LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

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\82X071619W.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	5.501	701	713	726	rBV	112747	326494	3.67%	0.591%
2	5.659	729	739	755	rVB	223533	629392	7.07%	1.139%
3	6.062	795	805	818	rBV	119514	308329	3.46%	0.558%
4	6.860	926	936	950	rBV	306240	718061	8.06%	1.299%
5	8.713	1234	1240	1249	rVB	628128	954759	10.72%	1.727%
6	10.110	1464	1469	1476	rBV	609601	822972	9.24%	1.489%
7	11.018	1612	1618	1622	rBV	219785	282320	3.17%	0.511%
8	11.134	1632	1637	1642	rBV	536624	670438	7.53%	1.213%
9	11.359	1668	1674	1679	rBV	419823	535419	6.01%	0.969%
10	11.804	1744	1747	1757	rVB5	45083	108650	1.22%	0.197%
11	11.945	1767	1770	1775	rVB	215426	257808	2.90%	0.466%
12	12.073	1787	1791	1797	rVB	675534	861424	9.68%	1.558%
13	12.298	1823	1828	1834	rVB	505678	620533	6.97%	1.123%
14	12.365	1835	1839	1841	rBV	196786	280337	3.15%	0.507%
15	12.384	1841	1842	1846	rVB	216595	173785	1.95%	0.314%
16	12.457	1846	1854	1861	rBV2	149051	255571	2.87%	0.462%
17	12.518	1861	1864	1870	rVB4	47358	91609	1.03%	0.166%
18	12.609	1870	1879	1881	rBV5	47291	111781	1.26%	0.202%
19	12.664	1884	1888	1892	rVV	776925	971405	10.91%	1.757%
20	12.701	1892	1894	1897	rVV	188777	185323	2.08%	0.335%
21	12.755	1897	1903	1910	rVV2	848198	1292091	14.51%	2.337%
22	12.871	1918	1922	1924	rBV2	95489	131228	1.47%	0.237%
23	12.944	1929	1934	1935	rBV3	98021	130491	1.47%	0.236%
24	12.975	1935	1939	1944	rVB	707334	849263	9.54%	1.536%
25	13.042	1944	1950	1952	rBV3	77688	107717	1.21%	0.195%
26	13.079	1952	1956	1962	rVB2	169329	264260	2.97%	0.478%
27	13.146	1963	1967	1971	rBV2	108391	180897	2.03%	0.327%
28	13.194	1971	1975	1977	rBV2	166234	198617	2.23%	0.359%
29	13.225	1977	1980	1983	rVB	386339	420391	4.72%	0.761%
30	13.316	1990	1995	1996	rBV2	101493	142530	1.60%	0.258%
31	13.347	1996	2000	2006	rVB2	959068	1436296	16.13%	2.598%
32	13.402	2006	2009	2010	rBV2	126904	137139	1.54%	0.248%
33	13.426	2010	2013	2018	rVV2	156701	268830	3.02%	0.486%
34	13.475	2018	2021	2030	rVB8	187383	392275	4.41%	0.710%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

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\82X071619W.M
Title : SW846 8260

35	13.560	2030	2035	2038	rBV	233345	338532	3.80%	0.612%
36	13.597	2038	2041	2044	rVV	398344	505566	5.68%	0.915%
37	13.639	2044	2048	2051	rVV2	397577	544274	6.11%	0.985%
38	13.694	2051	2057	2058	rVV2	222615	355279	3.99%	0.643%
39	13.719	2058	2061	2066	rVV2	451526	703958	7.91%	1.273%
40	13.804	2070	2075	2078	rBV2	101558	165201	1.86%	0.299%
41	13.853	2079	2083	2086	rBV3	157175	202293	2.27%	0.366%
42	13.908	2086	2092	2097	rVB2	697374	1070180	12.02%	1.936%
43	13.981	2097	2104	2108	rBV2	806843	1175676	13.20%	2.127%
44	14.024	2108	2111	2115	rBV	153902	231395	2.60%	0.419%
45	14.127	2124	2128	2132	rBV	621281	852783	9.58%	1.543%
46	14.267	2147	2151	2161	rVB2	625074	1088833	12.23%	1.970%
47	14.377	2165	2169	2174	rVV	389179	502744	5.65%	0.909%
48	14.426	2174	2177	2182	rVB2	560515	739114	8.30%	1.337%
49	14.475	2182	2185	2190	rVB2	129717	152888	1.72%	0.277%
50	14.536	2190	2195	2199	rBV2	775885	1062724	11.94%	1.923%
51	14.578	2199	2202	2205	rVB3	106914	124179	1.39%	0.225%
52	14.670	2212	2217	2223	rBV3	793726	1189025	13.35%	2.151%
53	14.731	2223	2227	2231	rVB2	378207	520163	5.84%	0.941%
54	14.779	2231	2235	2238	rBV2	196021	293211	3.29%	0.530%
55	14.840	2240	2245	2249	rVV	6974851	8903604	100.00%	16.107%
56	14.901	2253	2255	2260	rVV4	271056	380957	4.28%	0.689%
57	14.962	2262	2265	2272	rVV2	170013	388935	4.37%	0.704%
58	15.017	2272	2274	2282	rVB4	119736	251665	2.83%	0.455%
59	15.249	2306	2312	2314	rBV	507676	797181	8.95%	1.442%
60	15.273	2314	2316	2320	rVV	390159	440638	4.95%	0.797%
61	15.328	2320	2325	2331	rVB7	163315	377477	4.24%	0.683%
62	15.444	2339	2344	2346	rBV	268106	392280	4.41%	0.710%
63	15.474	2346	2349	2355	rVB	463519	679691	7.63%	1.230%
64	15.548	2355	2361	2366	rBV	1661578	2661571	29.89%	4.815%
65	15.688	2374	2384	2387	rBV	2765149	4626939	51.97%	8.370%
66	15.724	2387	2390	2398	rVB	1825473	2604274	29.25%	4.711%
67	15.810	2399	2404	2410	rBV4	134613	276835	3.11%	0.501%
68	15.913	2411	2421	2431	rVB	906294	2328727	26.15%	4.213%
69	16.066	2440	2446	2457	rVB	660240	1235104	13.87%	2.234%
70	16.163	2458	2462	2467	rBV5	57908	101382	1.14%	0.183%
71	16.273	2478	2480	2487	rVB	111964	174472	1.96%	0.316%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

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\82X071619W.M
Title : SW846 8260

72	16.346	2487	2492	2494	rBV	93562	157890	1.77%	0.286%
73	16.438	2503	2507	2515	rVB	213202	400330	4.50%	0.724%
74	16.505	2515	2518	2523	rVB	80031	124645	1.40%	0.225%
75	16.560	2523	2527	2532	rVB2	94959	155997	1.75%	0.282%
76	16.688	2544	2548	2553	rVB	255987	448162	5.03%	0.811%
77	16.749	2553	2558	2564	rBV2	233440	432975	4.86%	0.783%

Sum of corrected areas: 55278184

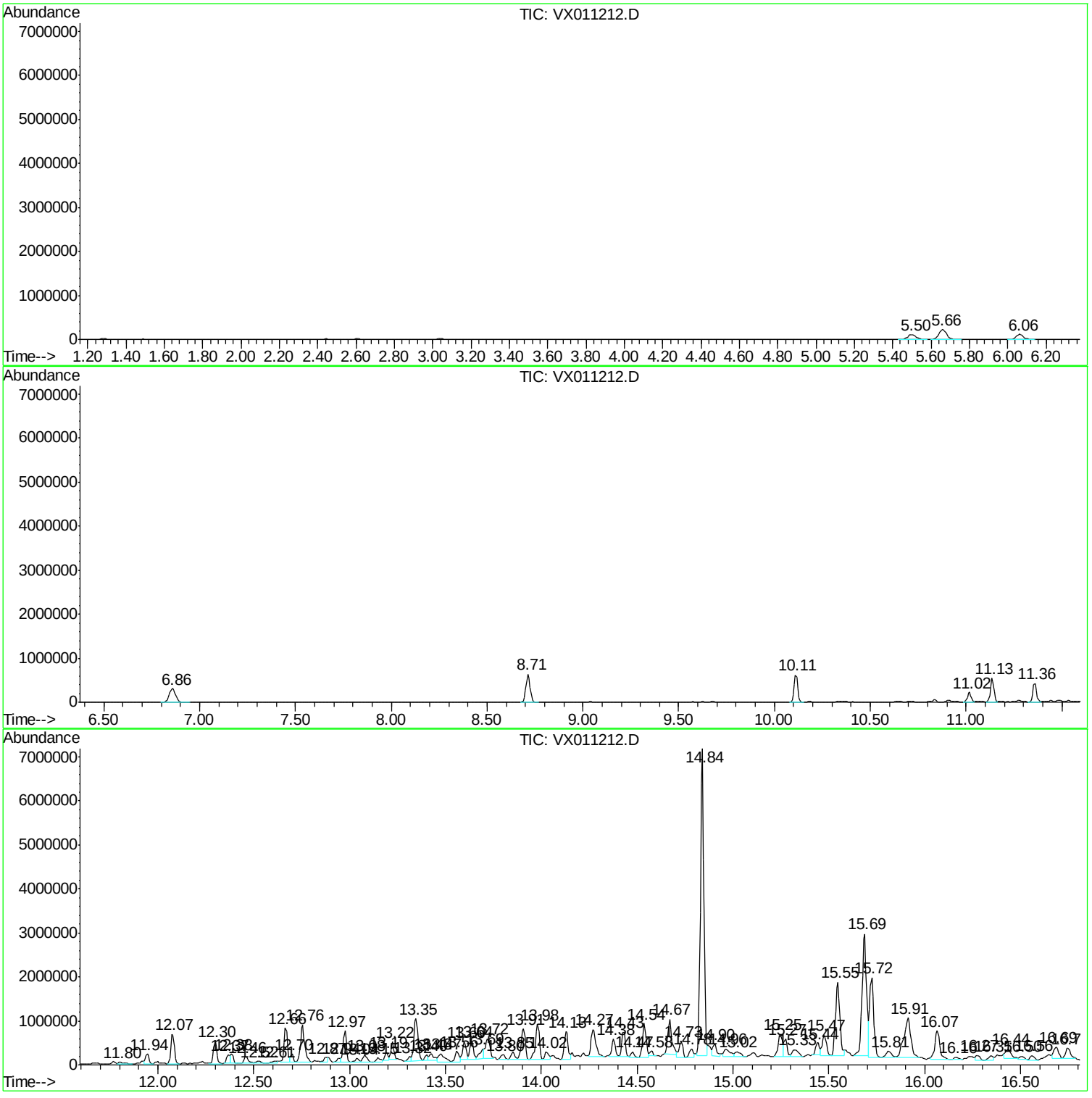
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

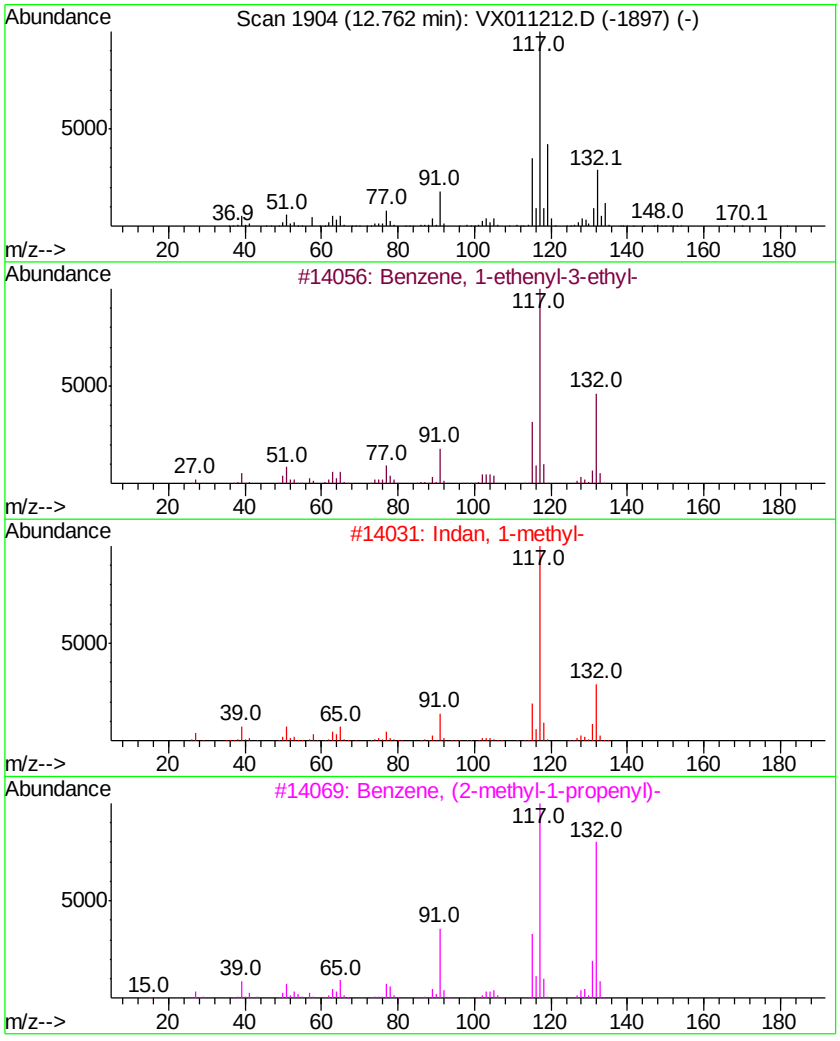
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

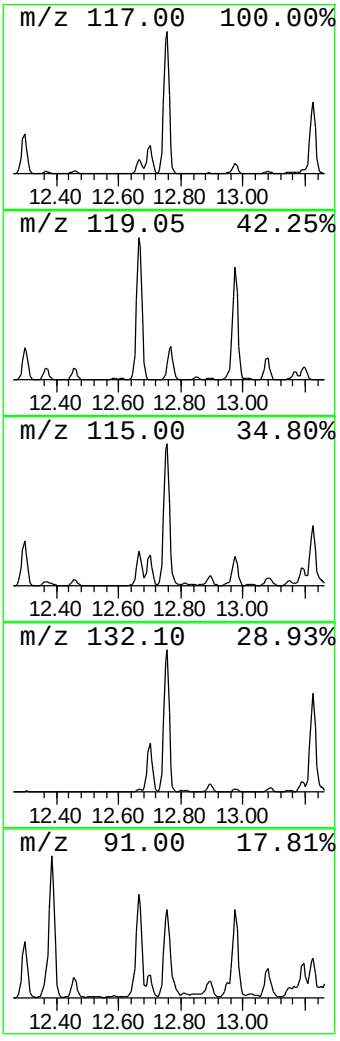
****

Peak Number 1 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	75.00 ug/l	1292090	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	87
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	74





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

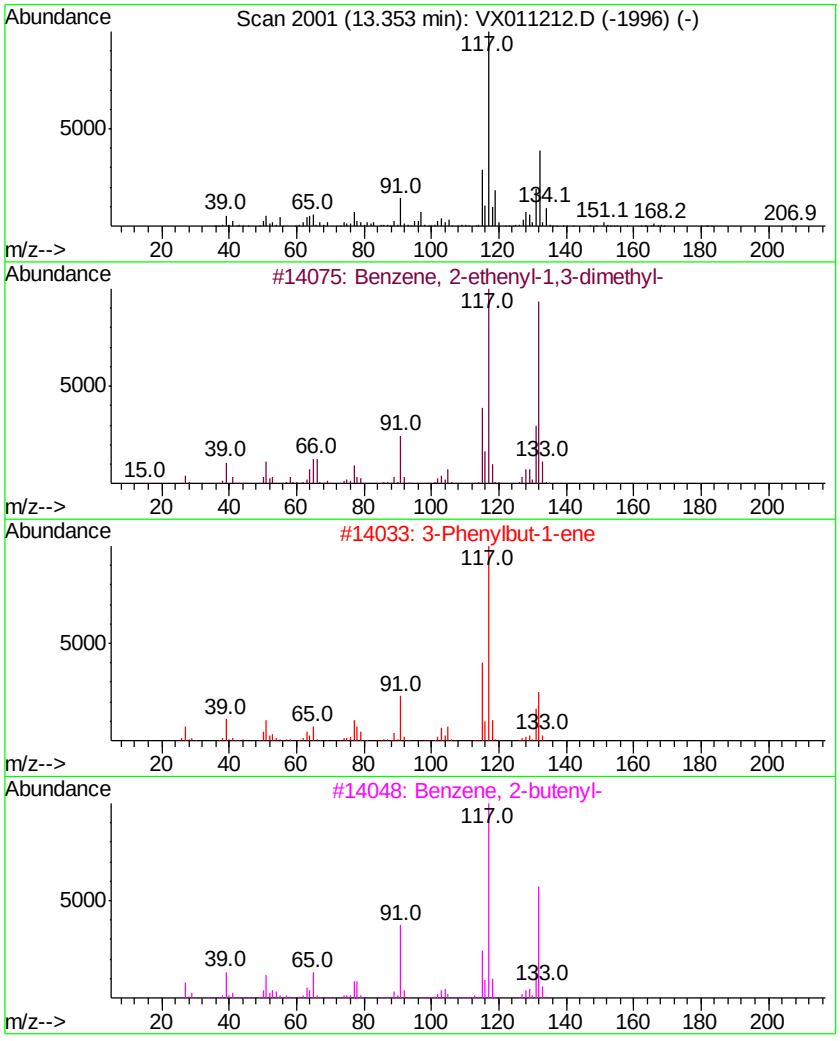
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

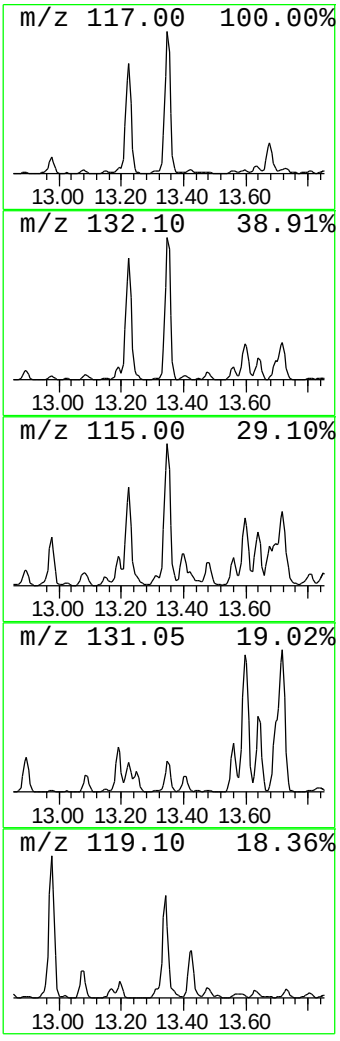
****

Peak Number 2 Benzene, 2-ethenyl-1,3-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	83.37 ug/l	1436300	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	86
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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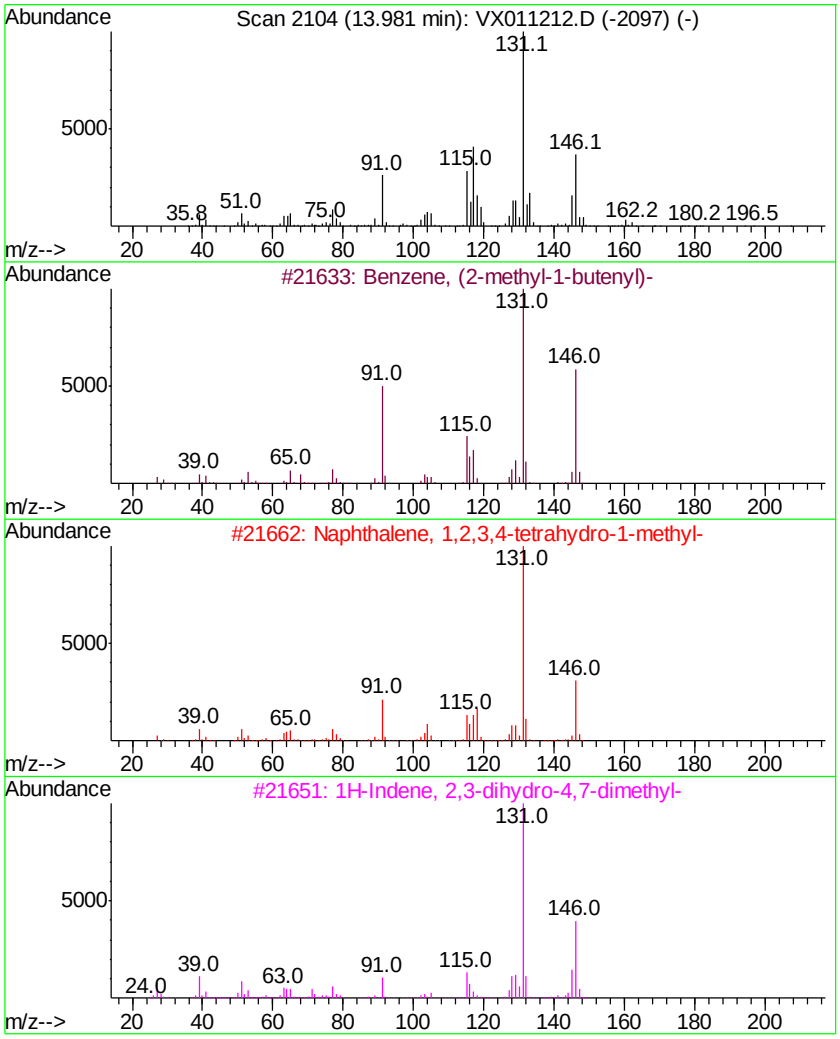
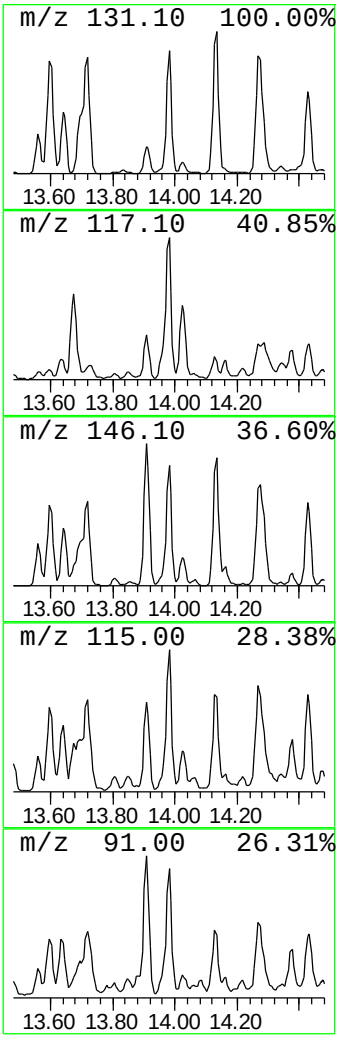
Peak Number 3 Benzene, (2-methyl-1-butenyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.98	68.24 ug/l	1175680	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	92
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	83
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	83
5	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	81





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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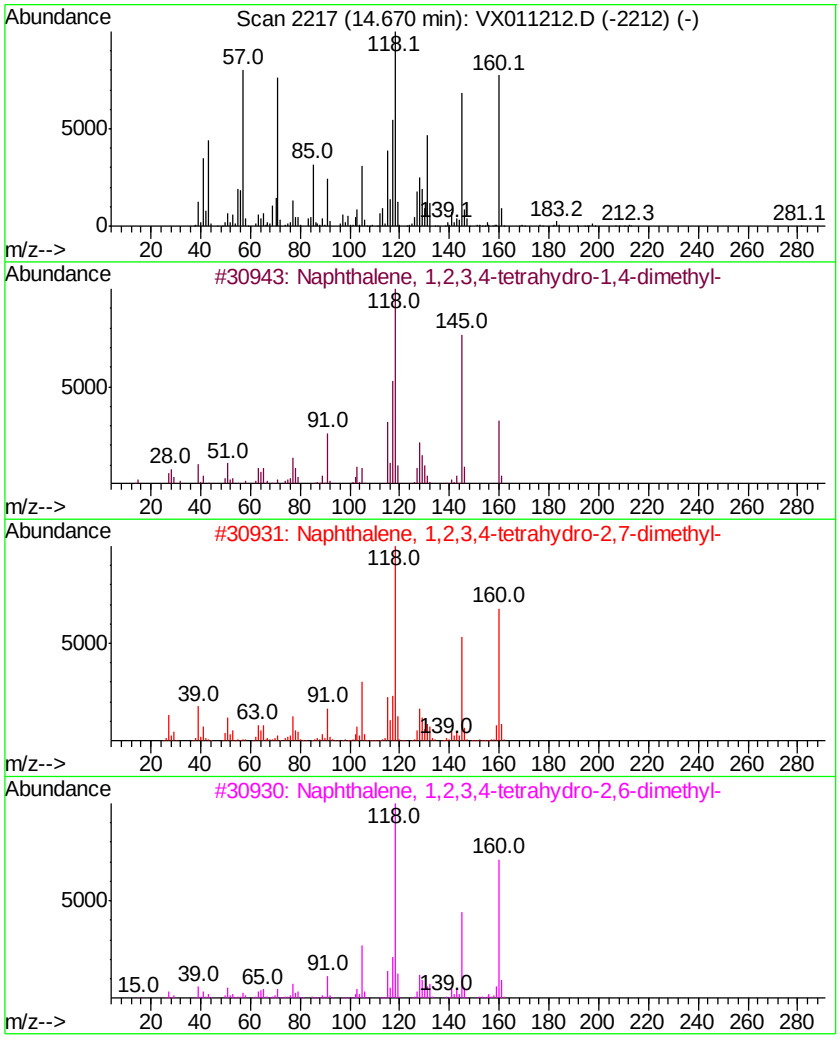
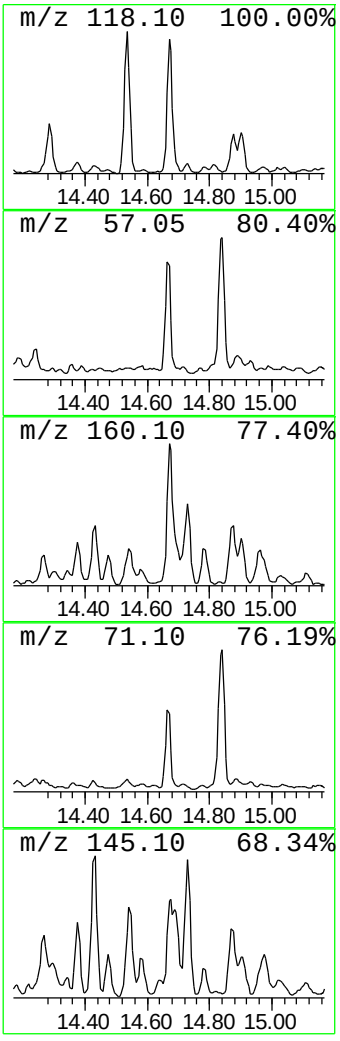
Peak Number 4 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	69.02 ug/l	1189030	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	91
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	55
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	46
4	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	40
5	2-Phenylpropyl isobutyrate	206	C13H18O2	065813-53-8	38





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

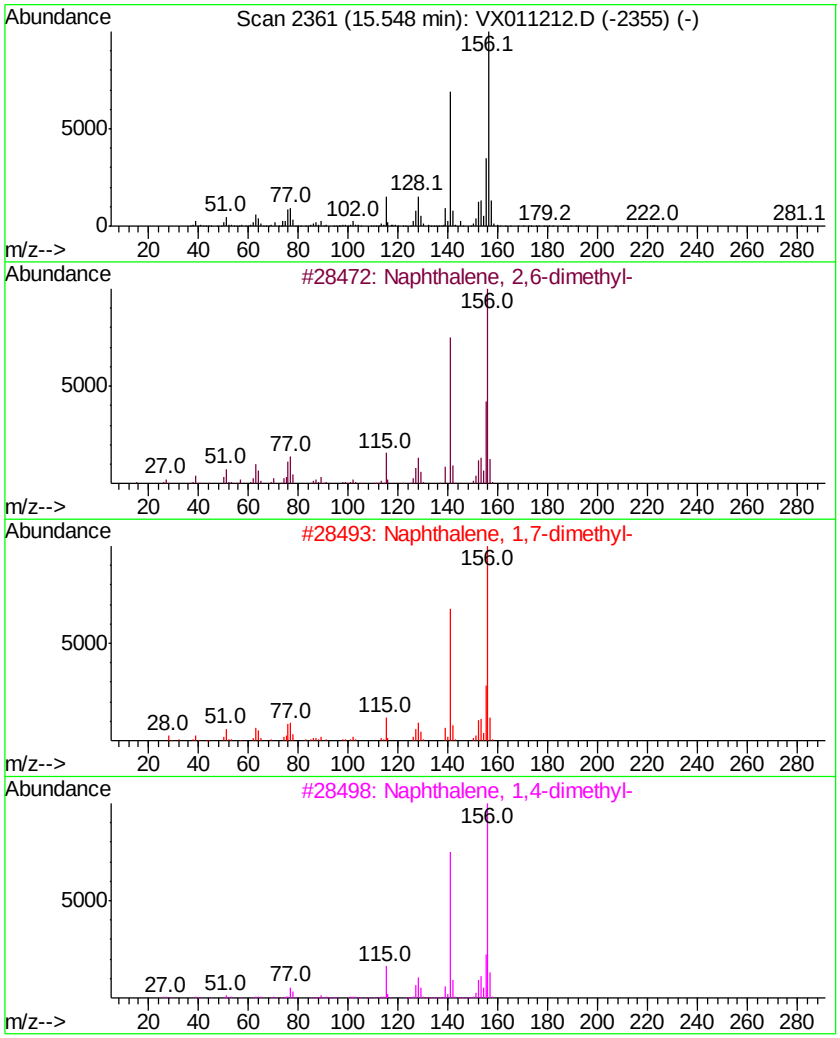
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

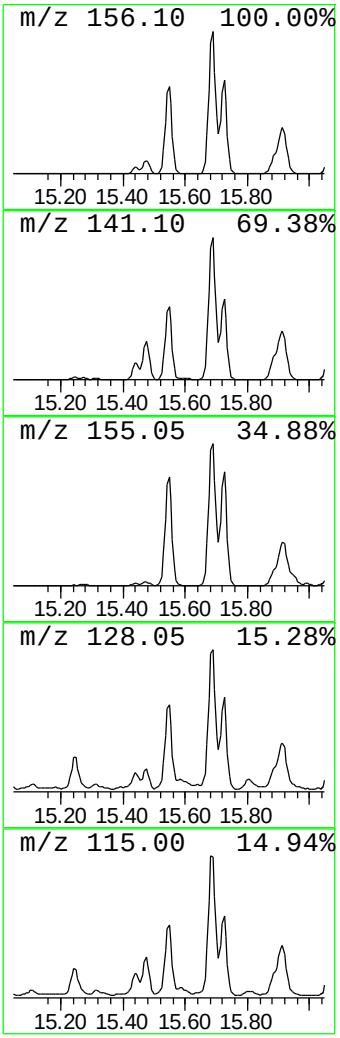
****

Peak Number 6 Naphthalene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.55	154.49 ug/l	2661570	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, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

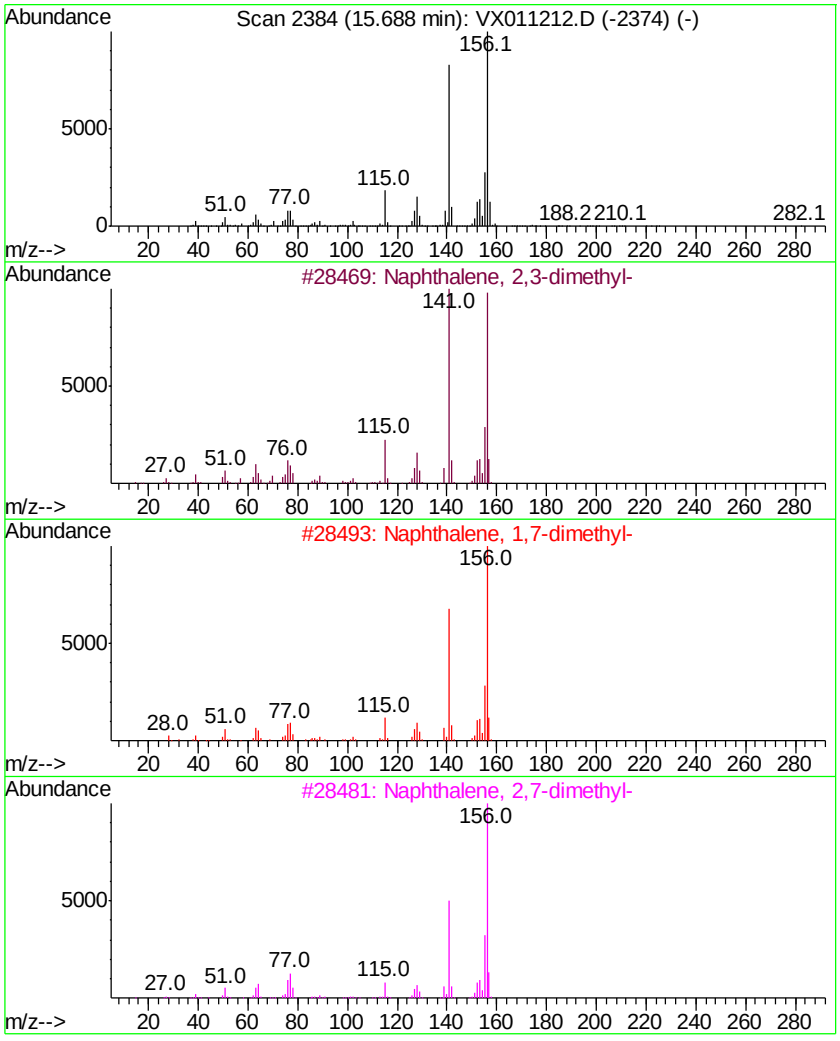
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

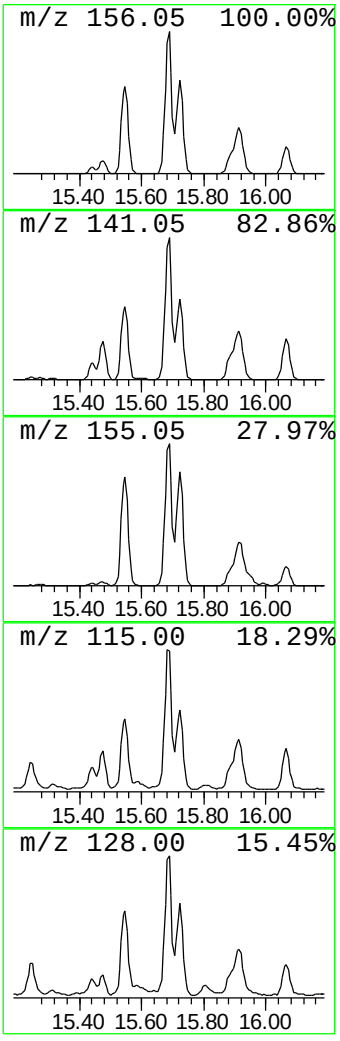
****

Peak Number 7 Naphthalene, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.69	268.56 ug/l	4626940	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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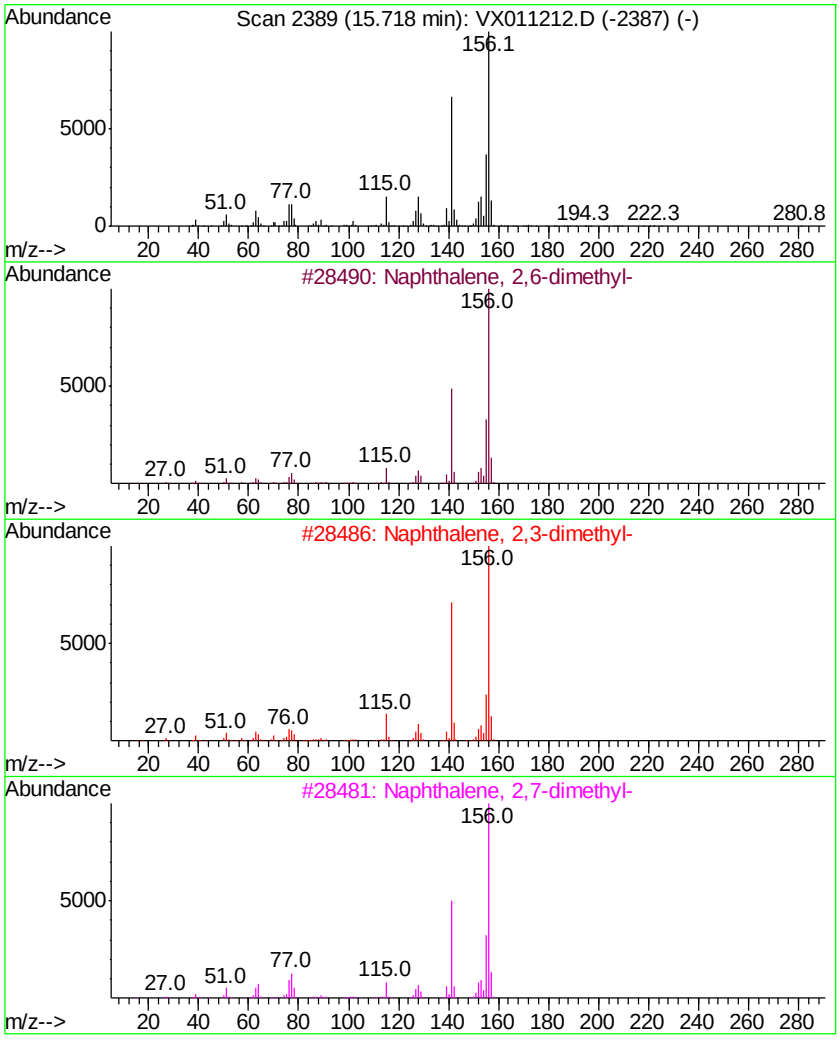
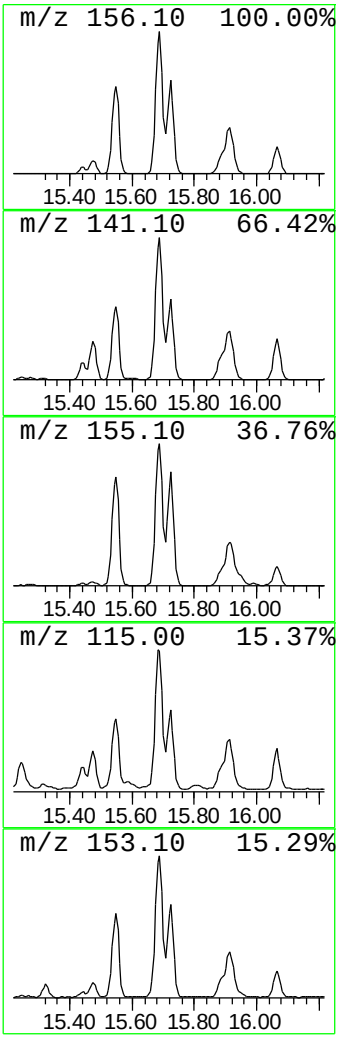
Peak Number 8 Naphthalene, 2,7-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.72	151.16 ug/l	2604270	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, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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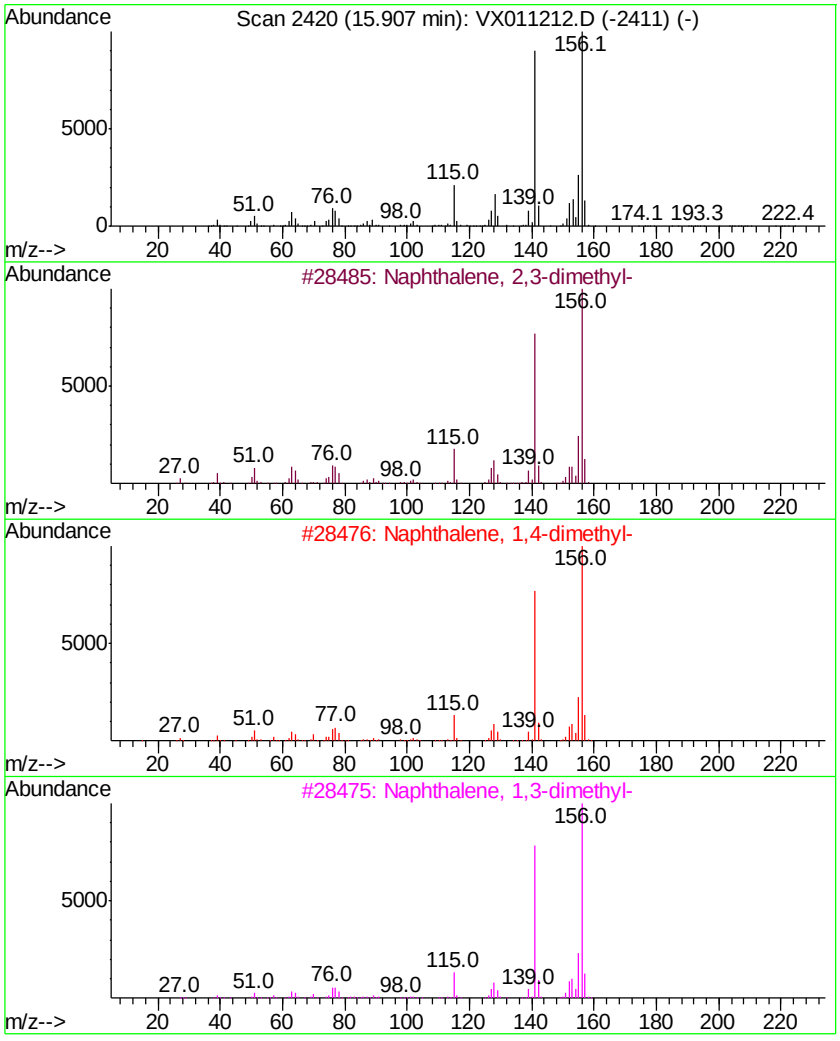
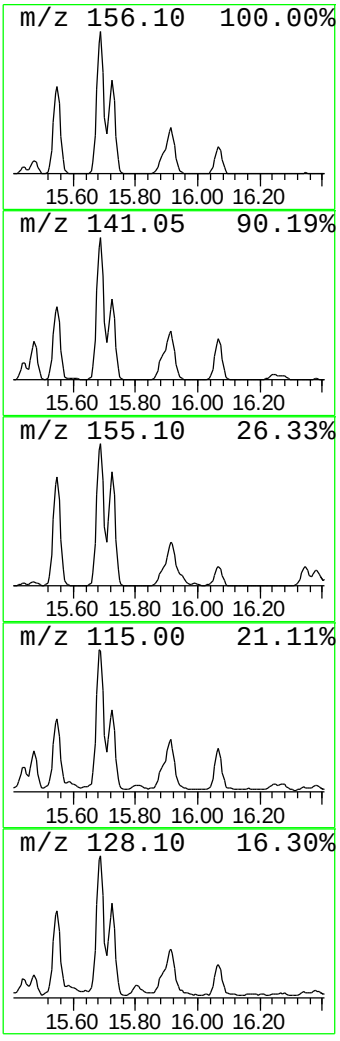
Peak Number 9 Naphthalene, 1,7-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	135.17 ug/l	2328730	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
3	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.M
Quant Title : SW846 8260

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

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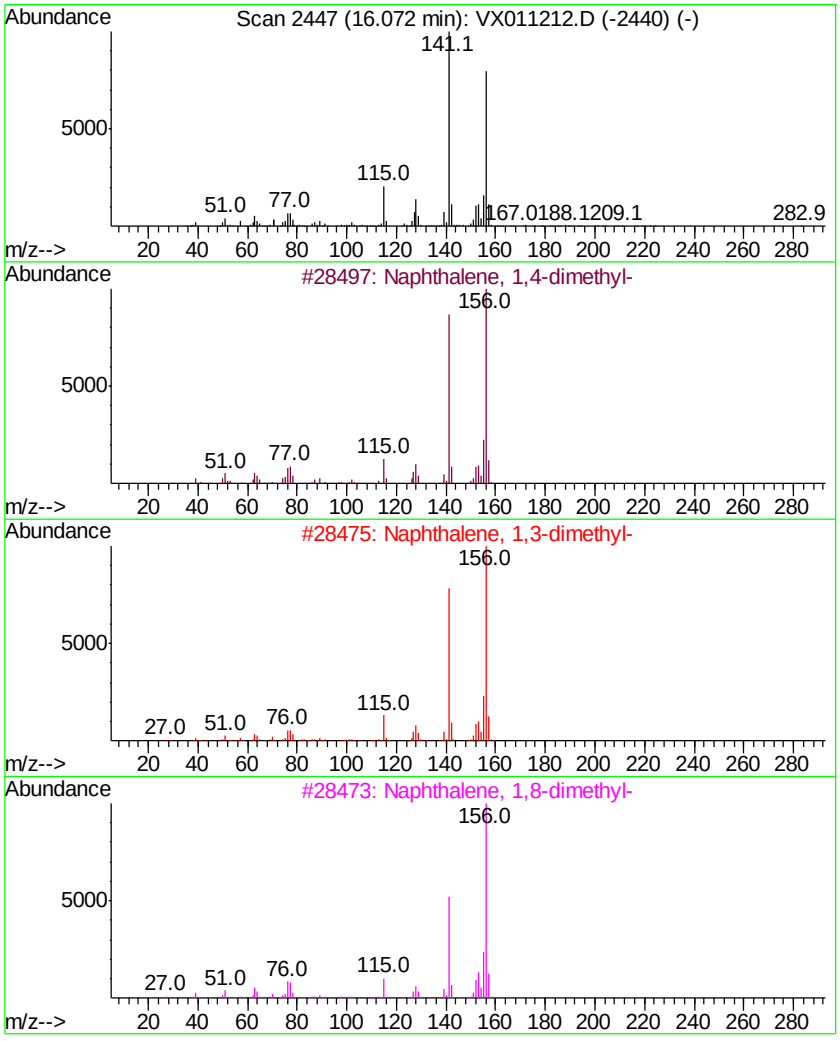
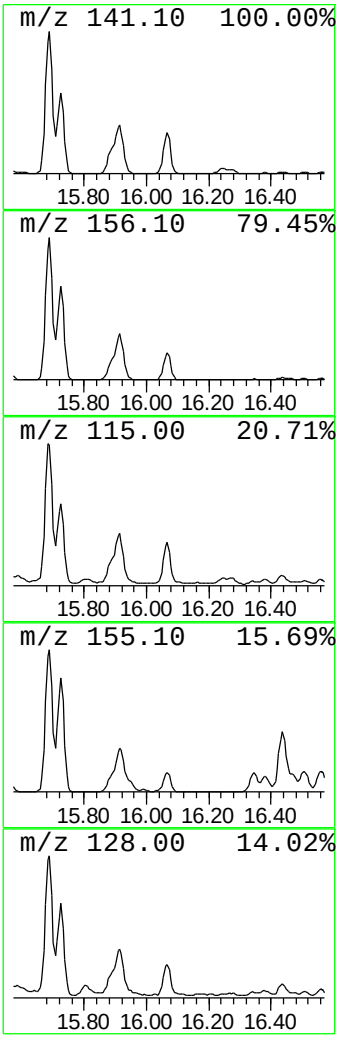
Peak Number 10 Naphthalene, 1,4-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	71.69 ug/l	1235100	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	96
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	95





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX073019\
Data File : VX011212.D
Acq On : 30 Jul 2019 21:57
Operator : JC/SP
Sample : K4024-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 28 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : TP-8

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X071619W.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-etheny...	12.76	75.0	ug/l	1292090	4	12.08	861424	50.0
Benzene, 2-etheny...	13.35	83.4	ug/l	1436300	4	12.08	861424	50.0
Benzene, (2-methy...	13.98	68.2	ug/l	1175680	4	12.08	861424	50.0
Naphthalene, 1,2,...	14.67	69.0	ug/l	1189030	4	12.08	861424	50.0
Naphthalene, 2,6-...	15.55	154.5	ug/l	2661570	4	12.08	861424	50.0
Naphthalene, 2,3-...	15.69	268.6	ug/l	4626940	4	12.08	861424	50.0
Naphthalene, 2,7-...	15.72	151.2	ug/l	2604270	4	12.08	861424	50.0
Naphthalene, 1,7-...	15.91	135.2	ug/l	2328730	4	12.08	861424	50.0
Naphthalene, 1,4-...	16.07	71.7	ug/l	1235100	4	12.08	861424	50.0