LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW101718\
Data File : VW006149.D
Acq On : 18 Oct 2018 03:13
Operator : VA/AP
Sample : J5579-01
Misc : 5.26G/5ML/MSVOA W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

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_W\METHOD\82W101518S.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	4.916	483	494	513	rVB	15865	53810	6.80%	0.857%
2	7.159	852	862	873	rBV	5749	18227	2.30%	0.290%
3	7.884	968	981	986	rBV	144273	340225	43.02%	5.418%
4	7.952	986	992	1003	rVB	285200	629504	79.60%	10.025%
5	8.311	1041	1051	1061	rBV	130744	292863	37.03%	4.664%
6	8.842	1129	1138	1151	rBV	400095	790872	100.00%	12.595%
7	10.329	1373	1382	1388	rBV	431313	775432	98.05%	12.349%
8	10.390	1388	1392	1398	rVB	15793	26704	3.38%	0.425%
9	11.634	1588	1596	1604	rBV	417571	686676	86.83%	10.935%
10	11.731	1606	1612	1620	rVV	46153	78538	9.93%	1.251%
11	11.841	1620	1630	1641	rVV	14801	33435	4.23%	0.532%
12	12.170	1672	1684	1690	rBV	18772	36327	4.59%	0.579%
13	12.371	1707	1717	1723	rBV3	20060	38525	4.87%	0.614%
14	12.469	1728	1733	1738	rBV2	12669	20397	2.58%	0.325%
15	12.524	1738	1742	1748	rVB2	29840	47156	5.96%	0.751%
16	12.621	1751	1758	1765	rBV	244399	399186	50.47%	6.357%
17	12.725	1765	1775	1781	rVB	36977	69513	8.79%	1.107%
18	12.810	1781	1789	1794	rVB	10145	21764	2.75%	0.347%
19	12.884	1794	1801	1804	rBV6	6750	16750	2.12%	0.267%
20	12.951	1804	1812	1823	rVB9	17397	58752	7.43%	0.936%
21	13.054	1823	1829	1834	rBV	47936	88535	11.19%	1.410%
22	13.109	1834	1838	1844	rVB	14322	24857	3.14%	0.396%
23	13.262	1858	1863	1869	rBV2	14272	28191	3.56%	0.449%
24	13.390	1878	1884	1889	rBV7	7102	15852	2.00%	0.252%
25	13.451	1889	1894	1899	rVV2	14869	27913	3.53%	0.445%
26	13.499	1899	1902	1906	rVV	9585	15352	1.94%	0.244%
27	13.560	1906	1912	1922	rVB	229703	411679	52.05%	6.556%
28	13.725	1931	1939	1942	rBV5	12617	26607	3.36%	0.424%
29	13.761	1942	1945	1949	rVV2	14471	21356	2.70%	0.340%
30	13.883	1961	1965	1970	rVB4	19647	34575	4.37%	0.551%
31	13.957	1972	1977	1982	rBV	17410	27083	3.42%	0.431%
32	14.018	1982	1987	1990	rBV3	9148	17307	2.19%	0.276%
33	14.048	1990	1992	1995	rVV2	8566	9925	1.25%	0.158%
34	14.103	1995	2001	2008	rVB	40038	67049	8.48%	1.068%

LSC Area Percent Report

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ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

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_W\METHOD\82W101518S.M
Title : SW846 8260

35	14.164	2008	2011	2014	rBV	6698	10534	1.33%	0.168%
36	14.213	2014	2019	2025	rVB2	32319	54943	6.95%	0.875%
37	14.292	2025	2032	2036	rVB8	6220	12887	1.63%	0.205%
38	14.347	2036	2041	2044	rBV2	13728	21967	2.78%	0.350%
39	14.396	2045	2049	2050	rBV3	7803	9232	1.17%	0.147%
40	14.426	2050	2054	2058	rVV2	19755	28987	3.67%	0.462%
41	14.463	2058	2060	2064	rVB	15036	20016	2.53%	0.319%
42	14.511	2064	2068	2071	rBV2	11960	16949	2.14%	0.270%
43	14.548	2071	2074	2081	rBV7	6217	9878	1.25%	0.157%
44	14.633	2081	2088	2089	rBV4	9338	14844	1.88%	0.236%
45	14.694	2094	2098	2102	rVV3	13656	24292	3.07%	0.387%
46	14.737	2102	2105	2110	rVB2	16462	24599	3.11%	0.392%
47	14.804	2110	2116	2123	rBV3	58170	144601	18.28%	2.303%
48	14.865	2123	2126	2132	rVB3	11681	19039	2.41%	0.303%
49	14.969	2139	2143	2148	rBV2	12889	22243	2.81%	0.354%
50	15.072	2156	2160	2164	rBV3	22701	36261	4.58%	0.577%
51	15.188	2174	2179	2184	rVB3	27694	59593	7.54%	0.949%
52	15.334	2198	2203	2206	rBV5	10226	21189	2.68%	0.337%
53	15.377	2206	2210	2214	rVB2	20413	34637	4.38%	0.552%
54	15.426	2214	2218	2223	rBV6	9175	23146	2.93%	0.369%
55	15.481	2223	2227	2230	rBV2	9734	13671	1.73%	0.218%
56	15.566	2238	2241	2250	rVB7	15654	36389	4.60%	0.579%
57	15.664	2250	2257	2265	rBV5	15005	41034	5.19%	0.653%
58	15.749	2267	2271	2276	rVV7	7114	14761	1.87%	0.235%
59	15.828	2276	2284	2288	rVV5	12864	37236	4.71%	0.593%
60	15.871	2288	2291	2296	rVB3	15039	27090	3.43%	0.431%
61	15.962	2303	2306	2311	rVB5	10971	18205	2.30%	0.290%
62	16.035	2312	2318	2323	rBV5	8900	22868	2.89%	0.364%
63	16.182	2335	2342	2354	rBV7	15791	47544	6.01%	0.757%
64	16.389	2371	2376	2381	rVB6	10841	23152	2.93%	0.369%
65	16.456	2381	2387	2401	rVB	24667	67465	8.53%	1.074%
66	16.657	2412	2420	2426	rBV2	22852	61060	7.72%	0.972%
67	16.767	2434	2438	2441	rBV4	4107	8227	1.04%	0.131%

Sum of corrected areas: 6279476

LSC Report - Integrated Chromatogram

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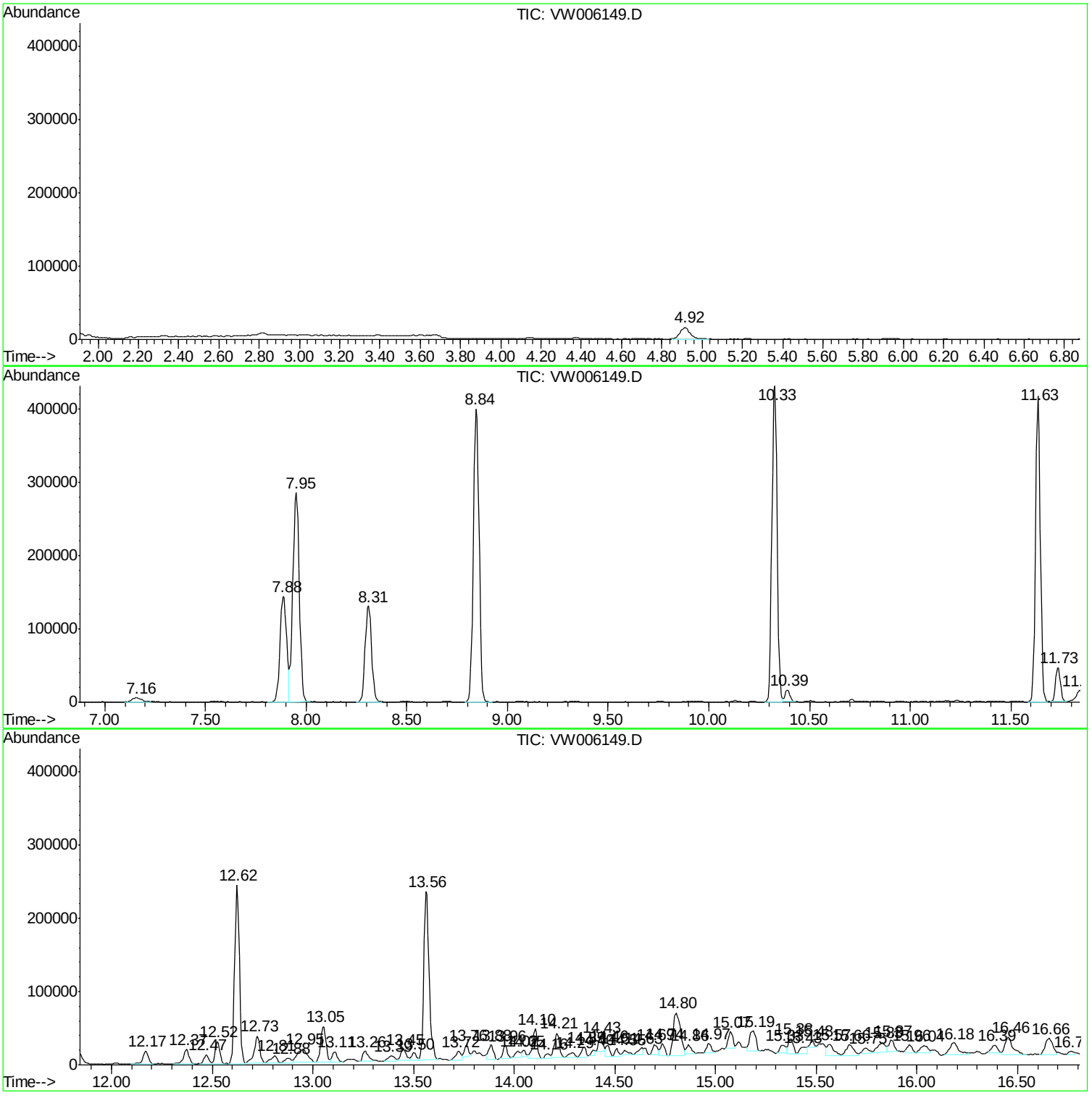
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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



Library Search Compound Report

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Quant Title : SW846 8260

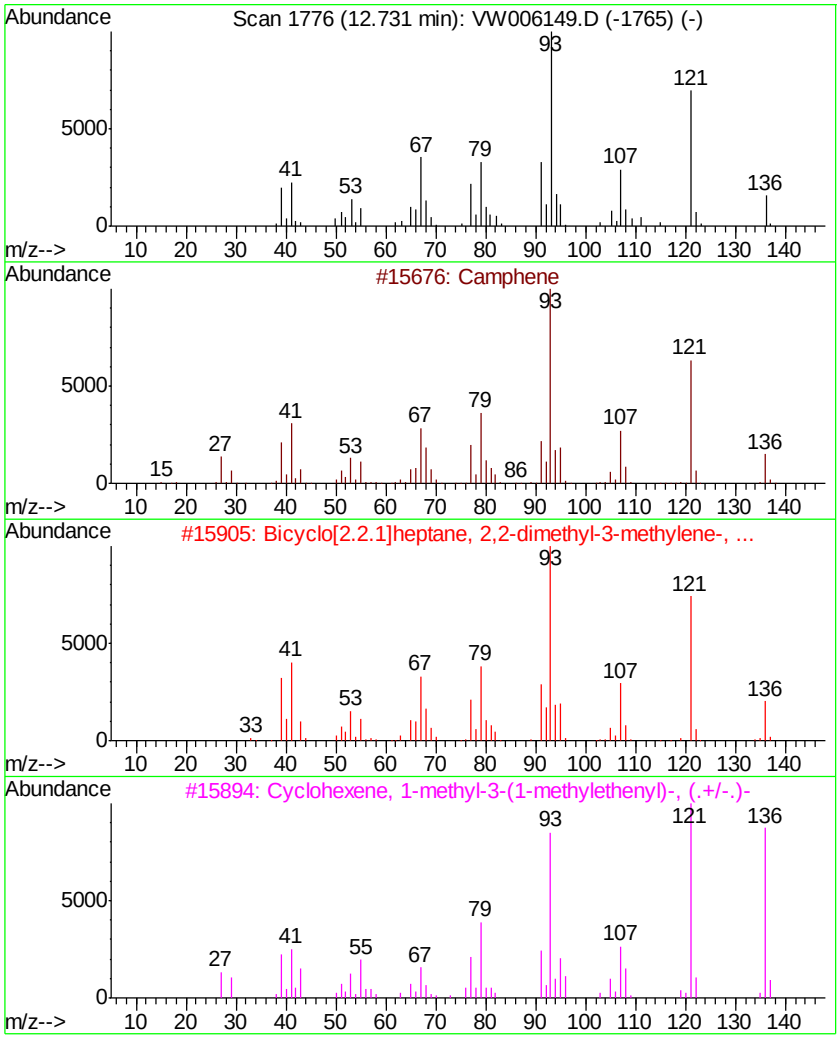
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

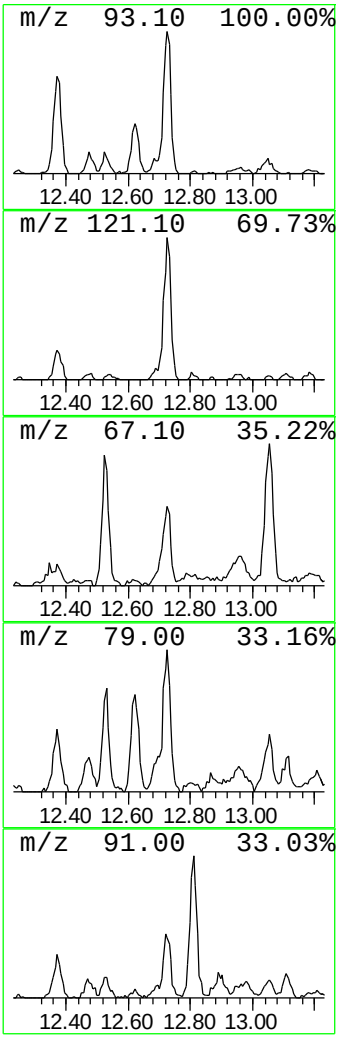
****

Peak Number 1 Camphene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.73	8.44 ug/l	69513	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphene	136	C10H16	000079-92-5	98
2	Bicyclo[2.2.1]heptane, 2,2-dimet...	136	C10H16	005794-04-7	95
3	Cyclohexene, 1-methyl-3-(1-methy...	136	C10H16	000499-03-6	94
4	1,3,6-Heptatriene, 2,5,5-trimethyl-	136	C10H16	029548-02-5	87
5	(+)-4-Carene	136	C10H16	029050-33-7	83





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW101718\
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Sample : J5579-01
Misc : 5.26G/5ML/MSVOA_W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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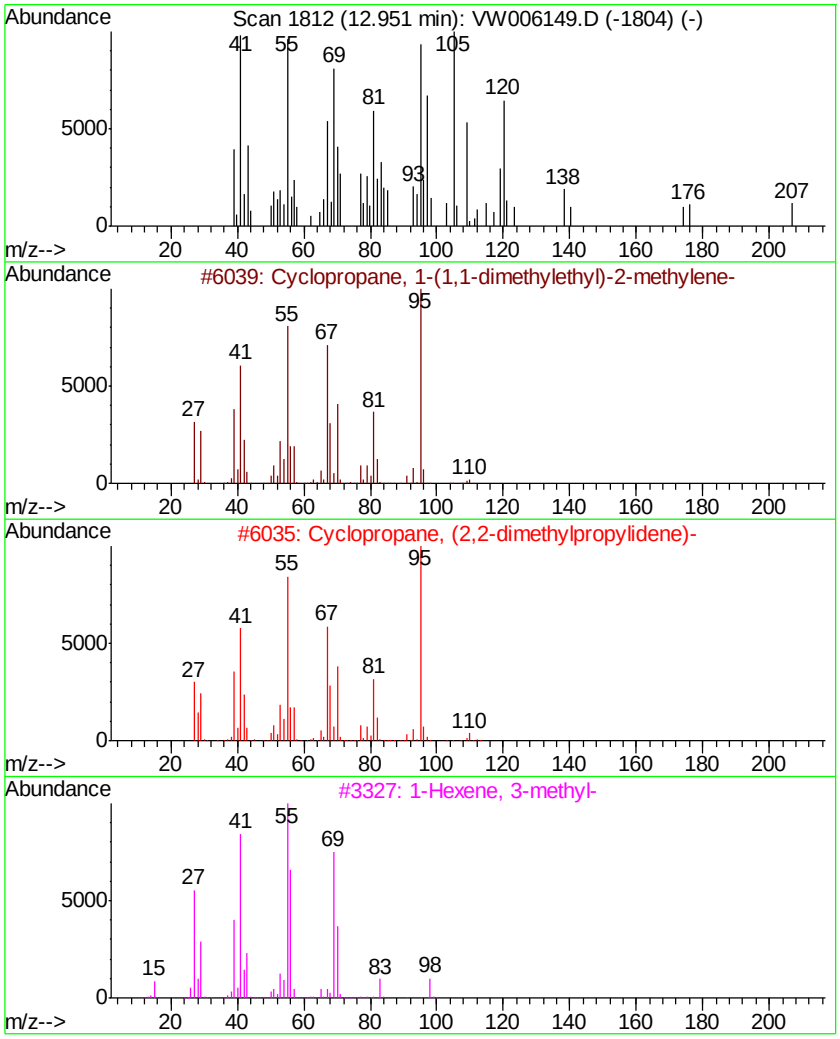
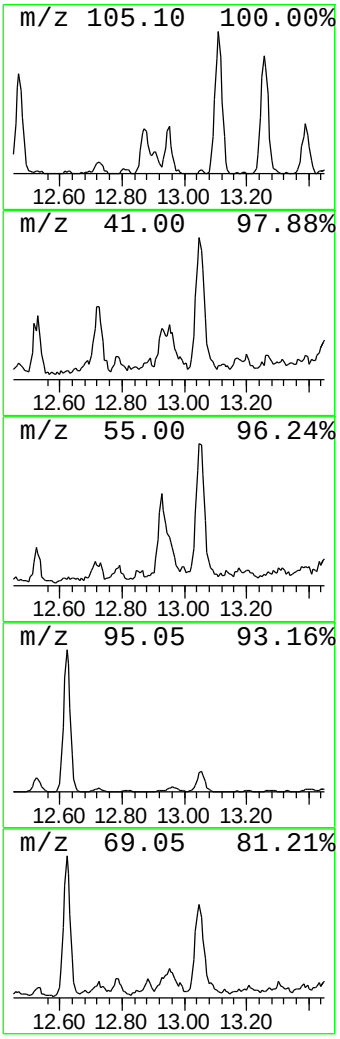
Peak Number 2 unknown12.95 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	7.14 ug/l	58752	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1-(1,1-dimethyleth...	110	C8H14	061142-25-4	42
2	Cyclopropane, (2,2-dimethylpropy...	110	C8H14	039647-71-7	38
3	1-Hexene, 3-methyl-	98	C7H14	003404-61-3	35
4	Cyclopentane, pentyl-	140	C10H20	003741-00-2	25
5	Cyclohexene, 1-methyl-4-(1-methy...	138	C10H18	001195-31-9	25





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Sample : J5579-01
Misc : 5.26G/5ML/MSVOA_W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

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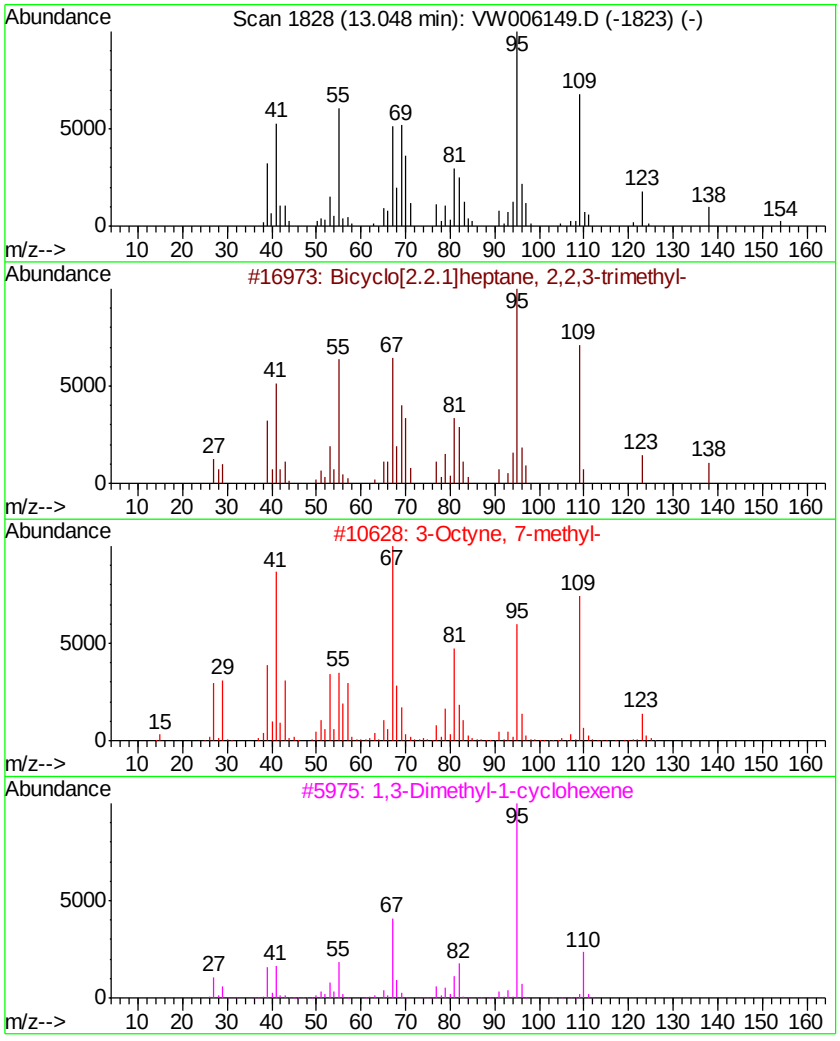
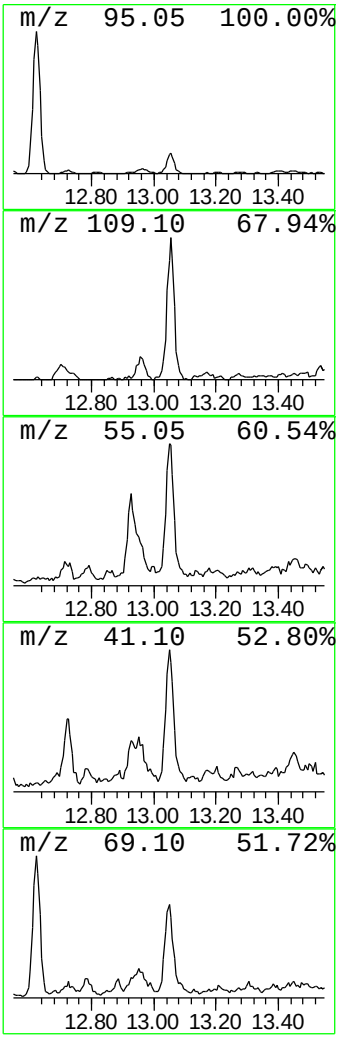
Peak Number 3 Bicyclo[2.2.1]heptane, 2,2,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.05	10.75 ug/l	88535	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	000473-19-8	96
2	3-Octyne, 7-methyl-	124	C9H16	037050-06-9	58
3	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	43
4	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	41
5	Cyclopentane, (2-methylbutylidene)-	138	C10H18	053366-54-4	38





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Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
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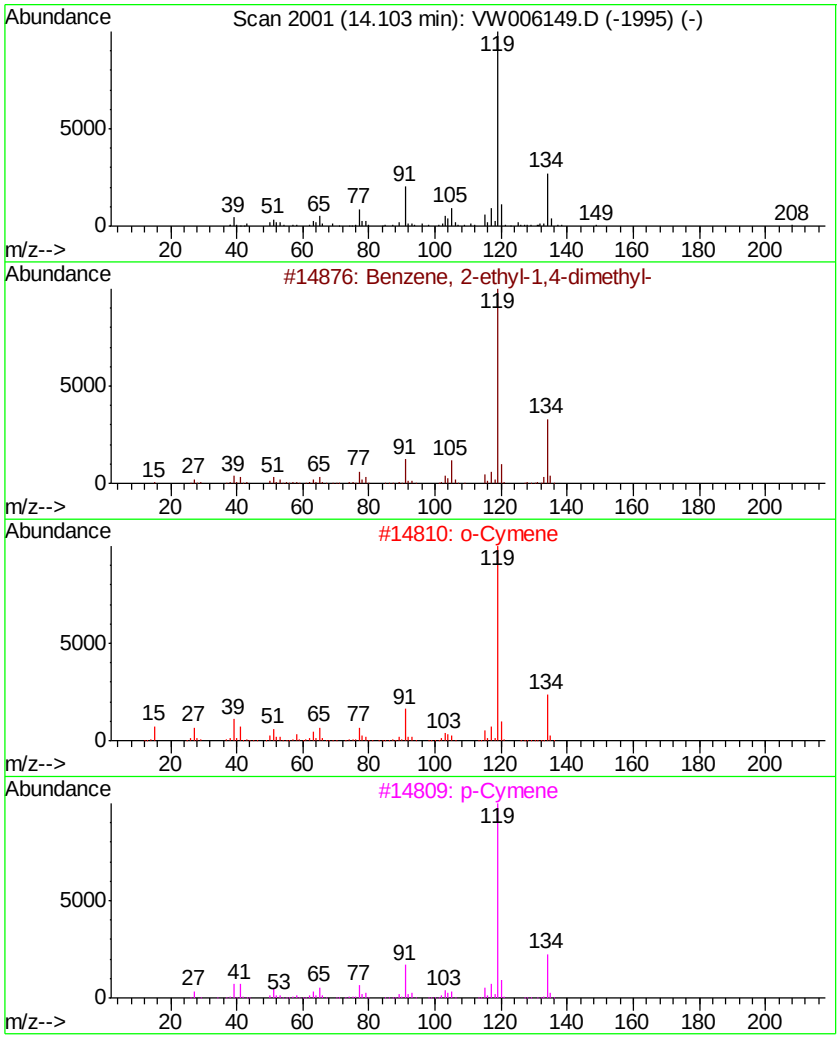
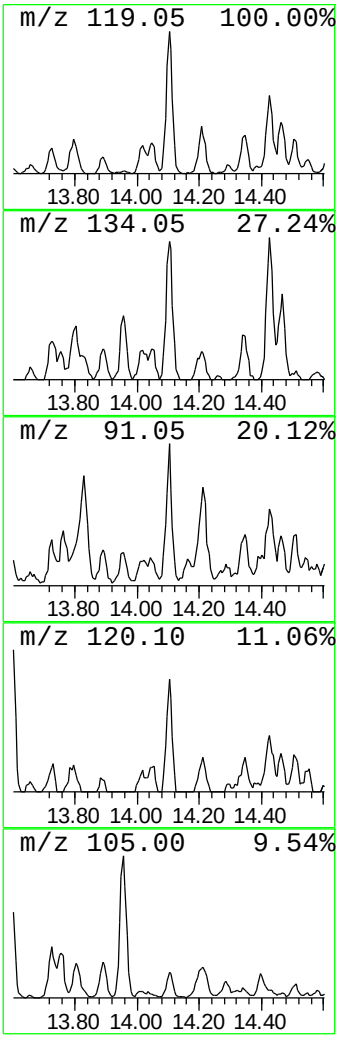
****

Peak Number 4 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.10	8.14 ug/l	67049	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
2	o-Cymene	134	C10H14	000527-84-4	94
3	p-Cymene	134	C10H14	000099-87-6	94
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91





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ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

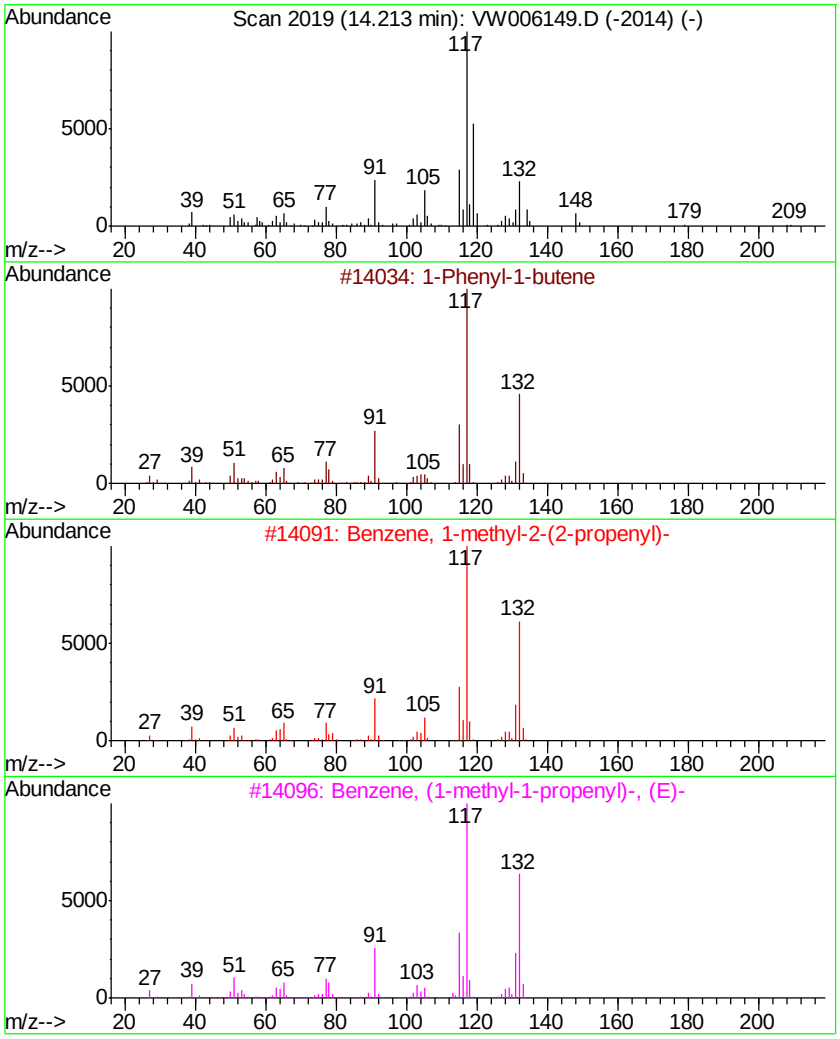
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

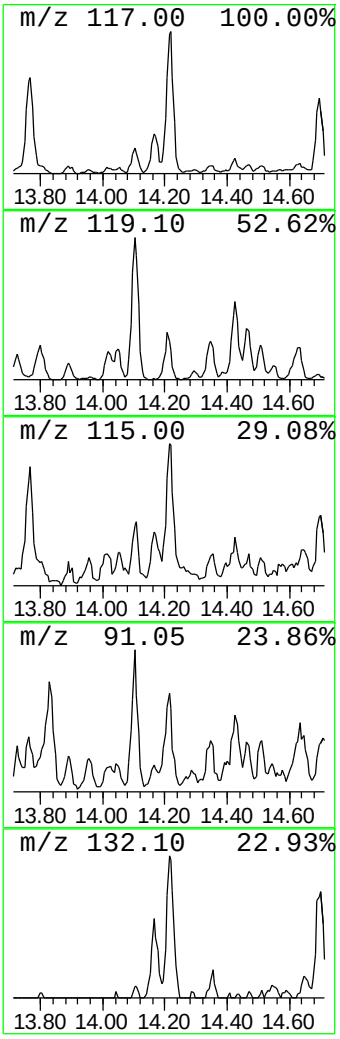
****

Peak Number 5 1-Phenyl-1-butene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	6.67 ug/l	54943	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	70
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	62
3	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	62
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	62
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	62





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Quant Title : SW846 8260

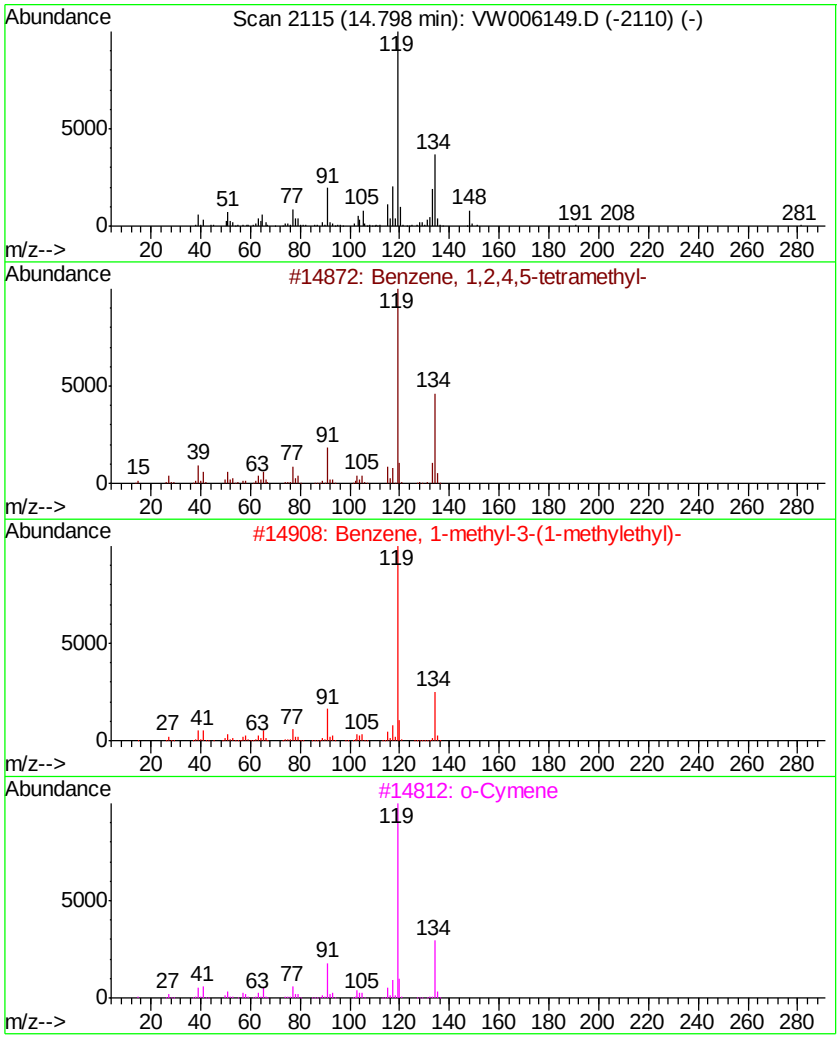
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

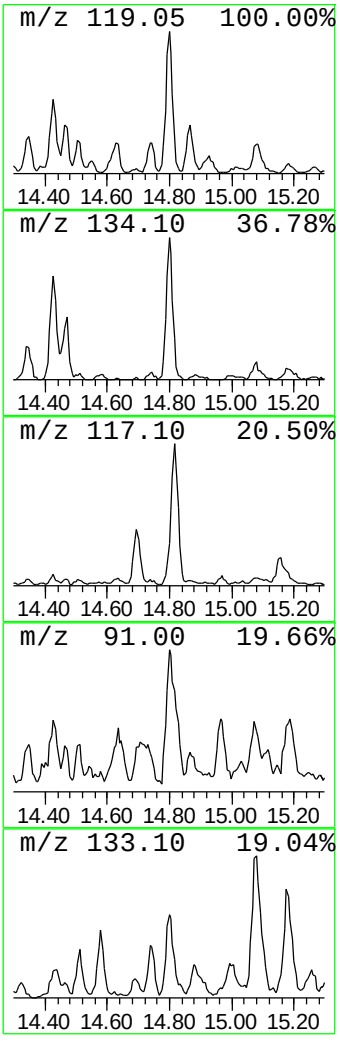
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	17.56 ug/l	144601	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
3	o-Cymene	134	C10H14	000527-84-4	64
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	60
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	60





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ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

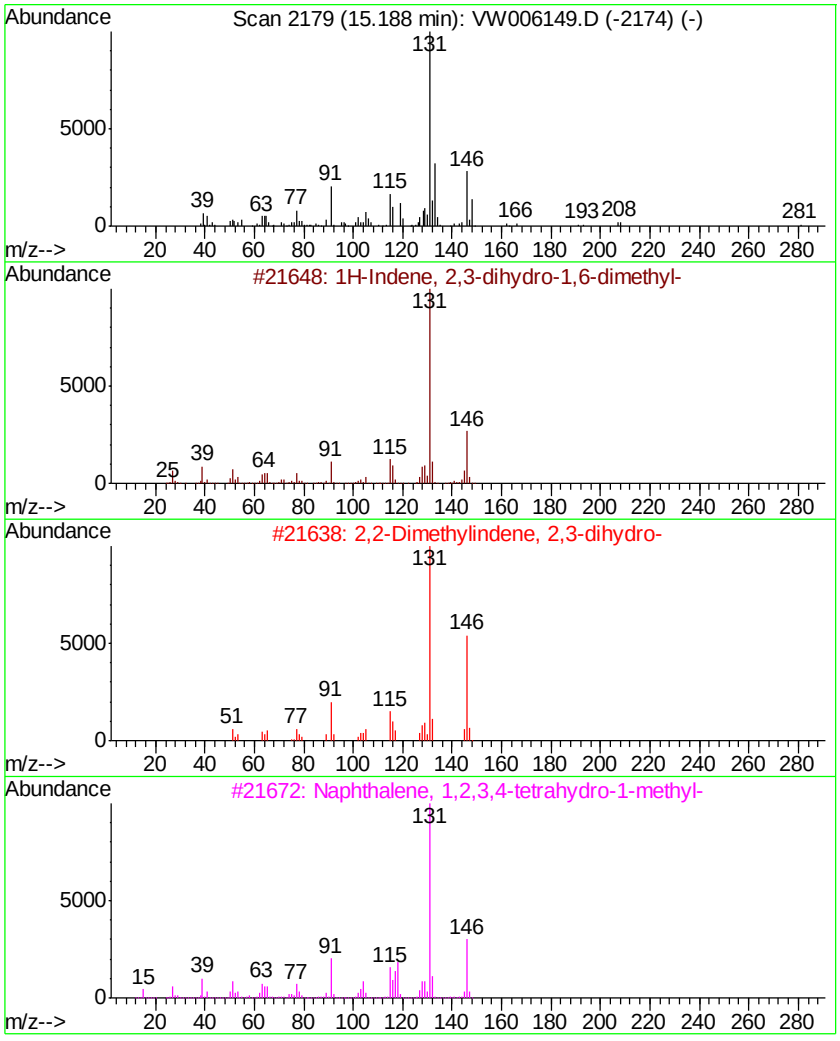
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

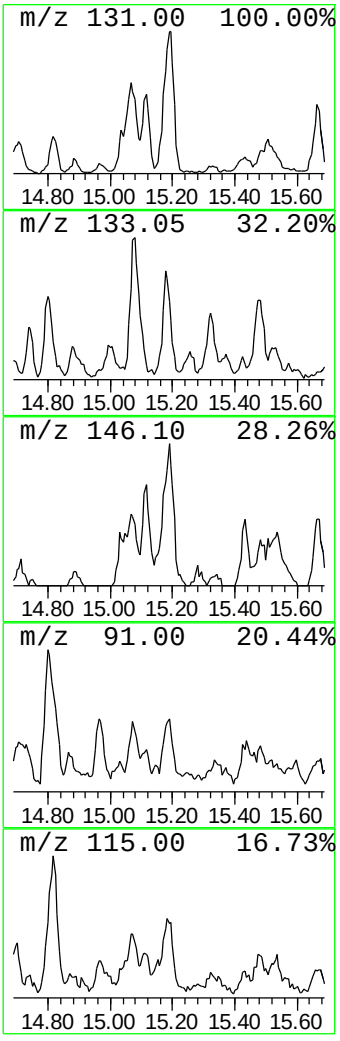
****

Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	7.24 ug/l	59593	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	92
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	70
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70
4	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	70
5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	62





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Data File : VW006149.D
Acq On : 18 Oct 2018 03:13
Operator : VA/AP
Sample : J5579-01
Misc : 5.26G/5ML/MSVOA_W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

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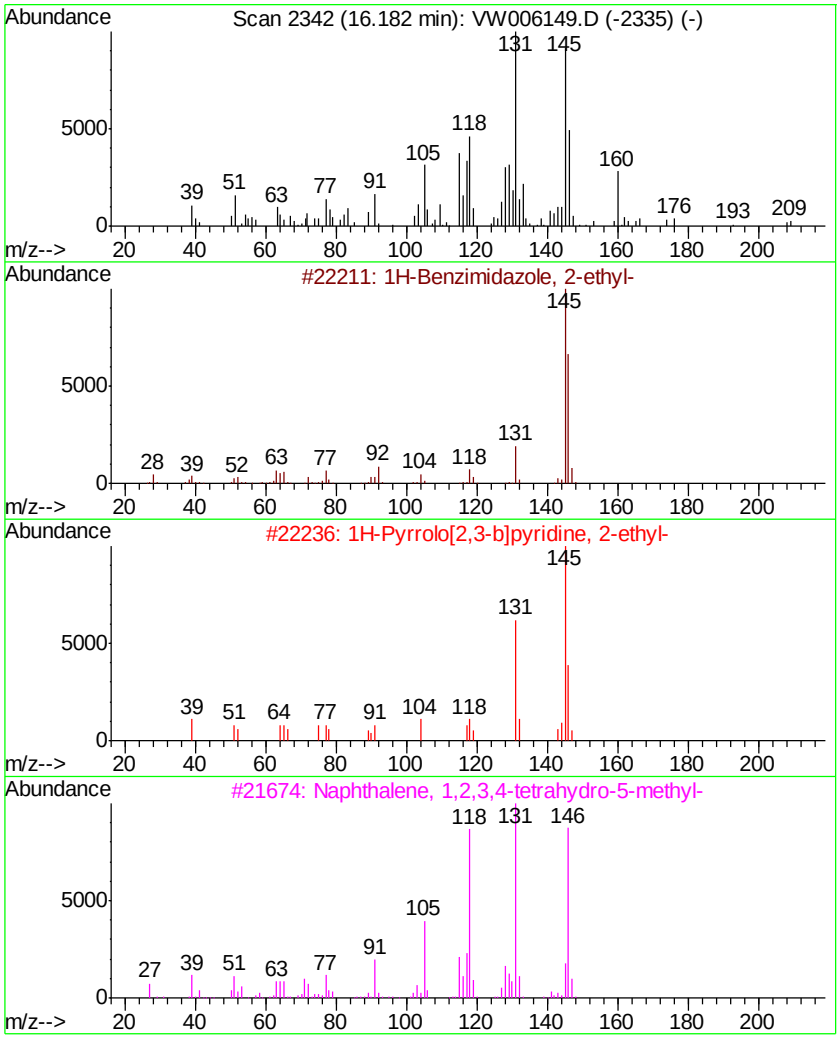
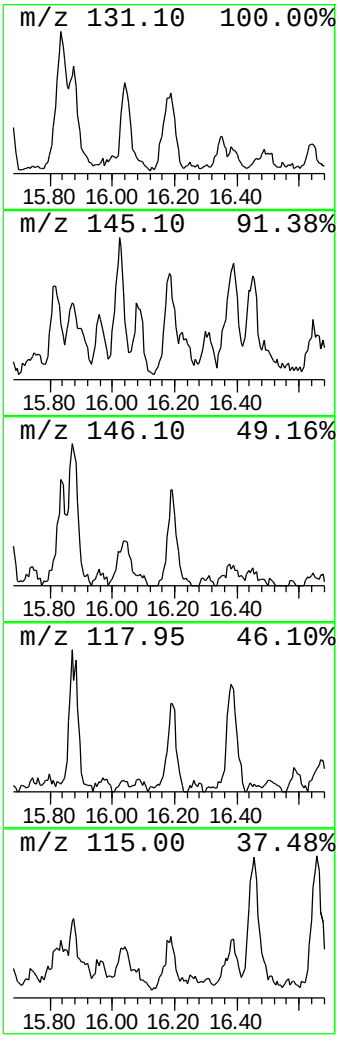
Peak Number 8 1H-Benzimidazole, 2-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.18	5.77 ug/l	47544	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	58
2	1H-Pyrrolo[2,3-b]pyridine, 2-ethyl-	146	C9H10N2	023612-49-9	52
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	46
4	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	45
5	4-Methyl-2H-benzopyrane	146	C10H10O	021776-94-3	41





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW101718\
Data File : VW006149.D
Acq On : 18 Oct 2018 03:13
Operator : VA/AP
Sample : J5579-01
Misc : 5.26G/5ML/MSVOA_W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

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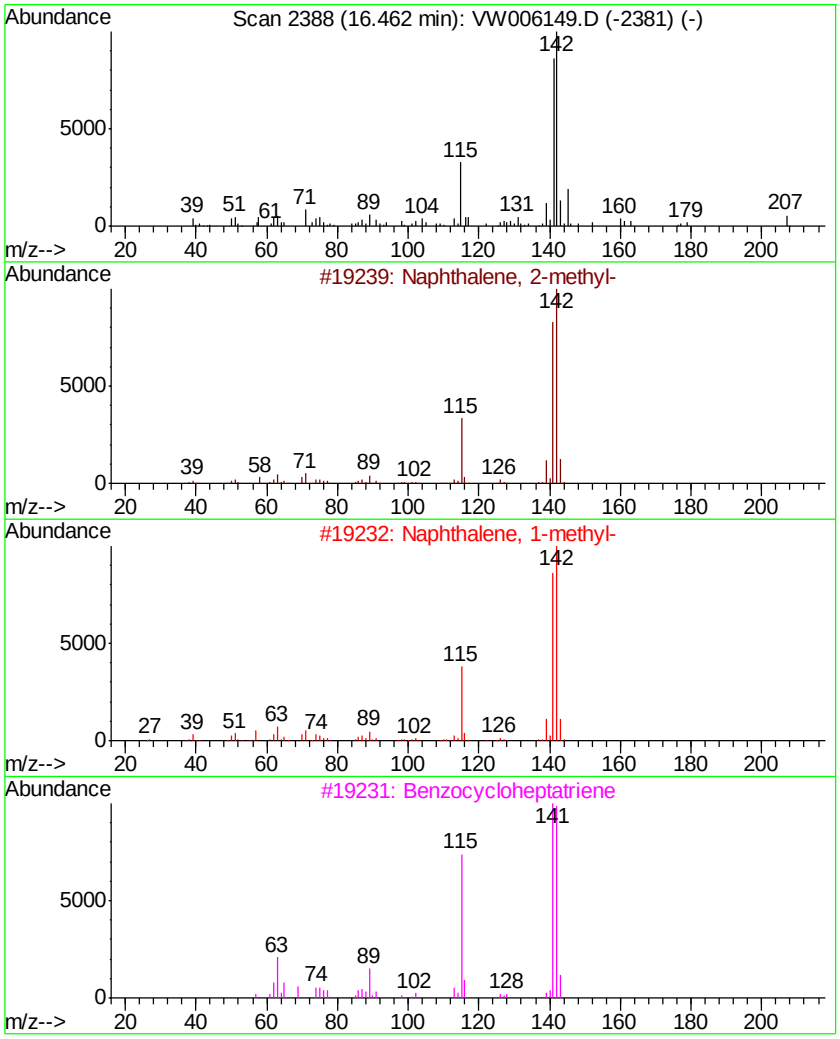
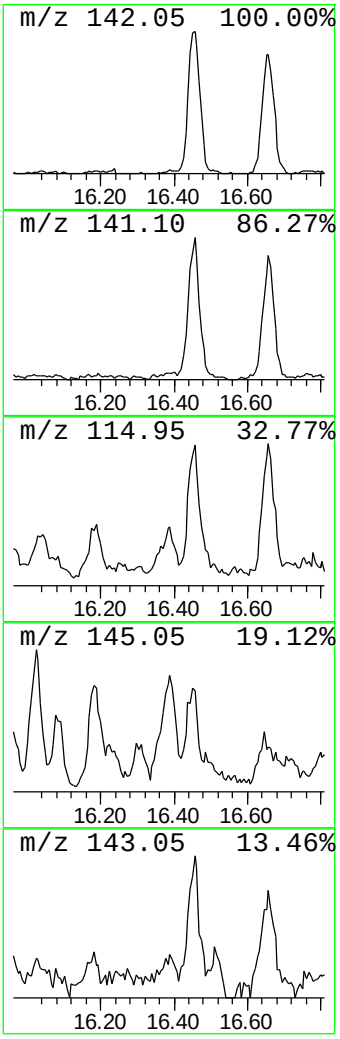
Peak Number 9 Naphthalene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.46	8.19 ug/l	67465	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
3	Benzocycloheptatriene	142	C11H10	000264-09-5	81
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	72
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW101718\
Data File : VW006149.D
Acq On : 18 Oct 2018 03:13
Operator : VA/AP
Sample : J5579-01
Misc : 5.26G/5ML/MSVOA_W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.M
Quant Title : SW846 8260

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

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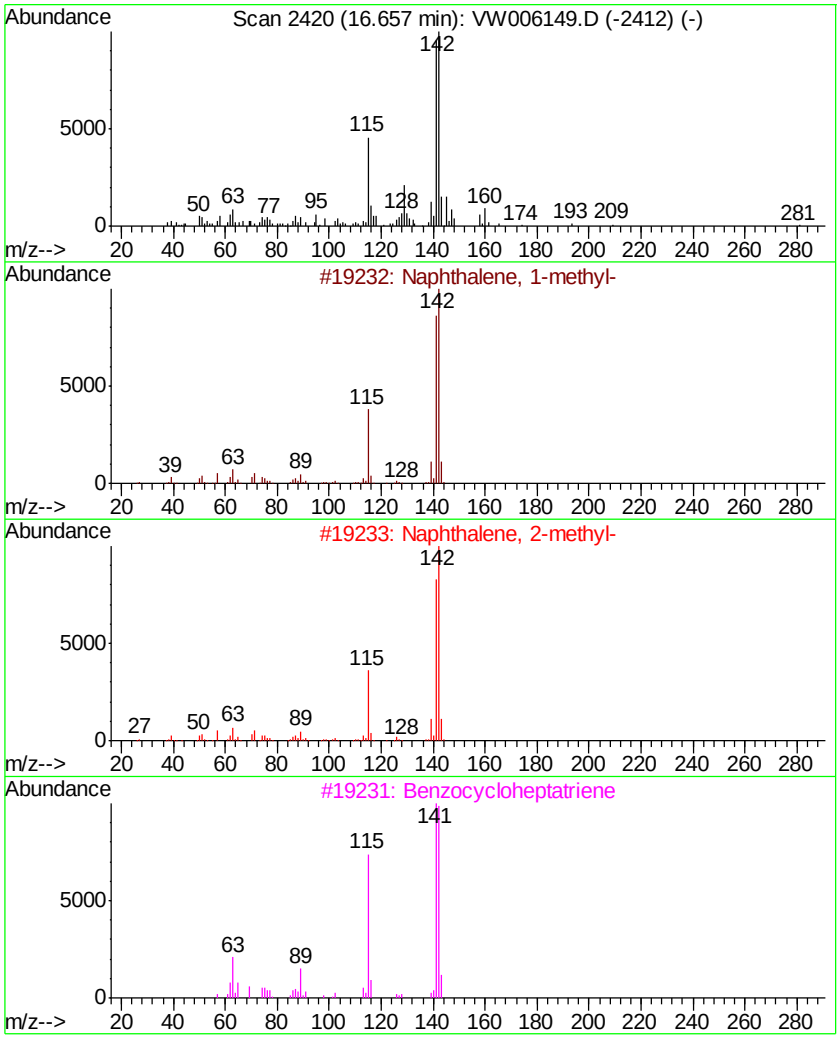
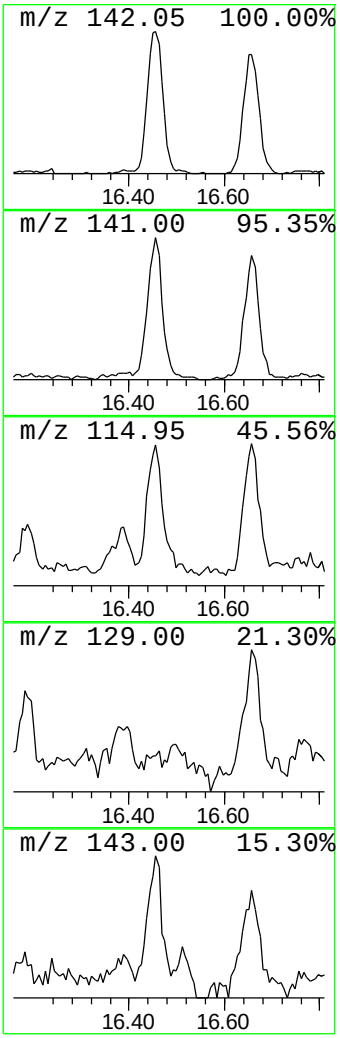
Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.66	7.42 ug/l	61060	1,4-Dichlorobenzene-d4	13.57

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	93
3	Benzocycloheptatriene	142	C11H10	000264-09-5	81
4	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	81
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_W\DATA\VW101718\
Data File : VW006149.D
Acq On : 18 Oct 2018 03:13
Operator : VA/AP
Sample : J5579-01
Misc : 5.26G/5ML/MSVOA_W/SOIL
ALS Vial : 36 Sample Multiplier: 1

Instrument : MSVOA_W
ClientSampleId : 7-1-ROLLOFFS

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_W\METHOD\82W101518S.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
Camphene	12.73	8.4	ug/l	69513	4	13.57	411679	50.0
unknown12.95	12.95	7.1	ug/l	58752	4	13.57	411679	50.0
Bicyclo[2.2.1]hep...	13.05	10.8	ug/l	88535	4	13.57	411679	50.0
Benzene, 2-ethyl-...	14.10	8.1	ug/l	67049	4	13.57	411679	50.0
1-Phenyl-1-butene	14.21	6.7	ug/l	54943	4	13.57	411679	50.0
Benzene, 1,2,4,5-...	14.80	17.6	ug/l	144601	4	13.57	411679	50.0
1H-Indene, 2,3-di...	15.19	7.2	ug/l	59593	4	13.57	411679	50.0
1H-Benzimidazole,...	16.18	5.8	ug/l	47544	4	13.57	411679	50.0
Naphthalene, 2-me...	16.46	8.2	ug/l	67465	4	13.57	411679	50.0
Naphthalene, 1-me...	16.66	7.4	ug/l	61060	4	13.57	411679	50.0