LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

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\82X061919W.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	2.849	270	278	289	rBV3	21433	50906	1.14%	0.165%
2	4.312	508	518	528	rBV2	15007	45811	1.03%	0.148%
3	5.501	699	713	722	rBV	132369	383720	8.61%	1.242%
4	5.665	729	740	748	rBV	224321	647299	14.52%	2.096%
5	6.061	795	805	820	rVB	127460	336941	7.56%	1.091%
6	6.348	844	852	864	rVB	75641	223884	5.02%	0.725%
7	6.860	925	936	950	rBV	337417	805842	18.07%	2.609%
8	7.457	1023	1034	1044	rVB4	22597	62984	1.41%	0.204%
9	8.055	1119	1132	1140	rBV2	53484	114897	2.58%	0.372%
10	8.177	1144	1152	1161	rBV	93719	183202	4.11%	0.593%
11	8.713	1223	1240	1256	rBV	740345	1176717	26.39%	3.810%
12	10.109	1463	1469	1478	rBV2	715557	1210071	27.14%	3.918%
13	10.183	1478	1481	1486	rVB	36721	51404	1.15%	0.166%
14	10.335	1501	1506	1510	rBV	38073	62449	1.40%	0.202%
15	10.658	1544	1559	1564	rBV3	63545	154987	3.48%	0.502%
16	10.835	1584	1588	1596	rVB2	82639	118621	2.66%	0.384%
17	10.914	1596	1601	1607	rBV5	42119	76869	1.72%	0.249%
18	11.018	1613	1618	1622	rBV	289032	359927	8.07%	1.165%
19	11.134	1632	1637	1642	rBV	627966	782533	17.55%	2.534%
20	11.182	1642	1645	1649	rVB	74598	91071	2.04%	0.295%
21	11.280	1657	1661	1665	rVB	79376	98942	2.22%	0.320%
22	11.359	1669	1674	1678	rBV	83797	103028	2.31%	0.334%
23	11.438	1683	1687	1692	rVB2	30274	48833	1.10%	0.158%
24	11.670	1719	1725	1732	rBV	128886	184370	4.14%	0.597%
25	11.768	1732	1741	1745	rBV	240622	310371	6.96%	1.005%
26	11.920	1758	1766	1768	rBV	528698	776705	17.42%	2.515%
27	11.944	1768	1770	1776	rVV	713456	752363	16.87%	2.436%
28	12.079	1786	1792	1803	rBV	774813	1074500	24.10%	3.479%
29	12.231	1812	1817	1823	rVB	882988	1046937	23.48%	3.390%
30	12.298	1823	1828	1833	rVV	625242	756306	16.96%	2.449%
31	12.365	1835	1839	1849	rVB2	128064	211852	4.75%	0.686%
32	12.457	1849	1854	1862	rBV2	113083	159089	3.57%	0.515%
33	12.670	1878	1889	1892	rBV	2954699	3752130	84.16%	12.148%
34	12.700	1892	1894	1898	rVV	575281	709219	15.91%	2.296%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

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\82X061919W.M
Title : SW846 8260

35	12.755	1898	1903	1910	rVV2	3254620	4458514	100.00%	14.435%
36	12.816	1910	1913	1918	rVV	142712	173824	3.90%	0.563%
37	12.896	1918	1926	1931	rVV	700811	960588	21.55%	3.110%
38	12.975	1931	1939	1944	rVV	1732118	2241038	50.26%	7.256%
39	13.085	1952	1957	1963	rVB2	155006	228307	5.12%	0.739%
40	13.152	1963	1968	1971	rBV	189595	238303	5.34%	0.772%
41	13.188	1971	1974	1977	rVV	223810	270453	6.07%	0.876%
42	13.225	1977	1980	1981	rVV	83354	91083	2.04%	0.295%
43	13.249	1981	1984	1988	rVB	137481	166689	3.74%	0.540%
44	13.310	1988	1994	1995	rBV2	86069	110061	2.47%	0.356%
45	13.347	1995	2000	2006	rVV2	2260359	3026721	67.89%	9.800%
46	13.402	2006	2009	2011	rVV2	85805	117600	2.64%	0.381%
47	13.420	2011	2012	2016	rVV	79278	96178	2.16%	0.311%
48	13.475	2018	2021	2027	rVB	55011	82525	1.85%	0.267%
49	13.597	2037	2041	2044	rVV	76544	108609	2.44%	0.352%
50	13.639	2044	2048	2052	rVV2	291226	406761	9.12%	1.317%
51	13.694	2052	2057	2059	rVV2	164988	301010	6.75%	0.975%
52	13.719	2059	2061	2070	rVB2	240029	290601	6.52%	0.941%
53	13.804	2070	2075	2079	rBV	59752	80132	1.80%	0.259%
54	13.847	2079	2082	2088	rVB	42117	59634	1.34%	0.193%
55	13.908	2088	2092	2097	rBV	52487	64272	1.44%	0.208%
56	13.981	2097	2104	2108	rBV3	53962	92221	2.07%	0.299%
57	14.029	2108	2112	2116	rVV2	36504	46321	1.04%	0.150%
58	14.267	2147	2151	2157	rBV	44450	66908	1.50%	0.217%
59	14.535	2191	2195	2200	rBV2	43487	57241	1.28%	0.185%
60	14.840	2238	2245	2253	rBV2	88573	126002	2.83%	0.408%

Sum of corrected areas: 30886376

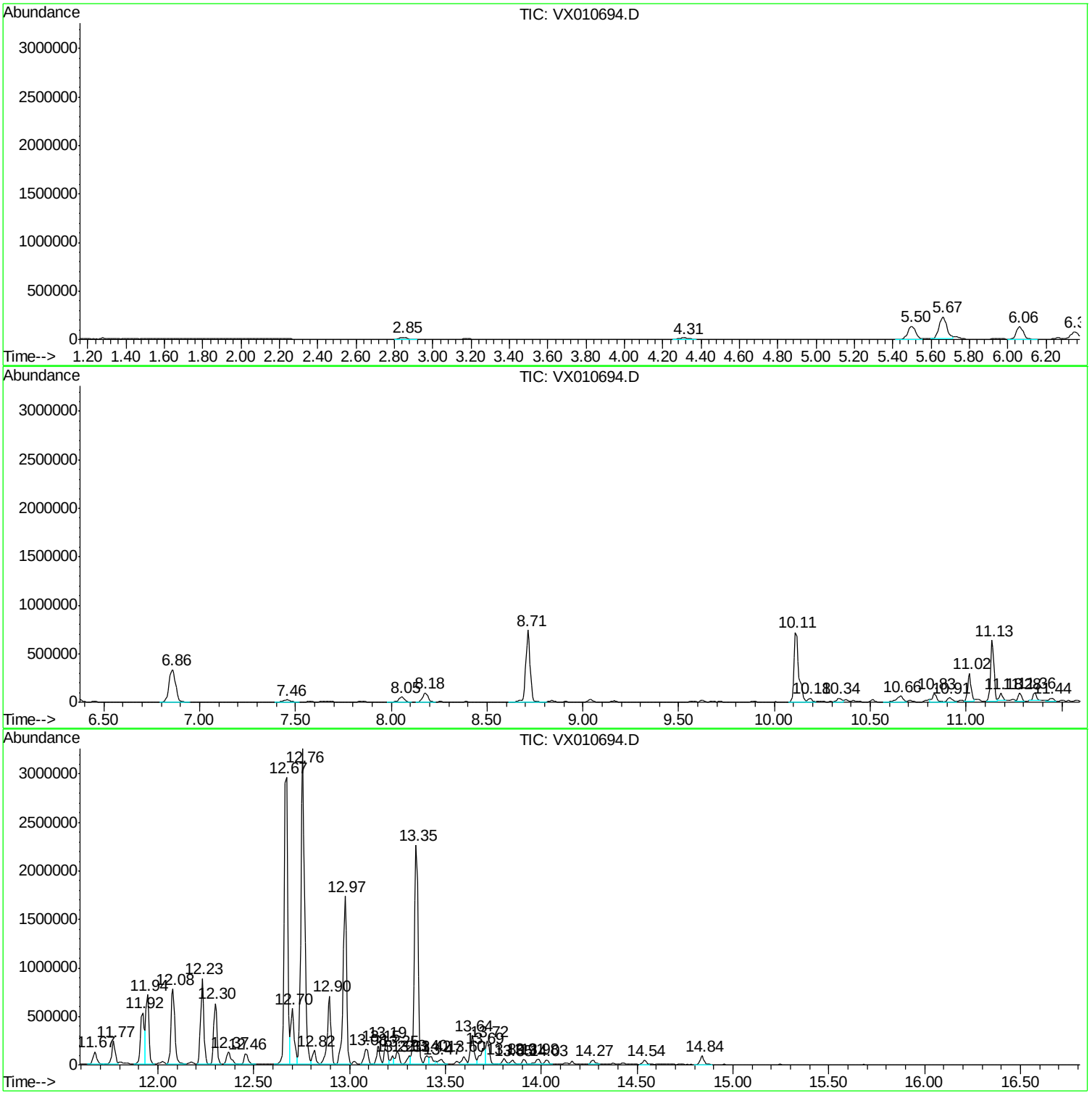
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

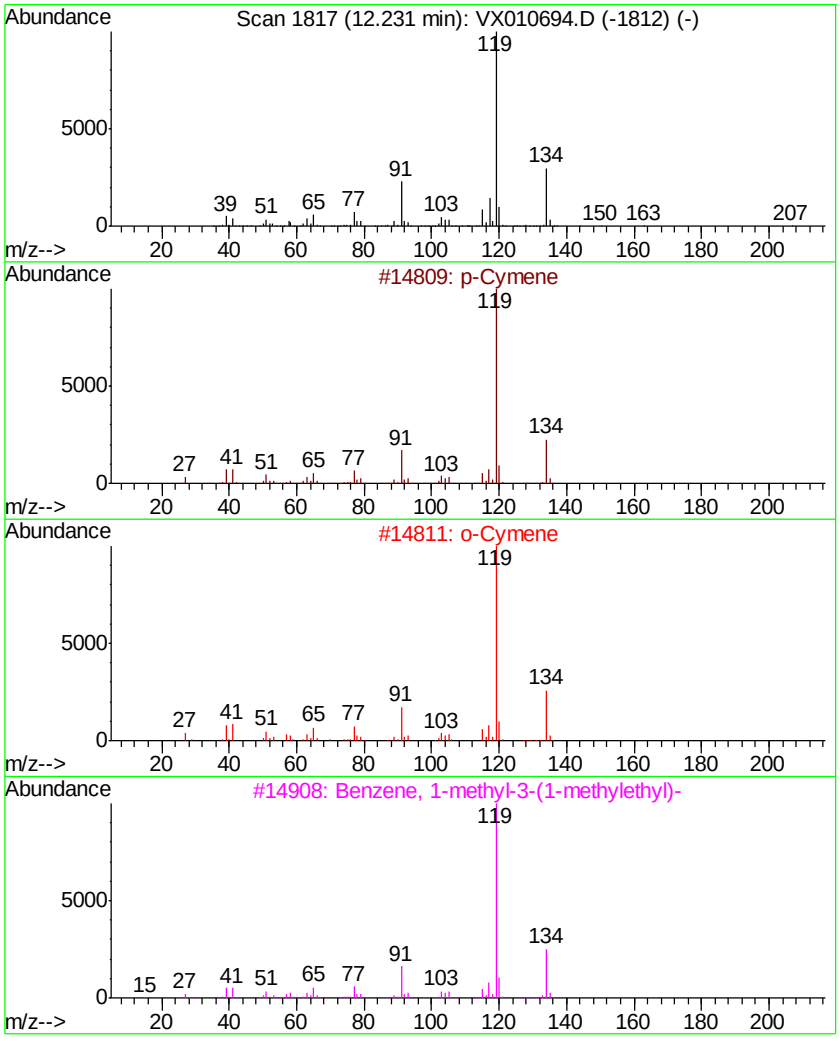
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

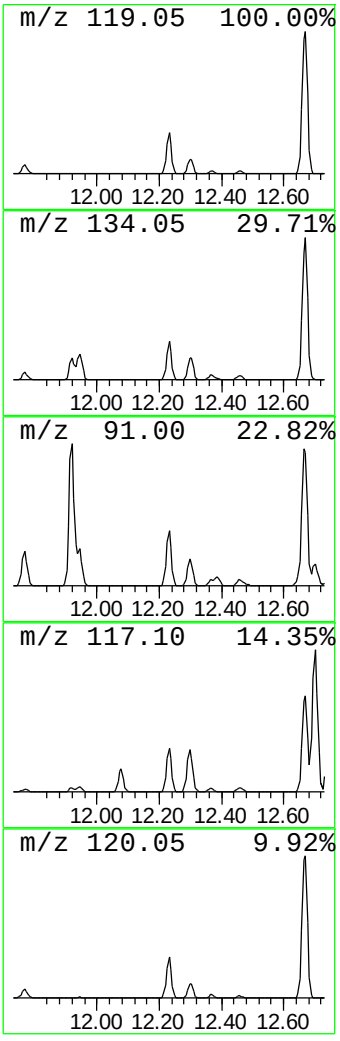
****

Peak Number 1 p-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	48.72 ug/l	1046940	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

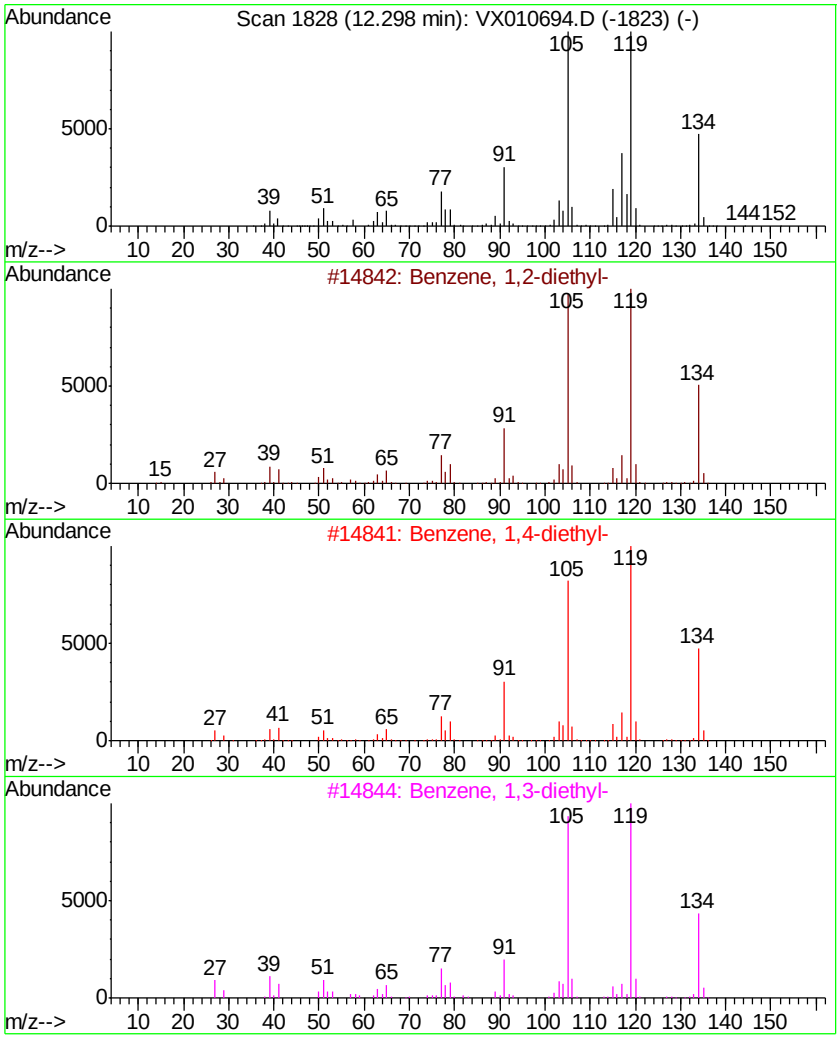
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

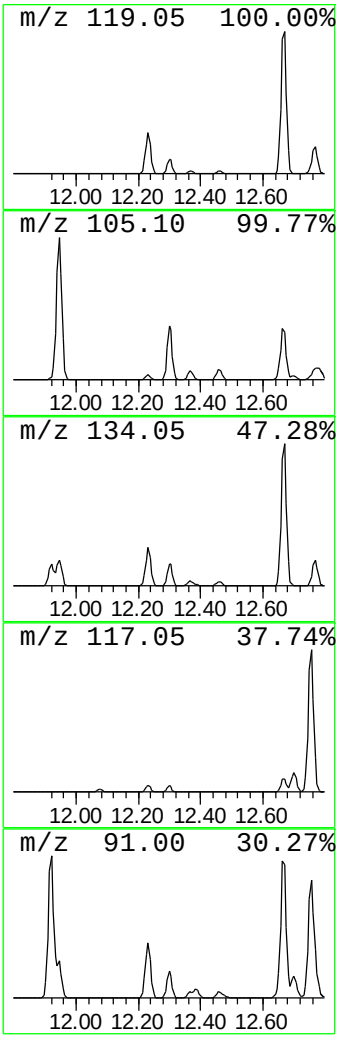
****

Peak Number 2 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	35.19 ug/l	756306	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	83
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

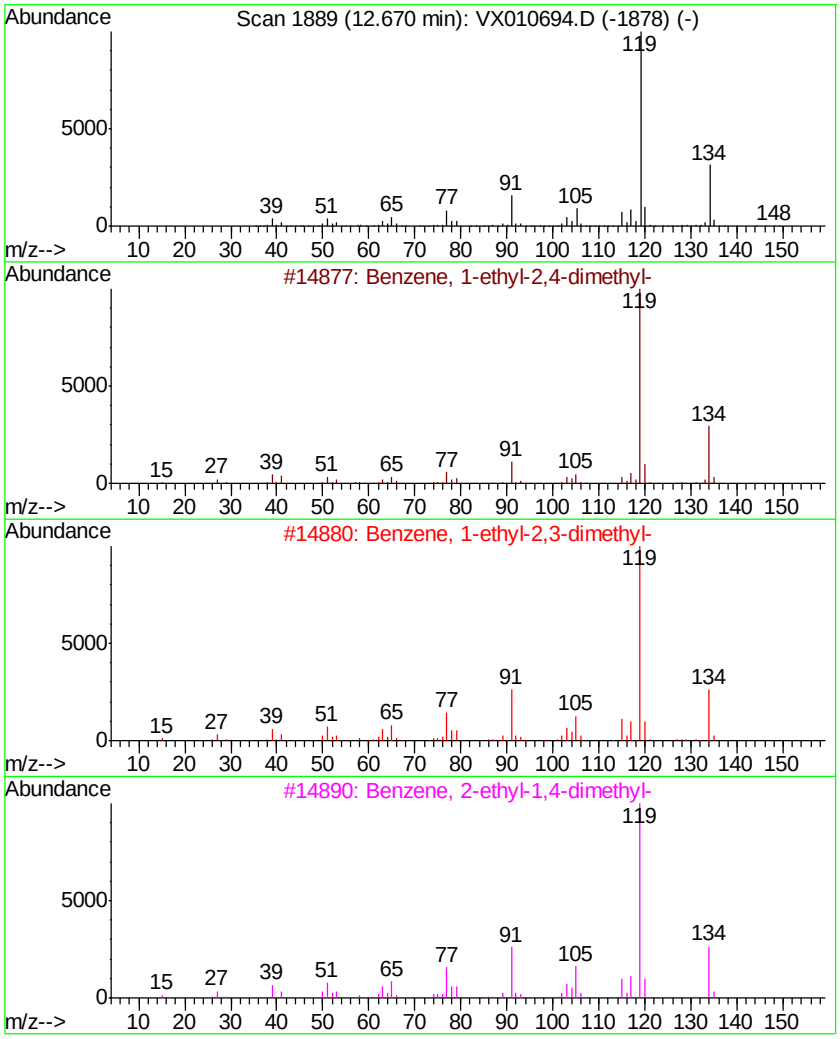
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

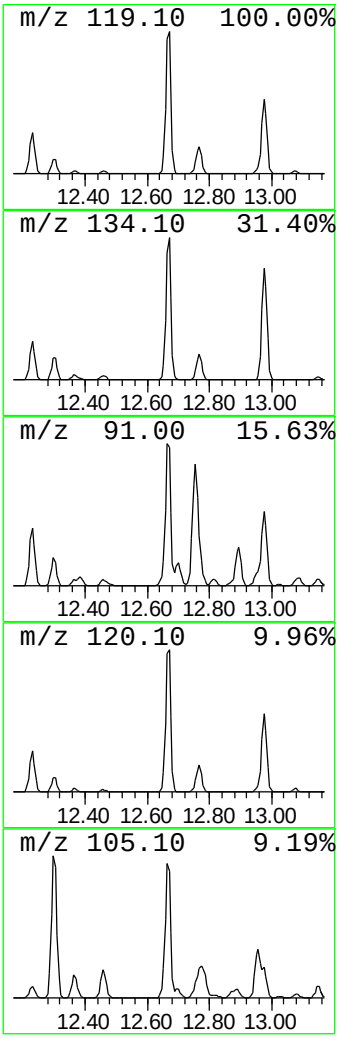
****

Peak Number 3 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	174.60 ug/l	3752130	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, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
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	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

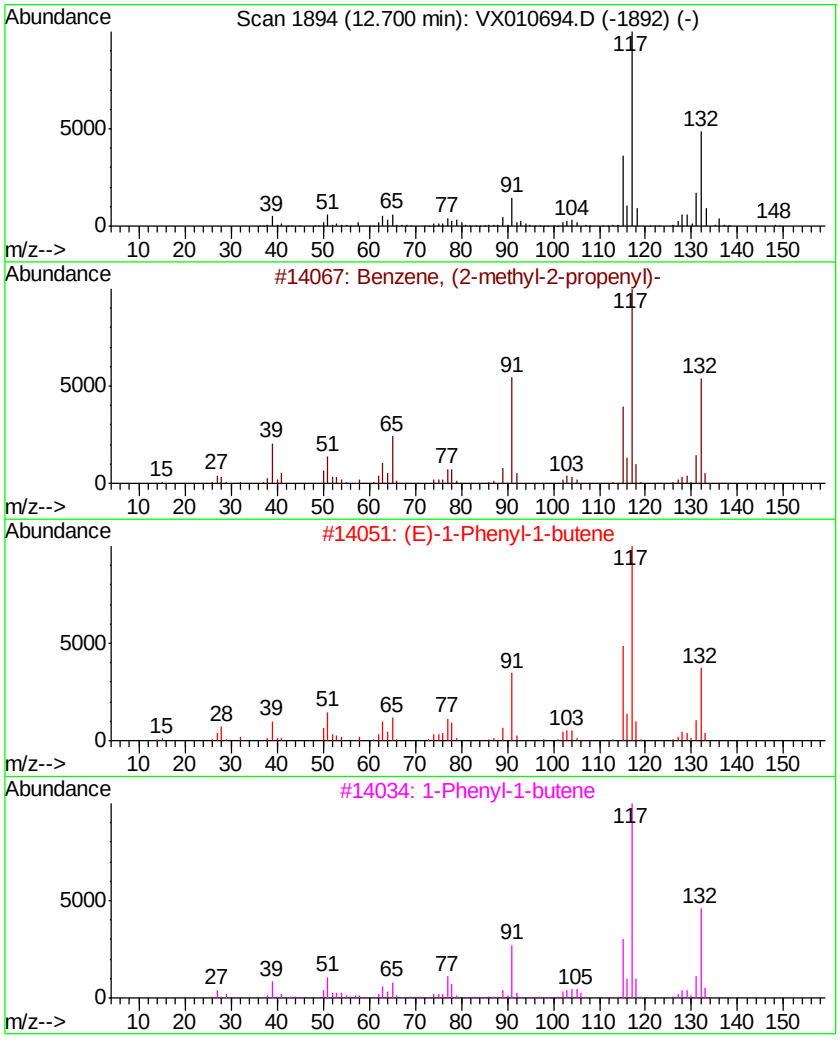
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

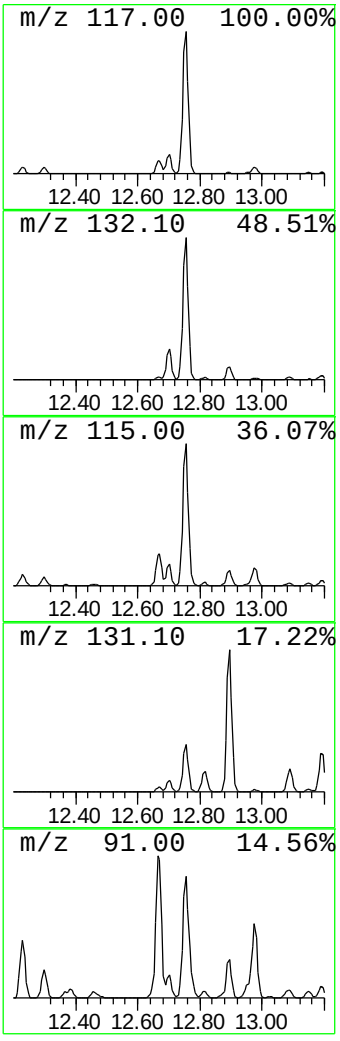
****

Peak Number 4 Benzene, (2-methyl-2-propen... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	33.00 ug/l	709219	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	91
2	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	91
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

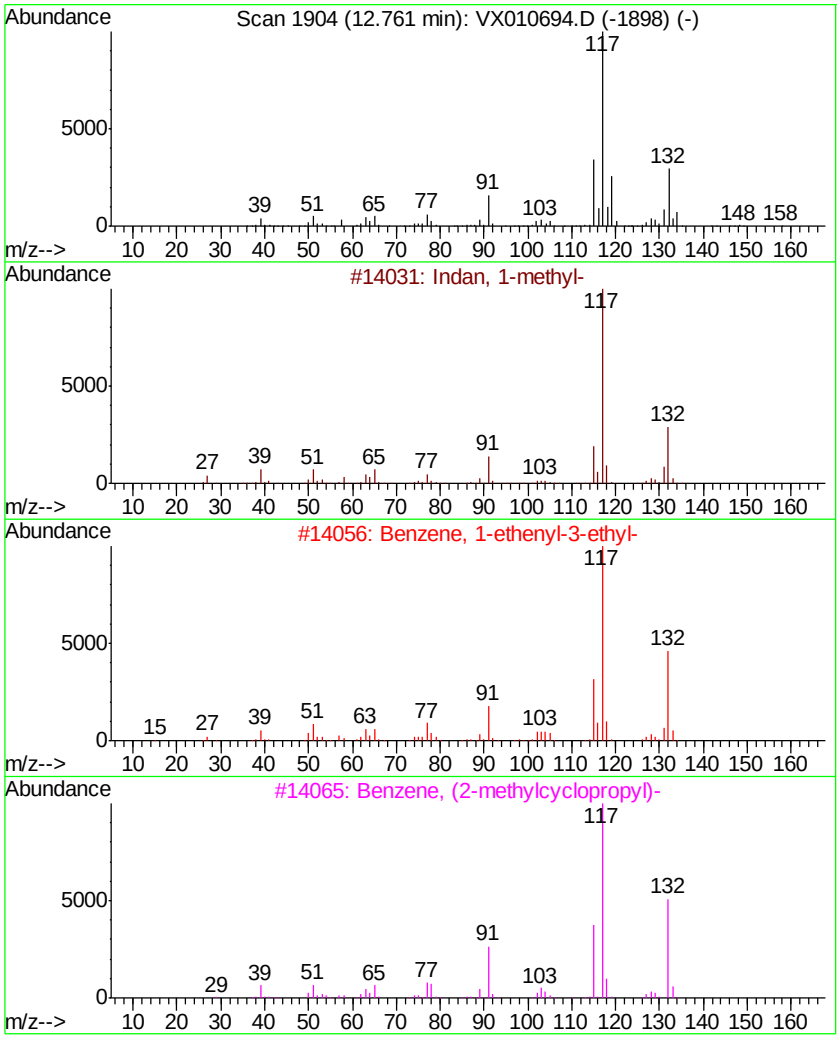
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

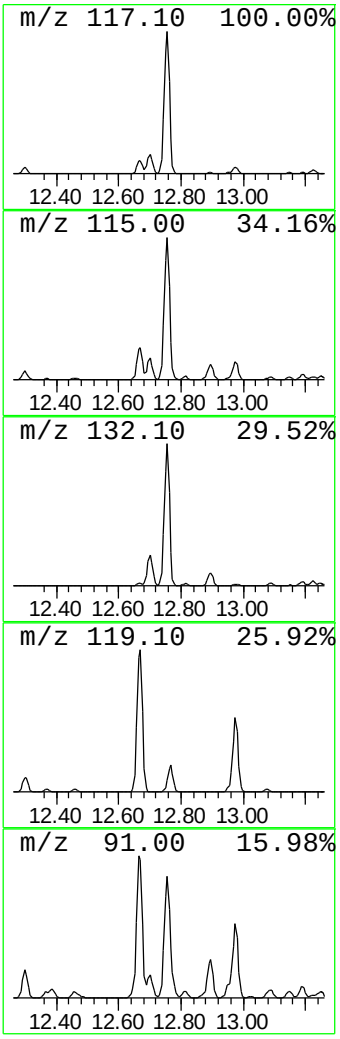
****

Peak Number 5 Indan, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	207.47 ug/l	4458510	1,4-Dichlorobenzene-d4	12.08

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	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

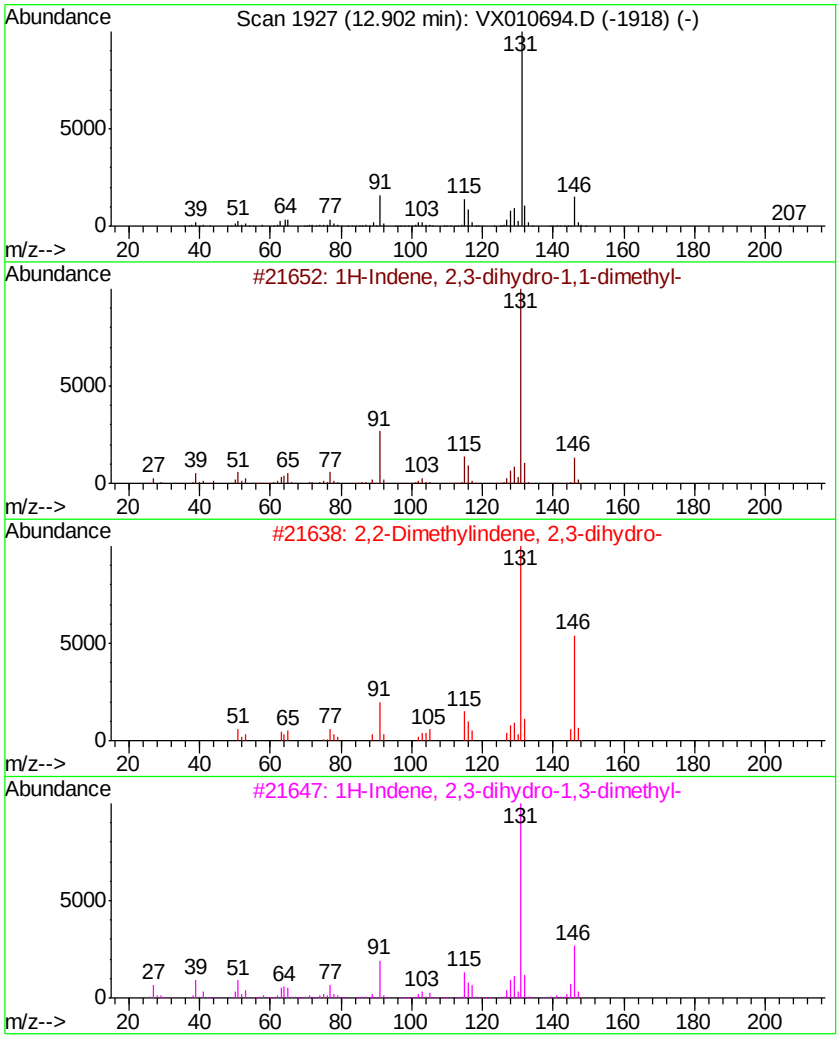
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

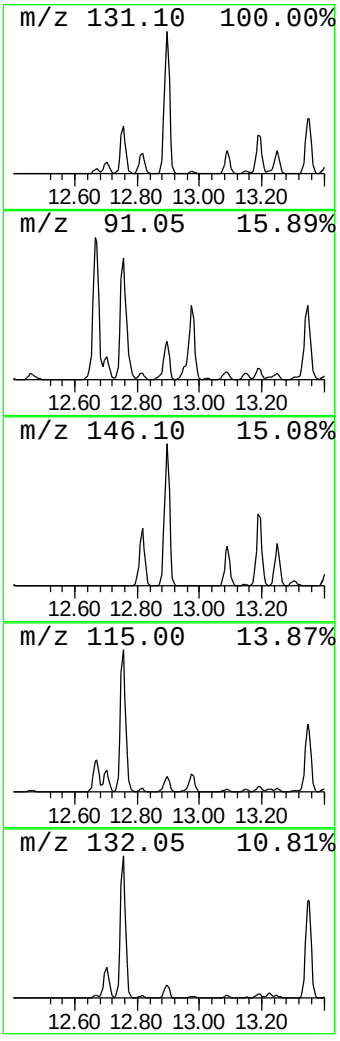
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	44.70 ug/l	960588	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	94
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	90
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

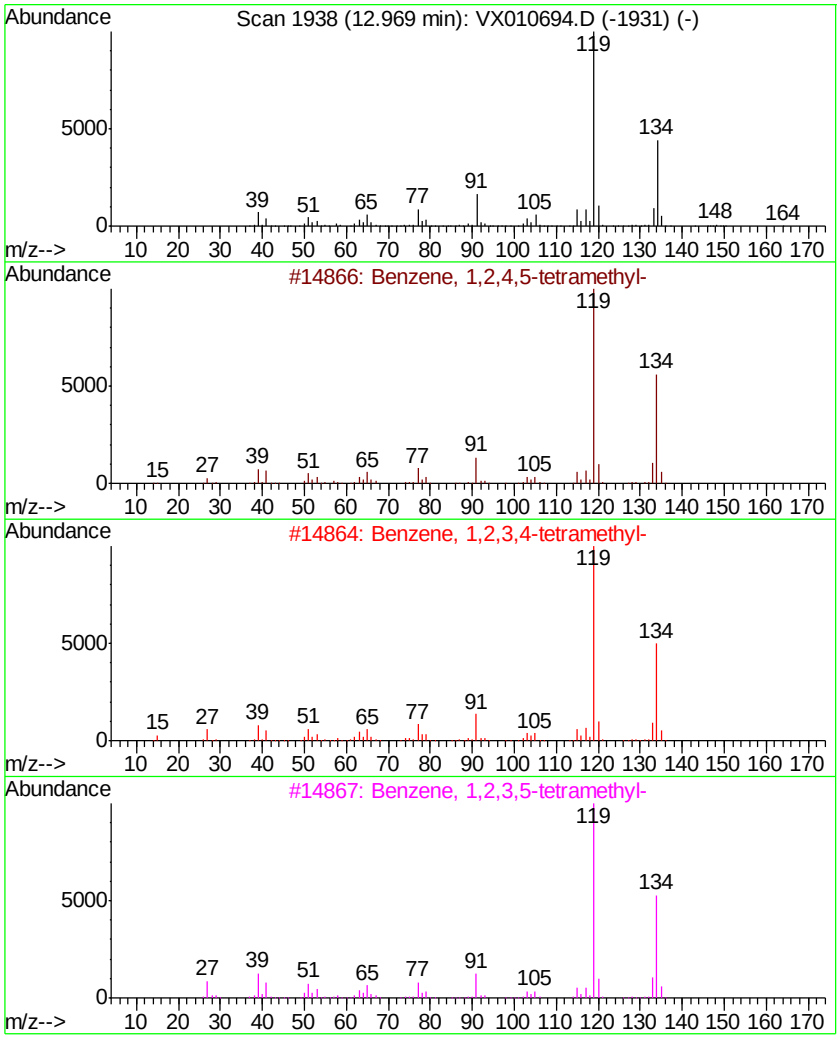
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

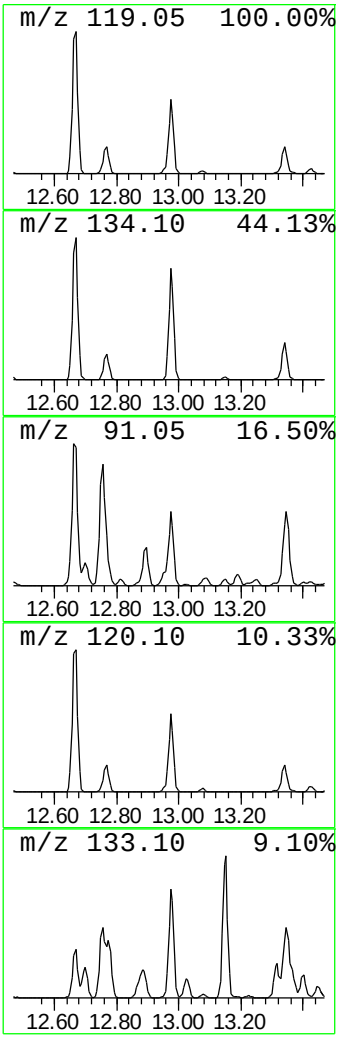
****

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	104.28 ug/l	2241040	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

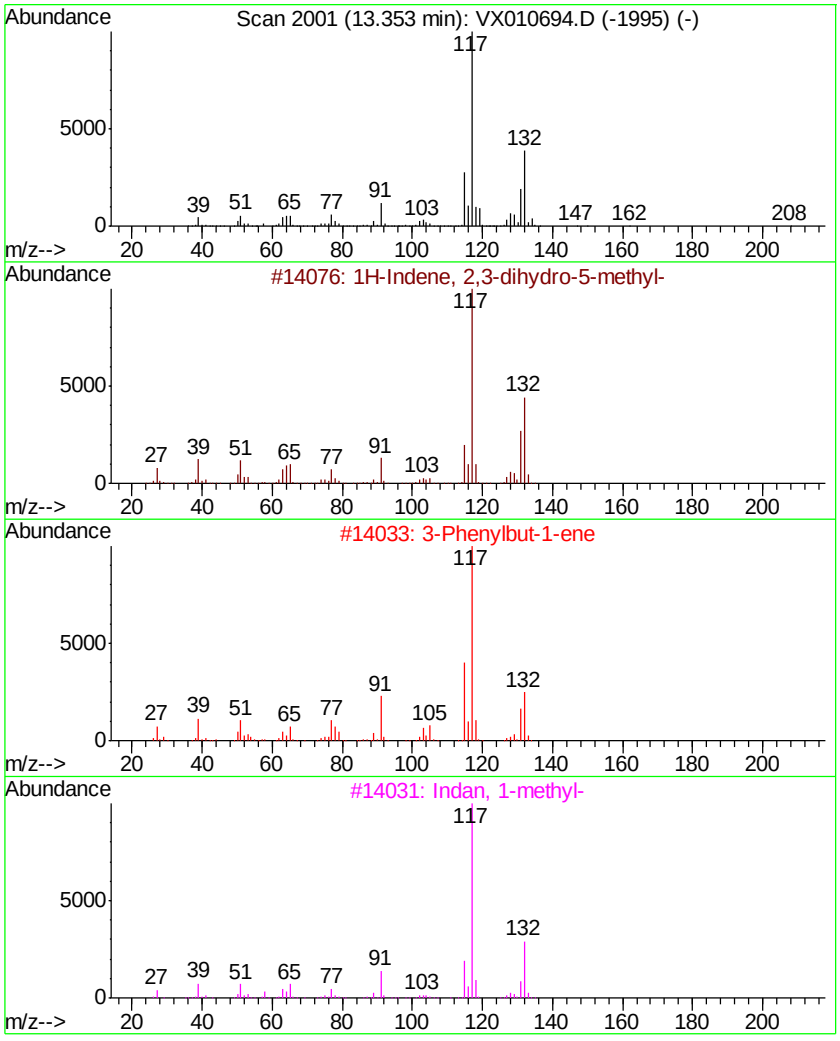
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

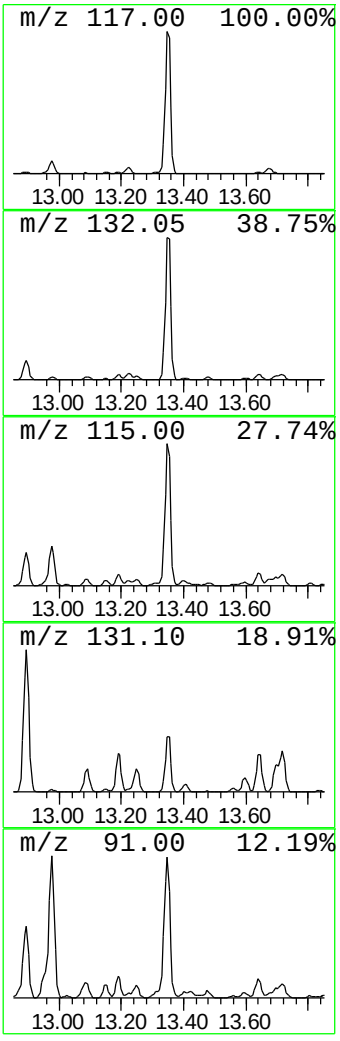
****

Peak Number 8 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	140.84 ug/l	3026720	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	81
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	81
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

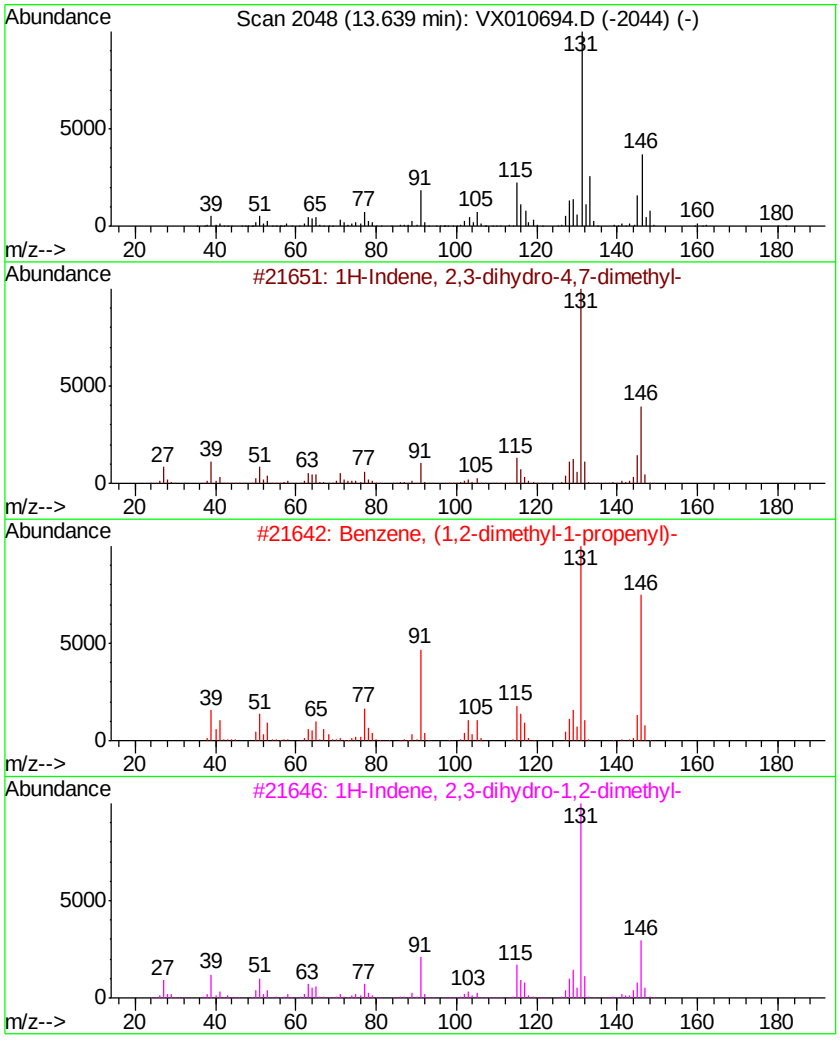
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

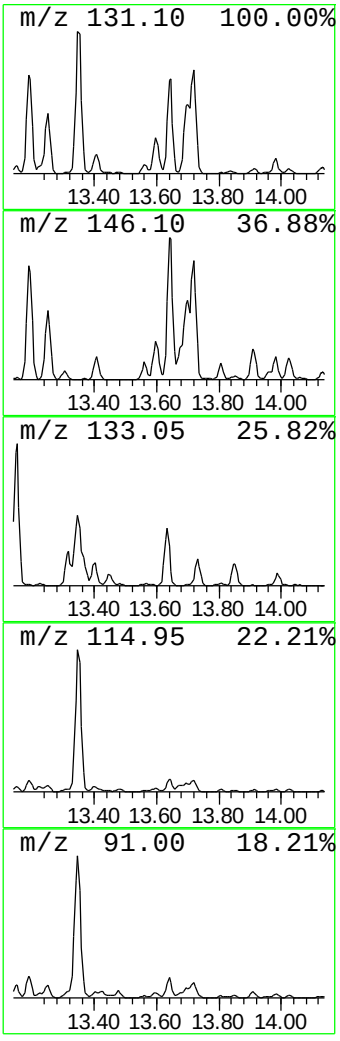
****

Peak Number 9 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	18.93 ug/l	406761	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	95
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	95
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

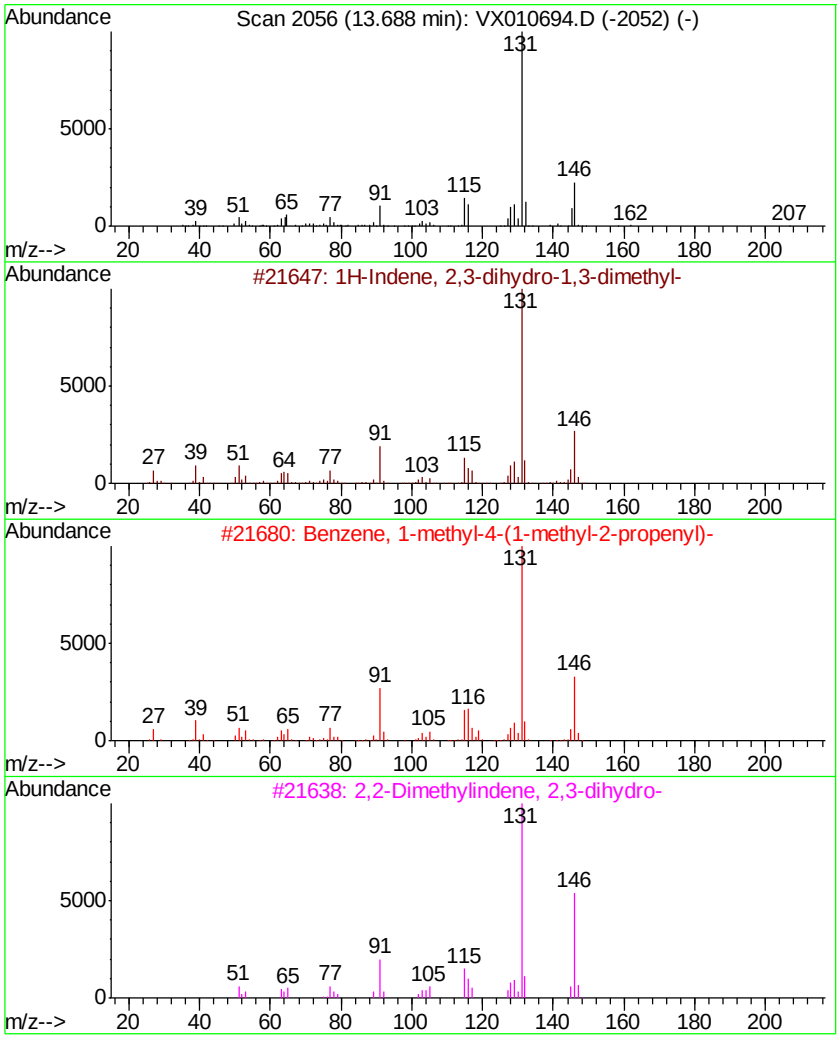
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

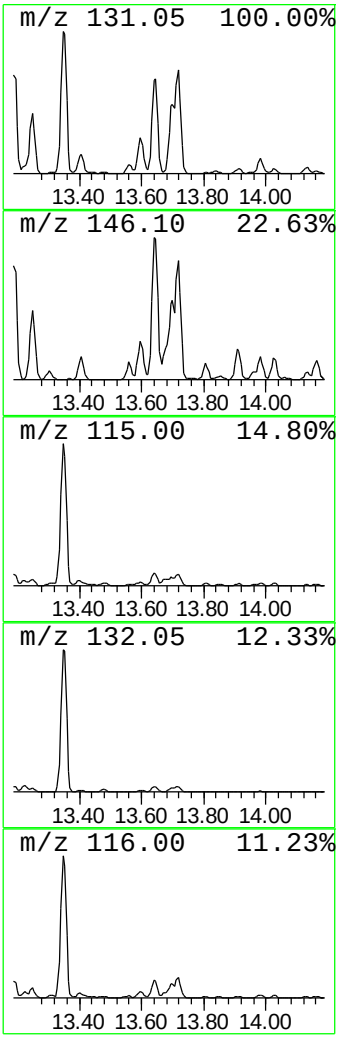
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	14.01 ug/l	301010	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
2	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	91
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX070819\
Data File : VX010694.D
Acq On : 08 Jul 2019 14:39
Operator : JC/SP
Sample : K3669-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : ST008MW037-190701

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.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
p-Cymene	12.23	48.7	ug/l	1046940	4	12.08	1074500	50.0
Benzene, 1,2-diet...	12.30	35.2	ug/l	756306	4	12.08	1074500	50.0
Benzene, 1-ethyl-...	12.67	174.6	ug/l	3752130	4	12.08	1074500	50.0
Benzene, (2-methy...	12.70	33.0	ug/l	709219	4	12.08	1074500	50.0
Indan, 1-methyl-	12.76	207.5	ug/l	4458510	4	12.08	1074500	50.0
1H-Indene, 2,3-di...	12.90	44.7	ug/l	960588	4	12.08	1074500	50.0
Benzene, 1,2,4,5-...	12.97	104.3	ug/l	2241040	4	12.08	1074500	50.0
1H-Indene, 2,3-di...	13.35	140.8	ug/l	3026720	4	12.08	1074500	50.0
1H-Indene, 2,3-di...	13.64	18.9	ug/l	406761	4	12.08	1074500	50.0
1H-Indene, 2,3-di...	13.69	14.0	ug/l	301010	4	12.08	1074500	50.0