87%	0.719%
63	13.853	2090	2094	2098	rBV2	33460	50476	7.47%	0.682%
64	13.926	2103	2106	2110	rVB2	41041	57271	8.48%	0.774%
65	13.975	2110	2114	2118	rVB	34350	49509	7.33%	0.669%
66	14.061	2126	2128	2133	rVB	9643	12112	1.79%	0.164%
67	14.140	2137	2141	2142	rBV2	12608	17028	2.52%	0.230%
68	14.158	2142	2144	2148	rVB2	10515	11323	1.68%	0.153%
69	14.225	2148	2155	2160	rBV5	32615	66880	9.90%	0.904%
70	14.317	2166	2170	2175	rBV	117278	154983	22.94%	2.095%
71	14.371	2175	2179	2184	rVV	84682	112568	16.66%	1.522%
72	14.420	2184	2187	2193	rVB5	6644	9763	1.45%	0.132%
73	14.487	2193	2198	2201	rBV3	25141	41347	6.12%	0.559%
74	14.524	2201	2204	2210	rVV2	14459	23033	3.41%	0.311%
75	14.609	2210	2218	2225	rVV4	40449	96105	14.23%	1.299%
76	14.676	2225	2229	2233	rVB	29064	34809	5.15%	0.471%

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Integrator: RTE
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77	14.756	2239	2242	2245	rVB3	9297	9007	1.33%	0.122%
78	14.792	2245	2248	2250	rBV2	15980	20719	3.07%	0.280%
79	14.823	2250	2253	2255	rVV2	14292	18885	2.80%	0.255%
80	14.847	2255	2257	2261	rVB3	11087	14310	2.12%	0.193%
81	14.914	2264	2268	2274	rBV3	25789	37952	5.62%	0.513%
82	15.182	2308	2312	2319	rBV	24976	43629	6.46%	0.590%
83	15.249	2319	2323	2330	rVB6	16857	26794	3.97%	0.362%
84	15.371	2339	2343	2346	rBV5	6869	12038	1.78%	0.163%
85	15.408	2346	2349	2358	rVB	17859	26591	3.94%	0.360%
86	15.487	2358	2362	2365	rBV	14699	22497	3.33%	0.304%
87	15.664	2388	2391	2394	rBV5	8178	10415	1.54%	0.141%
88	15.810	2411	2415	2418	rBV	11798	16127	2.39%	0.218%
89	15.981	2439	2443	2448	rBV2	9268	16743	2.48%	0.226%
90	16.371	2503	2507	2513	rBV	17142	29124	4.31%	0.394%
91	16.652	2550	2553	2558	rVB3	4377	7128	1.06%	0.096%

Sum of corrected areas: 7396320

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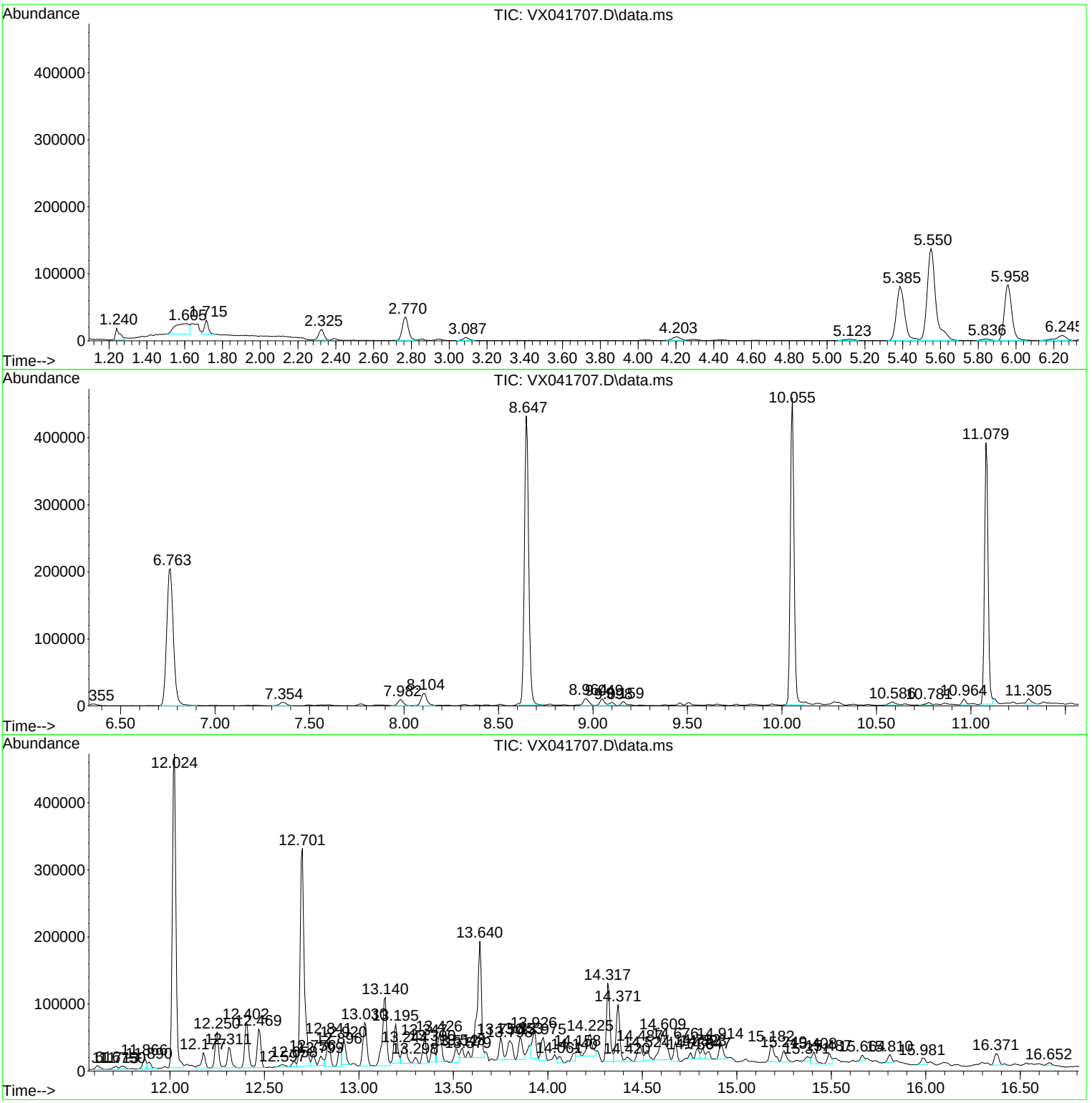
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060324\
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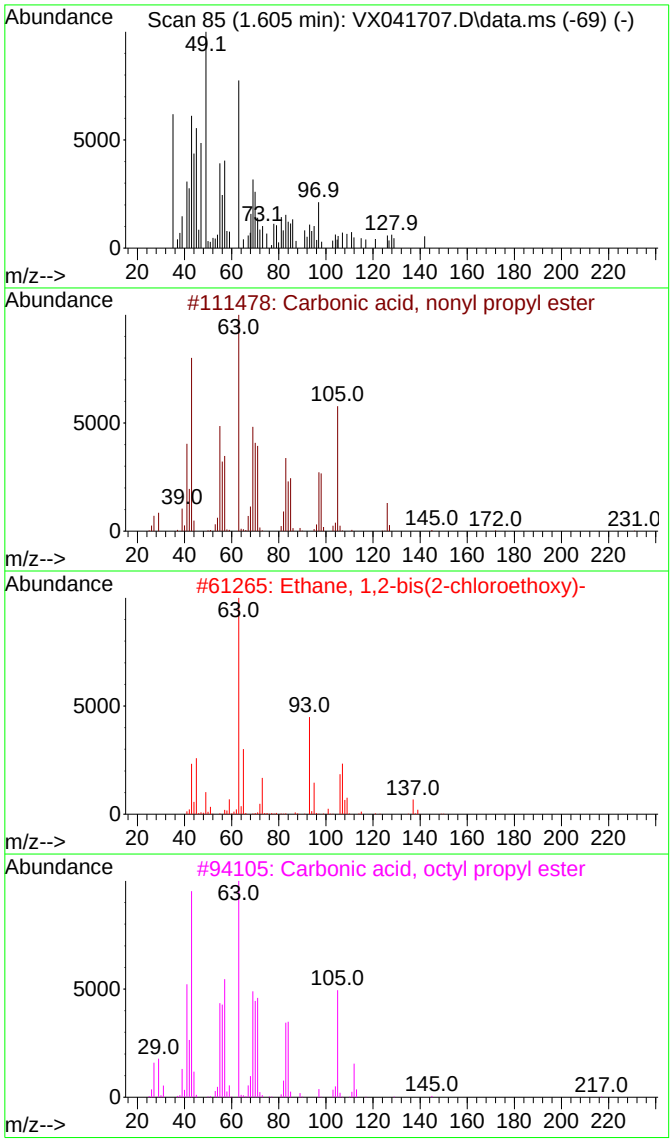
Quant Title : SW846 8260

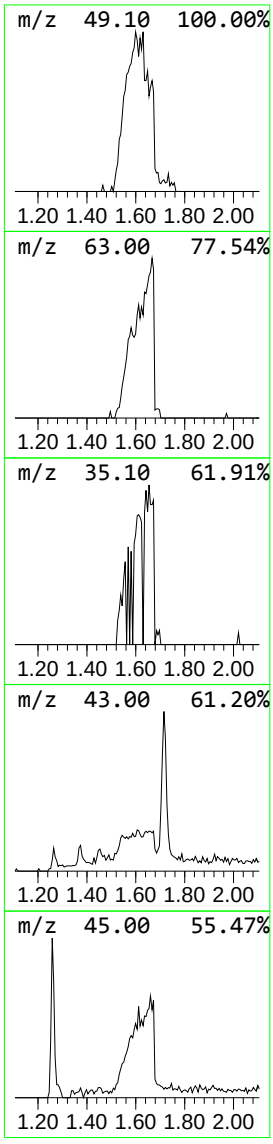
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 unknown1.605 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.605	9.92 ug/l	85788	Pentafluorobenzene	5.550

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, nonyl propyl ester	230	C13H26O3	1000314-53-6	22
2	Ethane, 1,2-bis(2-chloroethoxy)-	186	C6H12Cl2O2	000112-26-5	17
3	Carbonic acid, octyl propyl ester	216	C12H24O3	1000314-53-5	10
4	3,6,9,12-tetraoxatetradecane, 1,...	274	C10H20Cl2O4	1000401-80-9	10
5	Heptyl propyl carbonate	202	C11H22O3	959272-47-0	10





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ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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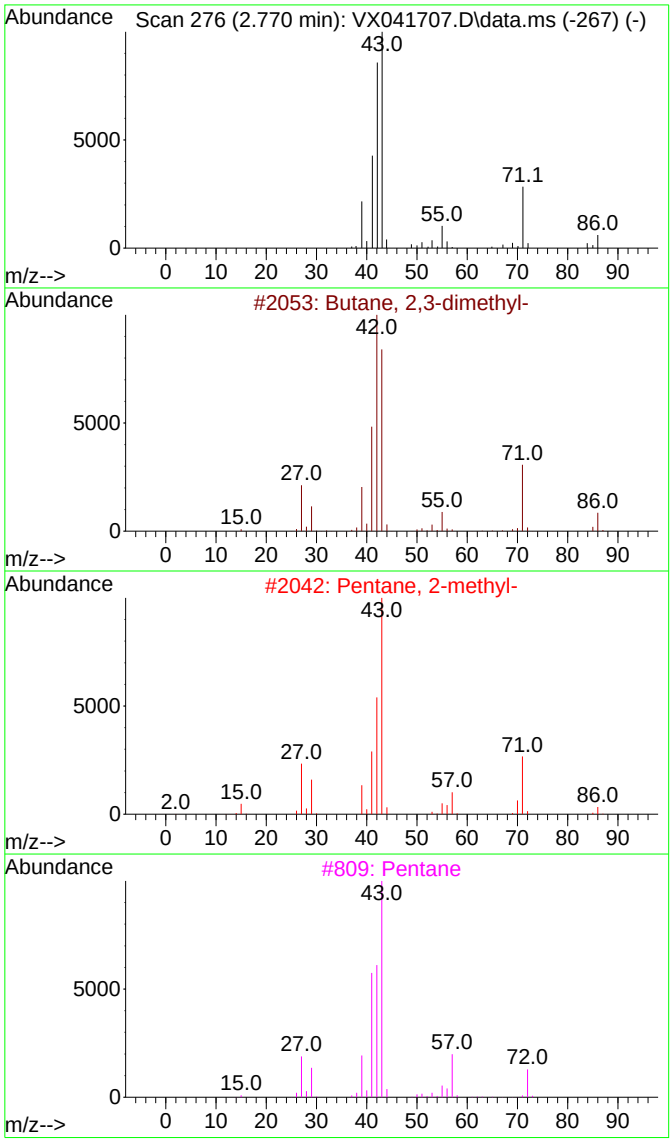
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

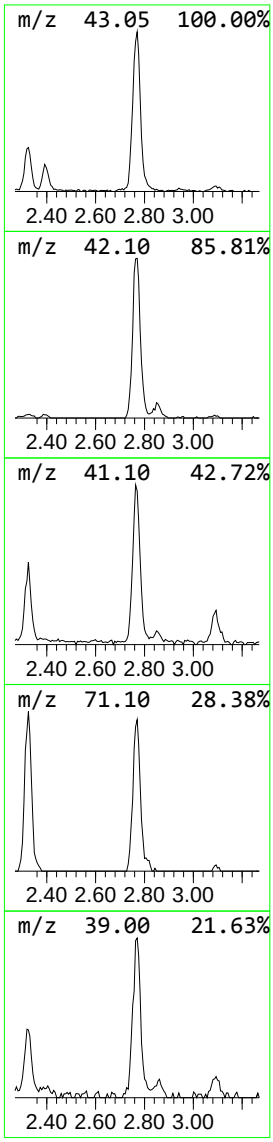
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.770	9.01 ug/l	77931	Pentafluorobenzene	5.550

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3	Pentane	72	C5H12	000109-66-0	38
4	C2H5C(CH3)2COCH3	114	C7H14O	020669-04-9	12
5	Butanal, 2,2-dimethyl-	100	C6H12O	002094-75-9	9





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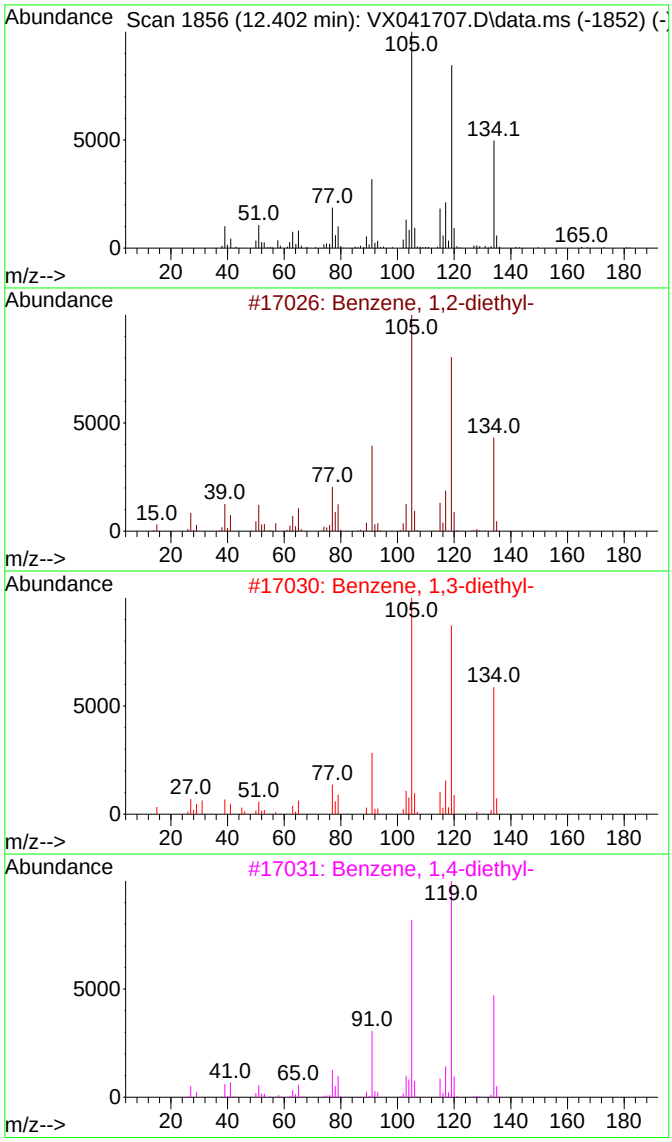
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

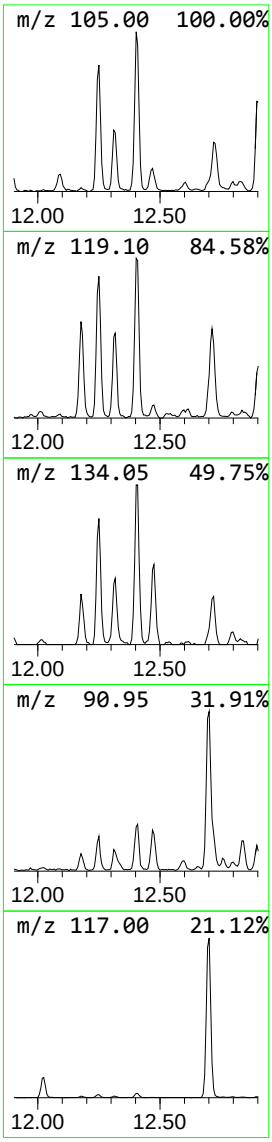
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.402	7.68 ug/l	93642	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	90
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
4	2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	87
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	80





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MSVOA_X
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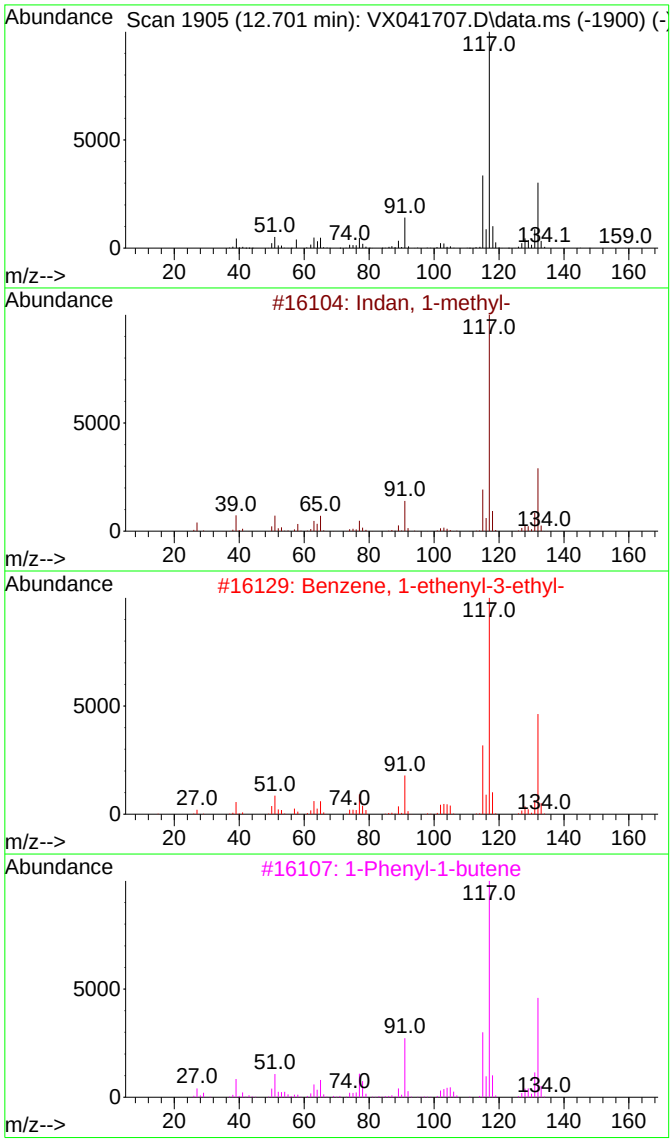
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

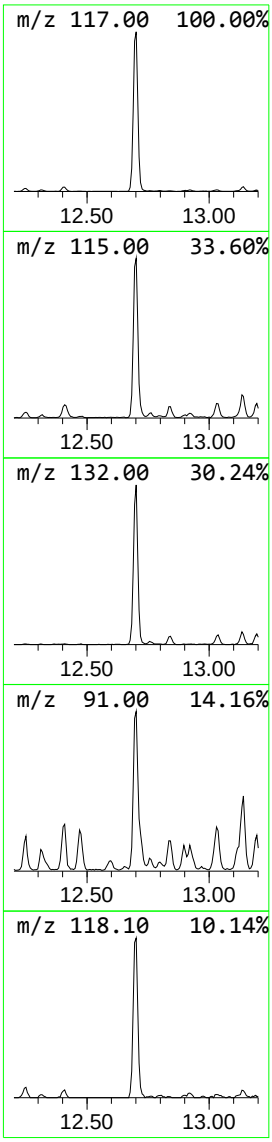
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Indan, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	40.44 ug/l	493310	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	91
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86





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ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

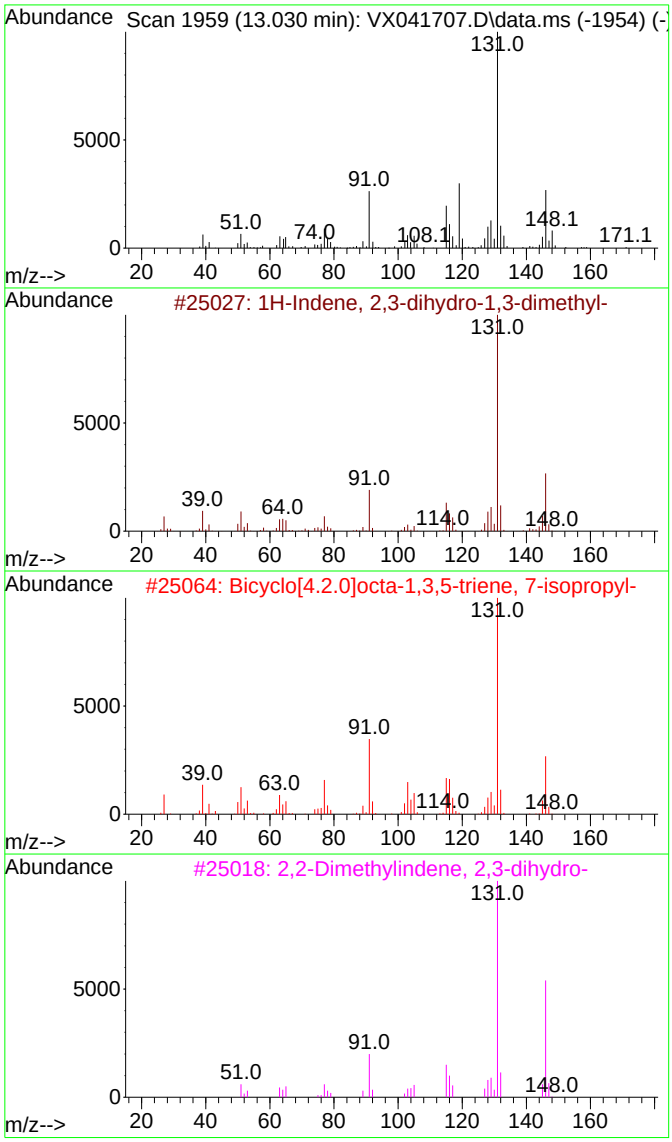
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

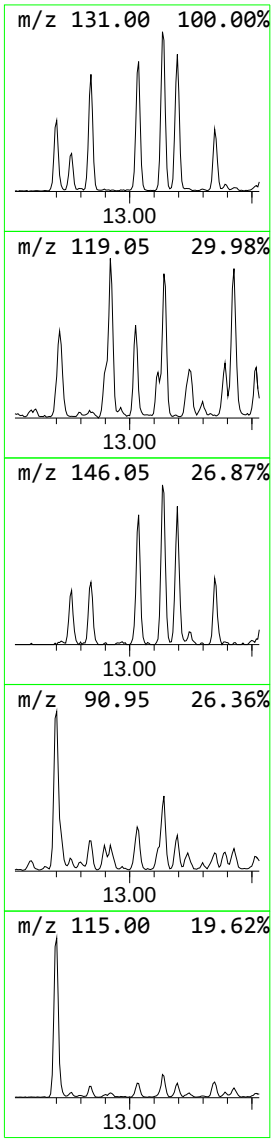
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.030	7.75 ug/l	94521	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	95
2	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	93
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	90





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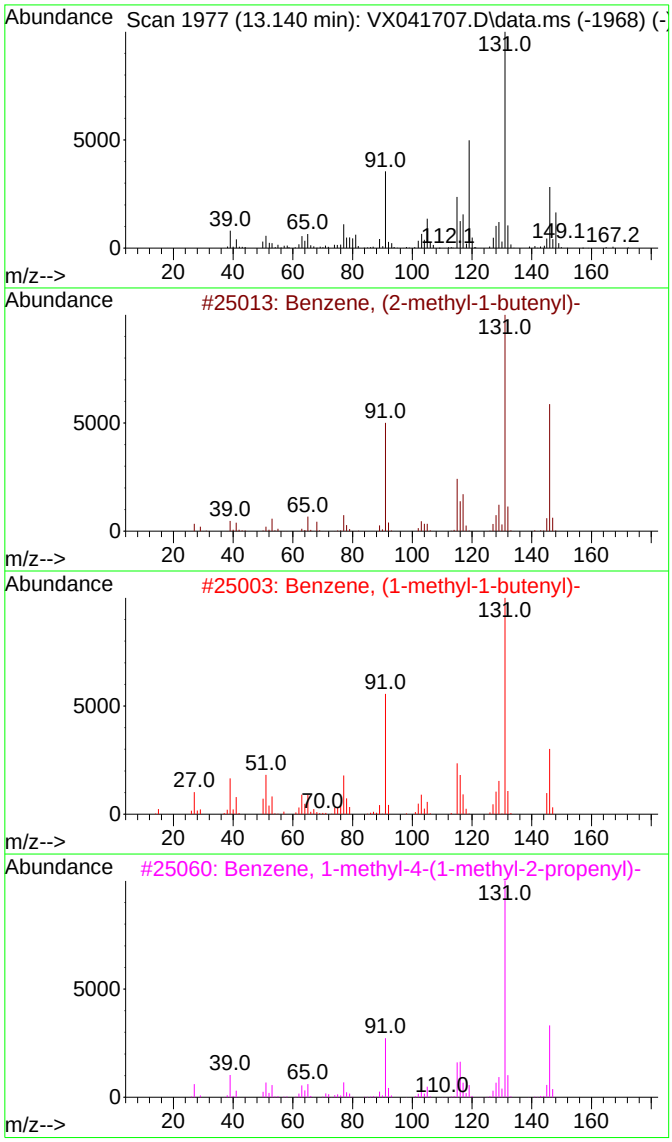
Quant Title : SW846 8260

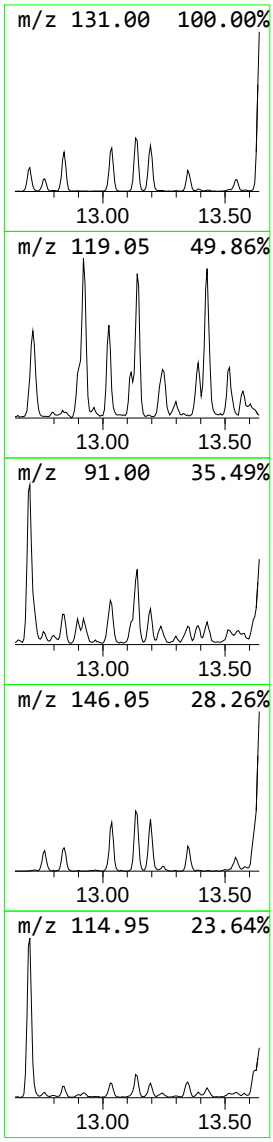
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (2-methyl-1-butenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.140	13.48 ug/l	164442	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	96
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
3	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70





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Data File : VX041707.D
Acq On : 03 Jun 2024 16:17
Operator : JC/MD
Sample : P2660-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

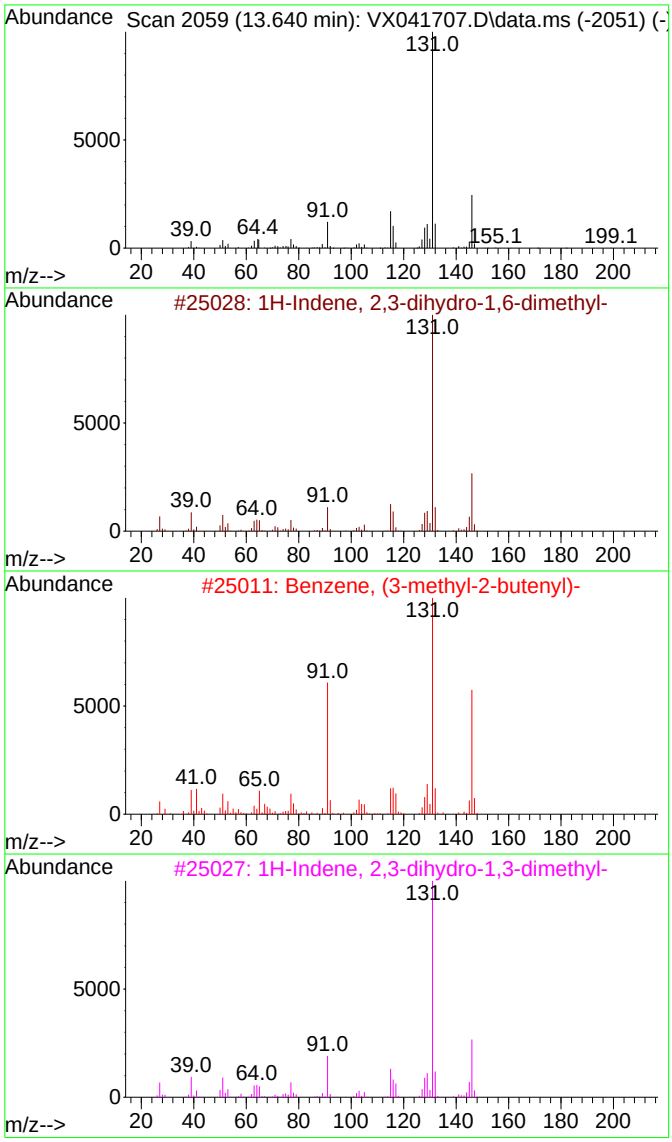
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

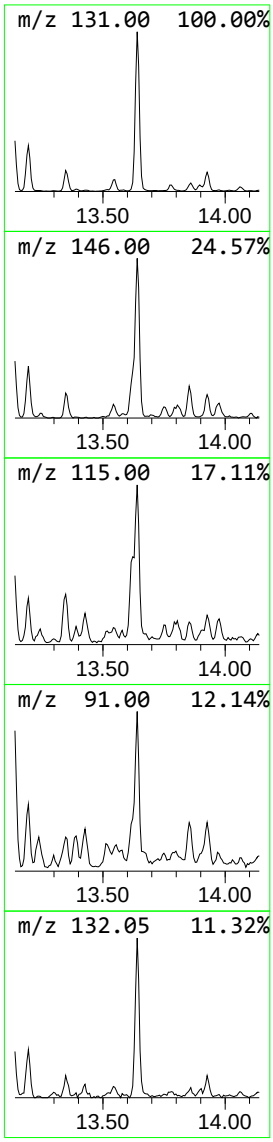
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.640	21.98 ug/l	268124	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060324\
Data File : VX041707.D
Acq On : 03 Jun 2024 16:17
Operator : JC/MD
Sample : P2660-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

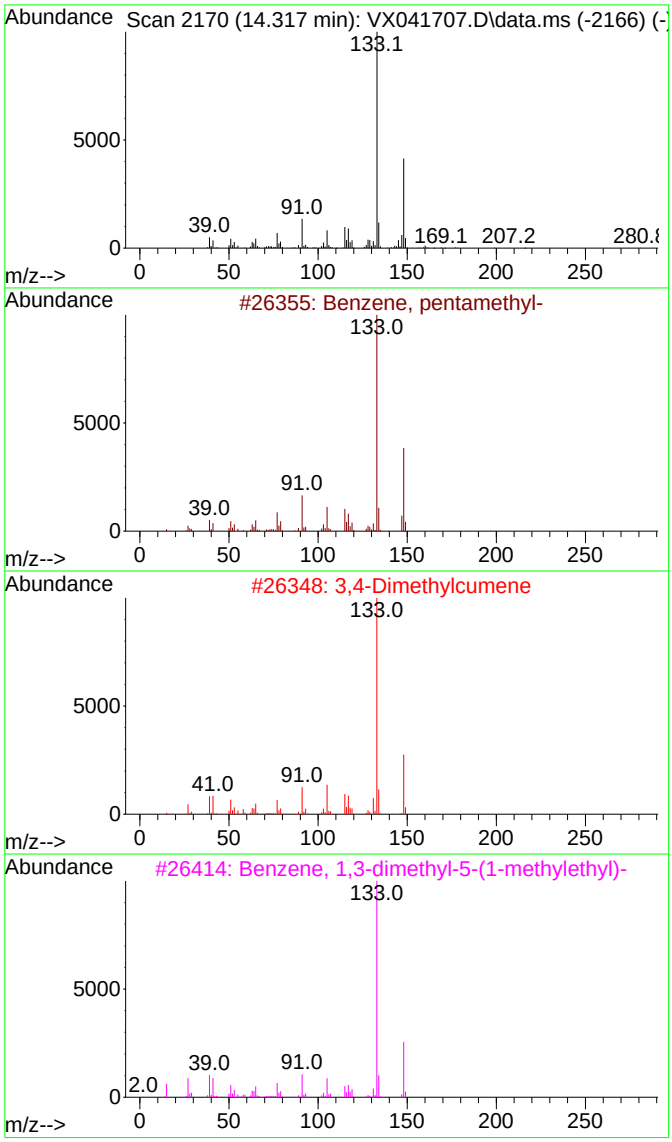
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

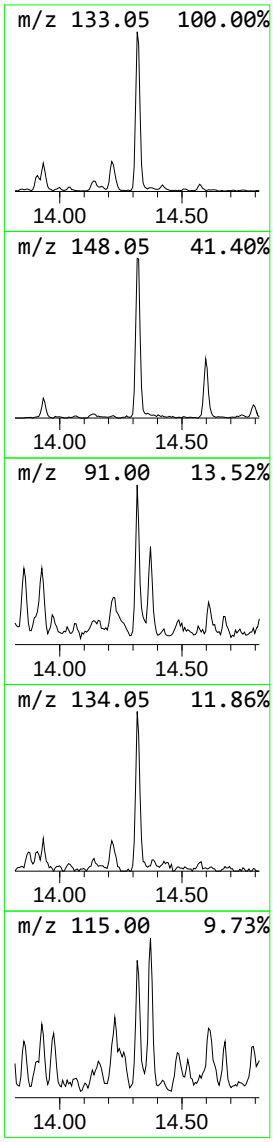
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, pentamethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.317	12.71 ug/l	154983	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	95
2	3,4-Dimethylcumene	148	C11H16	1000370-34-1	87
3	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	87
4	1,5,6,7-Tetramethylbicyclo[3.2.0...	148	C11H16	134329-46-7	83
5	Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060324\
Data File : VX041707.D
Acq On : 03 Jun 2024 16:17
Operator : JC/MD
Sample : P2660-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

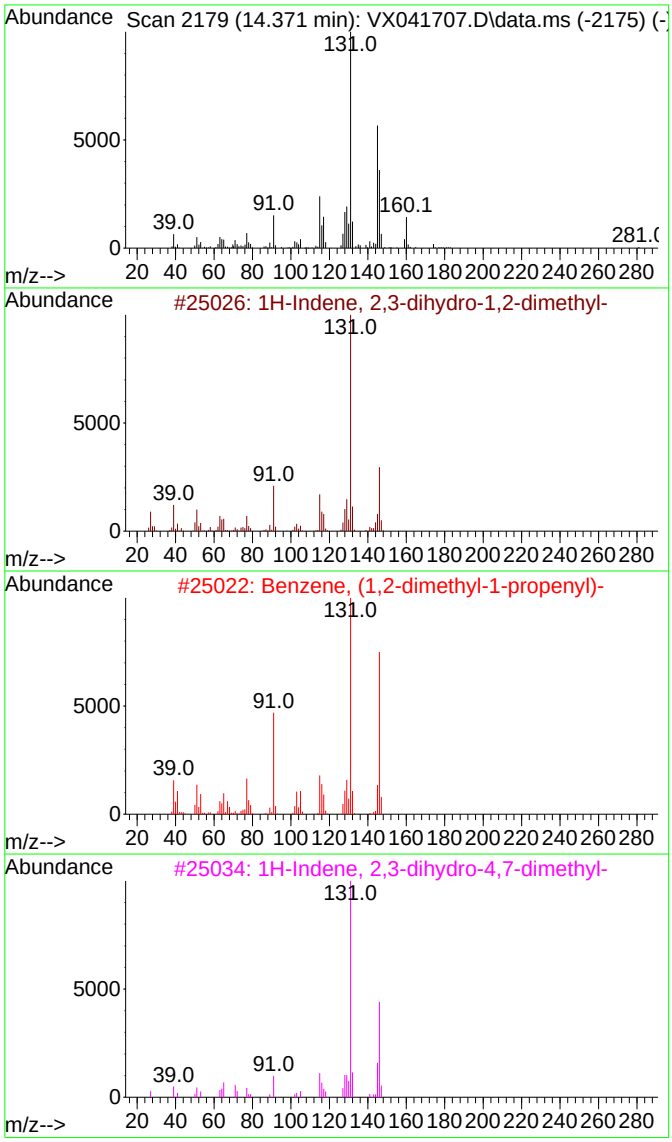
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

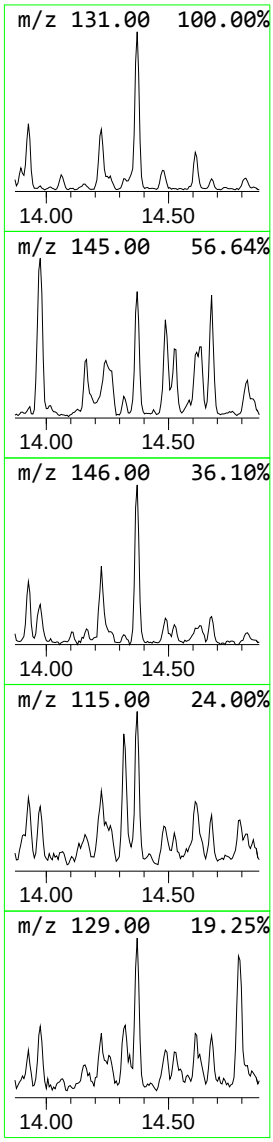
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.371	9.23 ug/l	112568	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	81
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	81
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	62
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	62





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060324\
Data File : VX041707.D
Acq On : 03 Jun 2024 16:17
Operator : JC/MD
Sample : P2660-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

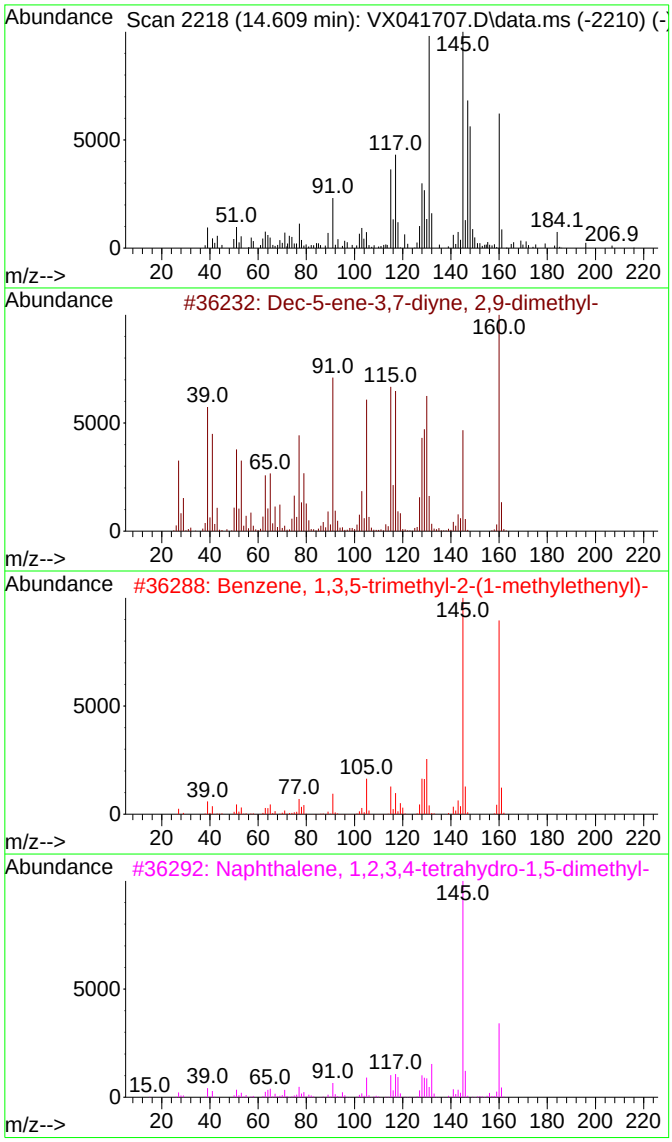
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

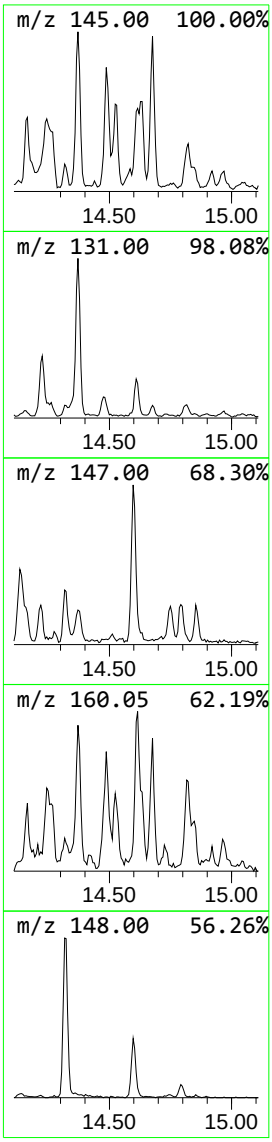
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Dec-5-ene-3,7-diyne, 2,9-di... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.609	7.88 ug/l	96105	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dec-5-ene-3,7-diyne, 2,9-dimethyl-	160	C12H16	1010211-23-3	51
2	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	50
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	50
4	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	50
5	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	49





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060324\
Data File : VX041707.D
Acq On : 03 Jun 2024 16:17
Operator : JC/MD
Sample : P2660-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X052924W.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown1.605	1.605	9.9	ug/l	85788	1	5.550	432445	50.0
Butane, 2,3-dim...	2.770	9.0	ug/l	77931	1	5.550	432445	50.0
Benzene, 1,2-di...	12.402	7.7	ug/l	93642	4	12.024	609930	50.0
Indan, 1-methyl-	12.701	40.4	ug/l	493310	4	12.024	609930	50.0
1H-Indene, 2,3-...	13.030	7.8	ug/l	94521	4	12.024	609930	50.0
Benzene, (2-met...	13.140	13.5	ug/l	164442	4	12.024	609930	50.0
1H-Indene, 2,3-...	13.640	22.0	ug/l	268124	4	12.024	609930	50.0
Benzene, pentam...	14.317	12.7	ug/l	154983	4	12.024	609930	50.0
1H-Indene, 2,3-...	14.371	9.2	ug/l	112568	4	12.024	609930	50.0
Dec-5-ene-3,7-d...	14.609	7.9	ug/l	96105	4	12.024	609930	50.0