LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

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\82X051222W.M
Title : SW846 8260

Signal : TIC: VX028714.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.075	154	162	172	rBV	68996	136068	6.22%	0.312%
2	2.380	203	212	222	rBV	147425	245538	11.22%	0.563%
3	4.562	561	570	582	rBV	20127	55659	2.54%	0.128%
4	5.391	694	706	721	rBV2	163246	482139	22.03%	1.106%
5	5.556	721	733	749	rVB	254762	722485	33.01%	1.657%
6	5.958	788	799	807	rBV	192080	482581	22.05%	1.107%
7	6.044	807	813	823	rVB	39572	105189	4.81%	0.241%
8	6.763	921	931	945	rBV	437163	1029433	47.04%	2.362%
9	7.458	1039	1045	1055	rBV	19663	42063	1.92%	0.096%
10	8.653	1234	1241	1246	rBV	900216	1414228	64.62%	3.244%
11	8.720	1246	1252	1263	rVB	970039	1455030	66.48%	3.338%
12	9.378	1355	1360	1365	rBV	19456	28106	1.28%	0.064%
13	10.055	1465	1471	1482	rBV	979452	1279569	58.47%	2.935%
14	10.195	1488	1494	1500	rBV	300196	384998	17.59%	0.883%
15	10.299	1505	1511	1518	rVV	1139120	1489059	68.04%	3.416%
16	10.640	1562	1567	1572	rBV	788149	997220	45.57%	2.288%
17	10.775	1582	1589	1595	rBV	31685	48961	2.24%	0.112%
18	10.860	1598	1603	1614	rVB7	11539	26669	1.22%	0.061%
19	10.964	1614	1620	1625	rBV	54281	69069	3.16%	0.158%
20	11.079	1634	1639	1651	rVB	796843	1027164	46.93%	2.356%
21	11.305	1669	1676	1682	rBV	190179	231697	10.59%	0.532%
22	11.378	1683	1688	1696	rVV	809528	1356520	61.98%	3.112%
23	11.451	1696	1700	1711	rVB	403570	543347	24.83%	1.247%
24	11.616	1721	1727	1732	rBV	436913	570581	26.07%	1.309%
25	11.750	1744	1749	1760	rVB	1788523	2176642	99.46%	4.993%
26	11.841	1760	1764	1766	rBV	19107	27586	1.26%	0.063%
27	11.890	1769	1772	1776	rVB	54225	66885	3.06%	0.153%
28	11.969	1781	1785	1789	rVV	93727	113535	5.19%	0.260%
29	12.024	1789	1794	1799	rVB	1188992	1552171	70.92%	3.561%
30	12.085	1799	1804	1810	rBV	839187	1038880	47.47%	2.383%
31	12.244	1822	1830	1833	rBV2	814086	1225250	55.98%	2.811%
32	12.317	1838	1842	1851	rVB3	535890	826683	37.77%	1.897%
33	12.408	1852	1857	1862	rBV3	68723	144525	6.60%	0.332%
34	12.463	1862	1866	1872	rVB	241244	285906	13.06%	0.656%
35	12.530	1872	1877	1879	rBV	375738	462221	21.12%	1.060%
36	12.555	1879	1881	1886	rVV	341208	368888	16.86%	0.846%

LSC Area Percent Report

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Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

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\82X051222W.M
Title : SW846 8260

37	12.616	1886	1891	1894	rVV	487789	620727	28.36%	1.424%
38	12.646	1894	1896	1900	rVB	160842	161453	7.38%	0.370%
39	12.701	1900	1905	1913	rBV2	439891	747704	34.16%	1.715%
40	12.768	1913	1916	1918	rBV3	17495	23122	1.06%	0.053%
41	12.799	1918	1921	1925	rVV2	43038	46956	2.15%	0.108%
42	12.847	1925	1929	1933	rVB2	226951	280709	12.83%	0.644%
43	12.921	1933	1941	1944	rBV	391282	488194	22.31%	1.120%
44	12.963	1944	1948	1954	rVB	652406	760080	34.73%	1.744%
45	13.024	1954	1958	1959	rBV	95783	103993	4.75%	0.239%
46	13.042	1959	1961	1964	rVV3	86982	110130	5.03%	0.253%
47	13.073	1964	1966	1968	rVB	54200	47119	2.15%	0.108%
48	13.170	1974	1982	1986	rBV	724364	987110	45.10%	2.265%
49	13.213	1986	1989	1992	rVB	109252	117798	5.38%	0.270%
50	13.292	1997	2002	2007	rVB2	1650224	2188540	100.00%	5.021%
51	13.372	2012	2015	2019	rVV2	71395	89998	4.11%	0.206%
52	13.420	2019	2023	2032	rVB	1191654	1532558	70.03%	3.516%
53	13.506	2032	2037	2040	rBV2	144460	203149	9.28%	0.466%
54	13.542	2040	2043	2046	rVV	250330	318448	14.55%	0.731%
55	13.585	2046	2050	2055	rVV3	314800	582936	26.64%	1.337%
56	13.640	2056	2059	2060	rVV	144982	174793	7.99%	0.401%
57	13.664	2060	2063	2069	rVB2	317093	473215	21.62%	1.086%
58	13.774	2077	2081	2088	rVB	472048	650724	29.73%	1.493%
59	13.853	2088	2094	2099	rVB	419684	561324	25.65%	1.288%
60	13.926	2099	2106	2110	rBV2	454448	652132	29.80%	1.496%
61	13.969	2110	2113	2117	rVB	156528	166574	7.61%	0.382%
62	14.036	2121	2124	2127	rVB3	123348	131935	6.03%	0.303%
63	14.073	2127	2130	2134	rBV	365757	429406	19.62%	0.985%
64	14.164	2142	2145	2147	rBV	26599	30344	1.39%	0.070%
65	14.231	2147	2156	2163	rVV	1115147	1808100	82.62%	4.148%
66	14.323	2167	2171	2175	rVV	156672	187911	8.59%	0.431%
67	14.371	2175	2179	2183	rVV	358409	434075	19.83%	0.996%
68	14.414	2183	2186	2192	rVB3	63655	103108	4.71%	0.237%
69	14.481	2192	2197	2207	rBV2	719315	984329	44.98%	2.258%
70	14.567	2207	2211	2215	rVB6	17835	29012	1.33%	0.067%
71	14.634	2215	2222	2226	rBV2	805270	1375766	62.86%	3.156%
72	14.670	2226	2228	2233	rVB	170456	212931	9.73%	0.488%
73	14.731	2233	2238	2239	rBV	64355	80702	3.69%	0.185%
74	14.780	2239	2246	2250	rVV2	460565	710859	32.48%	1.631%
75	14.817	2250	2252	2255	rVV	109842	151328	6.91%	0.347%
76	14.847	2255	2257	2261	rVV2	103849	105311	4.81%	0.242%

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

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

77	14.902	2262	2266	2273	rVB3	101093	174626	7.98%	0.401%
78	15.048	2284	2290	2294	rBV	51117	85334	3.90%	0.196%
79	15.109	2294	2300	2302	rBV6	15359	26288	1.20%	0.060%
80	15.182	2307	2312	2314	rBV	234959	321042	14.67%	0.737%
81	15.255	2320	2324	2331	rVB5	39137	102610	4.69%	0.235%
82	15.323	2331	2335	2339	rVB6	15907	23917	1.09%	0.055%
83	15.377	2339	2344	2347	rBV	87467	132912	6.07%	0.305%
84	15.481	2355	2361	2366	rBV2	212536	390590	17.85%	0.896%
85	15.524	2366	2368	2373	rVV3	52153	82993	3.79%	0.190%
86	15.615	2377	2383	2386	rVV	197844	343834	15.71%	0.789%
87	15.652	2386	2389	2397	rVB	116039	184647	8.44%	0.424%
88	15.841	2416	2420	2425	rVB3	40538	67232	3.07%	0.154%
89	15.987	2440	2444	2448	rBV2	30133	49718	2.27%	0.114%
90	16.085	2455	2460	2469	rBV	51309	103818	4.74%	0.238%
91	16.377	2501	2508	2513	rVB2	39500	76937	3.52%	0.177%
92	16.597	2541	2544	2550	rVB3	14300	23817	1.09%	0.055%
93	16.658	2550	2554	2559	rBV2	13570	22062	1.01%	0.051%

Sum of corrected areas: 43589525

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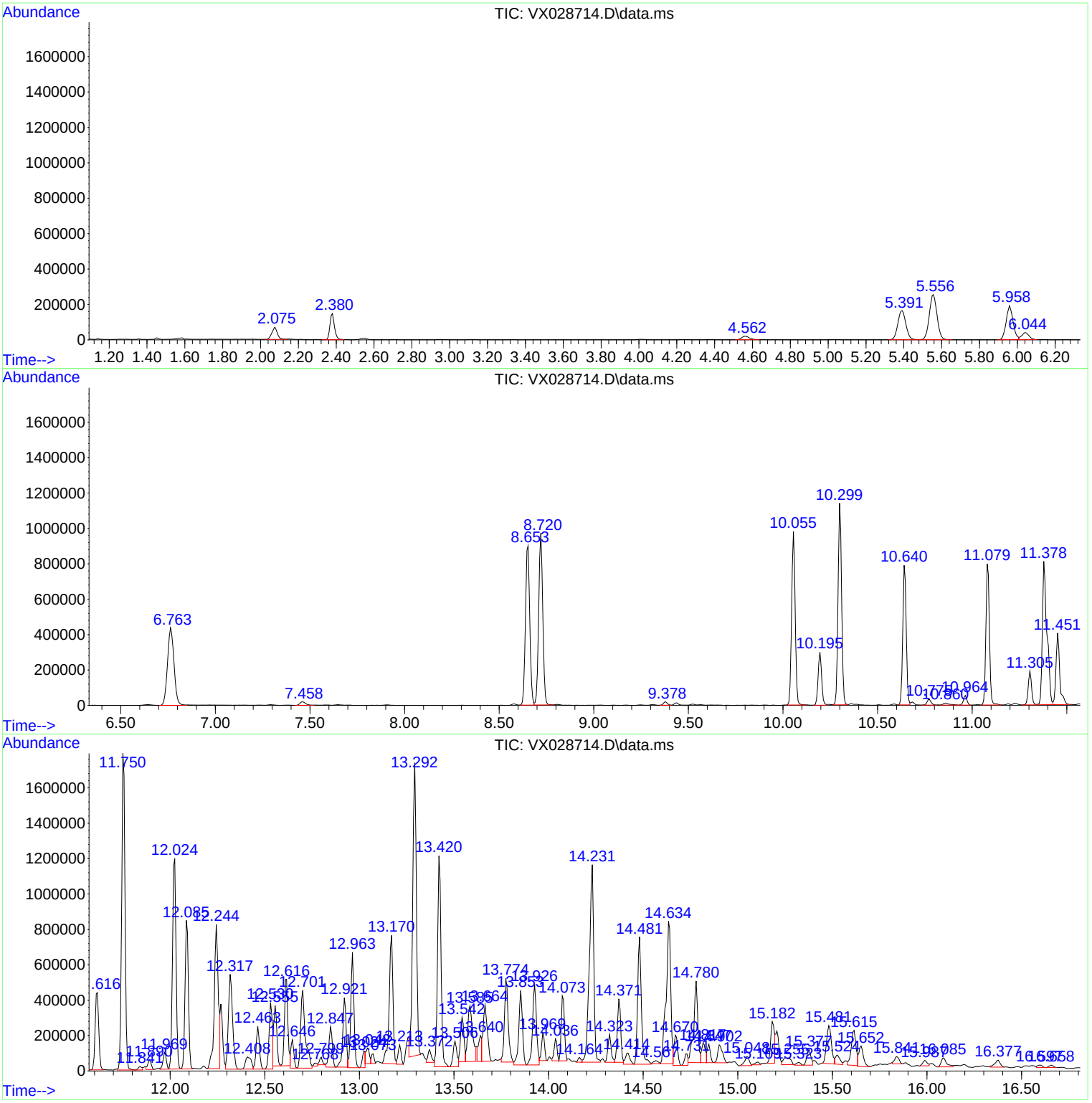
Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
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Instrument :
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ClientSampleId :
BUR-1144

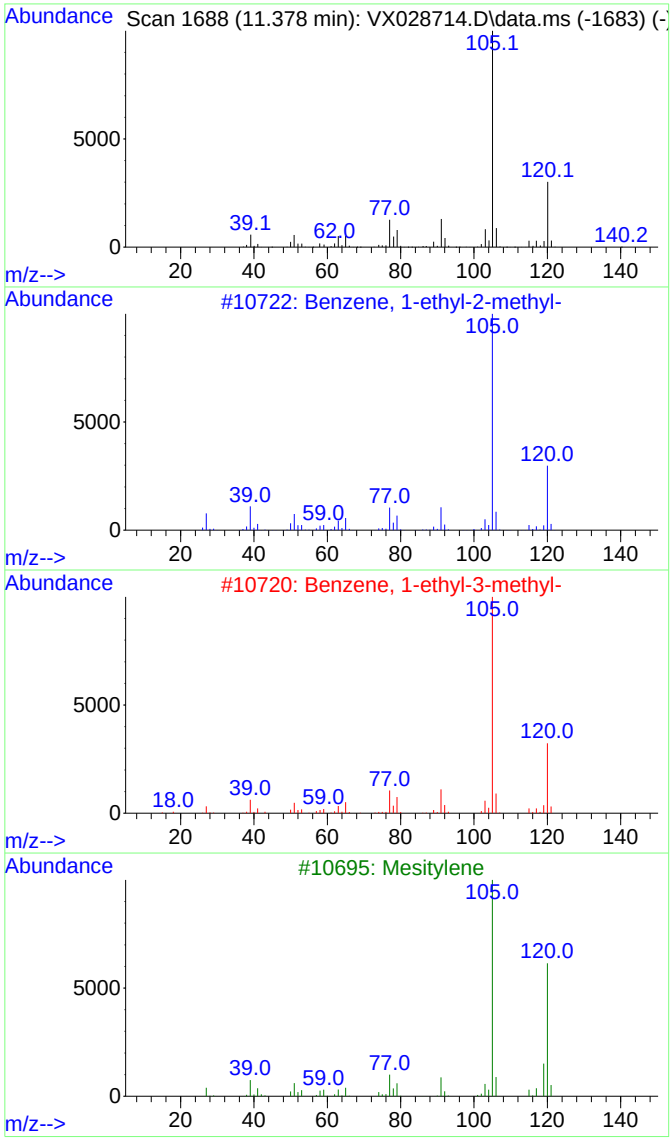
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

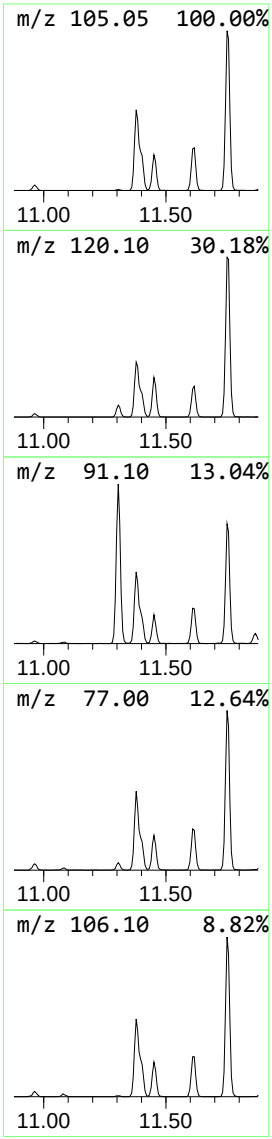
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	43.70 ug/l	1356520	1,4-Dichlorobenzene-d4	12.024

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	94
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

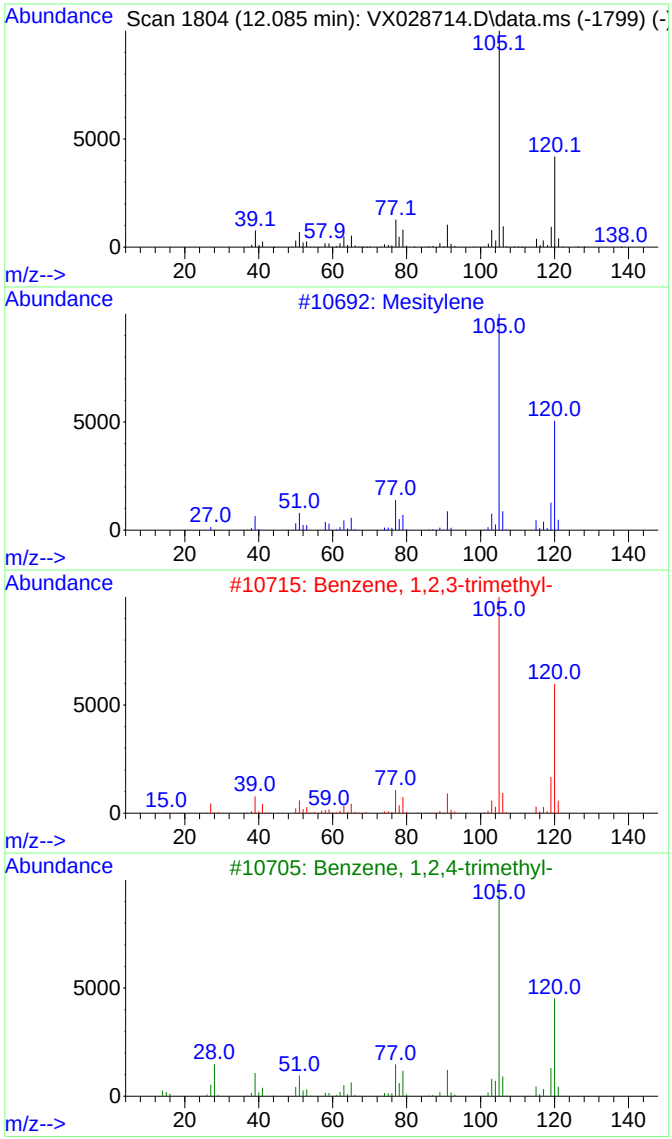
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

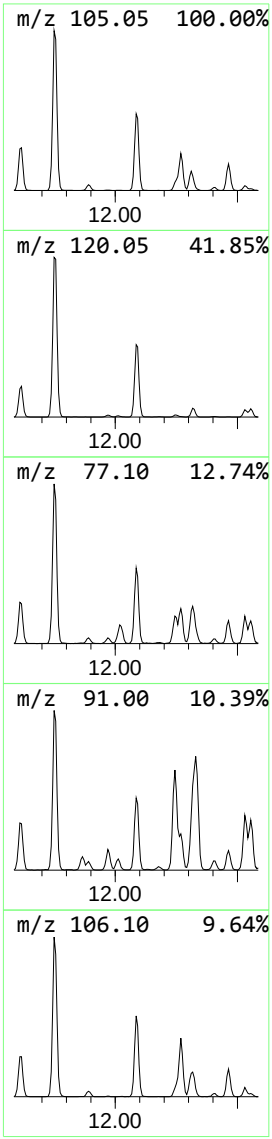
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	33.47 ug/l	1038880	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





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Instrument :
MSVOA_X
ClientSampleId :
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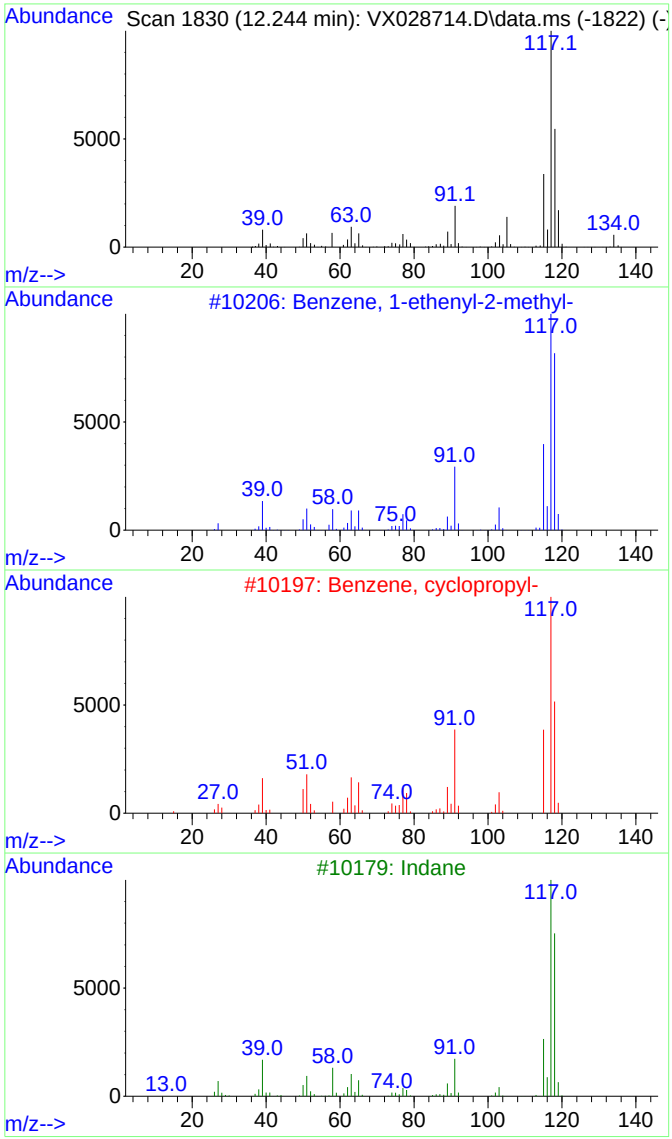
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

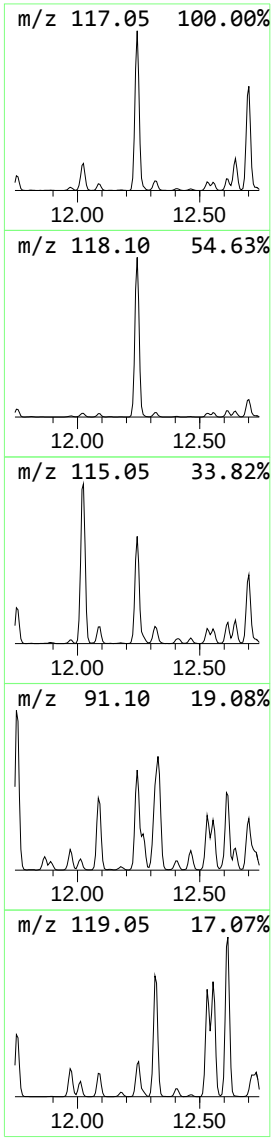
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethenyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	39.47 ug/l	1225250	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	87
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3	Indane	118	C9H10	000496-11-7	76
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	68





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
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ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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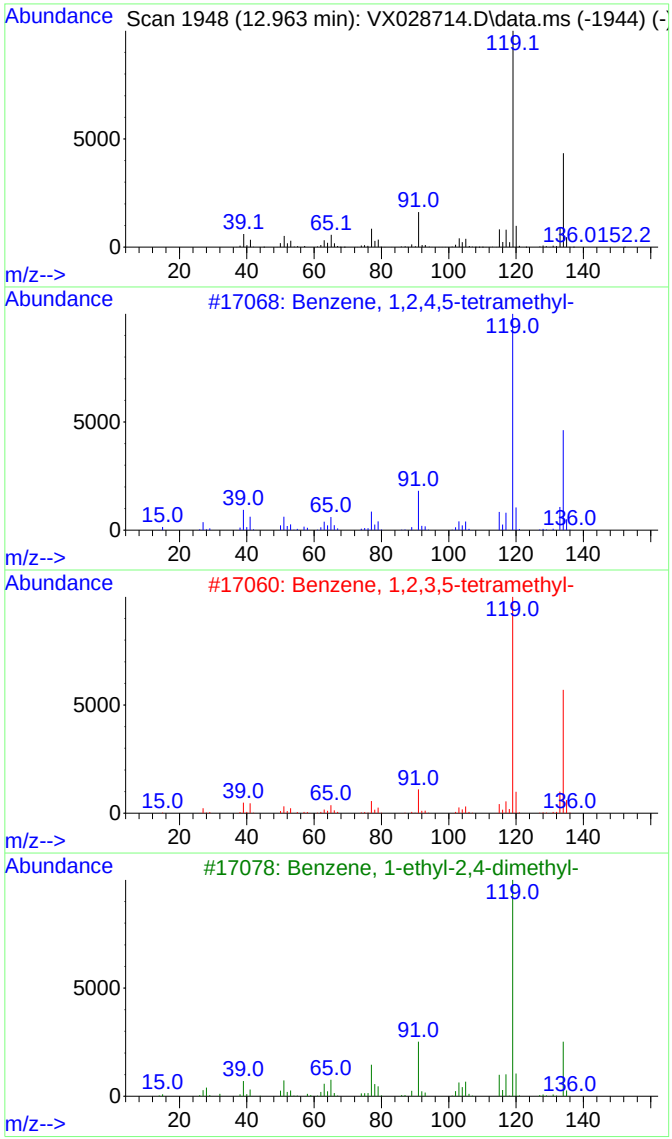
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

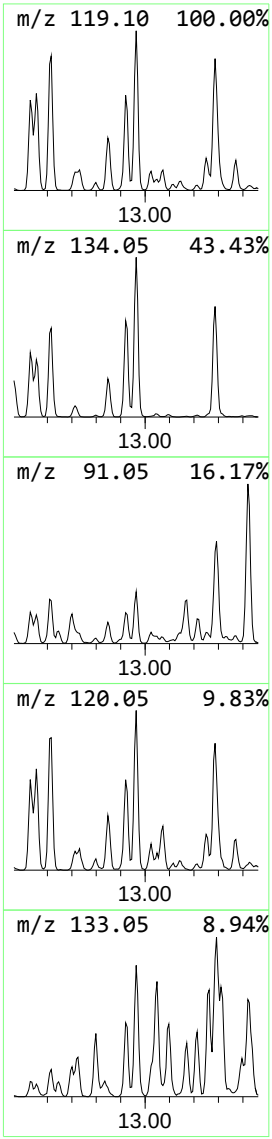
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	24.48 ug/l	760080	1,4-Dichlorobenzene-d4	12.024

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	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
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	94





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Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

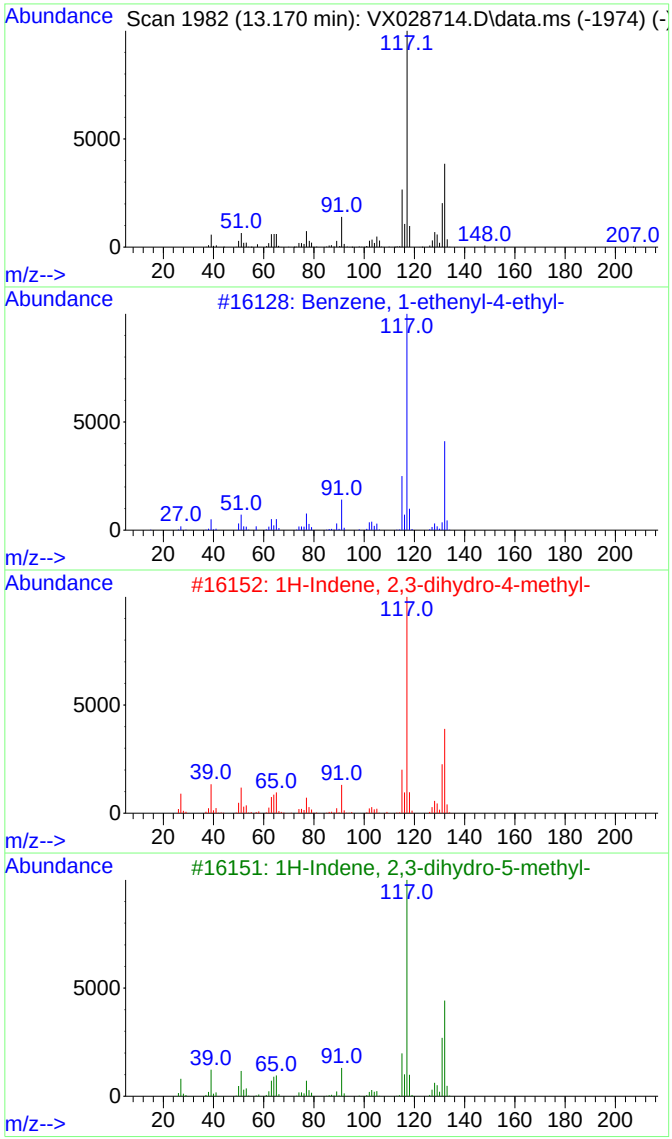
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

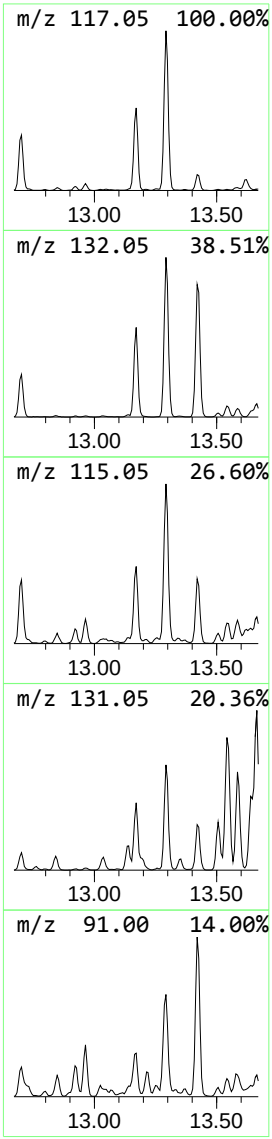
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.171	31.80 ug/l	987110	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

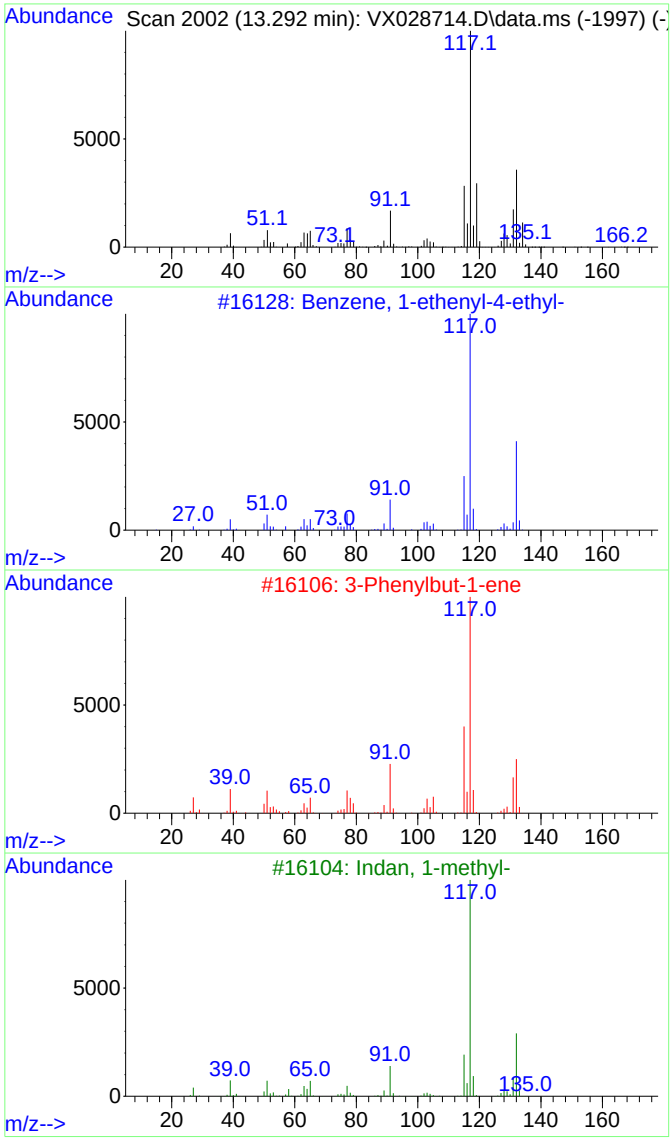
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

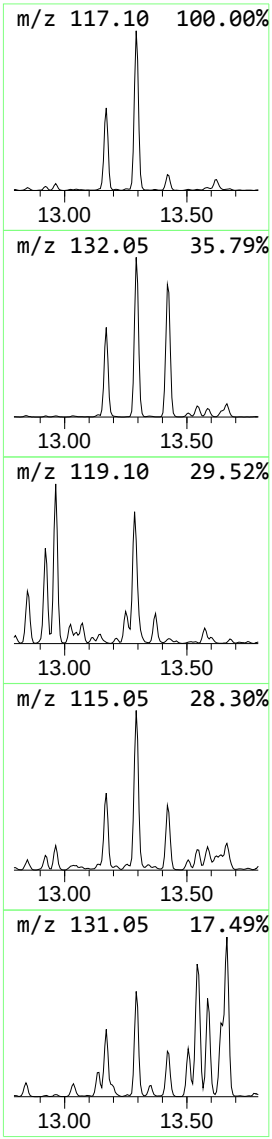
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 3-Phenylbut-1-ene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	70.50 ug/l	2188540	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	78
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

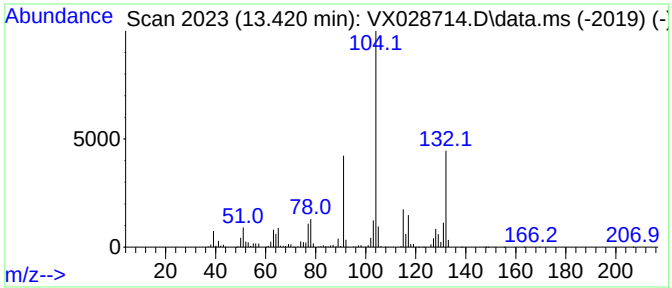
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

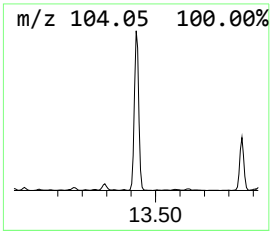
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

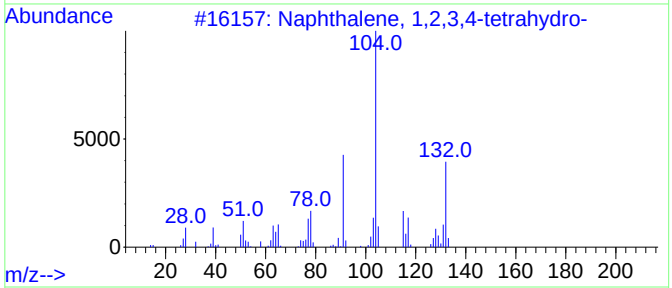
****
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 3

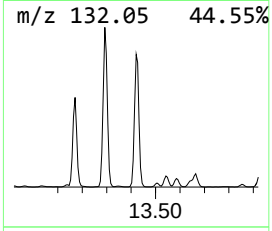
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	49.37 ug/l	1532560	1,4-Dichlorobenzene-d4	12.024

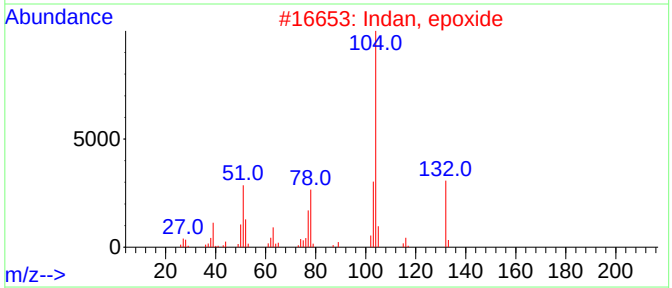
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Indan, epoxide	132	C9H8O	000768-22-9	58
3	Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
4	N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40
5	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	38

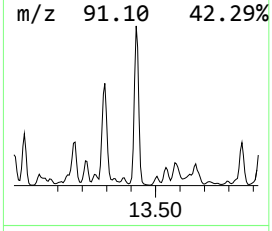


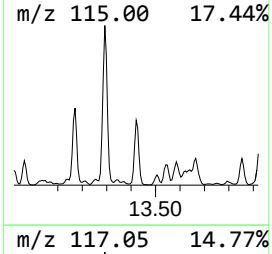


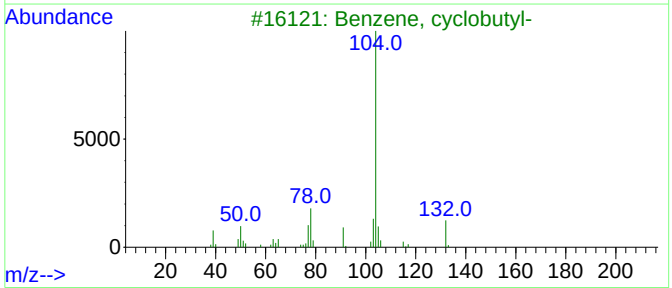


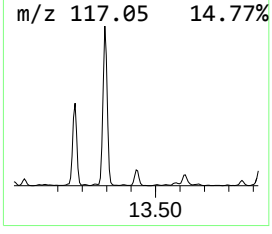












Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

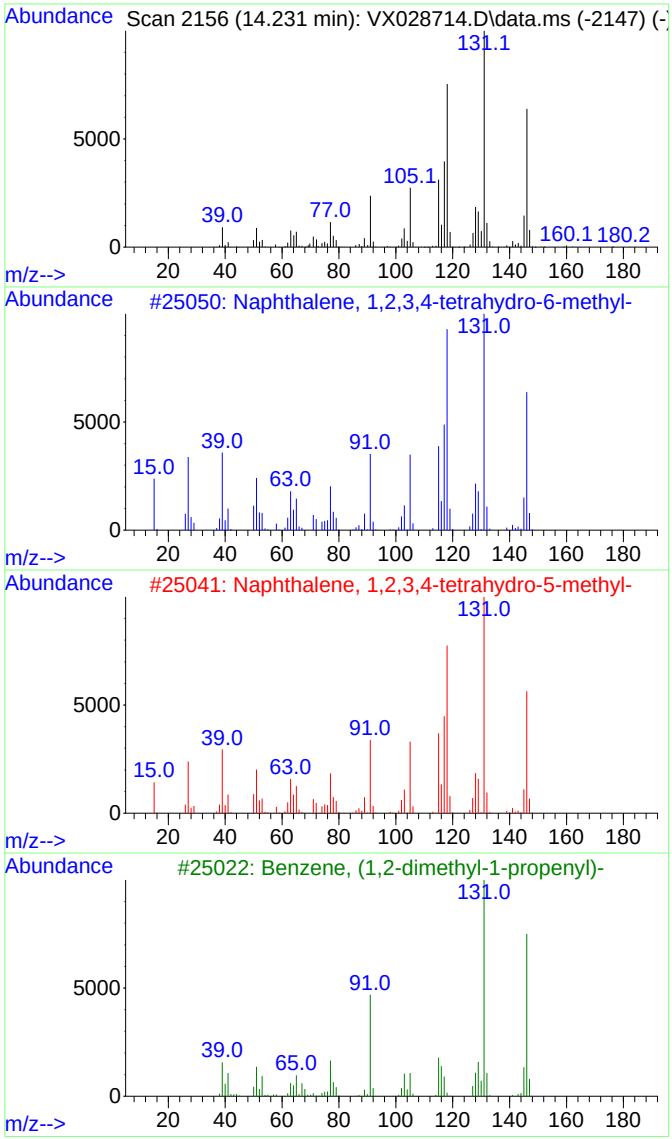
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

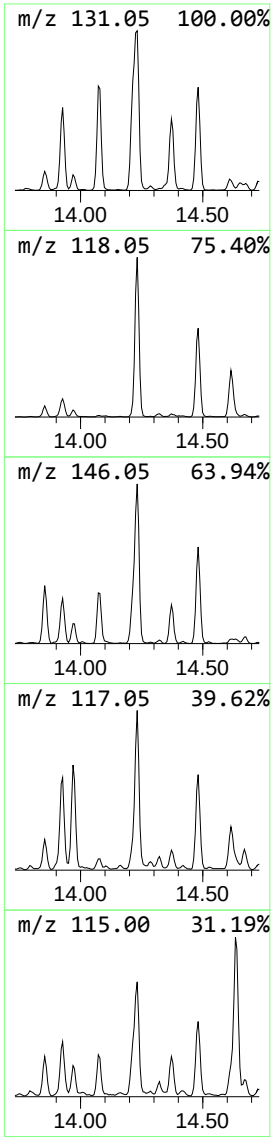
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.231	58.24 ug/l	1808100	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	97
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	60
5	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

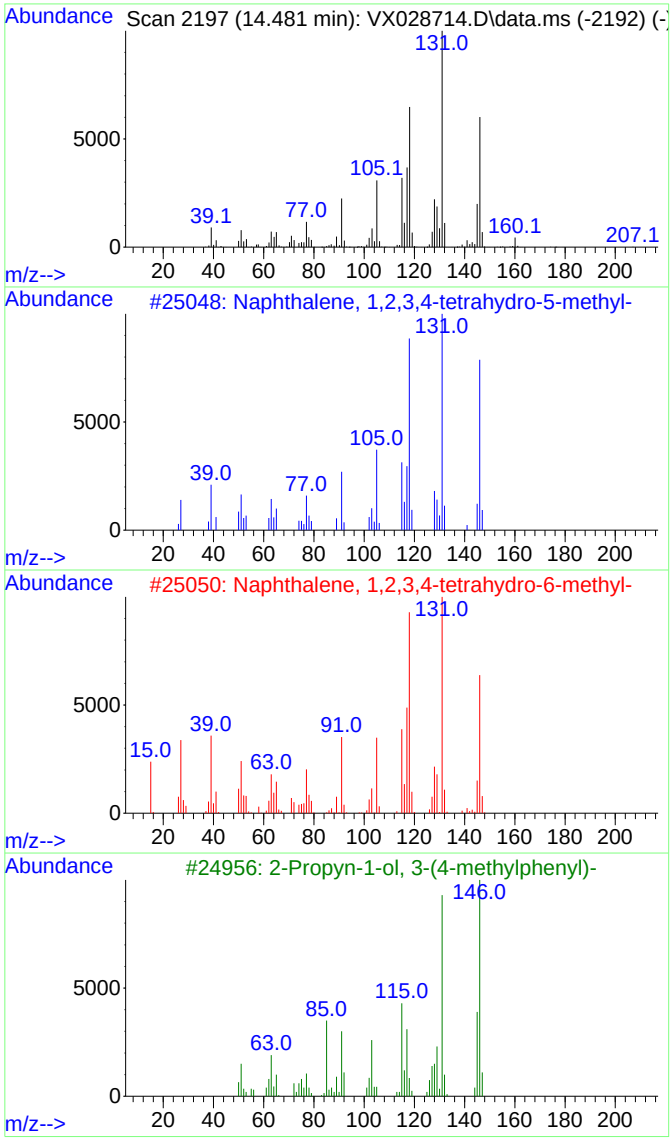
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

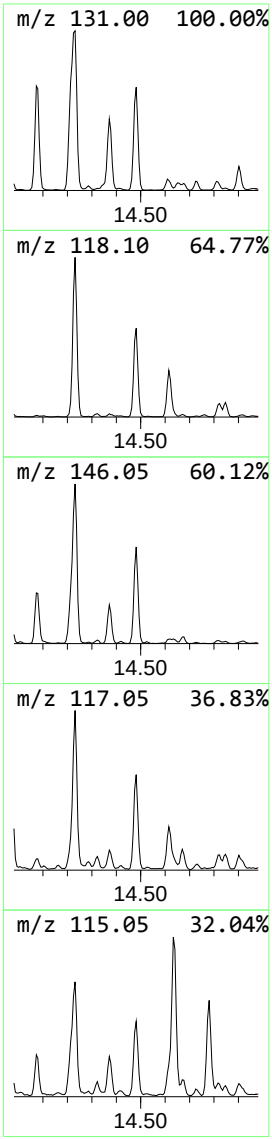
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.481	31.71 ug/l	984329	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	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	95
3	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	72
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

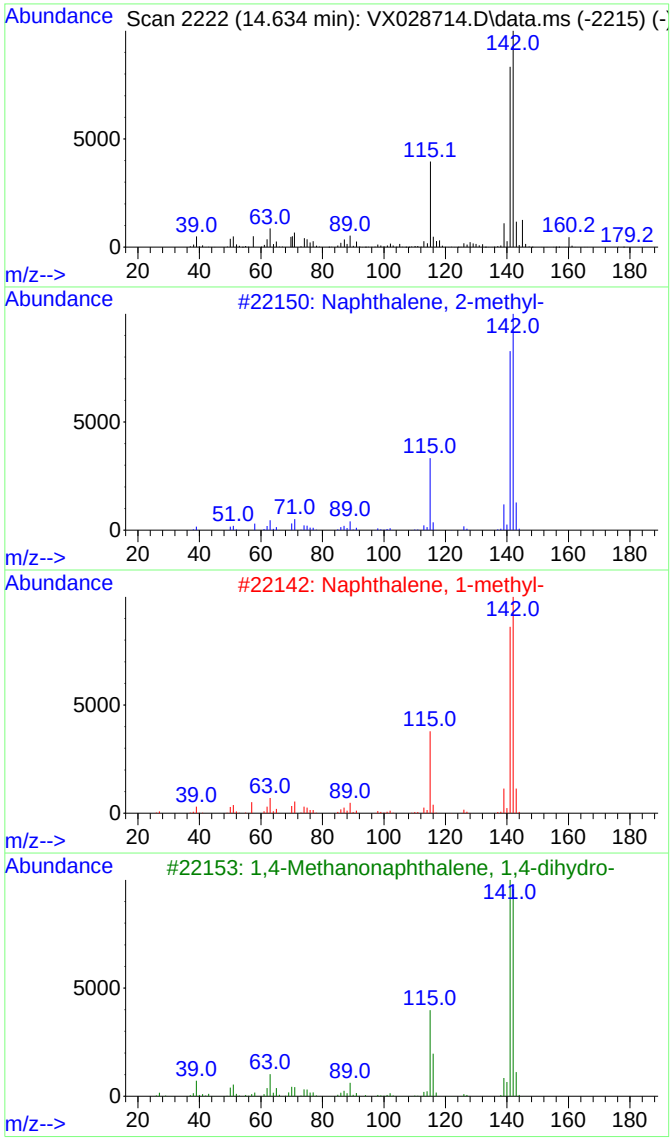
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
Quant Title : SW846 8260

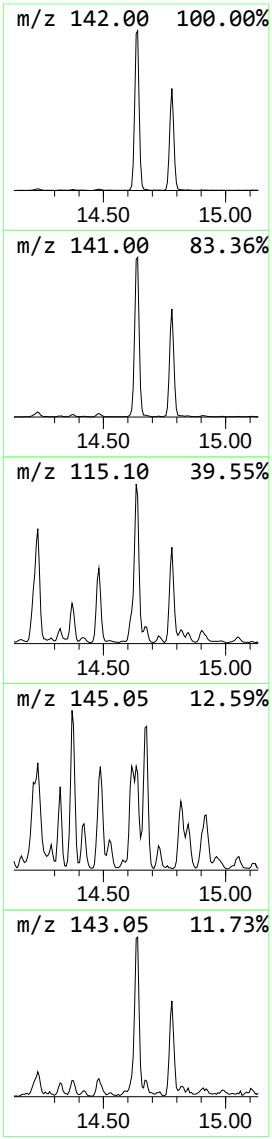
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.634	44.32 ug/l	1375770	1,4-Dichlorobenzene-d4	12.024

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	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051322\
Data File : VX028714.D
Acq On : 14 May 2022 06:04
Operator : JC/MD
Sample : N2813-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 51 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BUR-1144

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.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
Benzene, 1-ethy...	11.378	43.7	ug/l	1356520	4	12.024	1552170	50.0
Mesitylene	12.085	33.5	ug/l	1038880	4	12.024	1552170	50.0
Benzene, 1-ethe...	12.244	39.5	ug/l	1225250	4	12.024	1552170	50.0
Benzene, 1,2,4,...	12.963	24.5	ug/l	760080	4	12.024	1552170	50.0
Benzene, 1-ethe...	13.171	31.8	ug/l	987110	4	12.024	1552170	50.0
3-Phenylbut-1-ene	13.292	70.5	ug/l	2188540	4	12.024	1552170	50.0
Naphthalene, 1,...	13.420	49.4	ug/l	1532560	4	12.024	1552170	50.0
Naphthalene, 1,...	14.231	58.2	ug/l	1808100	4	12.024	1552170	50.0
Naphthalene, 1,...	14.481	31.7	ug/l	984329	4	12.024	1552170	50.0
Naphthalene, 1-...	14.634	44.3	ug/l	1375770	4	12.024	1552170	50.0