LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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	3.190	322	334	346	rBV	69915	188322	1.95%	0.352%
2	4.403	523	533	544	rVB2	42257	124522	1.29%	0.233%
3	5.501	701	713	721	rBV2	150803	438972	4.56%	0.820%
4	5.665	731	740	757	rVB	254476	737498	7.65%	1.378%
5	6.068	796	806	817	rBV	140948	375204	3.89%	0.701%
6	6.860	925	936	946	rBV	379898	993586	10.31%	1.856%
7	7.256	991	1001	1010	rBV2	57235	130622	1.36%	0.244%
8	7.464	1022	1035	1049	rVB2	95893	244208	2.53%	0.456%
9	8.213	1145	1158	1163	rBV	162646	298578	3.10%	0.558%
10	8.573	1210	1217	1226	rVB	135963	223983	2.32%	0.418%
11	8.713	1234	1240	1247	rVB	804674	1230080	12.76%	2.298%
12	9.219	1318	1323	1333	rVB2	72431	121876	1.26%	0.228%
13	10.109	1464	1469	1474	rBV	758898	1014819	10.53%	1.896%
14	11.018	1613	1618	1626	rBV	997992	1245542	12.92%	2.327%
15	11.134	1631	1637	1643	rBV	640972	827169	8.58%	1.545%
16	11.359	1668	1674	1680	rBV	2977666	3642495	37.80%	6.805%
17	11.664	1715	1724	1733	rBV	134015	202227	2.10%	0.378%
18	11.920	1760	1766	1767	rBV	118702	143666	1.49%	0.268%
19	11.945	1767	1770	1777	rVB	223251	272648	2.83%	0.509%
20	12.079	1786	1792	1798	rBV	739439	973033	10.10%	1.818%
21	12.231	1812	1817	1822	rBV	92767	112864	1.17%	0.211%
22	12.298	1822	1828	1833	rVV	8097104	9636738	100.00%	18.003%
23	12.365	1835	1839	1849	rVV2	395430	788906	8.19%	1.474%
24	12.457	1849	1854	1859	rVV	296627	369926	3.84%	0.691%
25	12.670	1881	1889	1892	rBV	2914999	3643395	37.81%	6.806%
26	12.700	1892	1894	1898	rVV	771865	782593	8.12%	1.462%
27	12.755	1898	1903	1910	rVV2	2432483	3330789	34.56%	6.222%
28	12.896	1921	1926	1931	rVB	130477	180482	1.87%	0.337%
29	12.975	1931	1939	1944	rBV	1929514	2329988	24.18%	4.353%
30	13.078	1951	1956	1962	rBV3	133902	199662	2.07%	0.373%
31	13.194	1971	1975	1976	rVV	152391	189688	1.97%	0.354%
32	13.225	1976	1980	1990	rVB	1934599	2369870	24.59%	4.427%
33	13.347	1995	2000	2006	rVV2	4765997	6664900	69.16%	12.451%
34	13.402	2006	2009	2011	rVV2	130380	188831	1.96%	0.353%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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	13.426	2011	2013	2018	rVV	121546	145963	1.51%	0.273%
36	13.481	2018	2022	2028	rVB	288903	371190	3.85%	0.693%
37	13.560	2028	2035	2038	rBV2	159833	231814	2.41%	0.433%
38	13.597	2038	2041	2044	rVV	336917	463709	4.81%	0.866%
39	13.639	2044	2048	2052	rVV3	443380	701580	7.28%	1.311%
40	13.694	2052	2057	2059	rVV3	253067	442356	4.59%	0.826%
41	13.719	2059	2061	2071	rVB3	446497	698089	7.24%	1.304%
42	13.853	2079	2083	2088	rVB	101020	136814	1.42%	0.256%
43	13.908	2088	2092	2097	rBV	151218	189176	1.96%	0.353%
44	13.981	2097	2104	2108	rBV2	303578	438755	4.55%	0.820%
45	14.023	2108	2111	2116	rVB	155282	195609	2.03%	0.365%
46	14.133	2124	2129	2132	rBV	258787	333791	3.46%	0.624%
47	14.286	2146	2154	2161	rBV2	409544	833015	8.64%	1.556%
48	14.377	2161	2169	2173	rBV	150258	214597	2.23%	0.401%
49	14.432	2173	2178	2183	rVB2	327262	474288	4.92%	0.886%
50	14.536	2190	2195	2200	rBV2	378184	500745	5.20%	0.935%
51	14.657	2209	2215	2223	rBV3	83689	199745	2.07%	0.373%
52	14.846	2240	2246	2253	rBV3	106682	231507	2.40%	0.432%
53	15.249	2305	2312	2318	rBV2	69247	119317	1.24%	0.223%
54	15.444	2331	2344	2347	rBV	142061	237135	2.46%	0.443%
55	15.548	2354	2361	2366	rBV	402158	640333	6.64%	1.196%
56	15.688	2374	2384	2388	rBV	551686	931680	9.67%	1.740%
57	15.913	2410	2421	2437	rVB2	150838	429735	4.46%	0.803%
58	16.066	2440	2446	2459	rVB	80144	151247	1.57%	0.283%

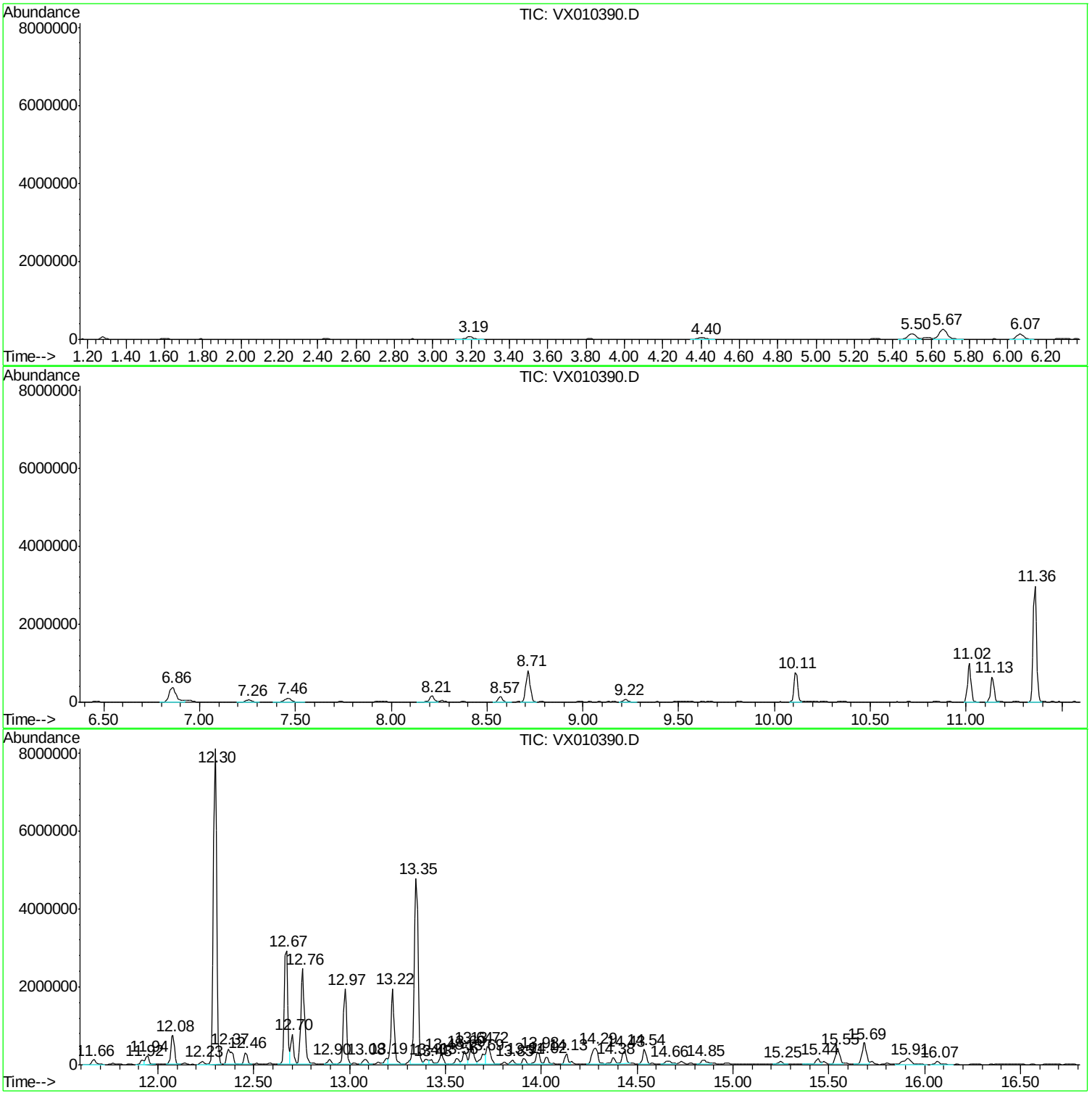
Sum of corrected areas: 53529872

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
T11-MW-5-9.0-062019

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

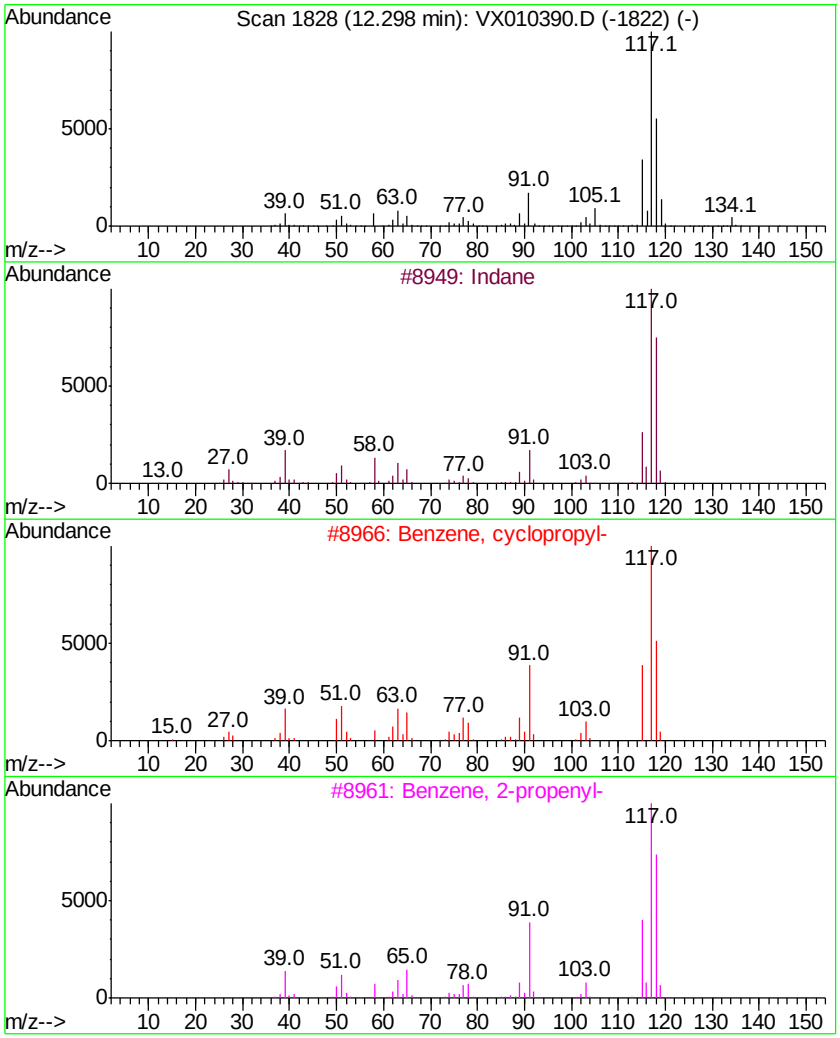
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

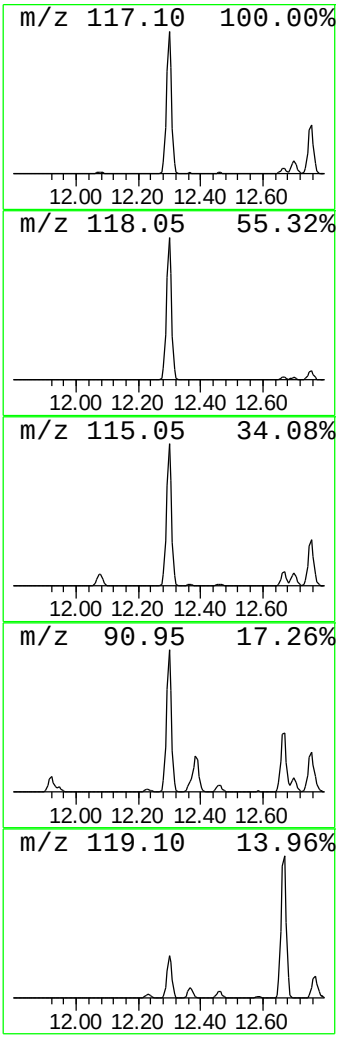
****

Peak Number 1 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	495.19 ug/l	9636740	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
4	1,4:5,8-Dimethanobiphenylene, 1,...	184	C14H16	001624-13-1	64
5	Deltacyclene	118	C9H10	007785-10-6	58





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

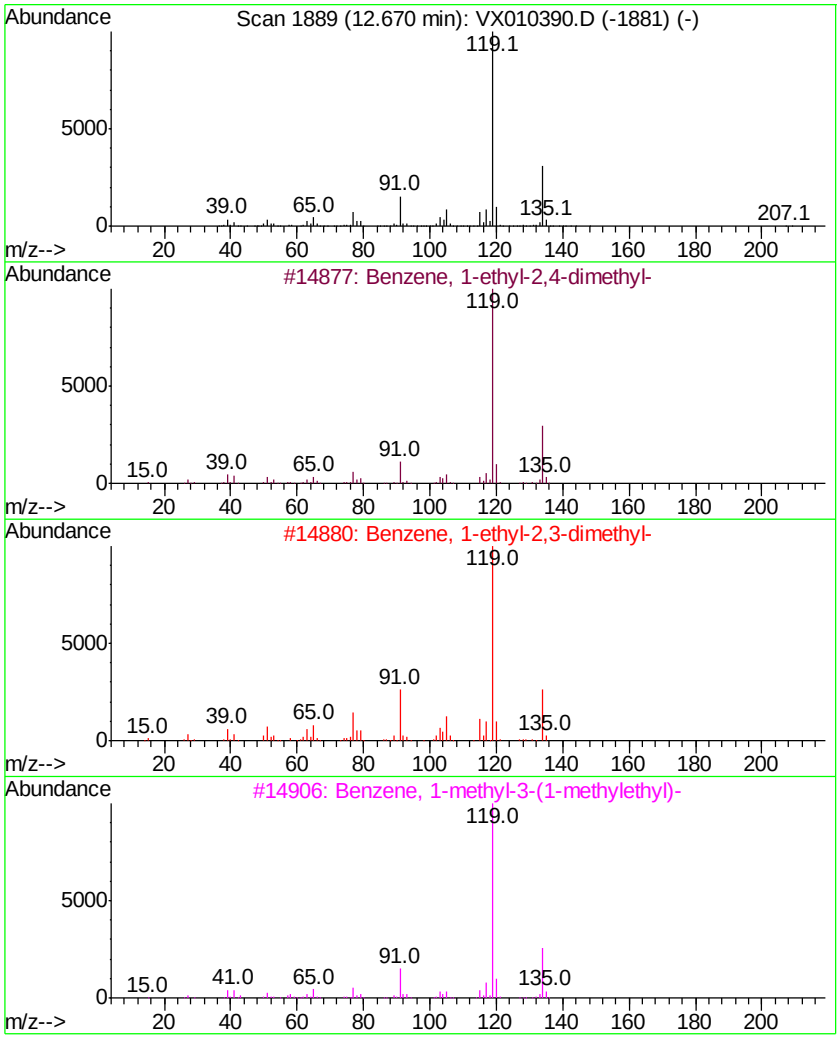
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Quant Title : SW846 8260

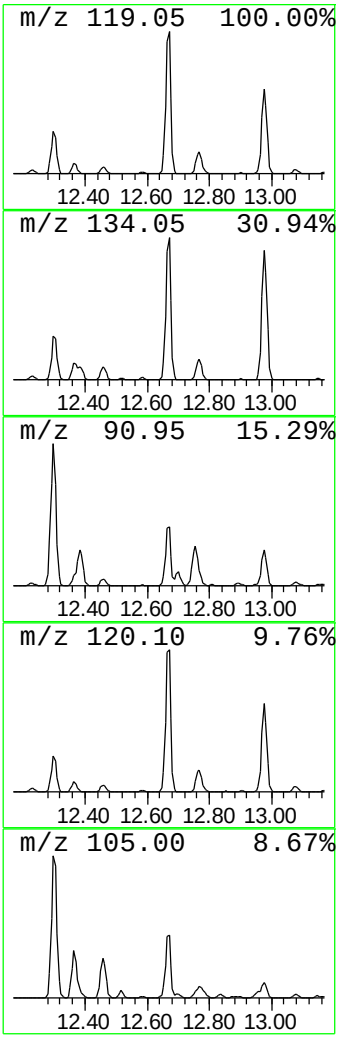
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	187.22 ug/l	3643400	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, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

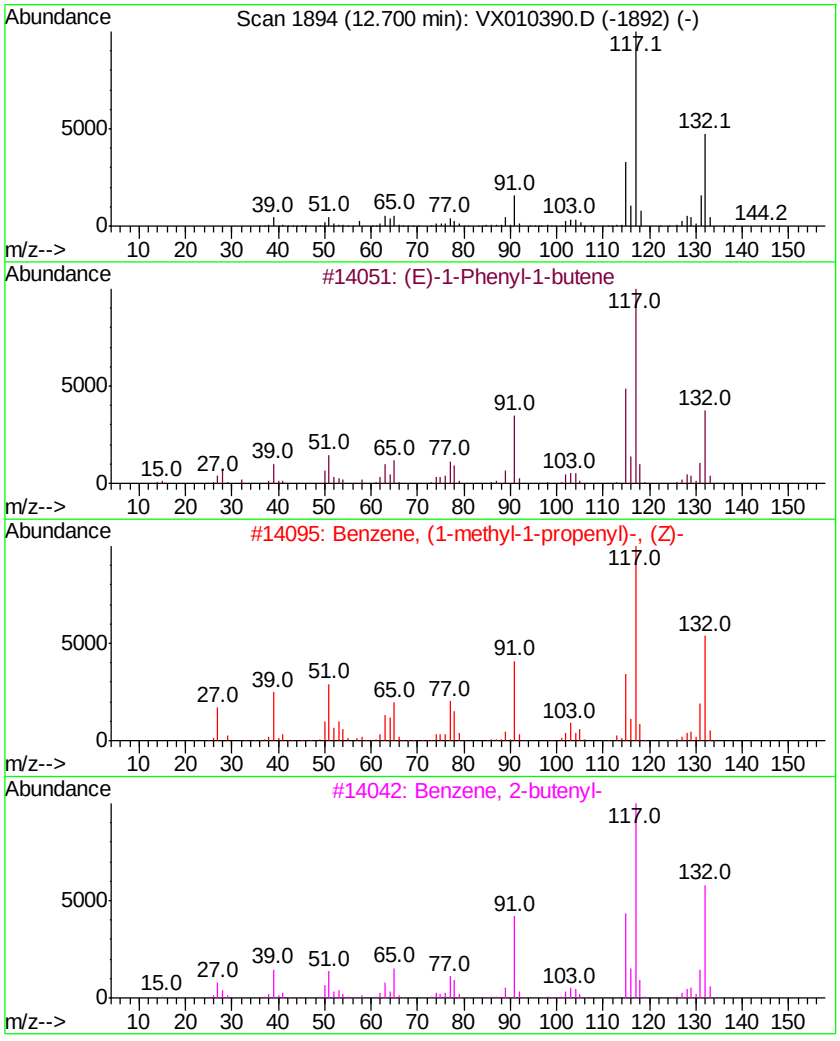
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

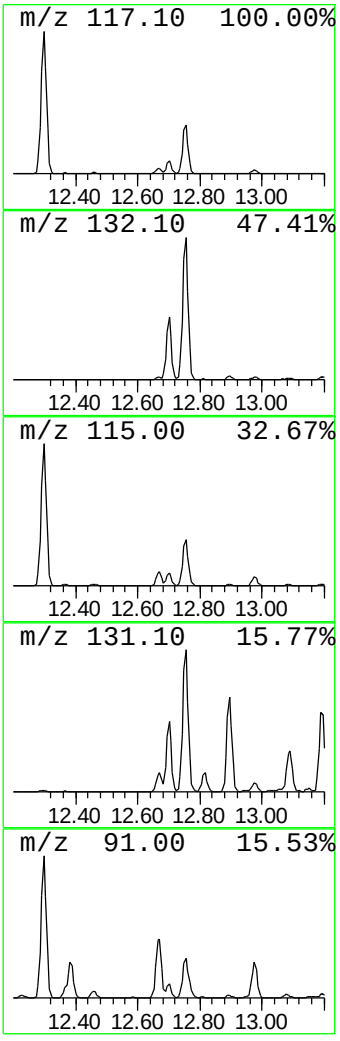
****

Peak Number 3 (E)-1-Phenyl-1-butene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	40.21 ug/l	782593	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	94
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
4	Benzene, (1-methylenepropyl)-	132	C10H12	002039-93-2	91
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

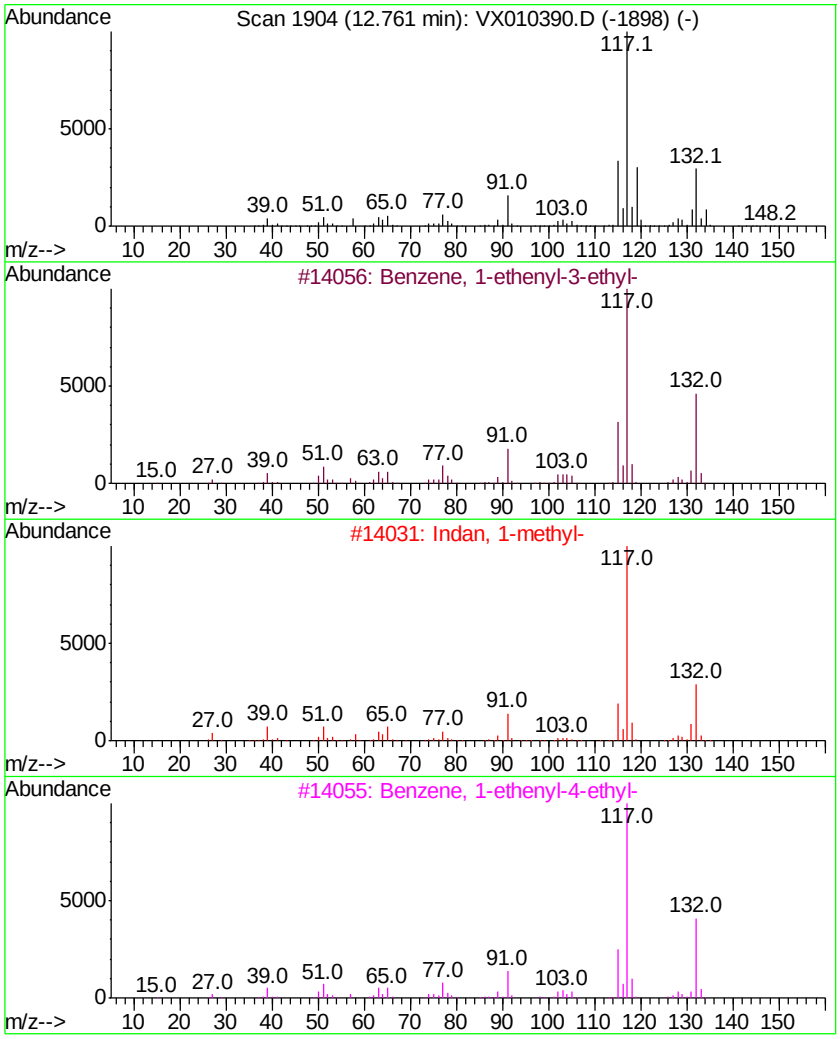
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

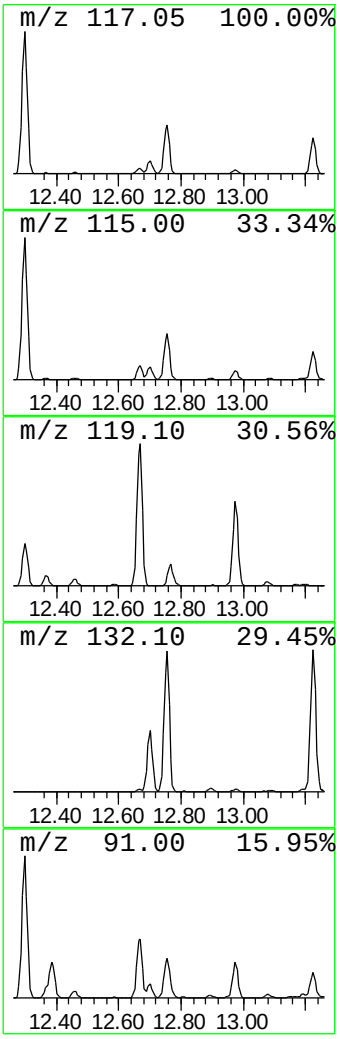
****

Peak Number 4 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	171.16 ug/l	3330790	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	74





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

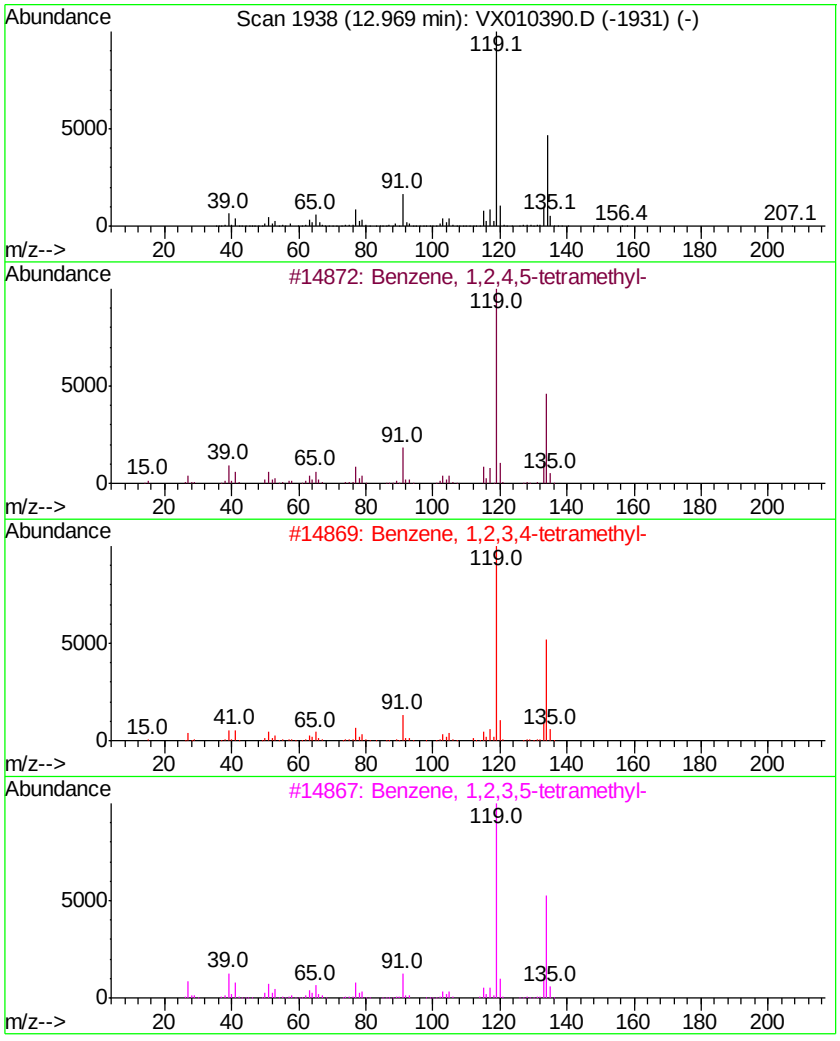
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

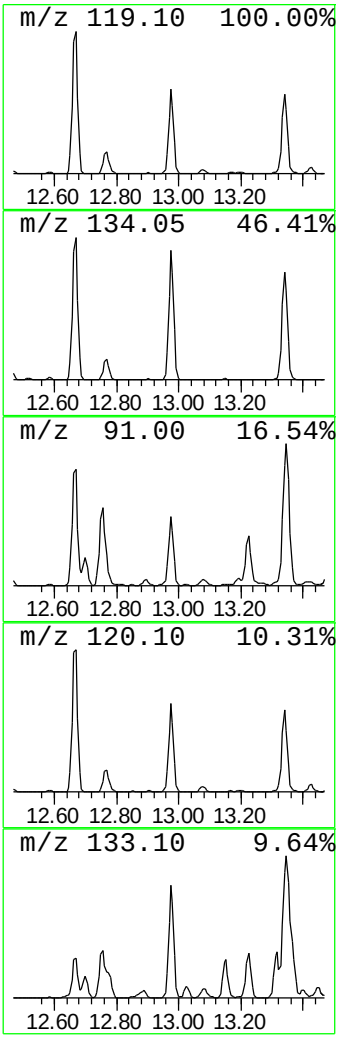
****

Peak Number 5 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	119.73 ug/l	2329990	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	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

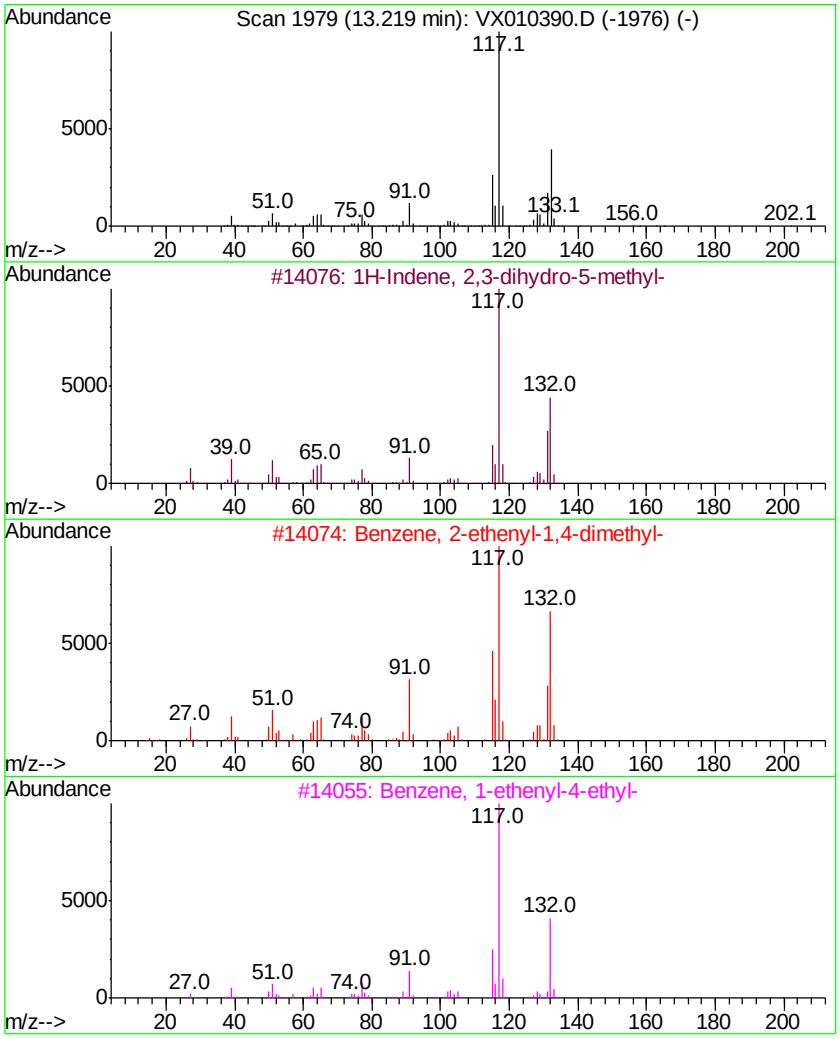
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

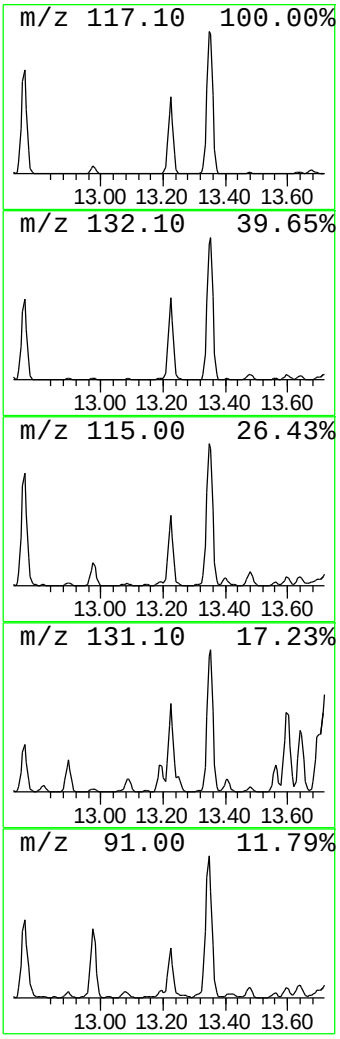
****

Peak Number 6 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	121.78 ug/l	2369870	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	93
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

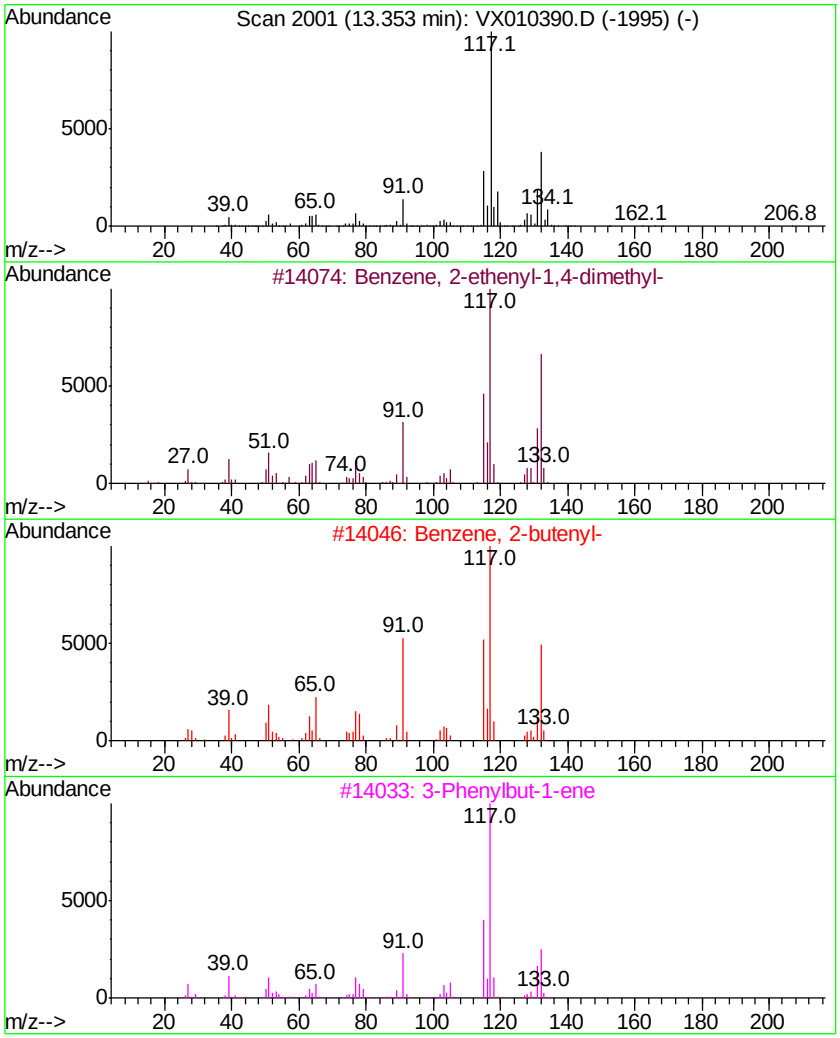
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

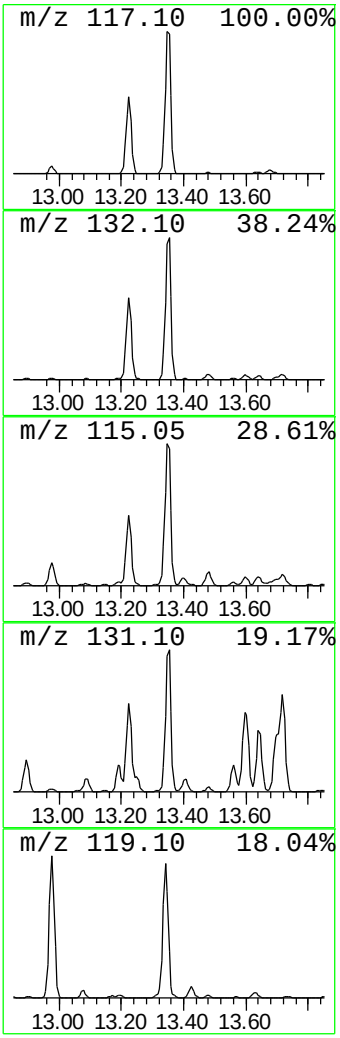
****

Peak Number 7 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	342.48 ug/l	6664900	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-butenyl-	132	C10H12	001560-06-1	70
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

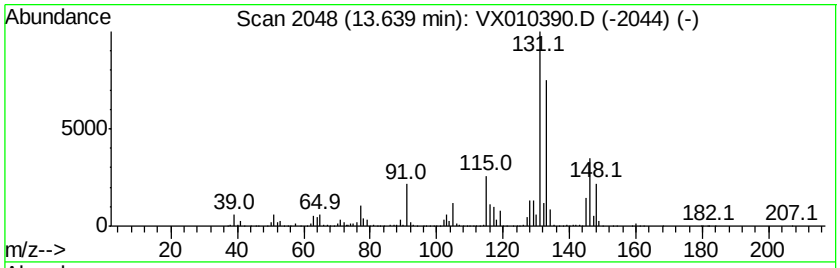
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

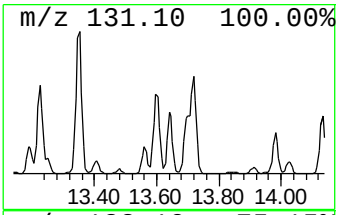
****

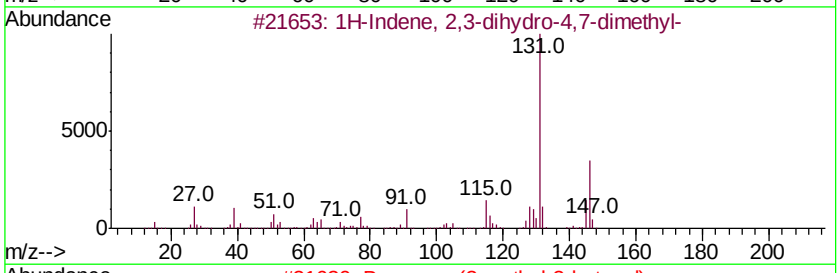
Peak Number 8 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

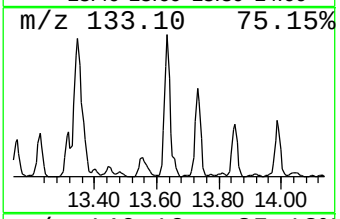
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	36.05 ug/l	701580	1,4-Dichlorobenzene-d4	12.08

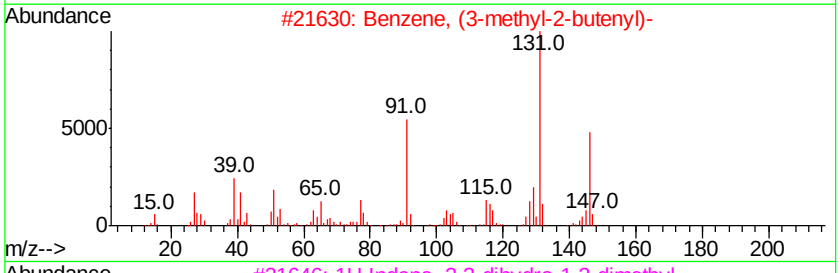
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87
2	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	78
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	78
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	60
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	60

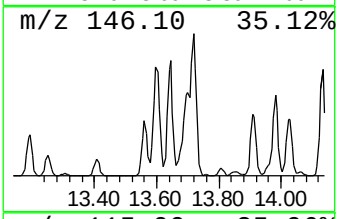


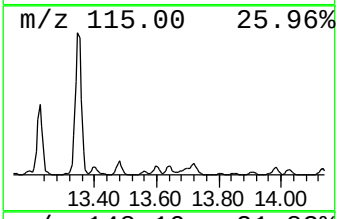


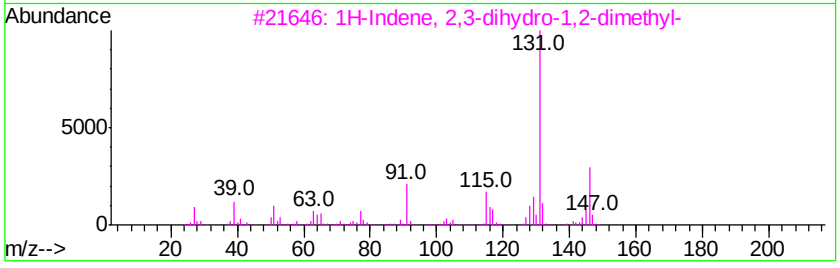


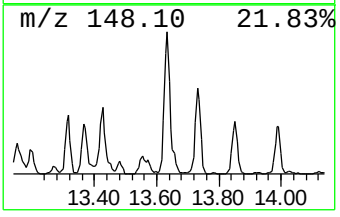












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Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

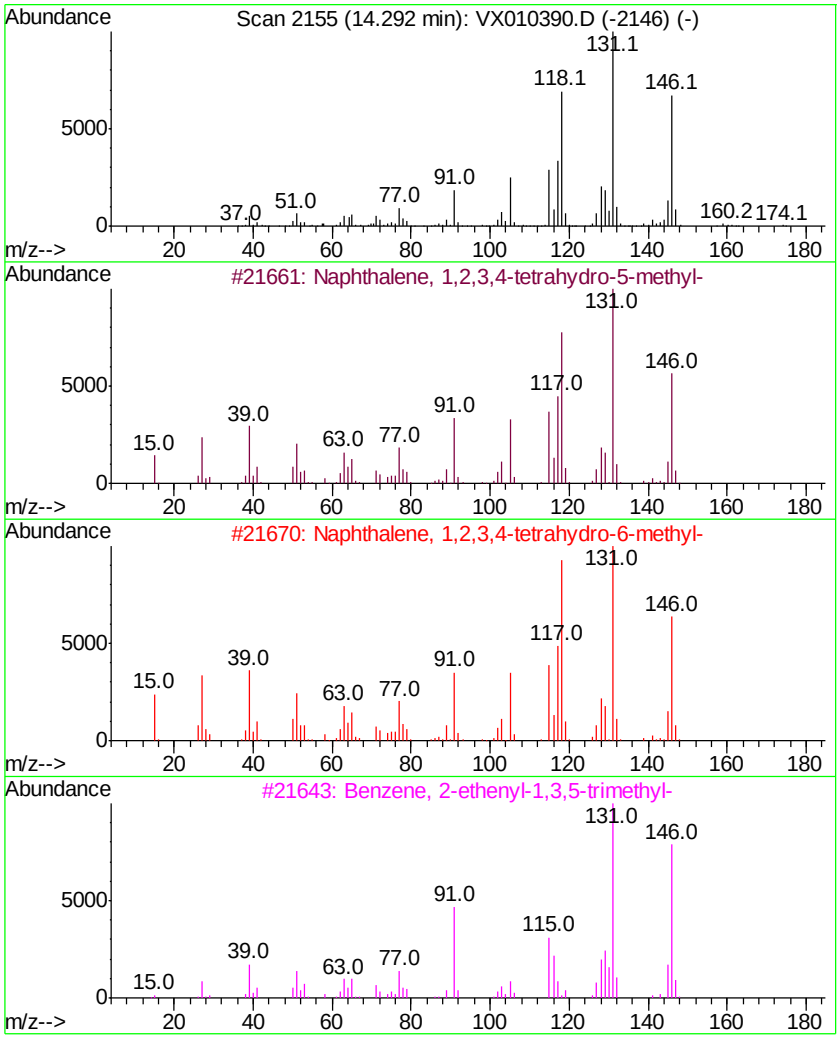
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

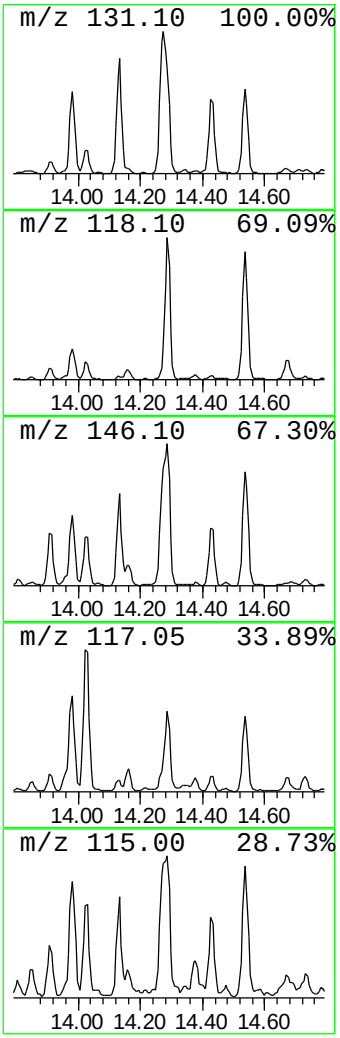
****

Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	42.81 ug/l	833015	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	97
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	93
3	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	89
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	70





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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

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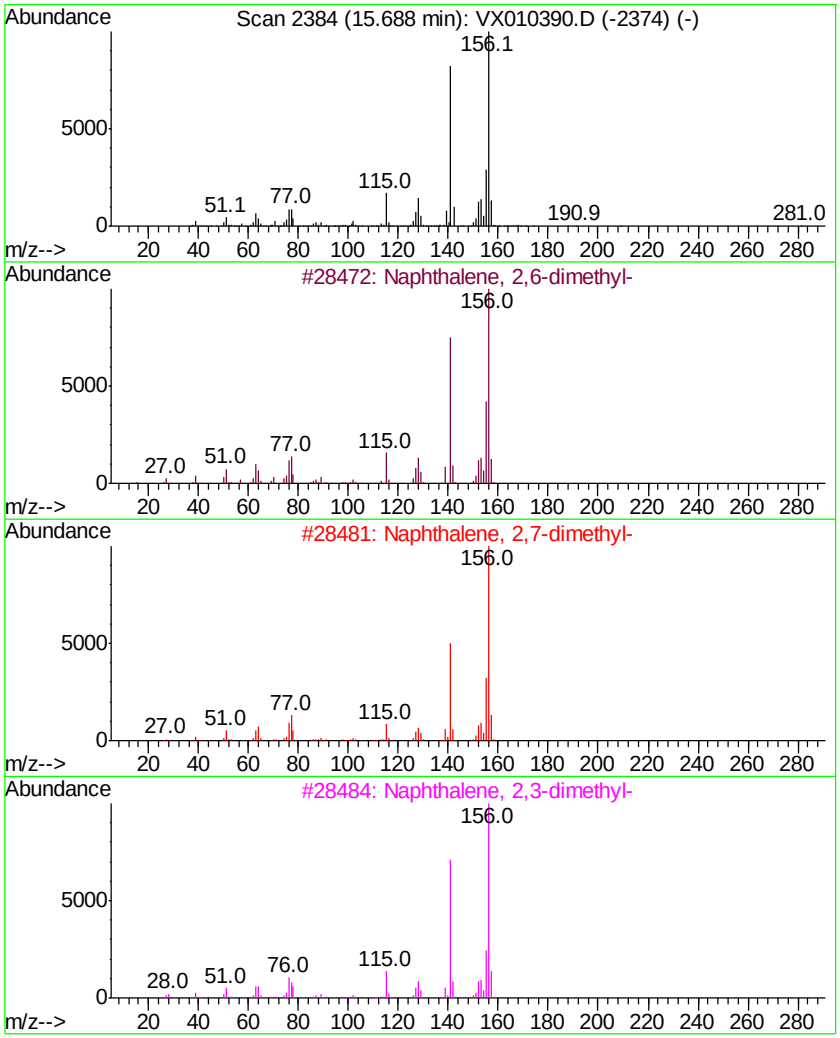
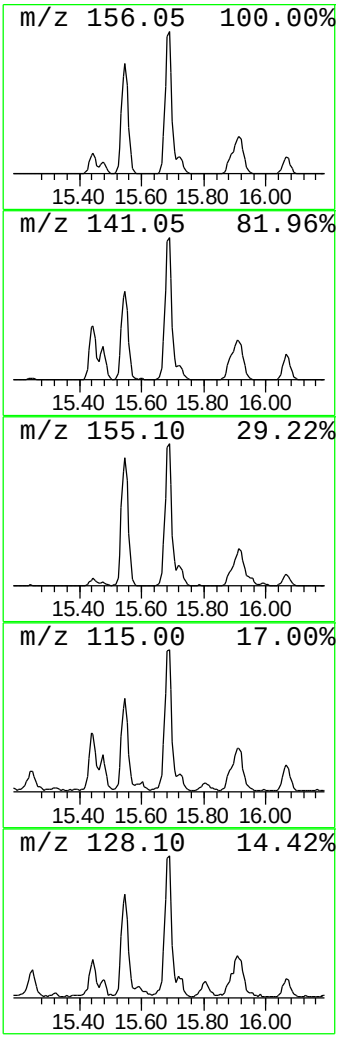
Peak Number 10 Naphthalene, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	47.88 ug/l	931680	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062119\
Data File : VX010390.D
Acq On : 21 Jun 2019 17:22
Operator : JC/SP
Sample : K3469-19
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T11-MW-5-9.0-062019

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
Indane	12.30	495.2	ug/l	9636740	4	12.08	973033	50.0
Benzene, 1-ethyl-...	12.67	187.2	ug/l	3643400	4	12.08	973033	50.0
(E)-1-Phenyl-1-bu...	12.70	40.2	ug/l	782593	4	12.08	973033	50.0
Benzene, 1-etheny...	12.76	171.2	ug/l	3330790	4	12.08	973033	50.0
Benzene, 1,2,4,5-...	12.97	119.7	ug/l	2329990	4	12.08	973033	50.0
1H-Indene, 2,3-di...	13.22	121.8	ug/l	2369870	4	12.08	973033	50.0
Benzene, 2-etheny...	13.35	342.5	ug/l	6664900	4	12.08	973033	50.0
1H-Indene, 2,3-di...	13.64	36.0	ug/l	701580	4	12.08	973033	50.0
Naphthalene, 1,2,...	14.29	42.8	ug/l	833015	4	12.08	973033	50.0
Naphthalene, 2,6-...	15.69	47.9	ug/l	931680	4	12.08	973033	50.0