LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060618\
Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

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\82X060418W.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	1.075	76	80	93	rBV	1269943	1589544	46.65%	4.767%
2	1.270	108	112	118	rBV	63534	85074	2.50%	0.255%
3	1.355	118	126	127	rBV	30962	66228	1.94%	0.199%
4	2.928	379	384	397	rVB	84717	190950	5.60%	0.573%
5	4.397	614	625	637	rVB	86625	254754	7.48%	0.764%
6	5.513	794	808	813	rBV	136661	409760	12.03%	1.229%
7	5.580	813	819	826	rVV	168664	507186	14.89%	1.521%
8	5.665	826	833	857	rVB	215753	661461	19.41%	1.984%
9	5.940	867	878	890	rVB5	22877	82572	2.42%	0.248%
10	6.074	890	900	909	rBV	157997	419374	12.31%	1.258%
11	6.306	920	938	944	rBV2	72190	272655	8.00%	0.818%
12	6.452	955	962	975	rVB3	45720	131914	3.87%	0.396%
13	6.866	1020	1030	1040	rBV	317073	763685	22.41%	2.290%
14	7.464	1115	1128	1141	rBV3	151730	391497	11.49%	1.174%
15	7.927	1198	1204	1206	rBV	41083	68085	2.00%	0.204%
16	7.958	1207	1209	1216	rVV2	46632	74368	2.18%	0.223%
17	8.214	1238	1251	1257	rBV	52297	106911	3.14%	0.321%
18	8.573	1304	1310	1319	rVB3	86513	165030	4.84%	0.495%
19	8.671	1319	1326	1329	rBV	72114	136928	4.02%	0.411%
20	8.720	1329	1334	1343	rVB	790567	1244294	36.52%	3.732%
21	8.842	1349	1354	1358	rBV	37879	59221	1.74%	0.178%
22	8.915	1363	1366	1373	rVB5	30473	58088	1.70%	0.174%
23	9.043	1382	1387	1394	rVB	76727	121618	3.57%	0.365%
24	9.165	1399	1407	1412	rVV	50455	89590	2.63%	0.269%
25	9.226	1412	1417	1429	rVB	59525	110011	3.23%	0.330%
26	9.573	1469	1474	1479	rVB3	41638	77761	2.28%	0.233%
27	9.628	1479	1483	1487	rBV2	60945	91030	2.67%	0.273%
28	9.683	1487	1492	1496	rVB2	42744	74144	2.18%	0.222%
29	10.116	1558	1563	1568	rBV	669530	882858	25.91%	2.648%
30	10.189	1568	1575	1578	rBV	57704	90511	2.66%	0.271%
31	10.341	1590	1600	1601	rBV4	41818	83749	2.46%	0.251%
32	10.366	1601	1604	1609	rVB2	68114	105709	3.10%	0.317%
33	10.646	1642	1650	1657	rVV5	42131	104879	3.08%	0.315%
34	10.841	1674	1682	1690	rBV2	73225	113790	3.34%	0.341%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX060618\
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Sample : J3311-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

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\82X060418W.M
Title : SW846 8260

35	10.915	1690	1694	1698	rBV3	53014	91332	2.68%	0.274%
36	10.957	1698	1701	1706	rVB5	33270	60025	1.76%	0.180%
37	11.024	1706	1712	1716	rBV	780142	1017236	29.86%	3.051%
38	11.140	1726	1731	1736	rVB	707359	865640	25.41%	2.596%
39	11.280	1751	1754	1759	rVB3	45261	61113	1.79%	0.183%
40	11.366	1763	1768	1778	rVB	730866	915554	26.87%	2.746%
41	11.457	1778	1783	1786	rBV2	60190	84602	2.48%	0.254%
42	11.670	1814	1818	1826	rBV	129981	208224	6.11%	0.624%
43	11.774	1826	1835	1838	rBV	74482	127405	3.74%	0.382%
44	11.927	1853	1860	1861	rBV	64298	99541	2.92%	0.299%
45	11.951	1861	1864	1868	rVB	169598	211267	6.20%	0.634%
46	12.085	1881	1886	1891	rBV	826369	1063179	31.20%	3.188%
47	12.237	1907	1911	1916	rVB	109876	151775	4.45%	0.455%
48	12.305	1917	1922	1927	rBV	2851135	3407104	100.00%	10.218%
49	12.372	1929	1933	1941	rVB2	193676	361034	10.60%	1.083%
50	12.463	1944	1948	1955	rVB	196412	264555	7.76%	0.793%
51	12.524	1955	1958	1964	rVB5	38294	71639	2.10%	0.215%
52	12.591	1965	1969	1972	rBV	50652	76134	2.23%	0.228%
53	12.670	1978	1982	1985	rBV	573947	678473	19.91%	2.035%
54	12.707	1985	1988	1992	rVB	246998	277782	8.15%	0.833%
55	12.762	1992	1997	2004	rVB2	1456238	2001036	58.73%	6.001%
56	12.823	2004	2007	2012	rVB5	43312	57496	1.69%	0.172%
57	12.902	2012	2020	2024	rVB2	133827	268358	7.88%	0.805%
58	12.981	2024	2033	2037	rBV	602629	833044	24.45%	2.498%
59	13.042	2037	2043	2047	rBV2	69763	116153	3.41%	0.348%
60	13.091	2047	2051	2057	rVB2	133719	225579	6.62%	0.677%
61	13.158	2057	2062	2065	rBV	114336	169081	4.96%	0.507%
62	13.201	2065	2069	2071	rBV	170472	199652	5.86%	0.599%
63	13.231	2071	2074	2077	rVB	383828	391888	11.50%	1.175%
64	13.317	2084	2088	2089	rBV2	70941	81533	2.39%	0.245%
65	13.353	2089	2094	2100	rVB2	1221413	1711186	50.22%	5.132%
66	13.408	2100	2103	2105	rBV2	85772	102038	2.99%	0.306%
67	13.487	2111	2116	2123	rVB2	244957	406247	11.92%	1.218%
68	13.566	2126	2129	2132	rBV4	72179	97862	2.87%	0.293%
69	13.603	2132	2135	2138	rBV	205159	263271	7.73%	0.790%
70	13.646	2138	2142	2146	rVB3	208071	279207	8.19%	0.837%
71	13.707	2146	2152	2153	rBV2	151073	240793	7.07%	0.722%

LSC Area Percent Report

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Sample : J3311-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

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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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Title : SW846 8260

72	13.725	2153	2155	2160	rVB2	289276	397812	11.68%	1.193%
73	13.810	2166	2169	2173	rVB2	74957	96226	2.82%	0.289%
74	13.859	2173	2177	2181	rVB3	137830	179832	5.28%	0.539%
75	13.914	2181	2186	2190	rBV2	208689	318168	9.34%	0.954%
76	13.987	2191	2198	2202	rVB2	270111	441991	12.97%	1.326%
77	14.036	2202	2206	2211	rBV4	89045	166873	4.90%	0.500%
78	14.133	2218	2222	2225	rBV2	99872	148317	4.35%	0.445%
79	14.164	2225	2227	2230	rVV3	76335	86028	2.52%	0.258%
80	14.280	2242	2246	2253	rVB4	216167	452029	13.27%	1.356%
81	14.383	2259	2263	2268	rBV	116326	165362	4.85%	0.496%
82	14.438	2268	2272	2276	rVB2	142538	191157	5.61%	0.573%
83	14.481	2276	2279	2285	rVB5	69626	95959	2.82%	0.288%
84	14.548	2285	2290	2294	rBV2	263469	392127	11.51%	1.176%
85	14.676	2307	2311	2318	rVV4	151634	228110	6.70%	0.684%
86	14.737	2318	2321	2325	rVB	97194	118688	3.48%	0.356%
87	14.786	2326	2329	2332	rBV2	48179	64206	1.88%	0.193%
88	14.847	2334	2339	2344	rVV	641617	891180	26.16%	2.673%
89	14.914	2347	2350	2353	rVB3	61723	78449	2.30%	0.235%
90	15.030	2367	2369	2377	rVB5	64123	117615	3.45%	0.353%
91	15.255	2402	2406	2408	rBV	88717	127709	3.75%	0.383%
92	15.280	2408	2410	2414	rVB	95308	106763	3.13%	0.320%
93	15.450	2435	2438	2441	rBV2	46814	58557	1.72%	0.176%
94	15.493	2441	2445	2449	rVB2	82627	129054	3.79%	0.387%
95	15.554	2451	2455	2460	rBV	151555	228636	6.71%	0.686%
96	15.694	2473	2478	2482	rBV	225296	357148	10.48%	1.071%
97	15.731	2482	2484	2492	rVB	72367	114594	3.36%	0.344%
98	15.901	2507	2512	2513	rBV	57912	88591	2.60%	0.266%
99	16.078	2537	2541	2553	rVB	66736	151920	4.46%	0.456%
100	16.426	2594	2598	2607	rVB2	64652	128101	3.76%	0.384%

Sum of corrected areas: 33344469

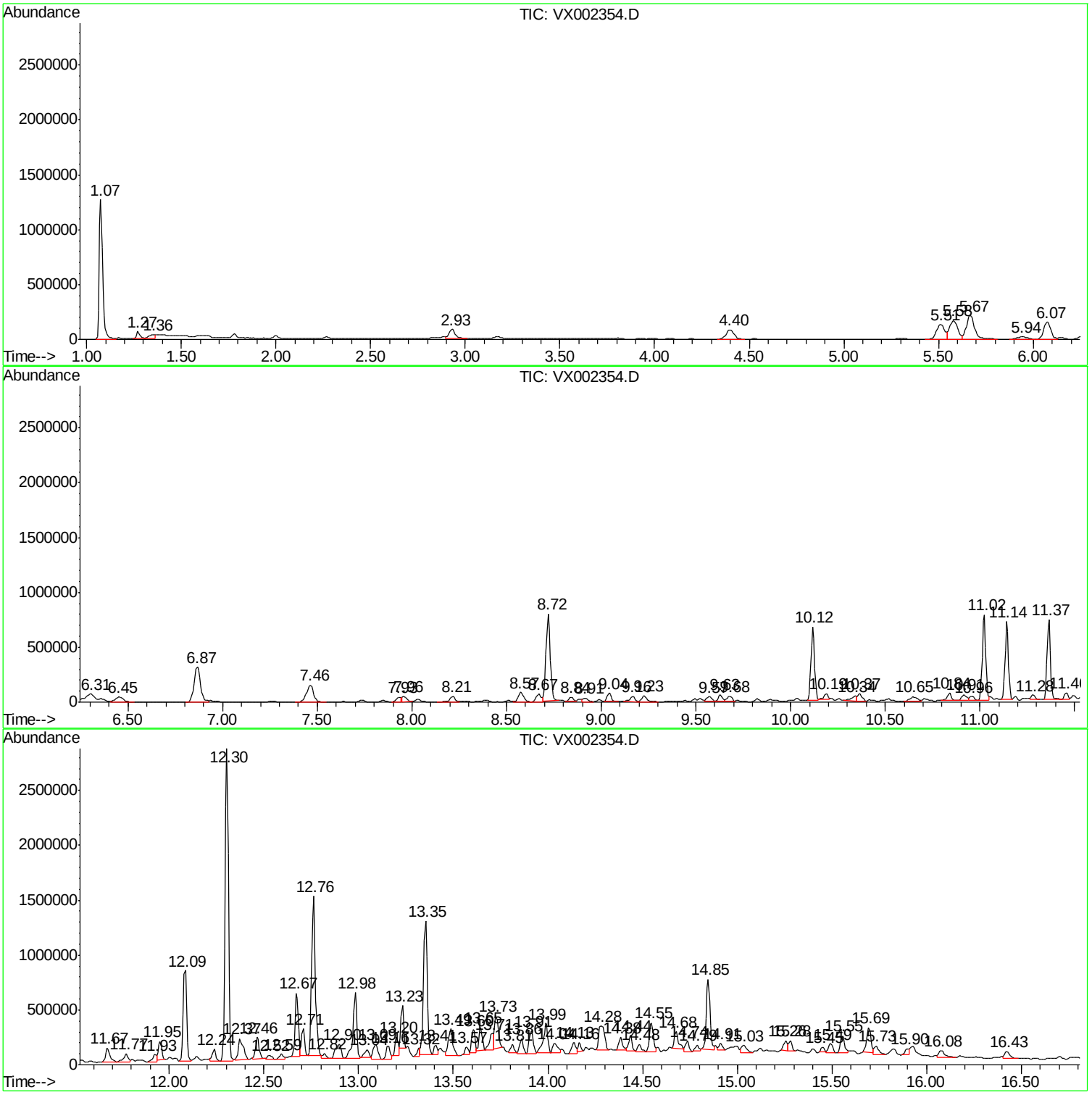
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.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\VX060618\
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Acq On : 06 Jun 2018 15:56
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Quant Title : SW846 8260

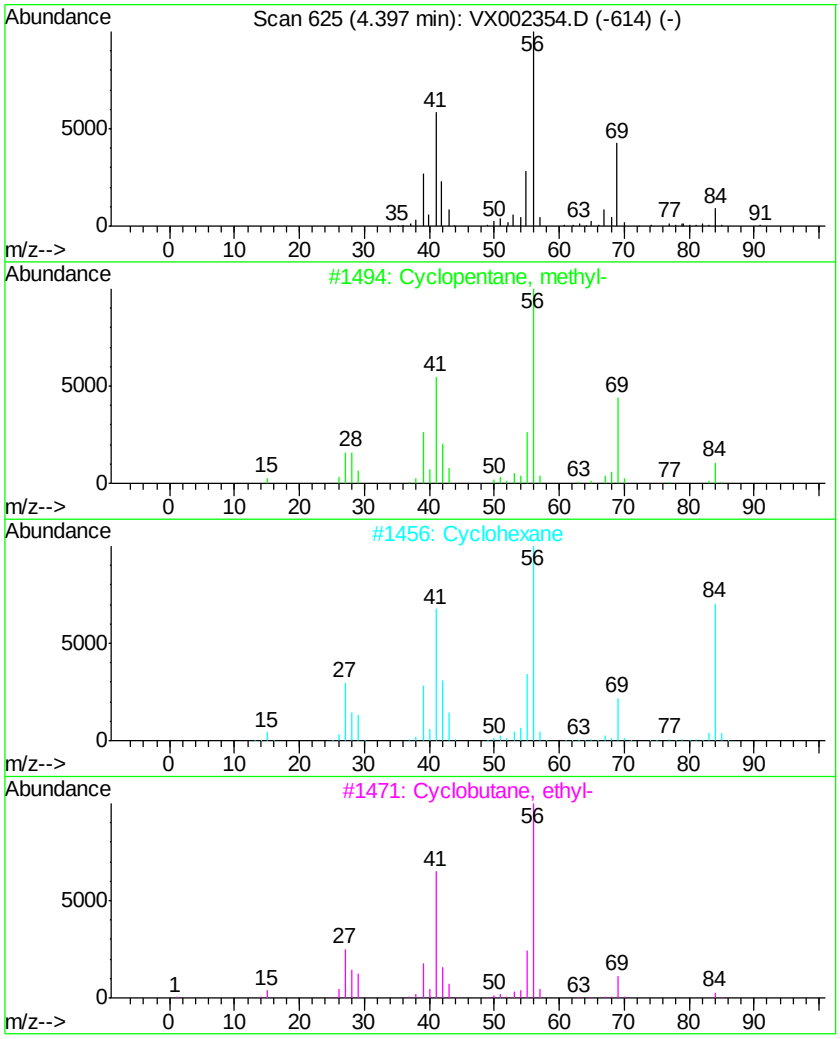
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

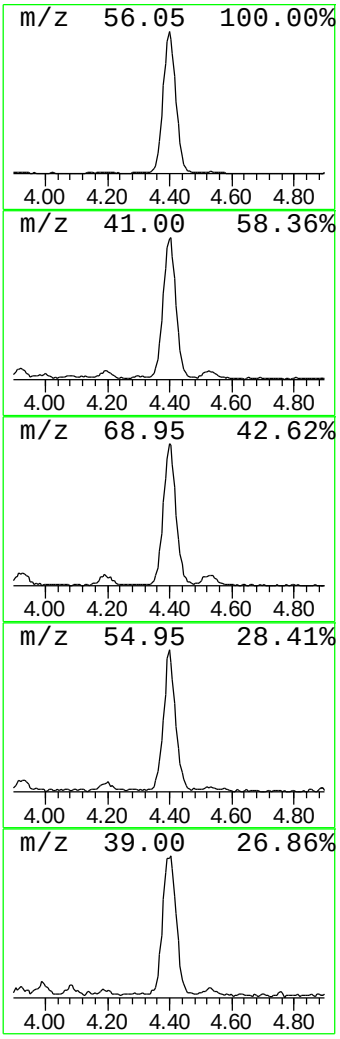
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	19.26 ug/l	254754	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclohexane	84	C6H12	000110-82-7	78
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	72
5	Cyclobutane	56	C4H8	000287-23-0	53





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

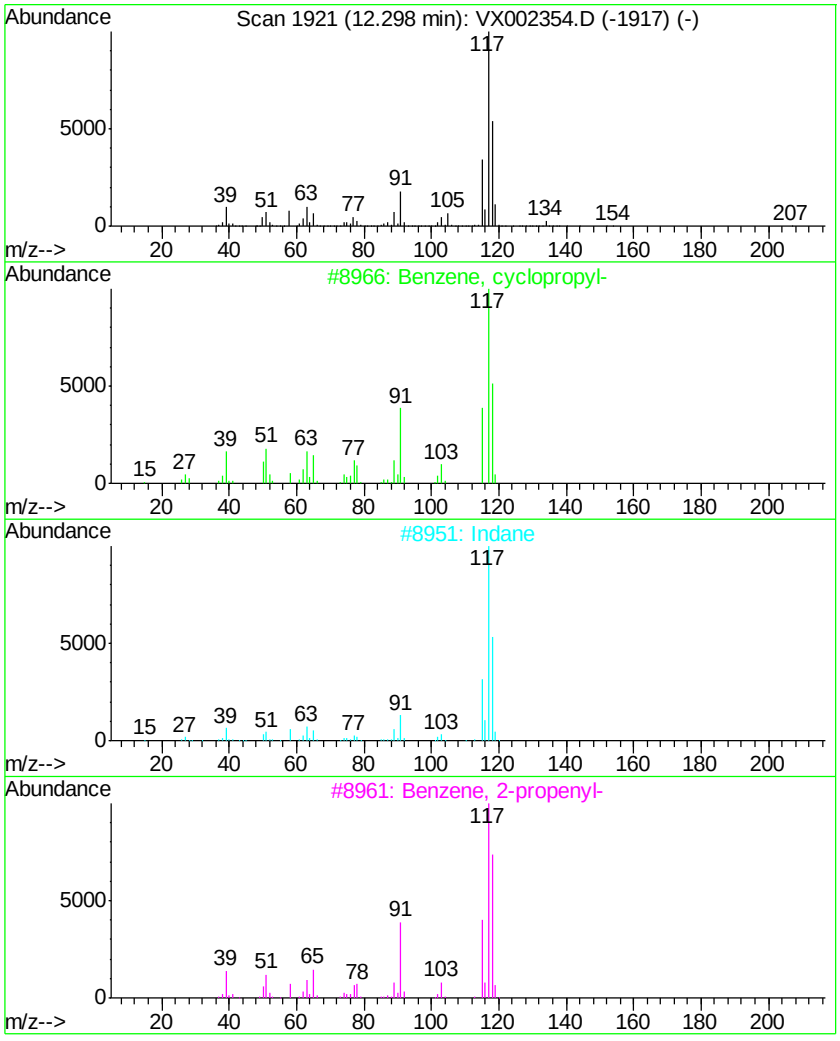
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

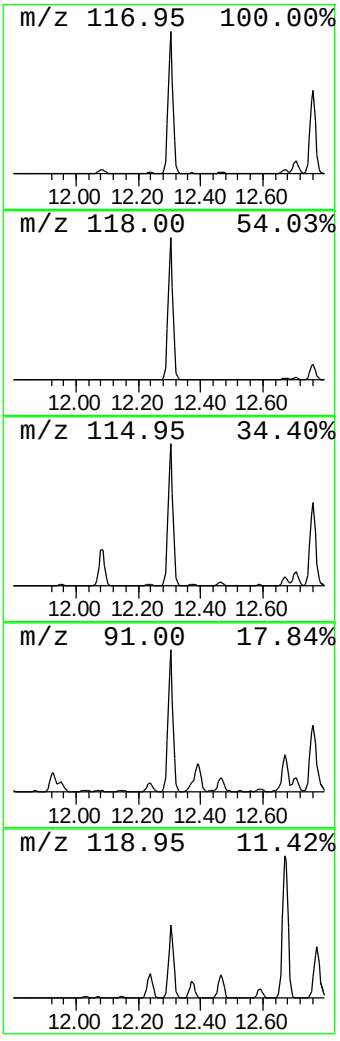
****

Peak Number 3 Benzene, cyclopropyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	160.23 ug/l	3407100	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
2	Indane	118	C9H10	000496-11-7	76
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4	Deltacyclene	118	C9H10	007785-10-6	58
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

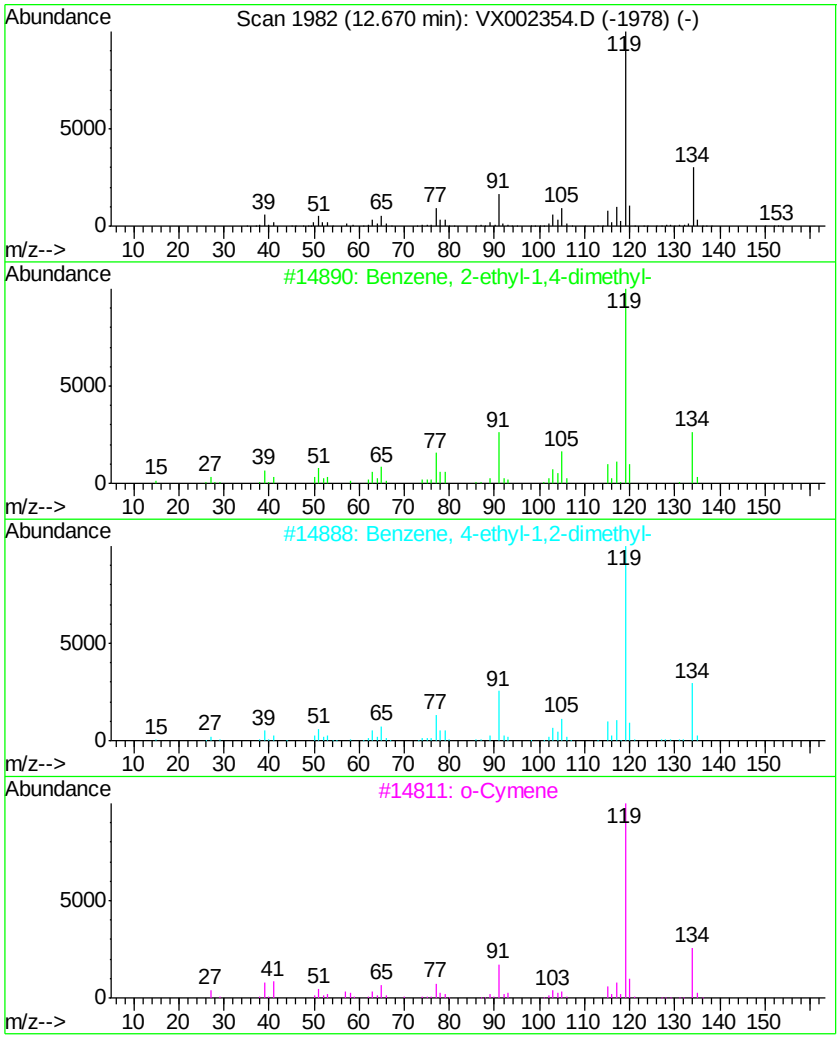
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

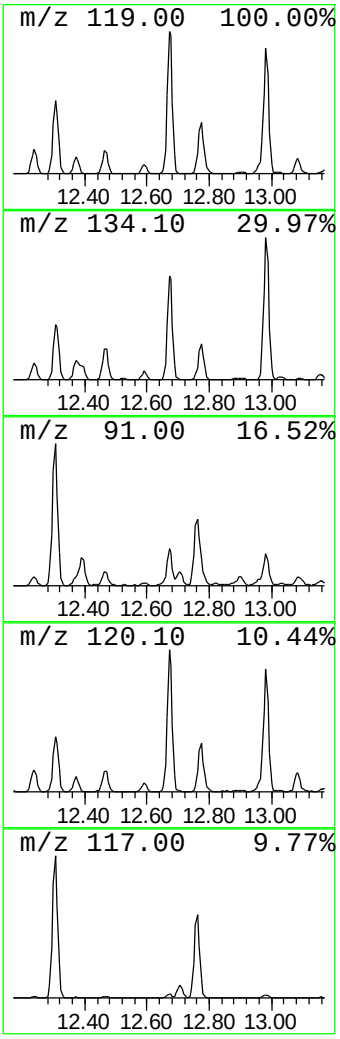
****

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.67	31.91 ug/l	678473	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	p-Cymene	134	C10H14	000099-87-6	94





Library Search Compound Report

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Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

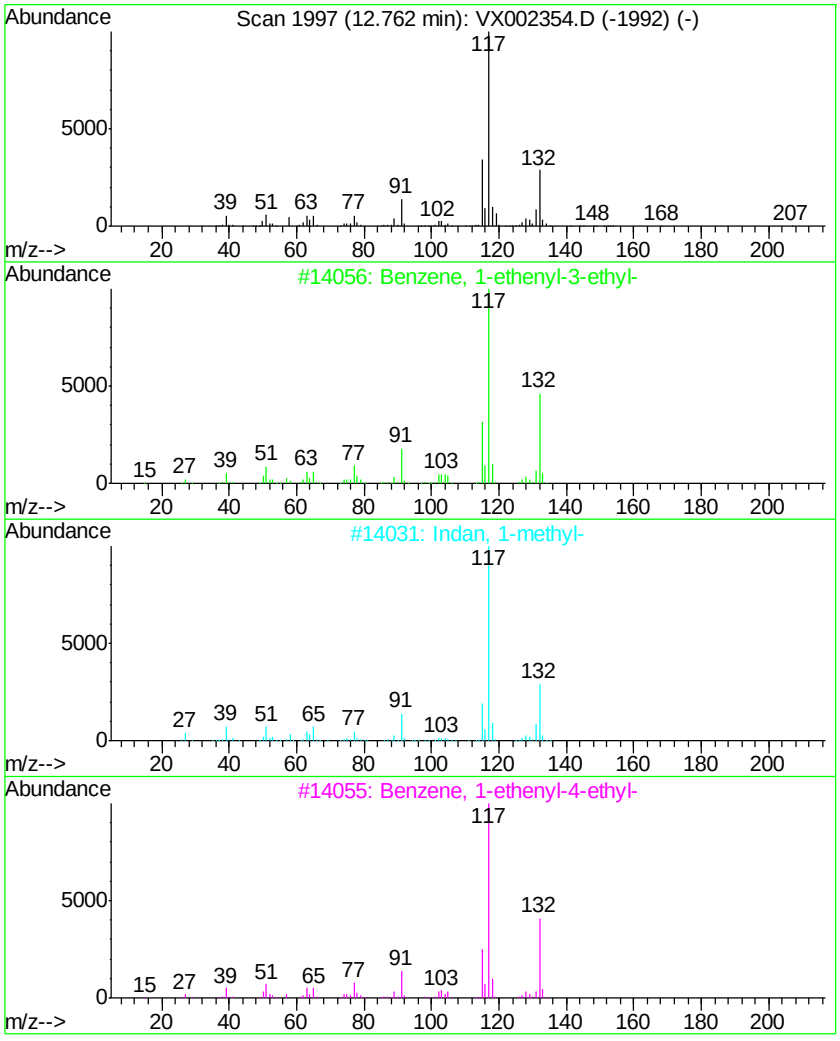
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

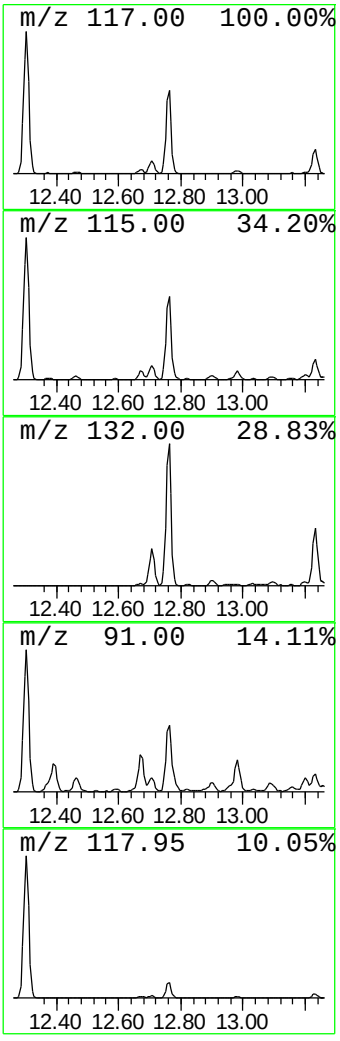
****

Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	94.11 ug/l	2001040	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	83
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	83





Library Search Compound Report

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Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

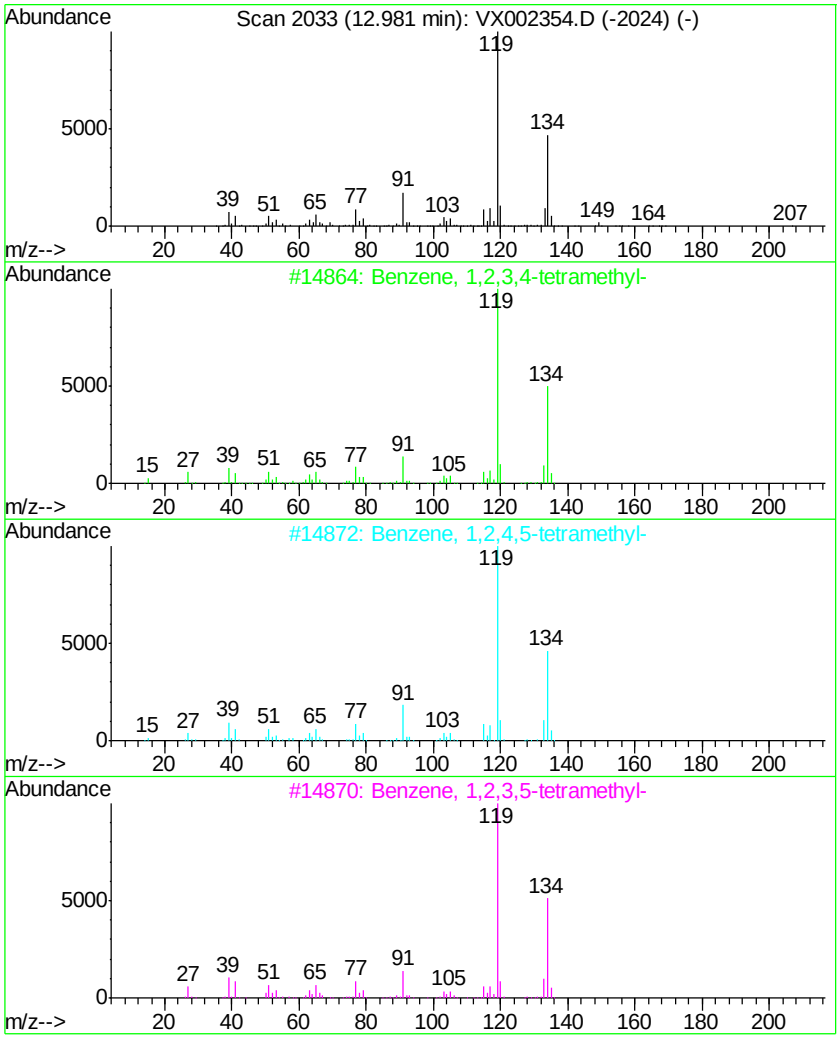
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

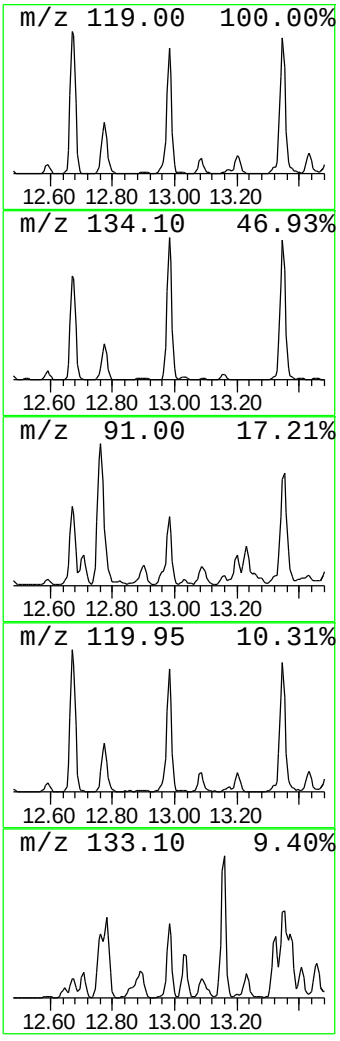
****

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.98	39.18 ug/l	833044	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	93
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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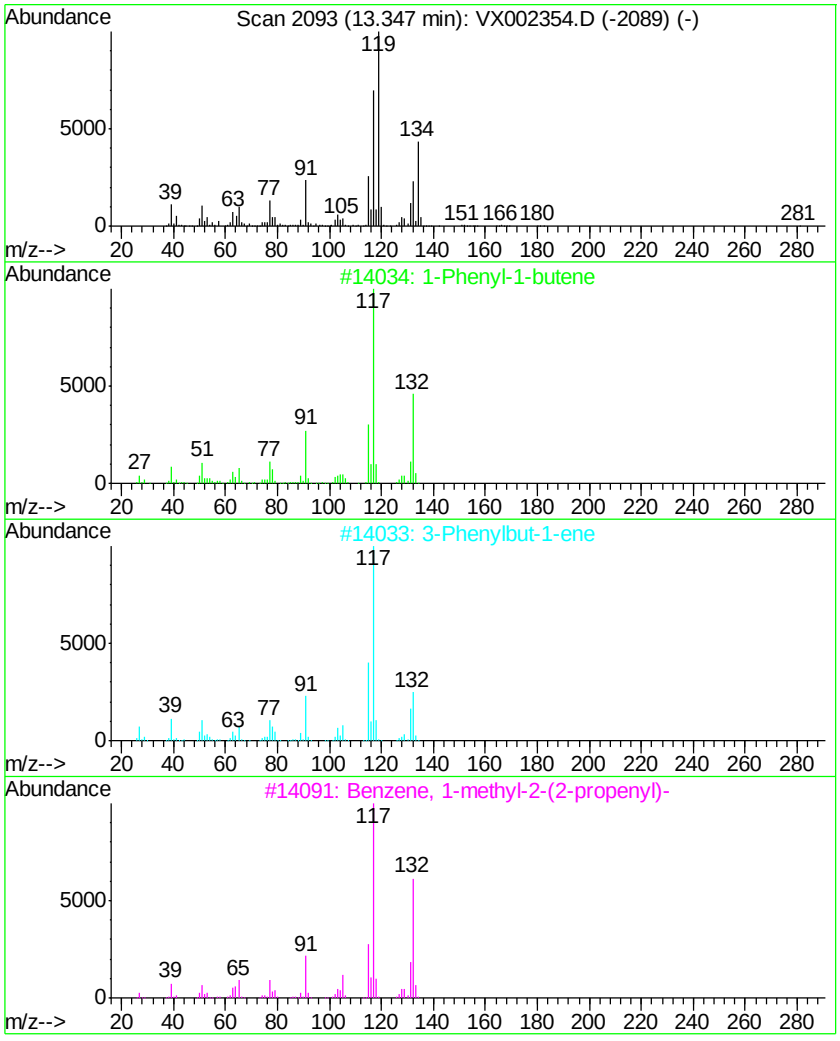
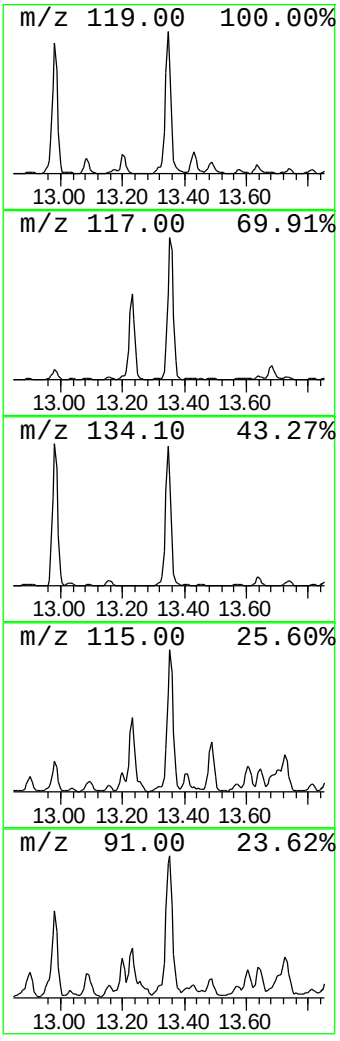
Peak Number 7 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	80.48 ug/l	1711190	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	64





Library Search Compound Report

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Data File : VX002354.D
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Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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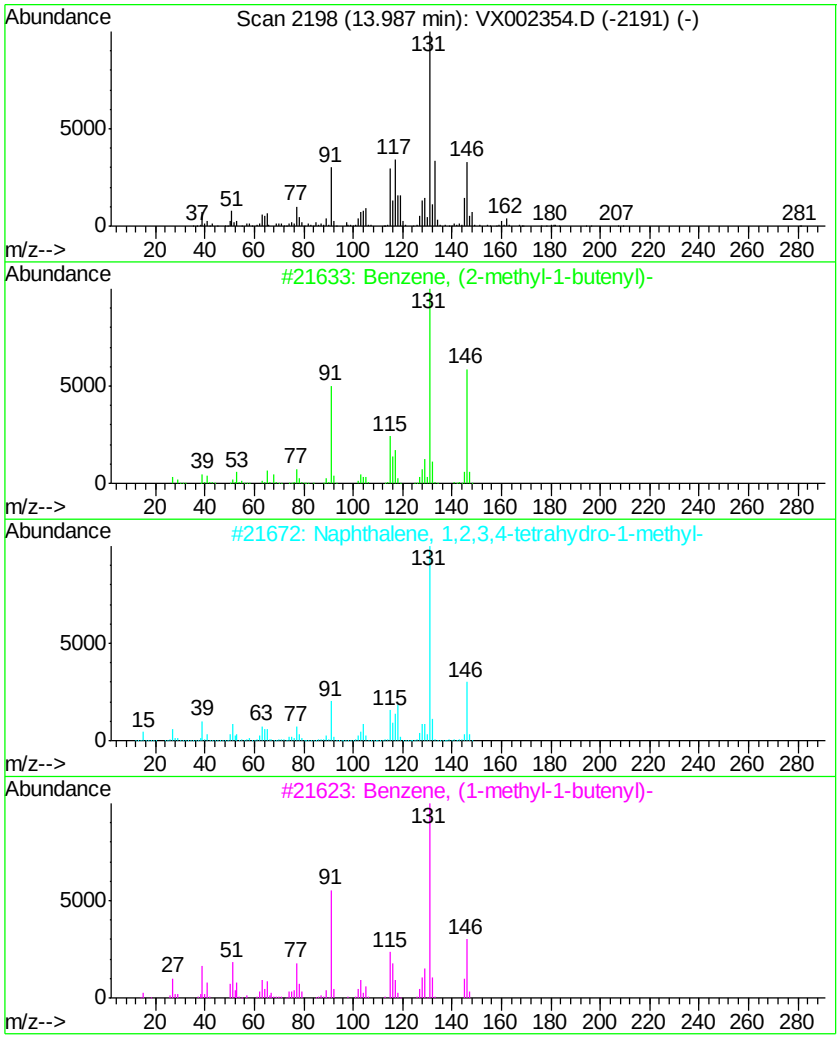
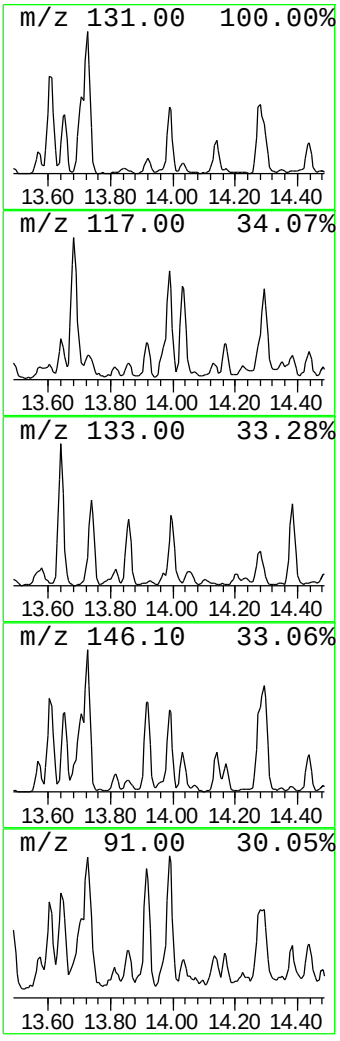
Peak Number 8 Benzene, (2-methyl-1-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	20.79 ug/l	441991	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	93
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	92
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	86
5	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	76





Library Search Compound Report

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Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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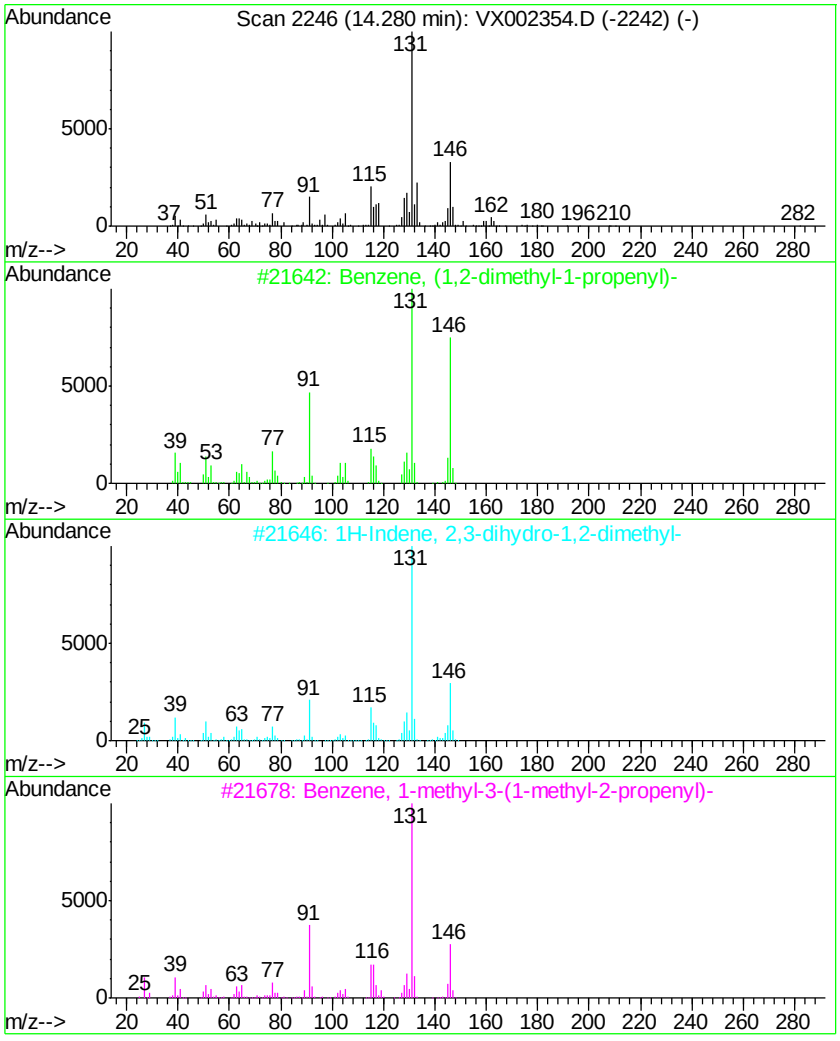
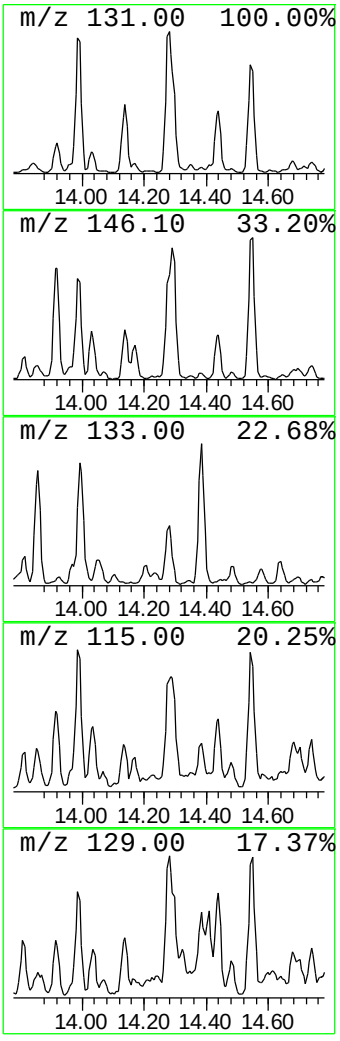
Peak Number 9 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.28	21.26 ug/l	452029	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	95
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
3	Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	87
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76





Library Search Compound Report

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ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.M
Quant Title : SW846 8260

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

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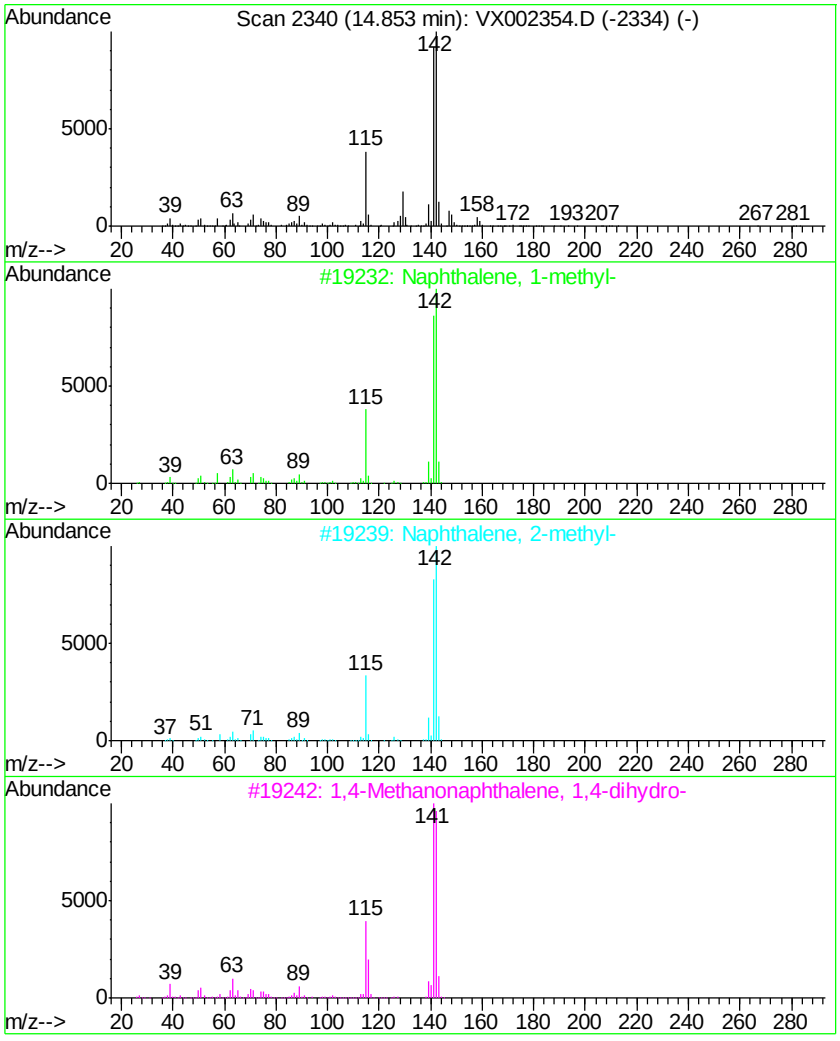
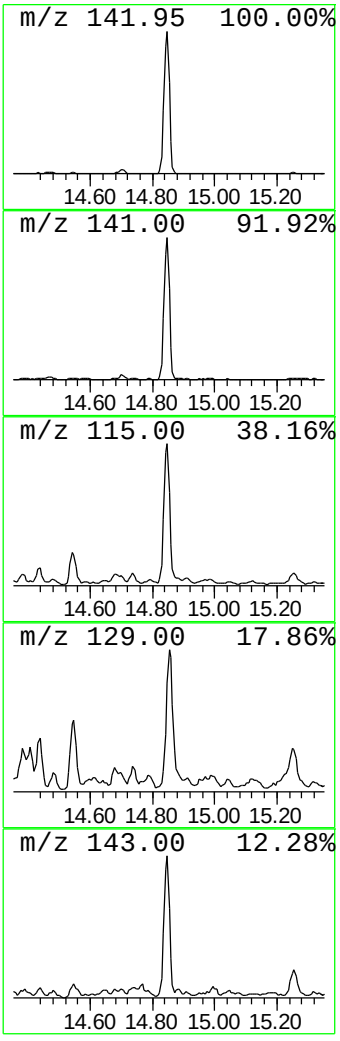
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	41.91 ug/l	891180	1,4-Dichlorobenzene-d4	12.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	Acetonitrile, 2-(3-cyanophenyl)-	142	C9H6N2	016532-78-8	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX060618\
Data File : VX002354.D
Acq On : 06 Jun 2018 15:56
Operator : JC/MD
Sample : J3311-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-35

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X060418W.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
Cyclopentane, met...	4.40	19.3	ug/l	254754	1	5.67	661461	50.0
Benzene, cyclopro...	12.30	160.2	ug/l	3407100	4	12.09	1063180	50.0
Benzene, 2-ethyl-...	12.67	31.9	ug/l	678473	4	12.09	1063180	50.0
Benzene, 1-etheny...	12.76	94.1	ug/l	2001040	4	12.09	1063180	50.0
Benzene, 1,2,3,4-...	12.98	39.2	ug/l	833044	4	12.09	1063180	50.0
1-Phenyl-1-butene	13.35	80.5	ug/l	1711190	4	12.09	1063180	50.0
Benzene, (2-methy...	13.99	20.8	ug/l	441991	4	12.09	1063180	50.0
Benzene, (1,2-dim...	14.28	21.3	ug/l	452029	4	12.09	1063180	50.0
Naphthalene, 1-me...	14.85	41.9	ug/l	891180	4	12.09	1063180	50.0