LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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	1.276	16	20	24	rBV	45147	63899	4.50%	0.387%
2	1.325	24	28	35	rVV	16140	45475	3.20%	0.275%
3	1.788	101	104	110	rVB	29538	43347	3.05%	0.263%
4	2.001	135	139	147	rVB2	19832	34904	2.46%	0.211%
5	2.269	177	183	193	rVB	26926	50287	3.54%	0.305%
6	2.446	207	212	216	rBV	28645	45367	3.19%	0.275%
7	2.605	229	238	249	rBV	27921	61223	4.31%	0.371%
8	2.843	269	277	280	rBV3	16902	37722	2.65%	0.229%
9	2.891	280	285	289	rVV2	47622	104489	7.35%	0.633%
10	2.934	289	292	305	rVB2	43302	100371	7.06%	0.608%
11	3.172	323	331	341	rBV	49489	119596	8.41%	0.725%
12	3.397	361	368	377	rBV3	14324	36663	2.58%	0.222%
13	3.525	382	389	399	rVB	19368	42947	3.02%	0.260%
14	3.818	425	437	447	rBV2	52078	130592	9.19%	0.791%
15	3.928	447	455	462	rVV2	34160	94731	6.66%	0.574%
16	3.995	462	466	475	rVV	16961	41504	2.92%	0.251%
17	4.092	475	482	492	rVV5	13200	42480	2.99%	0.257%
18	4.196	492	499	509	rVV	39172	107709	7.58%	0.652%
19	4.318	509	519	523	rVV4	9521	29936	2.11%	0.181%
20	4.403	523	533	546	rVV	158459	471025	33.14%	2.853%
21	4.531	546	554	566	rVB3	18562	55682	3.92%	0.337%
22	4.678	571	578	589	rVB	11117	33174	2.33%	0.201%
23	5.306	669	681	697	rVB2	69705	226844	15.96%	1.374%
24	5.501	700	713	721	rBV	173447	508579	35.78%	3.081%
25	5.586	721	727	731	rBV3	38750	97134	6.83%	0.588%
26	5.665	732	740	748	rVV	253704	663841	46.70%	4.022%
27	5.933	773	784	796	rBV3	30893	95594	6.73%	0.579%
28	6.061	796	805	813	rBV	155687	413219	29.07%	2.503%
29	6.141	813	818	827	rVB2	33536	81721	5.75%	0.495%
30	6.269	829	839	840	rBV3	29927	65339	4.60%	0.396%
31	6.354	849	853	862	rVB3	52010	134543	9.47%	0.815%
32	6.458	862	870	881	rVB2	31055	91085	6.41%	0.552%
33	6.702	902	910	919	rVB3	15180	41167	2.90%	0.249%
34	6.860	924	936	944	rBV	432542	1073719	75.54%	6.505%

LSC Area Percent Report

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Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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	6.939	946	949	960	rVB5	59413	147774	10.40%	0.895%
36	7.067	960	970	977	rBV3	11932	28007	1.97%	0.170%
37	7.147	977	983	991	rBV5	13697	29838	2.10%	0.181%
38	7.256	994	1001	1010	rVB	46620	98985	6.96%	0.600%
39	7.464	1024	1035	1045	rVB4	123719	337916	23.77%	2.047%
40	7.732	1073	1079	1091	rVB2	27919	58237	4.10%	0.353%
41	7.927	1104	1111	1113	rBV2	23737	43034	3.03%	0.261%
42	7.957	1113	1116	1123	rVV3	27436	52074	3.66%	0.315%
43	8.055	1123	1132	1144	rVB2	30470	82929	5.83%	0.502%
44	8.177	1144	1152	1154	rBV	47123	89884	6.32%	0.545%
45	8.214	1154	1158	1163	rVV	104128	184201	12.96%	1.116%
46	8.262	1163	1166	1176	rVV4	19546	38501	2.71%	0.233%
47	8.573	1210	1217	1225	rVB	75443	132222	9.30%	0.801%
48	8.671	1225	1233	1234	rBV3	22970	42948	3.02%	0.260%
49	8.713	1234	1240	1248	rVV	920965	1421414	100.00%	8.611%
50	8.902	1264	1271	1278	rVB4	17275	39726	2.79%	0.241%
51	9.219	1319	1323	1328	rVV	33407	51404	3.62%	0.311%
52	9.622	1385	1389	1392	rVV	22271	29538	2.08%	0.179%
53	10.110	1459	1469	1475	rBV	897517	1200546	84.46%	7.273%
54	10.195	1479	1483	1488	rVV4	19718	40181	2.83%	0.243%
55	10.359	1503	1510	1516	rVV2	64618	130062	9.15%	0.788%
56	10.512	1531	1535	1541	rVB2	29843	44261	3.11%	0.268%
57	10.652	1545	1558	1564	rBV5	48335	151006	10.62%	0.915%
58	10.707	1564	1567	1576	rVB2	16110	26888	1.89%	0.163%
59	10.835	1576	1588	1593	rBV2	90706	150077	10.56%	0.909%
60	10.908	1596	1600	1608	rBV4	44931	88345	6.22%	0.535%
61	11.018	1613	1618	1622	rBV	148064	186319	13.11%	1.129%
62	11.134	1632	1637	1642	rBV	754254	930192	65.44%	5.635%
63	11.183	1642	1645	1649	rVB	40917	47661	3.35%	0.289%
64	11.274	1657	1660	1665	rVB2	43484	56011	3.94%	0.339%
65	11.359	1669	1674	1683	rVB	247896	311652	21.93%	1.888%
66	11.451	1683	1689	1693	rBV2	32244	55019	3.87%	0.333%
67	11.670	1719	1725	1732	rBV2	64074	99031	6.97%	0.600%
68	11.768	1732	1741	1744	rBV2	21836	35620	2.51%	0.216%
69	11.804	1744	1747	1759	rVB2	18938	34383	2.42%	0.208%
70	11.914	1759	1765	1767	rBV3	34844	51348	3.61%	0.311%
71	11.945	1767	1770	1775	rVB	49324	62010	4.36%	0.376%

LSC Area Percent Report

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ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061919W.M
Title : SW846 8260

72	12.079	1786	1792	1797	rBV	898285	1152096	81.05%	6.979%
73	12.231	1812	1817	1820	rBV	36664	42889	3.02%	0.260%
74	12.298	1820	1828	1835	rVV	624249	835692	58.79%	5.063%
75	12.365	1835	1839	1848	rVB2	63327	126382	8.89%	0.766%
76	12.457	1849	1854	1859	rBV	48056	62633	4.41%	0.379%
77	12.518	1859	1864	1869	rVB2	27092	41221	2.90%	0.250%
78	12.664	1881	1888	1891	rVV	252170	312210	21.96%	1.891%
79	12.701	1891	1894	1899	rVV2	83688	121184	8.53%	0.734%
80	12.755	1899	1903	1910	rVV2	218431	355327	25.00%	2.153%
81	12.896	1923	1926	1931	rVB	41819	53274	3.75%	0.323%
82	12.975	1931	1939	1944	rBV	147056	196411	13.82%	1.190%
83	13.085	1952	1957	1962	rVB2	19250	32705	2.30%	0.198%
84	13.194	1971	1975	1977	rVV2	27022	36437	2.56%	0.221%
85	13.225	1977	1980	1988	rVB	51186	75773	5.33%	0.459%
86	13.341	1995	1999	2006	rVV2	280328	405154	28.50%	2.454%
87	13.426	2010	2013	2016	rVV2	22124	28768	2.02%	0.174%
88	13.481	2018	2022	2029	rVB3	20612	32612	2.29%	0.198%
89	13.597	2038	2041	2044	rVV	19183	27105	1.91%	0.164%
90	13.639	2044	2048	2052	rVV3	42084	67924	4.78%	0.411%
91	13.694	2052	2057	2059	rVV4	22647	39269	2.76%	0.238%
92	13.719	2059	2061	2071	rVB2	34940	47687	3.35%	0.289%
93	13.828	2072	2079	2088	rVB	77762	130684	9.19%	0.792%
94	13.908	2088	2092	2098	rBV2	24155	36323	2.56%	0.220%
95	13.981	2098	2104	2108	rBV2	23546	34443	2.42%	0.209%
96	14.286	2147	2154	2158	rBV3	29421	58154	4.09%	0.352%
97	14.536	2191	2195	2200	rBV2	32627	44452	3.13%	0.269%
98	14.840	2239	2245	2250	rBV	110973	148037	10.41%	0.897%
99	15.688	2374	2384	2387	rBV	19456	36661	2.58%	0.222%
100	16.413	2495	2503	2510	rBV4	14035	30617	2.15%	0.185%

Sum of corrected areas: 16507276

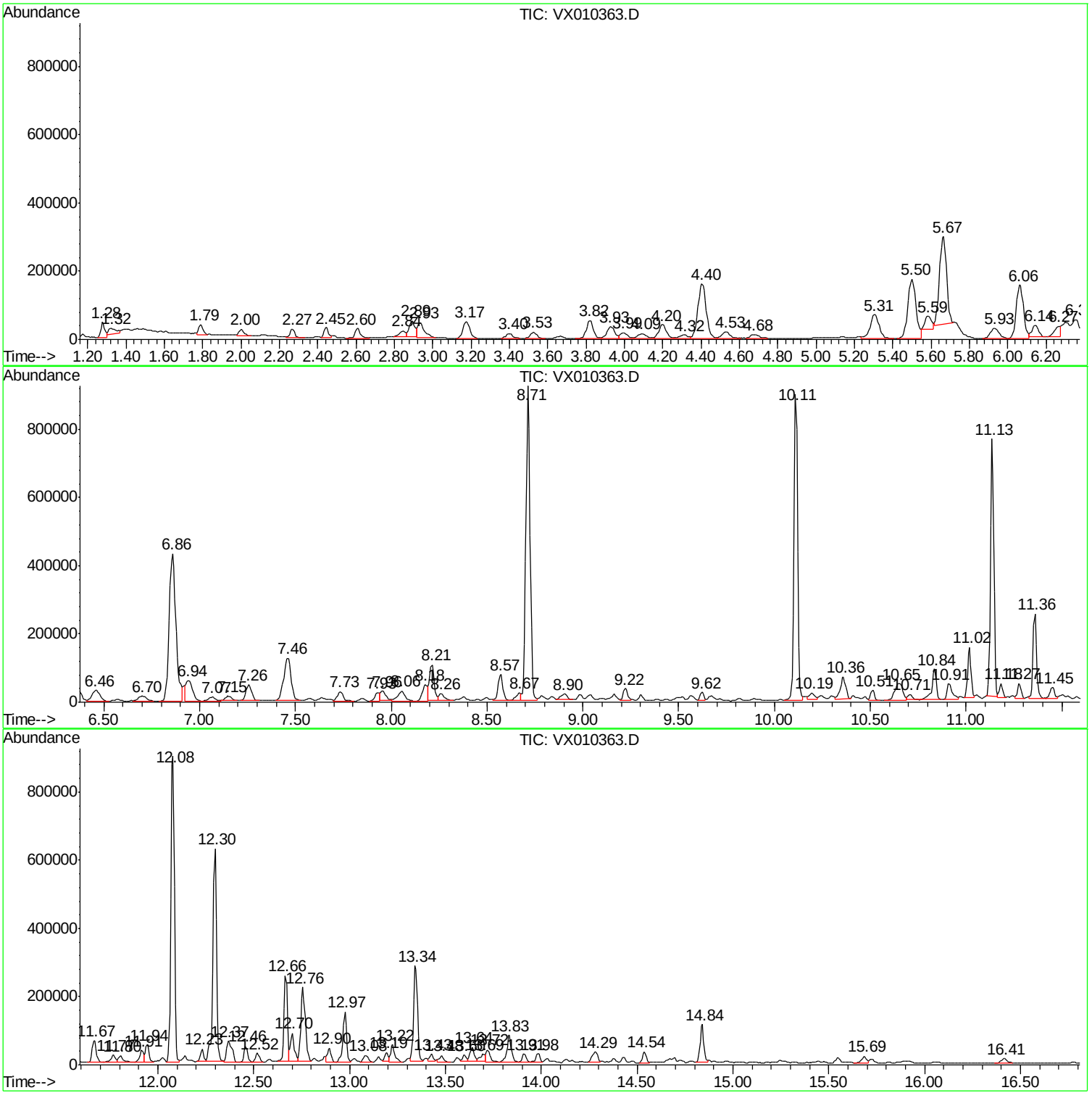
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
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ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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

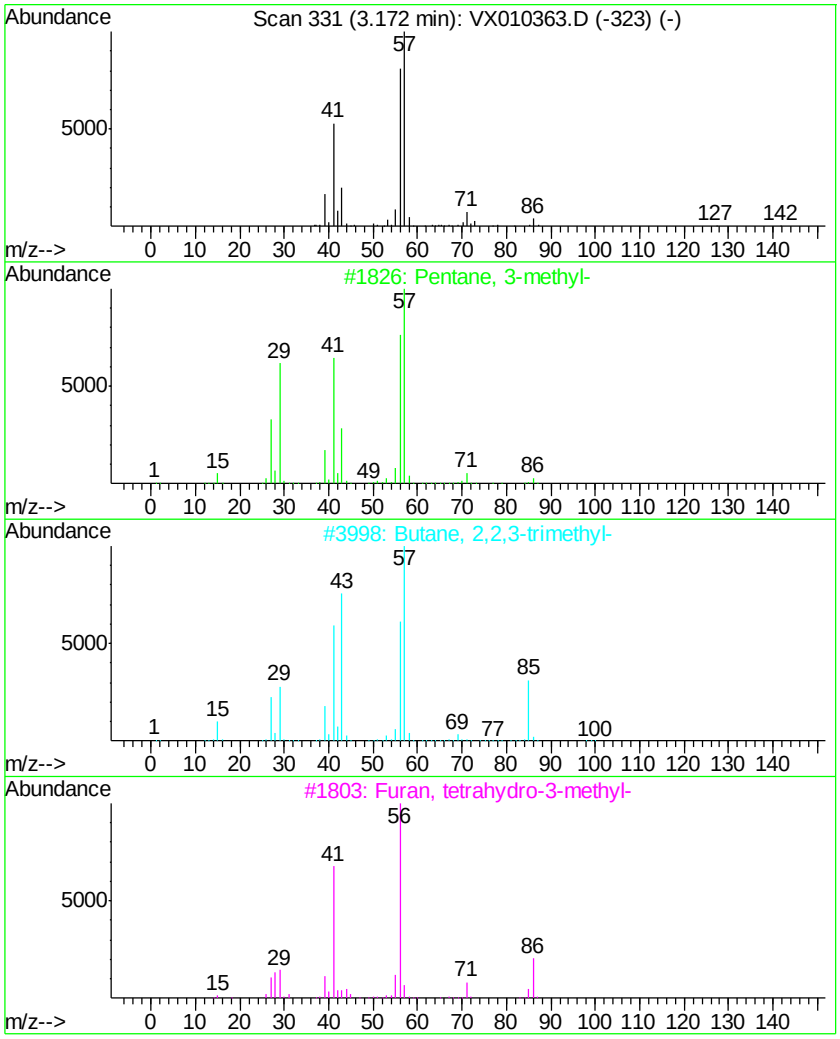
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

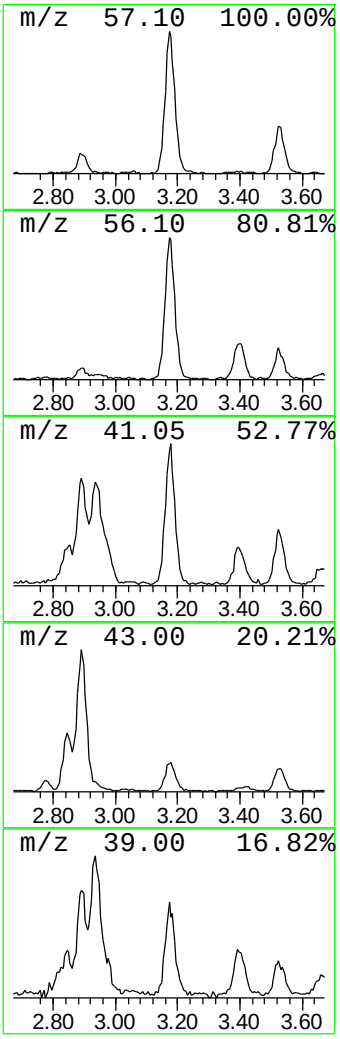
****

Peak Number 1 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.17	9.01 ug/l	119596	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	72
3	Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	64
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
5	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062019\
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Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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

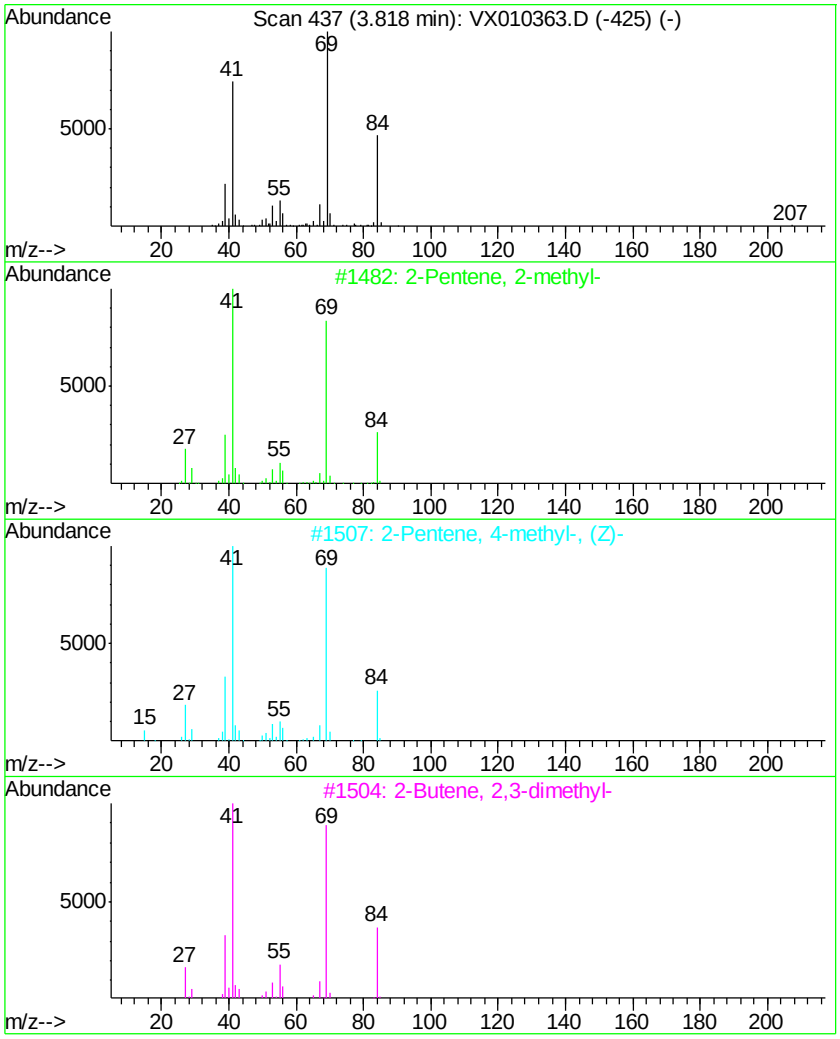
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

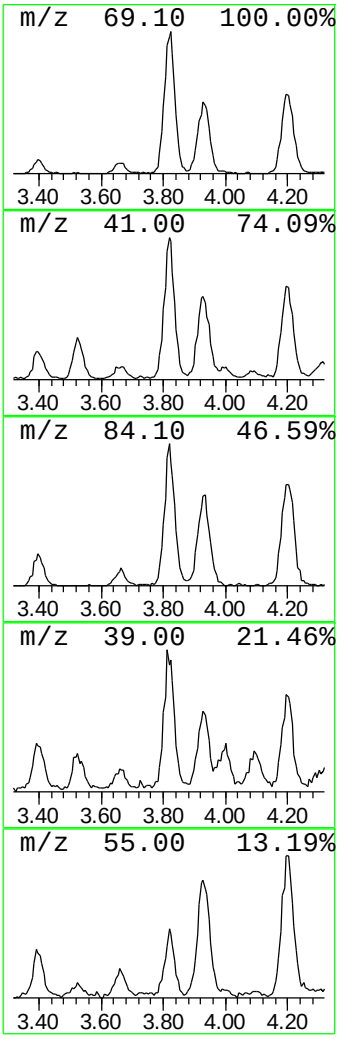
****

Peak Number 2 2-Pentene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.82	9.84 ug/l	130592	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 2-methyl-	84	C6H12	000625-27-4	91
2	2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	91
3	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	91
4	2-Pentene, 4-methyl-	84	C6H12	004461-48-7	91
5	Cyclopropane, 1,1,2-trimethyl-	84	C6H12	004127-45-1	86





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062019\
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Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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

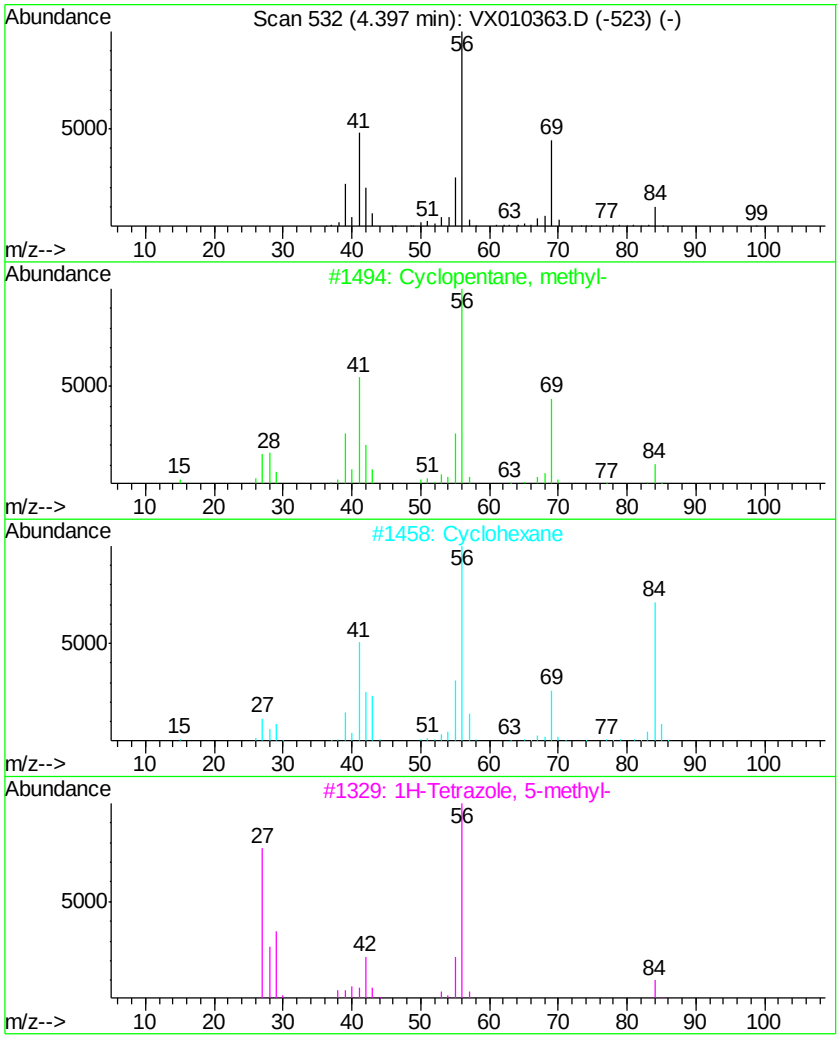
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

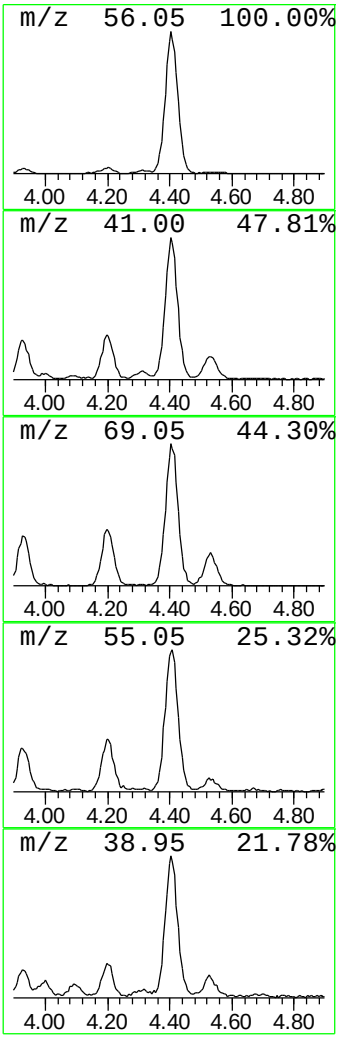
****

Peak Number 3 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.40	35.48 ug/l	471025	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclohexane	84	C6H12	000110-82-7	83
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	64
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062019\
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ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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

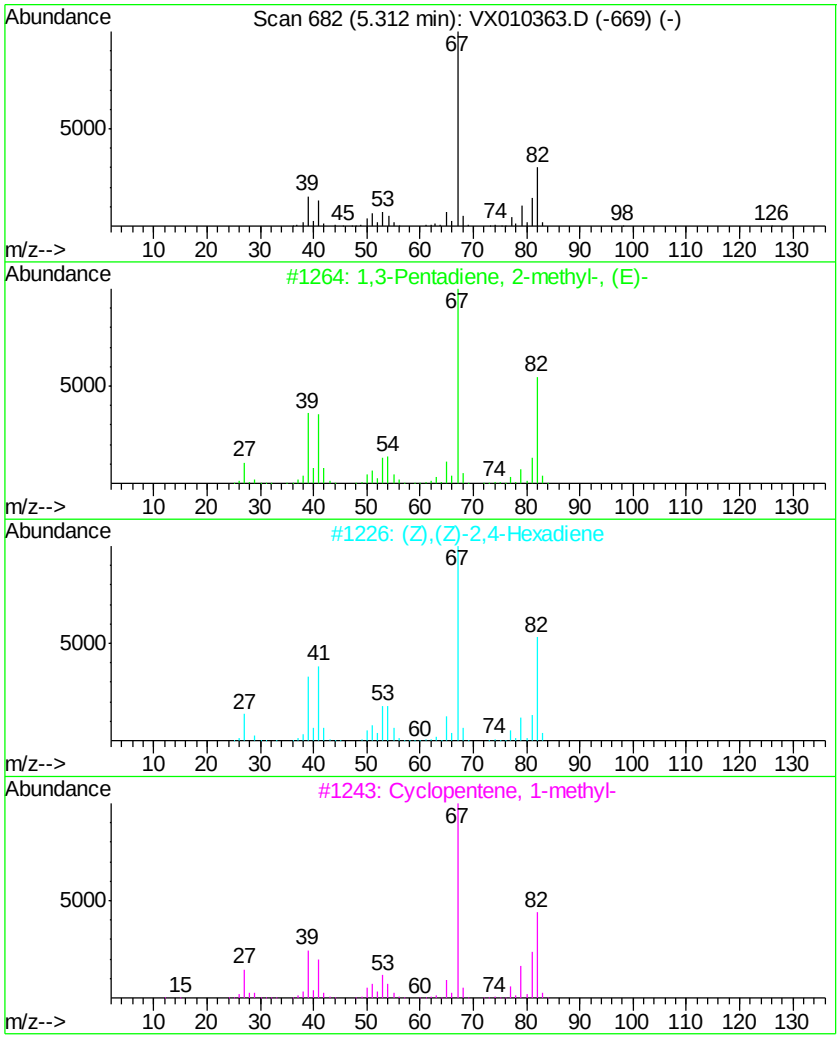
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

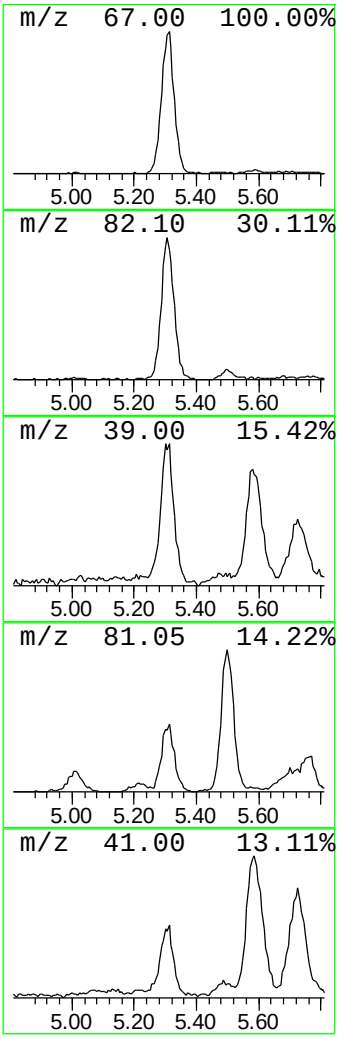
****

Peak Number 4 1,3-Pentadiene, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.31	17.09 ug/l	226844	Pentafluorobenzene	5.67

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Pentadiene, 2-methyl-, (E)-	82	C6H10	000926-54-5	83
2	(Z),(Z)-2,4-Hexadiene	82	C6H10	006108-61-8	83
3	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	81
4	Cyclopentene, 4-methyl-	82	C6H10	001759-81-5	80
5	2,4-Hexadiene, (E,E)-	82	C6H10	005194-51-4	80





Library Search Compound Report

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Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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

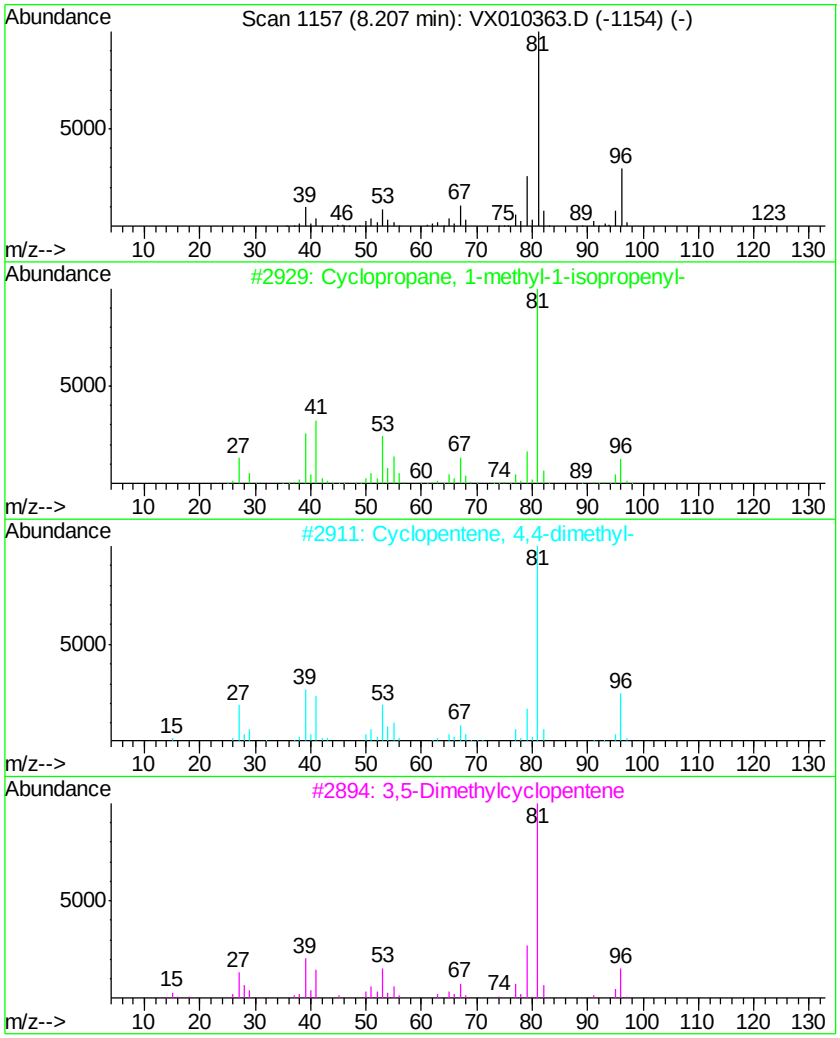
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

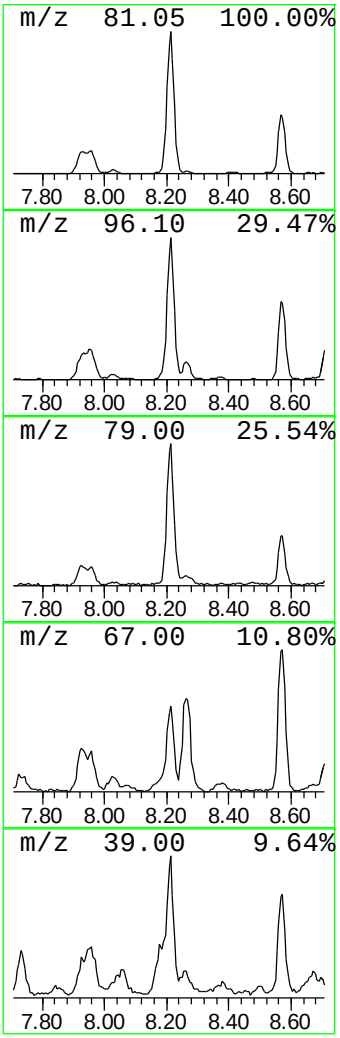
****

Peak Number 5 Cyclopropane, 1-methyl-1-is... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.21	8.58 ug/l	184201	1,4-Difluorobenzene	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-methyl-1-isoprop...	96	C7H12	003422-07-9	72
2	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	72
3	3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	72
4	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	72
5	Furan, 2-ethyl-	96	C6H8O	003208-16-0	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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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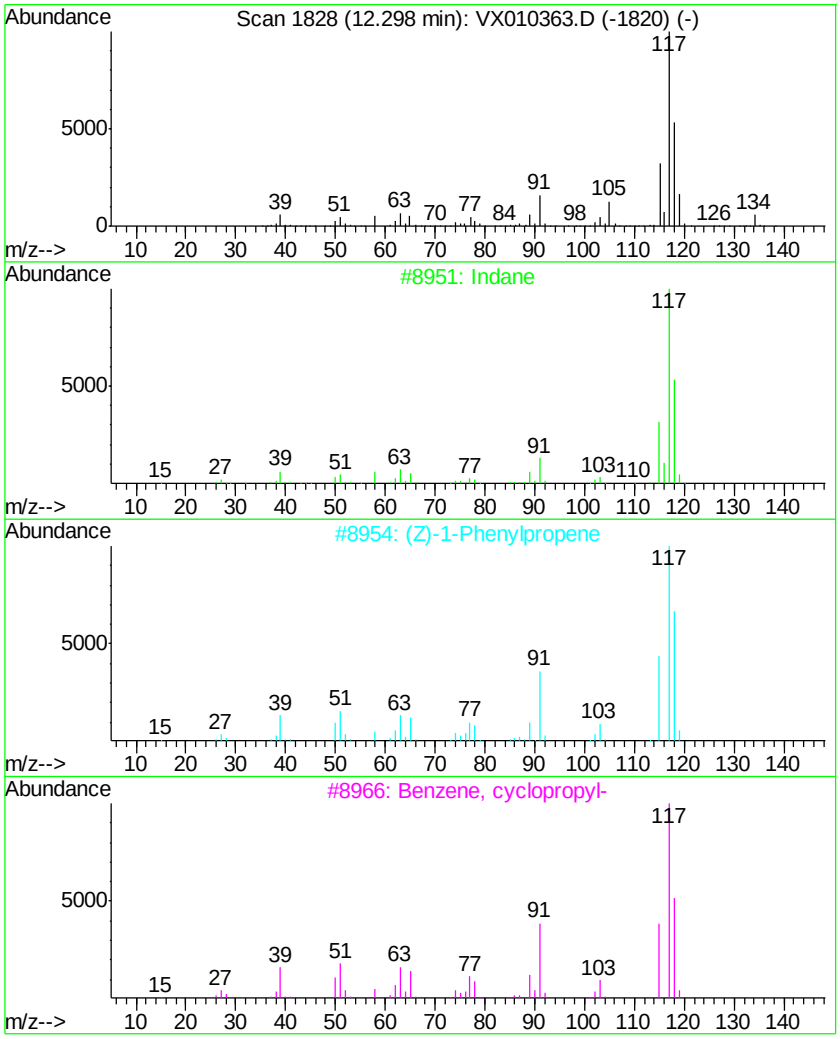
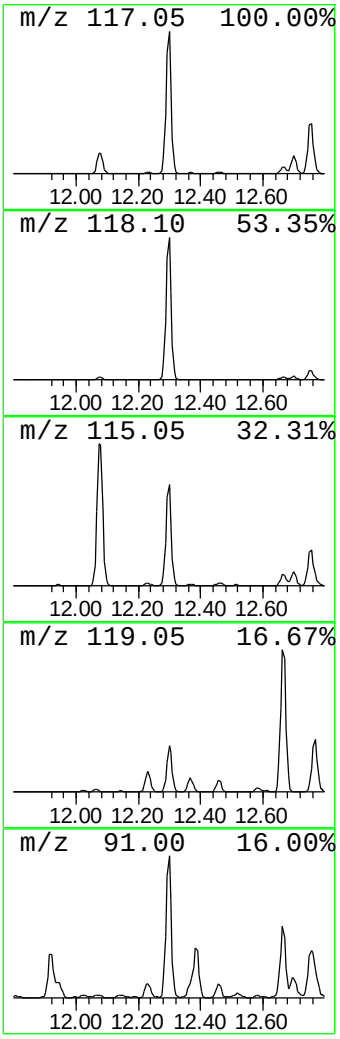
Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	36.27 ug/l	835692	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	76
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	68
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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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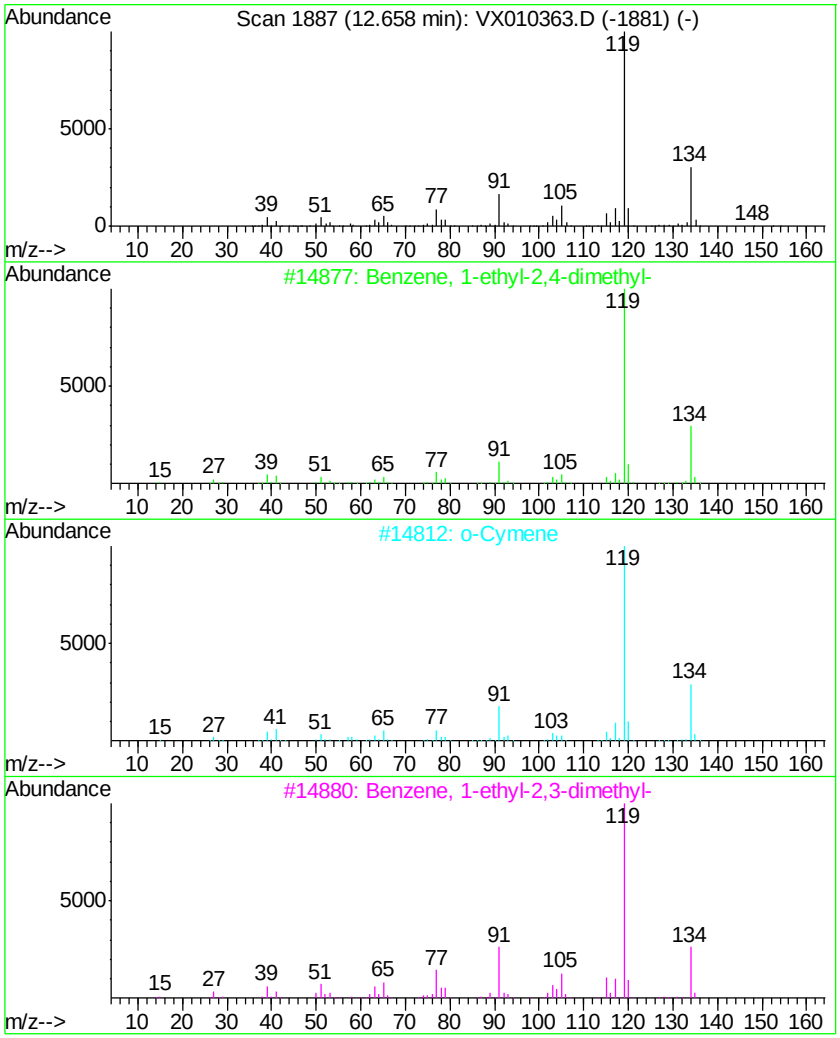
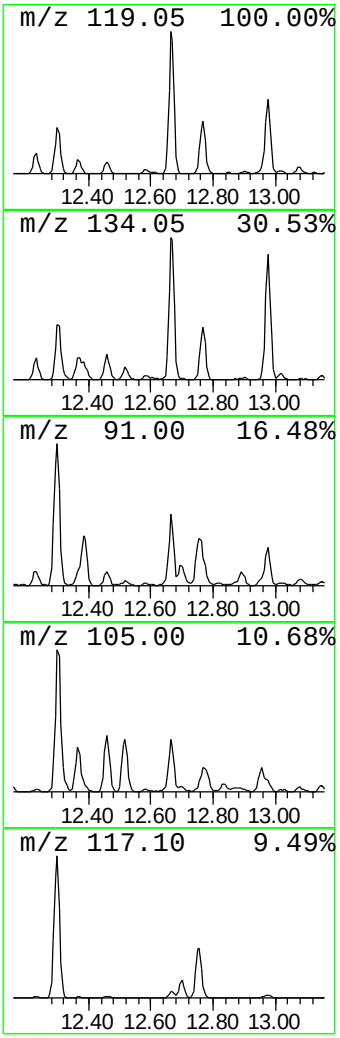
Peak Number 7 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	13.55 ug/l	312210	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	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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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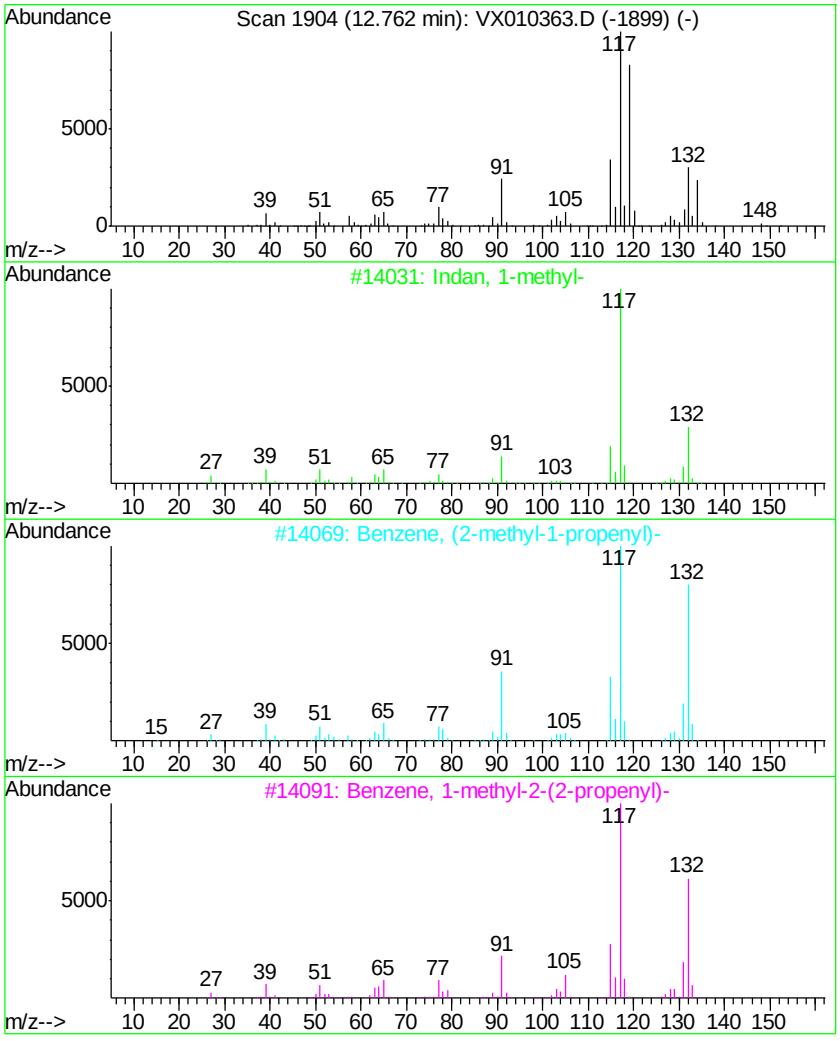
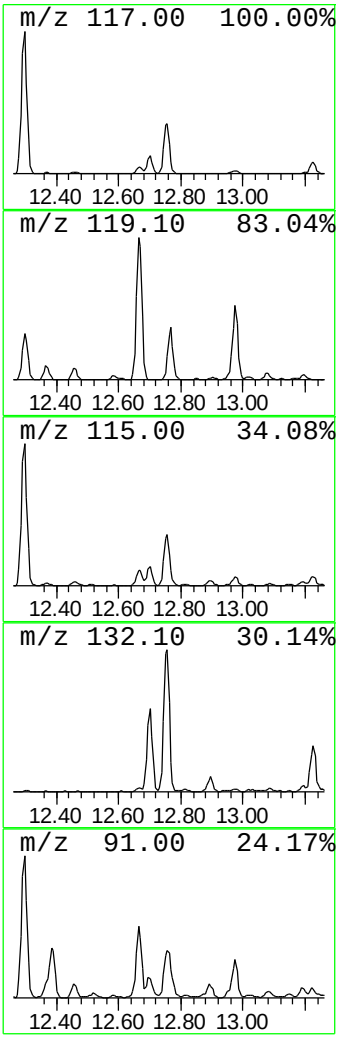
Peak Number 8 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	15.42 ug/l	355327	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, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	74
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	74





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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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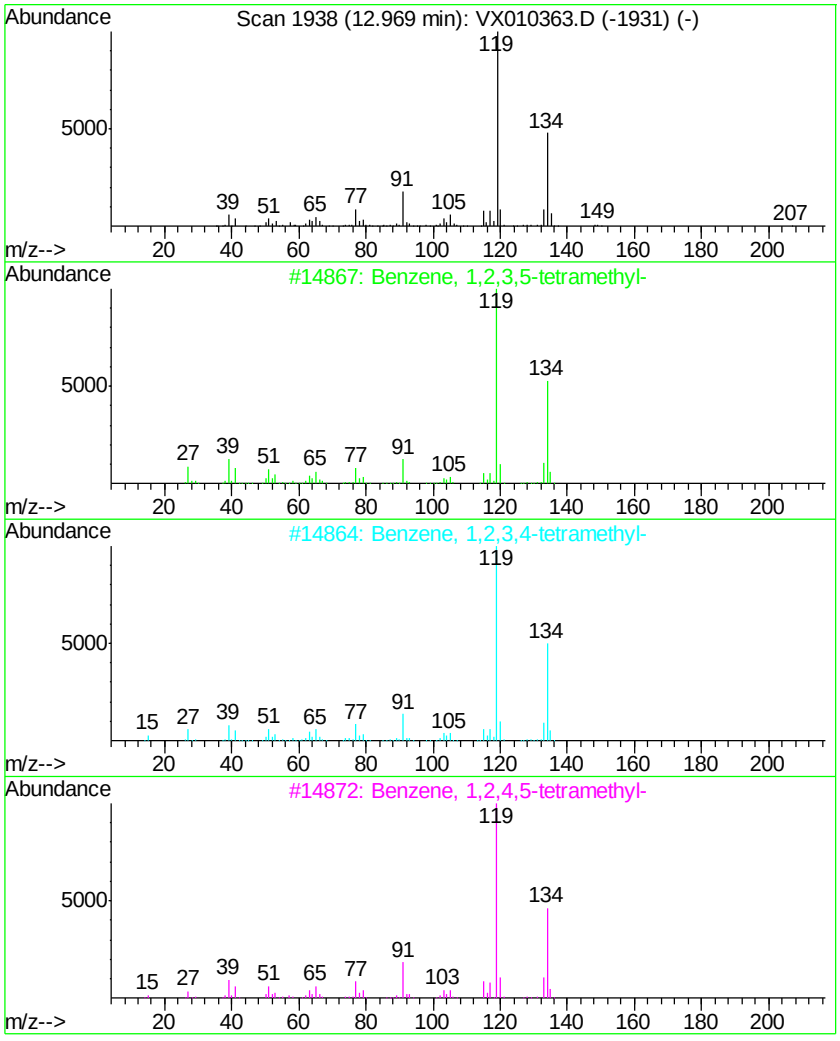
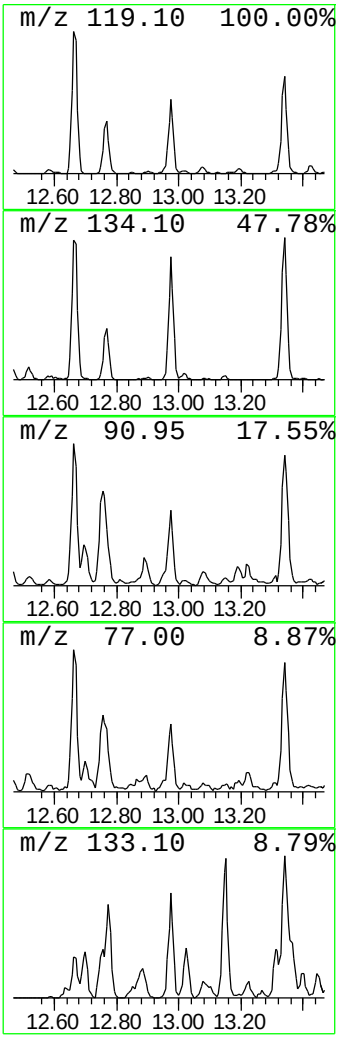
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	8.52 ug/l	196411	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX062019\
Data File : VX010363.D
Acq On : 20 Jun 2019 17:35
Operator : JC/SP
Sample : K3469-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : T3-MW-1-9.0-061919

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
Pentane, 3-methyl-	3.17	9.0	ug/l	119596	1	5.67	663841	50.0
2-Pentene, 2-methyl-	3.82	9.8	ug/l	130592	1	5.67	663841	50.0
Cyclopentane, met...	4.40	35.5	ug/l	471025	1	5.67	663841	50.0
1,3-Pentadiene, 2...	5.31	17.1	ug/l	226844	1	5.67	663841	50.0
Cyclopropane, 1-m...	8.21	8.6	ug/l	184201	2	6.86	1073720	50.0
Indane	12.30	36.3	ug/l	835692	4	12.08	1152100	50.0
Benzene, 1-ethyl-...	12.66	13.6	ug/l	312210	4	12.08	1152100	50.0
Indan, 1-methyl-	12.76	15.4	ug/l	355327	4	12.08	1152100	50.0
Benzene, 1,2,3,5-...	12.97	8.5	ug/l	196411	4	12.08	1152100	50.0