LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

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\82X031419W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.446	206	212	223	rVB	144779	241760	2.75%	0.336%
2	2.891	280	285	296	rVB2	49689	109112	1.24%	0.151%
3	3.172	322	331	345	rVB2	54479	145721	1.65%	0.202%
4	4.397	523	532	546	rVB	106650	310528	3.53%	0.431%
5	5.507	698	714	721	rBV	157863	461740	5.24%	0.641%
6	5.665	731	740	767	rVB	307360	1098934	12.48%	1.525%
7	5.927	772	783	796	rVB4	44267	136930	1.55%	0.190%
8	6.068	796	806	816	rBV2	157112	397135	4.51%	0.551%
9	6.257	823	837	845	rBV3	51727	171207	1.94%	0.238%
10	6.354	845	853	861	rVV3	90459	265468	3.01%	0.369%
11	6.458	861	870	882	rVB2	52855	149804	1.70%	0.208%
12	6.860	925	936	943	rBV	414057	990523	11.25%	1.375%
13	6.933	943	948	962	rVB3	81690	240114	2.73%	0.333%
14	7.256	993	1001	1009	rBV2	60600	131704	1.50%	0.183%
15	7.458	1021	1034	1048	rBV4	220527	630615	7.16%	0.875%
16	7.732	1073	1079	1087	rVB	67826	131936	1.50%	0.183%
17	8.055	1122	1132	1143	rVB	57607	146656	1.67%	0.204%
18	8.177	1143	1152	1154	rBV	98316	193254	2.19%	0.268%
19	8.214	1154	1158	1163	rVB	184991	306359	3.48%	0.425%
20	8.573	1210	1217	1226	rVB2	104631	203440	2.31%	0.282%
21	8.671	1226	1233	1235	rBV2	50645	95021	1.08%	0.132%
22	8.713	1235	1240	1248	rVB	823094	1296882	14.73%	1.800%
23	8.982	1279	1284	1289	rBV	65990	101975	1.16%	0.142%
24	9.165	1305	1314	1318	rBV	59540	109278	1.24%	0.152%
25	9.219	1318	1323	1334	rVB	186044	295979	3.36%	0.411%
26	9.561	1375	1379	1385	rVB3	46774	93324	1.06%	0.130%
27	9.658	1392	1395	1401	rVB2	64958	94032	1.07%	0.131%
28	10.110	1459	1469	1474	rBV	762273	1046291	11.88%	1.452%
29	11.018	1612	1618	1627	rBV	810437	1014276	11.52%	1.408%
30	11.134	1633	1637	1642	rBV	629570	797834	9.06%	1.108%
31	11.359	1665	1674	1680	rBV	3056068	3637155	41.30%	5.049%
32	11.664	1718	1724	1732	rBV	317079	411714	4.67%	0.572%
33	11.768	1733	1741	1745	rBV	66115	90376	1.03%	0.125%
34	11.920	1760	1766	1767	rBV	171185	209163	2.37%	0.290%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

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\82X031419W.M
Title : SW846 8260

35	11.945	1767	1770	1776	rVB	321181	388783	4.41%	0.540%
36	12.079	1786	1792	1797	rBV	771993	971394	11.03%	1.348%
37	12.231	1812	1817	1822	rVB	220055	262646	2.98%	0.365%
38	12.298	1822	1828	1834	rBV	4532316	5472429	62.14%	7.597%
39	12.365	1834	1839	1847	rVB2	1013814	1583611	17.98%	2.198%
40	12.457	1849	1854	1860	rVB	567351	690688	7.84%	0.959%
41	12.524	1860	1865	1870	rBV2	138074	218520	2.48%	0.303%
42	12.585	1870	1875	1882	rVB	676161	799614	9.08%	1.110%
43	12.670	1882	1889	1892	rBV	5616740	6890911	78.24%	9.566%
44	12.701	1892	1894	1898	rVV	1136281	1197358	13.60%	1.662%
45	12.755	1898	1903	1910	rVB2	3664148	5098054	57.89%	7.077%
46	12.896	1921	1926	1932	rVB2	254522	353777	4.02%	0.491%
47	12.975	1932	1939	1943	rBV	3386479	4104319	46.60%	5.697%
48	13.018	1943	1946	1952	rVB	1704869	1948924	22.13%	2.705%
49	13.085	1952	1957	1962	rBV3	221649	494134	5.61%	0.686%
50	13.194	1971	1975	1976	rBV2	191319	206548	2.35%	0.287%
51	13.225	1976	1980	1989	rVB	3058843	3719078	42.23%	5.163%
52	13.310	1989	1994	1996	rBV	248088	373223	4.24%	0.518%
53	13.347	1996	2000	2006	rVB2	6381532	8806901	100.00%	12.225%
54	13.402	2006	2009	2011	rBV2	144534	179371	2.04%	0.249%
55	13.426	2011	2013	2018	rVB	153836	159938	1.82%	0.222%
56	13.481	2018	2022	2028	rVB	254059	326354	3.71%	0.453%
57	13.560	2028	2035	2038	rBV2	277779	404952	4.60%	0.562%
58	13.597	2038	2041	2044	rBV	628273	803549	9.12%	1.115%
59	13.639	2044	2048	2053	rVB2	645200	1064856	12.09%	1.478%
60	13.719	2058	2061	2067	rVB2	727993	1131276	12.85%	1.570%
61	13.853	2078	2083	2088	rVB3	124768	188035	2.14%	0.261%
62	13.908	2088	2092	2097	rVB	147833	192204	2.18%	0.267%
63	13.981	2097	2104	2108	rBV2	383572	537426	6.10%	0.746%
64	14.023	2108	2111	2116	rVB	265902	328673	3.73%	0.456%
65	14.133	2124	2129	2133	rBV	463829	616739	7.00%	0.856%
66	14.273	2147	2152	2158	rBV2	499400	971631	11.03%	1.349%
67	14.377	2164	2169	2174	rBV	263928	386438	4.39%	0.536%
68	14.426	2174	2177	2182	rVB	291027	376317	4.27%	0.522%
69	14.536	2190	2195	2200	rBV2	294818	401385	4.56%	0.557%
70	14.657	2209	2215	2217	rBV	81118	133739	1.52%	0.186%
71	14.694	2217	2221	2226	rVB	924680	1108963	12.59%	1.539%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

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\82X031419W.M
Title : SW846 8260

72	14.834	2239	2244	2253	rBV	2108061	2774067	31.50%	3.851%
73	15.548	2354	2361	2366	rBV	96819	160118	1.82%	0.222%
74	15.688	2375	2384	2387	rBV	130737	227577	2.58%	0.316%
75	15.724	2387	2390	2396	rVB	87798	127254	1.44%	0.177%
76	15.913	2411	2421	2432	rVB2	35252	100365	1.14%	0.139%

Sum of corrected areas: 72038379

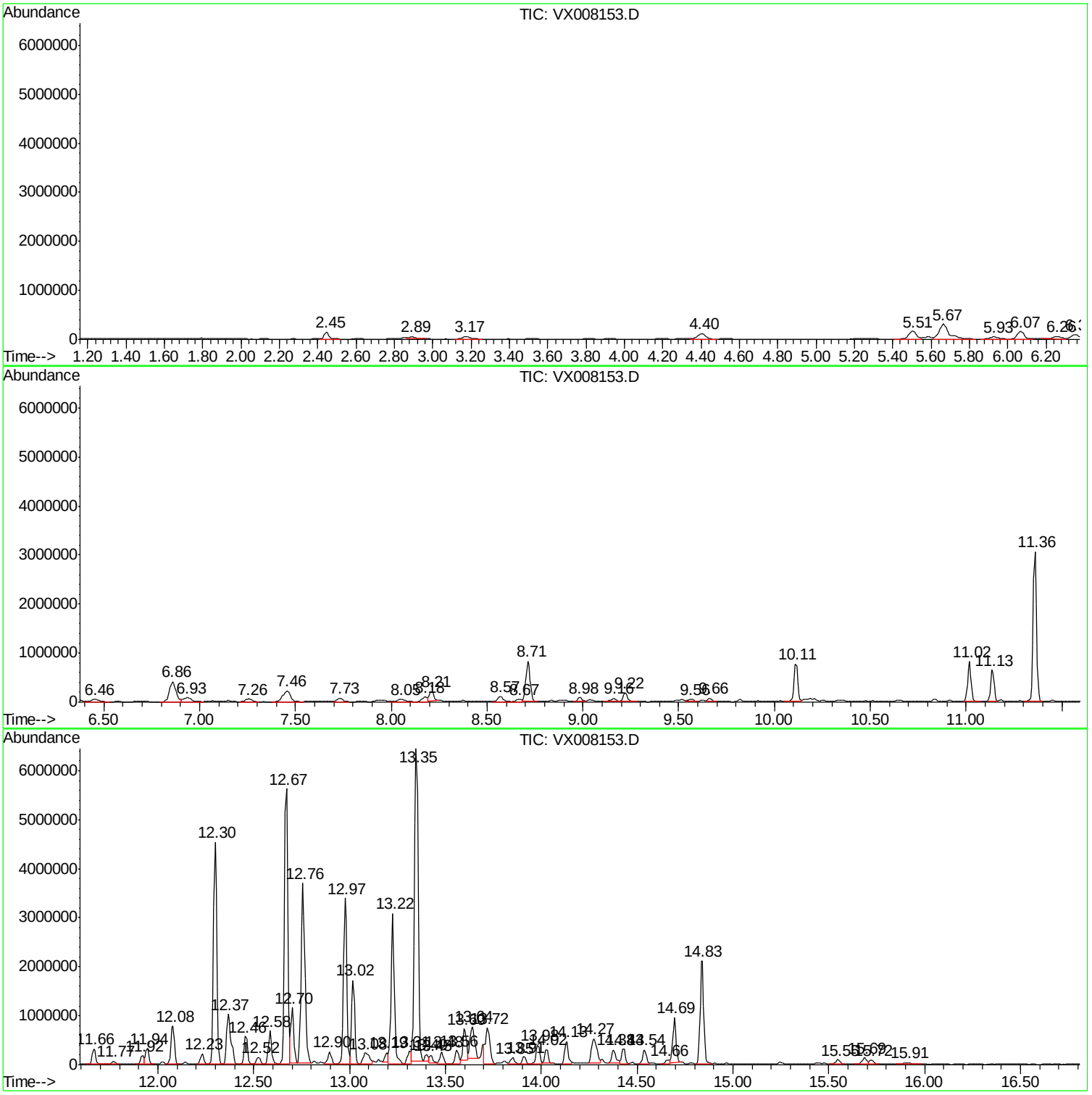
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

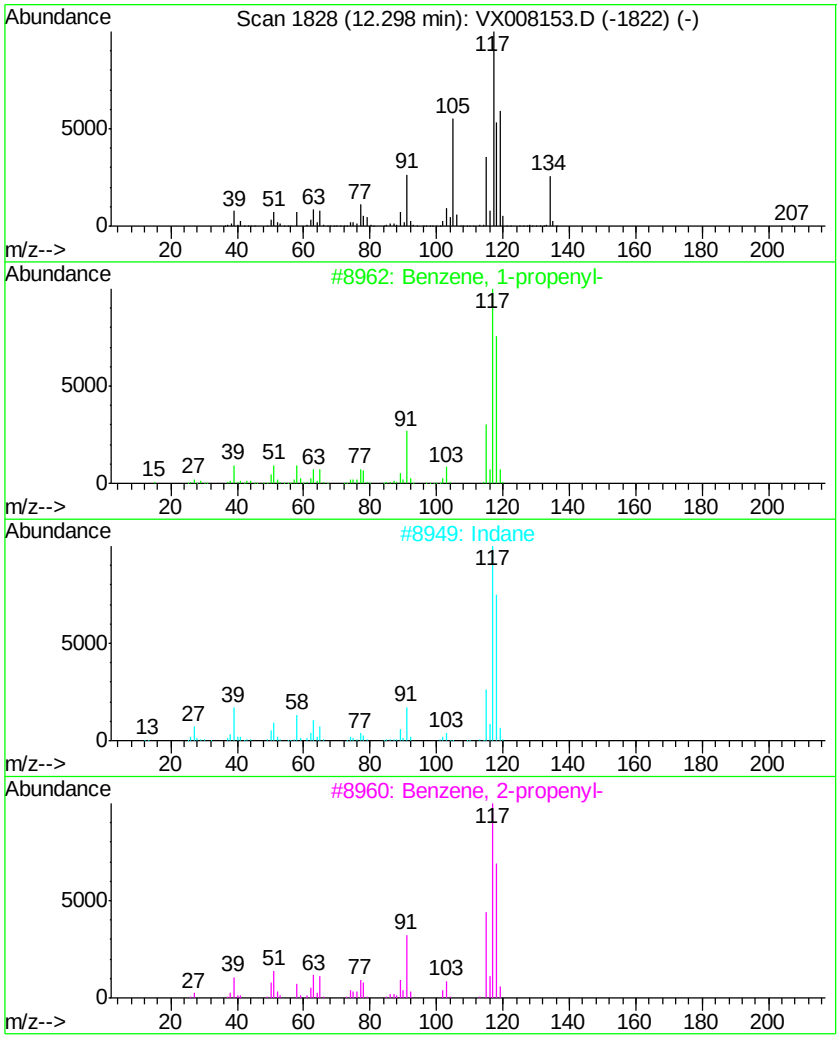
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

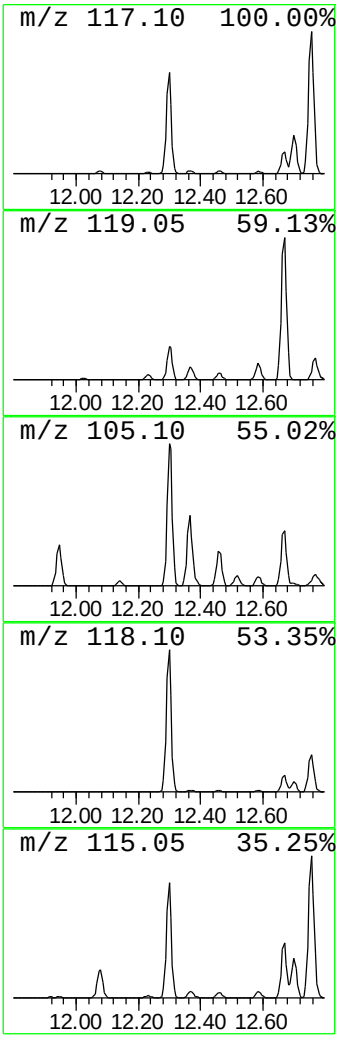
****

Peak Number 1 Benzene, 1-propenyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	281.68 ug/l	5472430	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
2	Indane	118	C9H10	000496-11-7	60
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

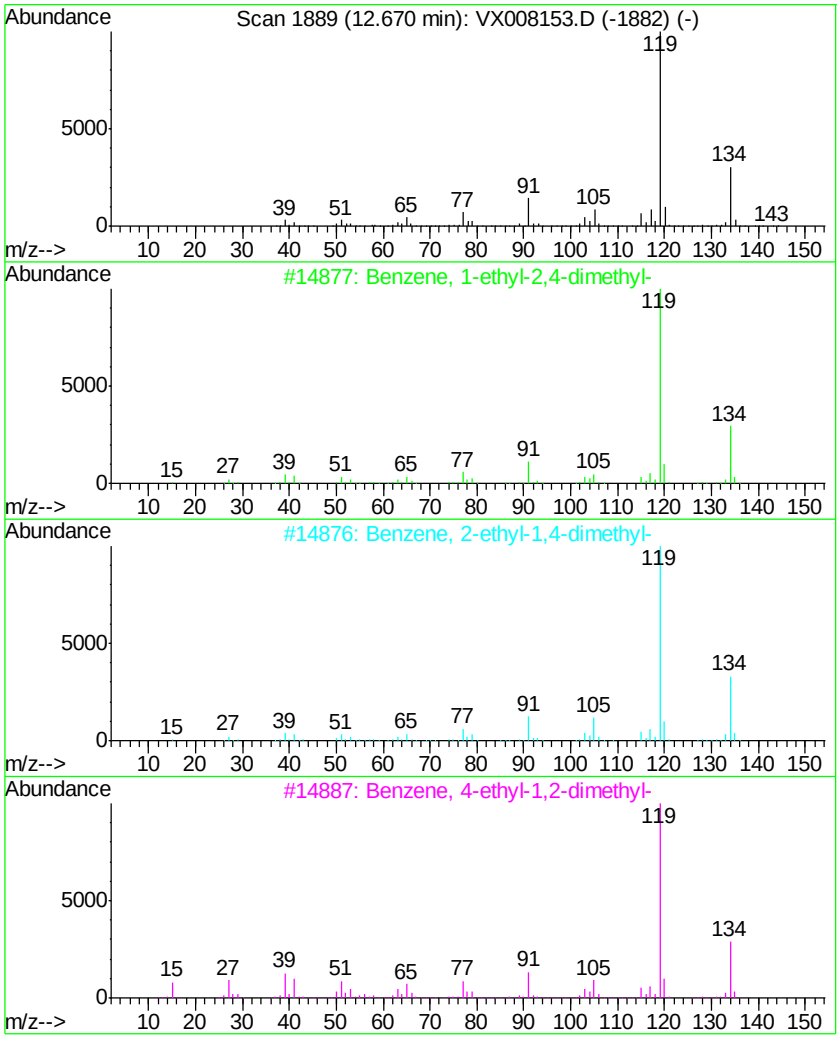
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

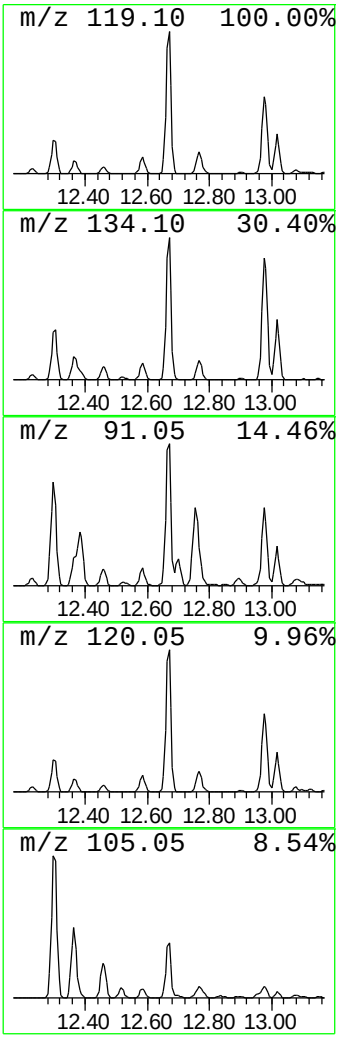
****

Peak Number 2 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	354.69 ug/l	6890910	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

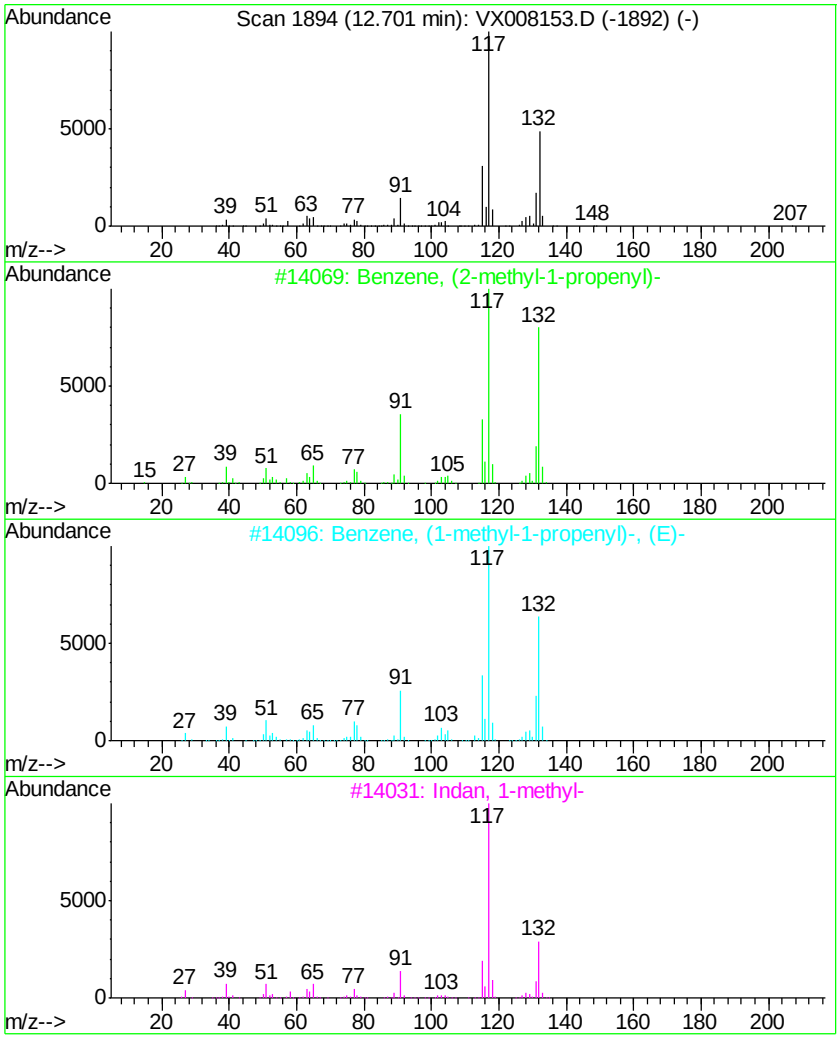
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

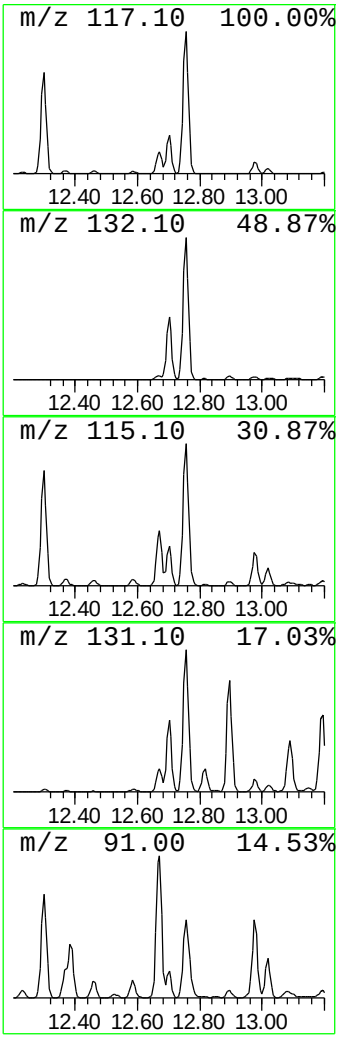
****

Peak Number 3 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	61.63 ug/l	1197360	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	90
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

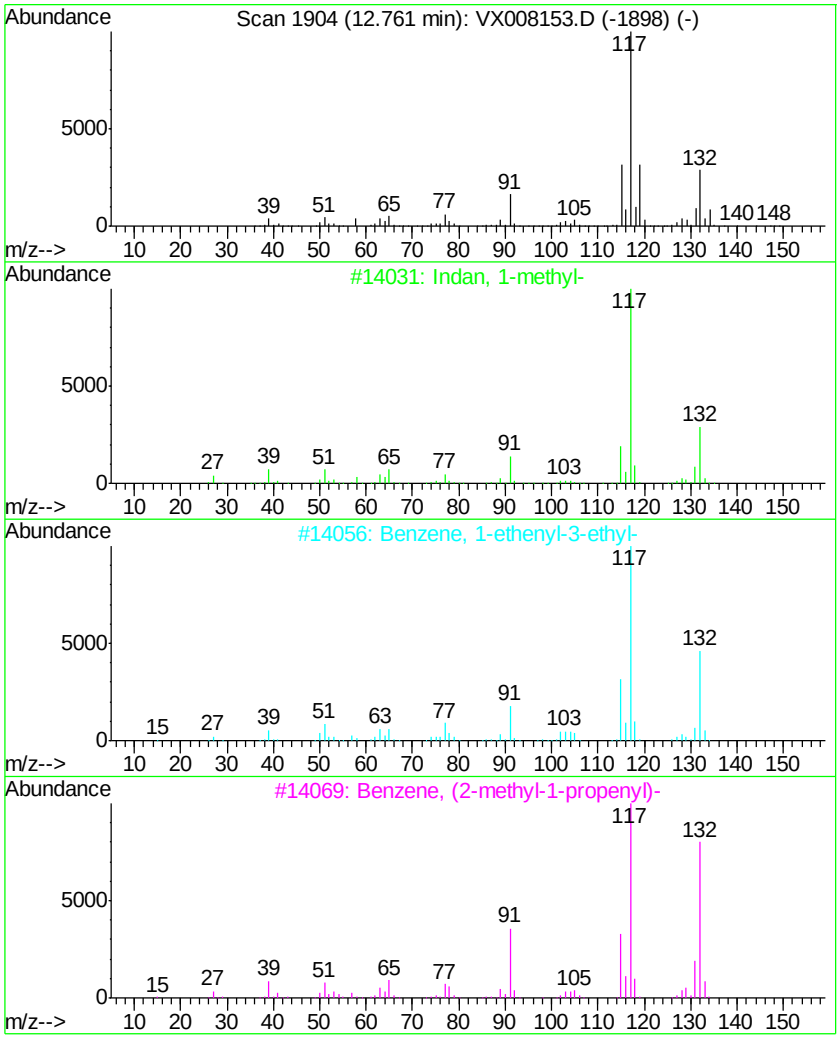
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

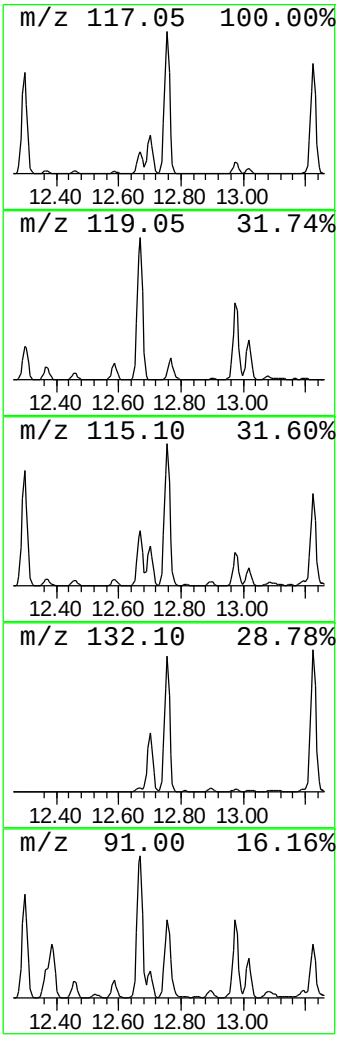
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	262.41 ug/l	5098050	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
5	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	81





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

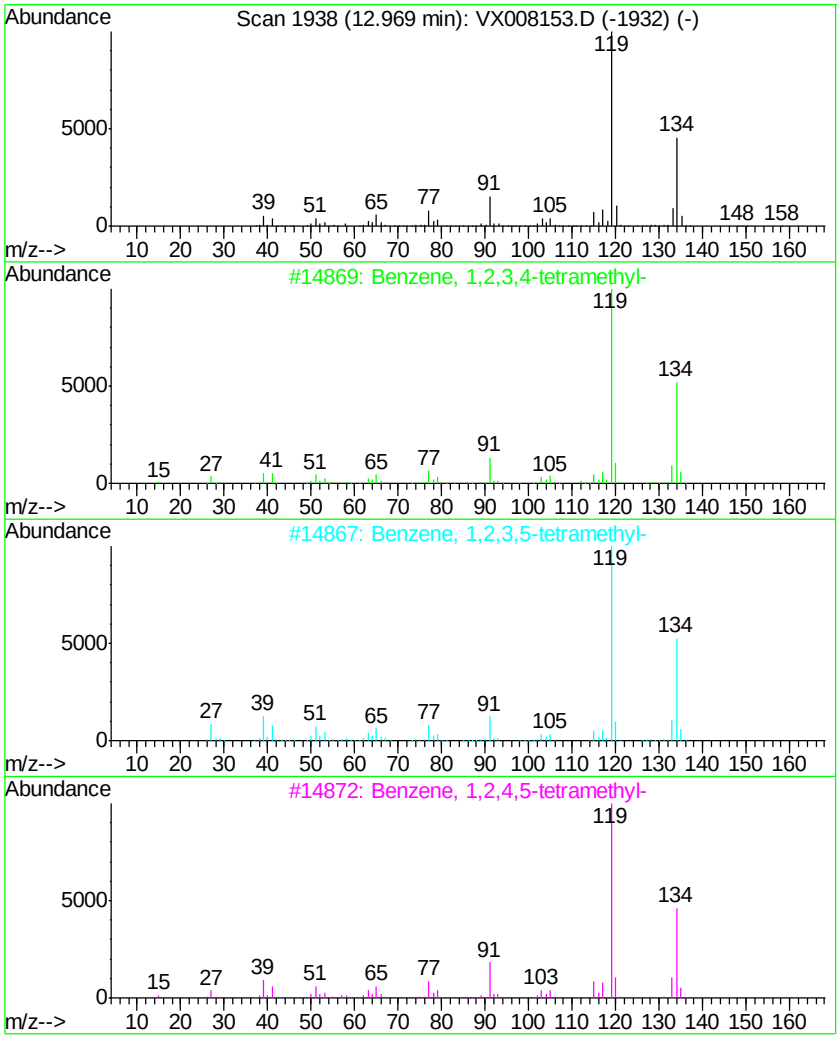
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

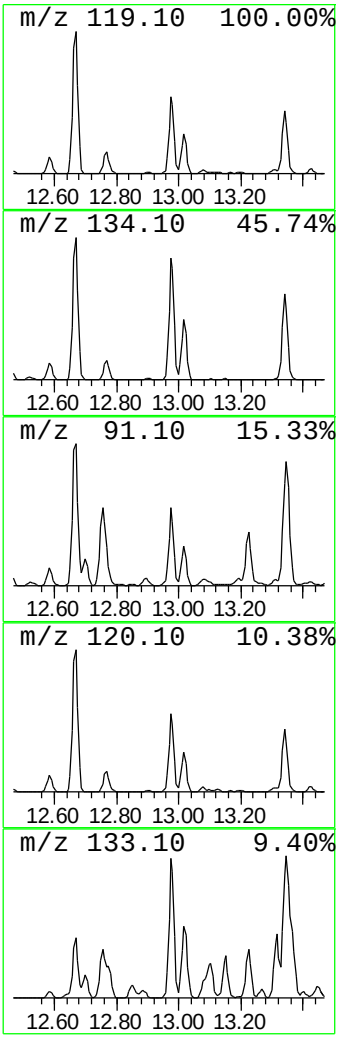
****

Peak Number 5 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	211.26 ug/l	4104320	1,4-Dichlorobenzene-d4	12.08

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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

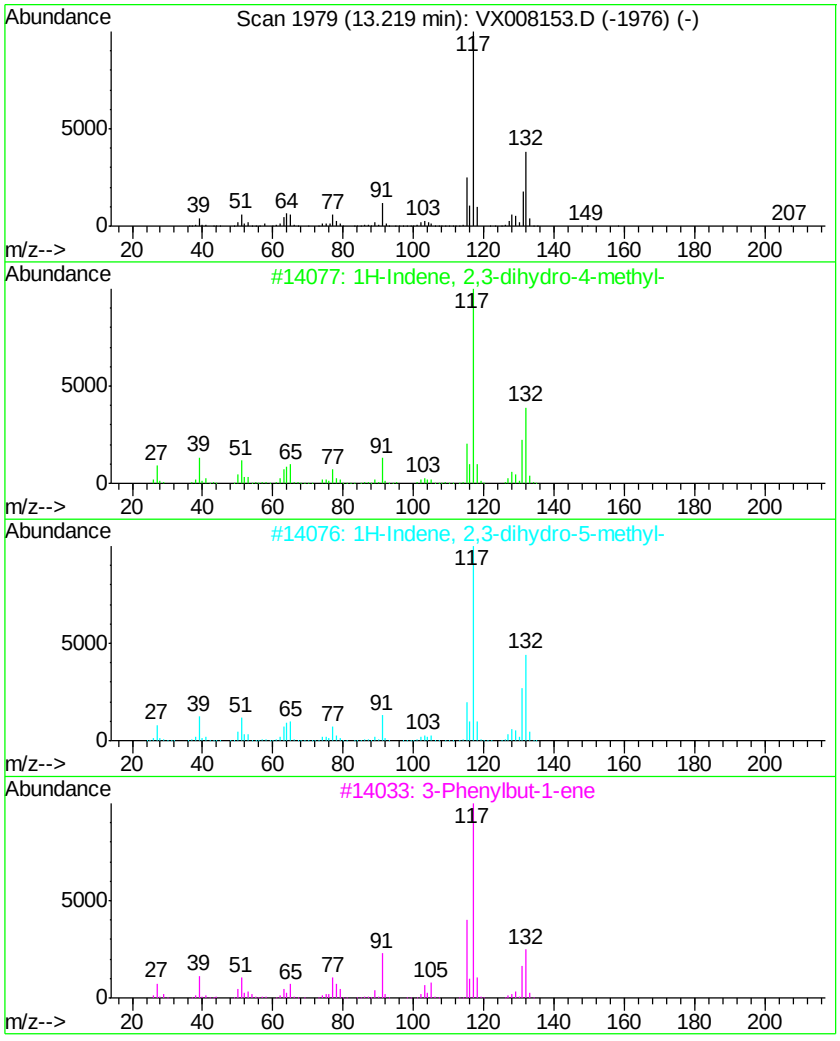
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

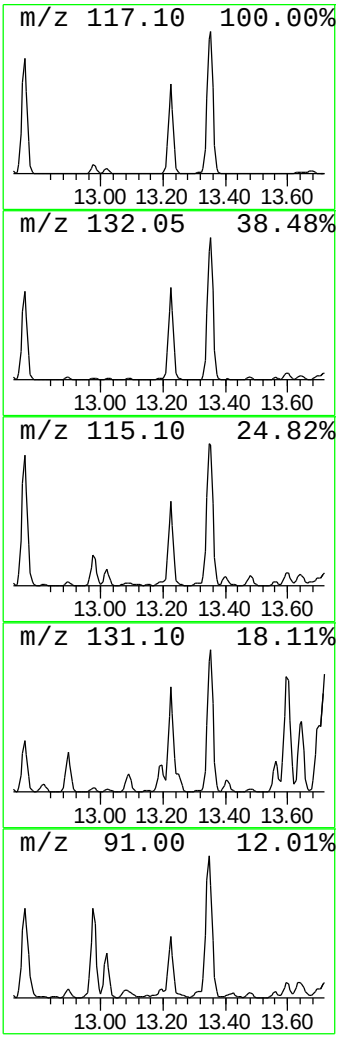
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	191.43 ug/l	3719080	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4	Indan, 1-methyl-	132	C10H12	000767-58-8	87
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

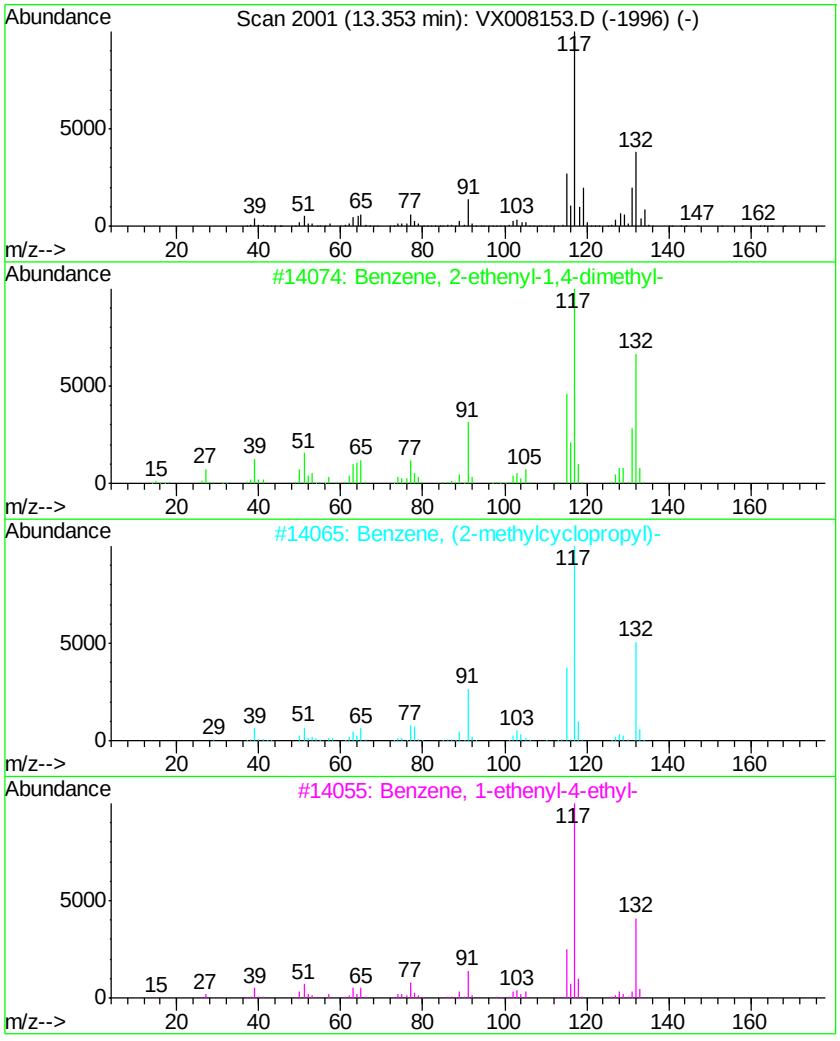
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

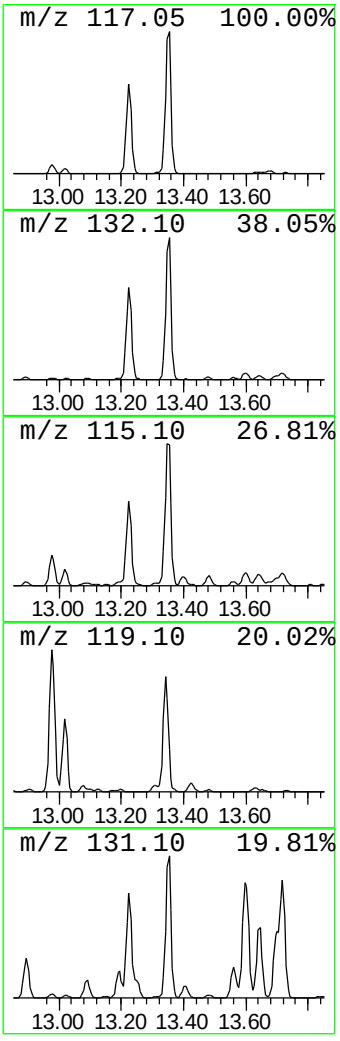
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	453.31 ug/l	8806900	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	90
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

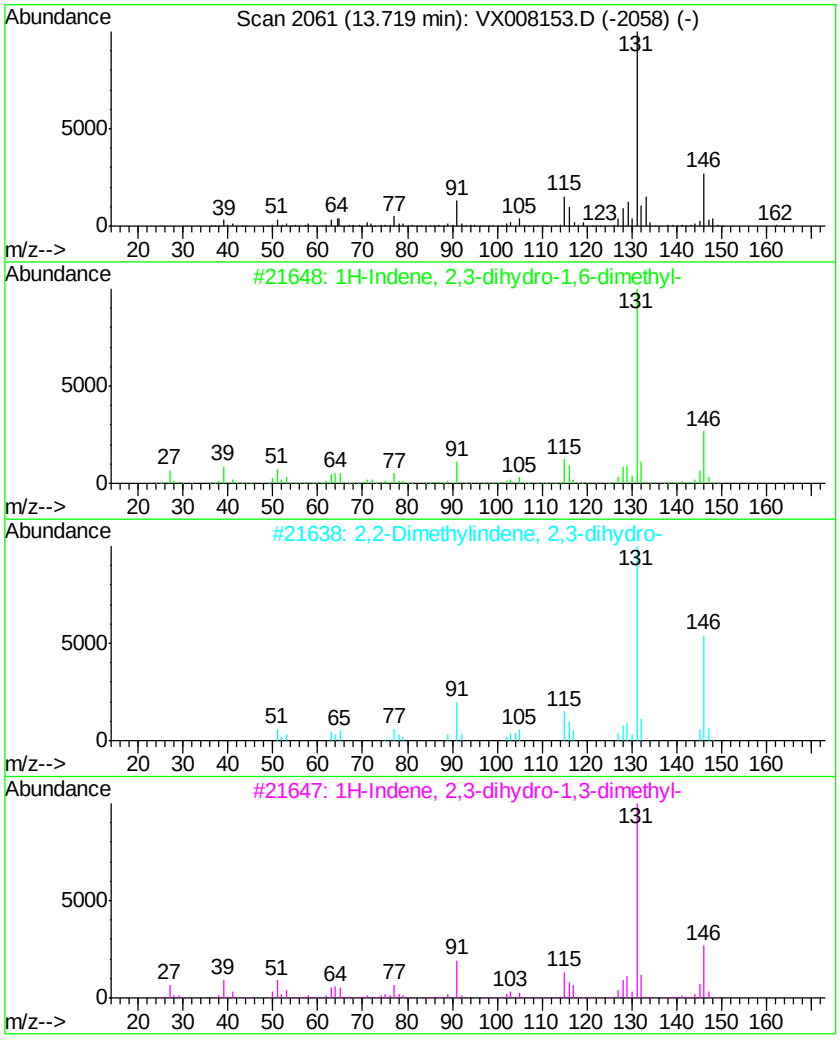
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

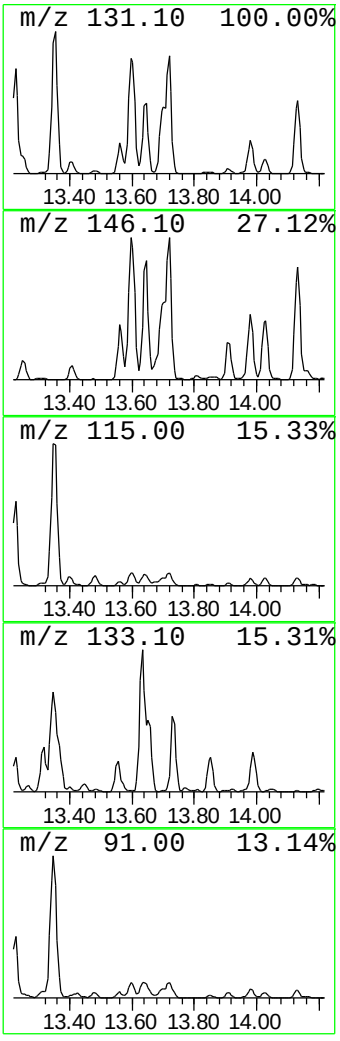
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	58.23 ug/l	1131280	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

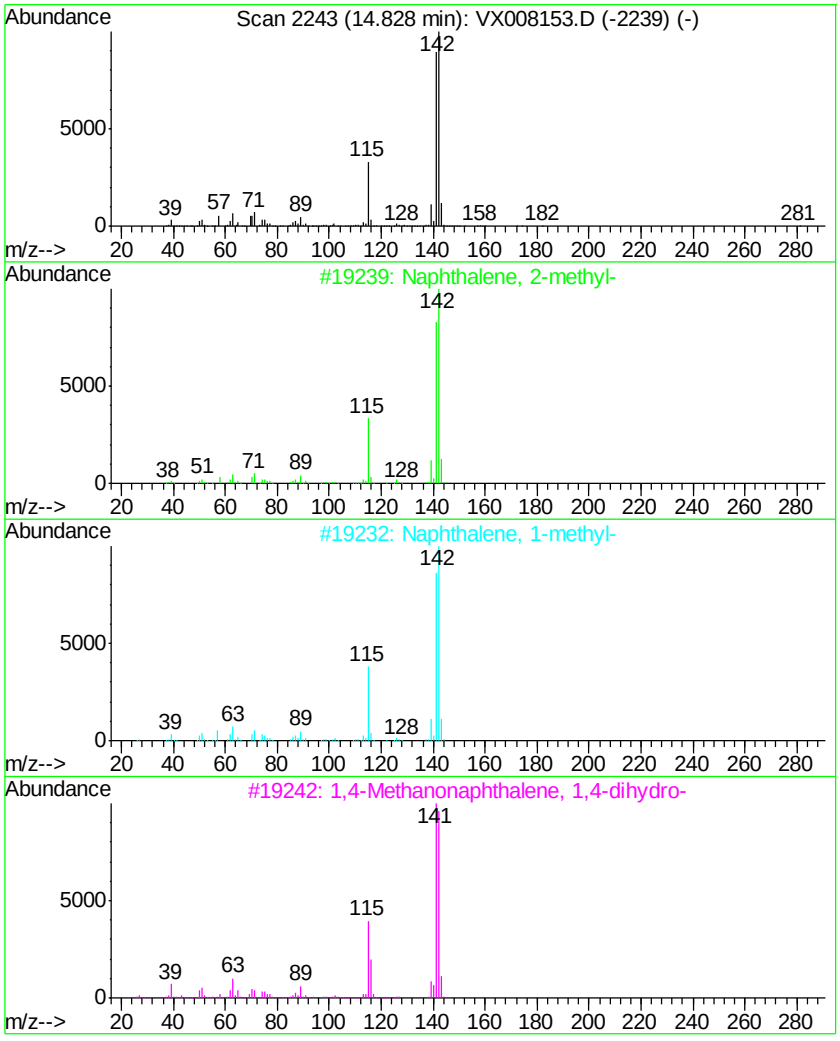
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

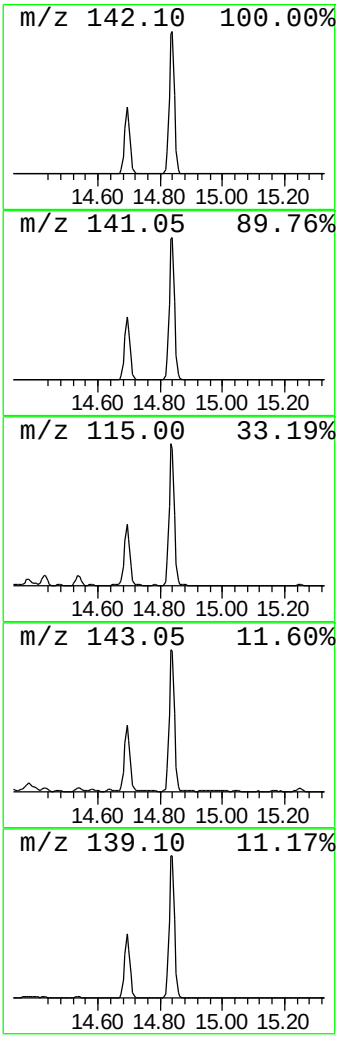
****

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	142.79 ug/l	2774070	1,4-Dichlorobenzene-d4	12.08

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	91
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX031619\
Data File : VX008153.D
Acq On : 16 Mar 2019 20:40
Operator : JC/SP
Sample : K1960-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-2-9.0-031419

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X031419W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-propenyl-	12.30	281.7	ug/l	5472430	4	12.08	971394	50.0
Benzene, 1-ethyl-...	12.67	354.7	ug/l	6890910	4	12.08	971394	50.0
Benzene, (2-methy...	12.70	61.6	ug/l	1197360	4	12.08	971394	50.0
Indan, 1-methyl-	12.76	262.4	ug/l	5098050	4	12.08	971394	50.0
Benzene, 1,2,3,4-...	12.97	211.3	ug/l	4104320	4	12.08	971394	50.0
1H-Indene, 2,3-di...	13.22	191.4	ug/l	3719080	4	12.08	971394	50.0
Benzene, 2-etheny...	13.35	453.3	ug/l	8806900	4	12.08	971394	50.0
1H-Indene, 2,3-di...	13.72	58.2	ug/l	1131280	4	12.08	971394	50.0
Naphthalene, 1-me...	14.83	142.8	ug/l	2774070	4	12.08	971394	50.0