LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

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_U\METHOD\82U061318W.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	1.091	148	156	174	rBV	678876	793154	26.04%	2.194%
2	2.747	664	671	680	rVB2	19296	34784	1.14%	0.096%
3	3.001	735	750	768	rBV3	27349	59381	1.95%	0.164%
4	4.104	1076	1093	1110	rBV2	52880	143169	4.70%	0.396%
5	4.889	1319	1337	1354	rBV	186090	429103	14.09%	1.187%
6	4.992	1354	1369	1387	rVV	249960	583504	19.16%	1.614%
7	5.085	1389	1398	1418	rVB8	15320	46835	1.54%	0.130%
8	5.316	1456	1470	1488	rVV	176869	377430	12.39%	1.044%
9	5.558	1534	1545	1556	rVV6	14276	34757	1.14%	0.096%
10	5.628	1556	1567	1581	rVB4	13922	32852	1.08%	0.091%
11	5.892	1635	1649	1662	rBV	355929	714368	23.45%	1.976%
12	6.226	1741	1753	1768	rVB4	19195	46390	1.52%	0.128%
13	6.419	1792	1813	1829	rBV4	57411	152672	5.01%	0.422%
14	6.847	1929	1946	1958	rBV7	14217	43758	1.44%	0.121%
15	7.072	1997	2016	2027	rBV2	59787	124276	4.08%	0.344%
16	7.432	2117	2128	2141	rBV2	30224	55929	1.84%	0.155%
17	7.570	2149	2171	2190	rBV	674595	1223024	40.15%	3.383%
18	9.094	2630	2645	2663	rBV	523239	884199	29.03%	2.446%
19	9.188	2667	2674	2686	rVB3	20072	40651	1.33%	0.112%
20	9.400	2720	2740	2749	rBV10	23382	78608	2.58%	0.217%
21	9.573	2783	2794	2808	rVB2	34533	59120	1.94%	0.164%
22	9.747	2821	2848	2859	rBV4	57343	190787	6.26%	0.528%
23	9.959	2887	2914	2931	rBV2	151006	316550	10.39%	0.876%
24	10.053	2931	2943	2958	rBV5	63888	129137	4.24%	0.357%
25	10.168	2969	2979	2990	rBV	452701	693310	22.76%	1.918%
26	10.310	3012	3023	3035	rBV	534353	853923	28.03%	2.362%
27	10.381	3036	3045	3054	rVB2	78523	130613	4.29%	0.361%
28	10.496	3073	3081	3092	rVB4	65458	104606	3.43%	0.289%
29	10.589	3092	3110	3131	rBV	1409587	2172304	71.31%	6.009%
30	10.705	3136	3146	3157	rVB5	30315	66162	2.17%	0.183%
31	10.979	3217	3231	3249	rBV4	150158	283471	9.31%	0.784%
32	11.104	3252	3270	3278	rBV	70990	120191	3.95%	0.332%
33	11.294	3314	3329	3333	rBV	244171	378657	12.43%	1.048%
34	11.326	3333	3339	3353	rVB	402077	630902	20.71%	1.745%

LSC Area Percent Report

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Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

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_U\METHOD\82U061318W.M
Title : SW846 8260

35	11.487	3378	3389	3403	rBV	687608	1082380	35.53%	2.994%
36	11.692	3442	3453	3464	rVB	163852	256934	8.43%	0.711%
37	11.773	3464	3478	3495	rBV3	1549276	3046152	100.00%	8.427%
38	11.869	3496	3508	3511	rBV3	284141	432856	14.21%	1.197%
39	11.985	3531	3544	3556	rBV	254561	402184	13.20%	1.113%
40	12.062	3560	3568	3578	rVB3	23209	37636	1.24%	0.104%
41	12.255	3605	3628	3633	rBV	570249	915643	30.06%	2.533%
42	12.290	3633	3639	3649	rVV	486943	753528	24.74%	2.085%
43	12.358	3649	3660	3680	rVV2	1065591	2430302	79.78%	6.723%
44	12.438	3680	3685	3692	rVB2	45887	59160	1.94%	0.164%
45	12.541	3708	3717	3732	rVB	246240	397164	13.04%	1.099%
46	12.654	3732	3752	3764	rVV	1168079	1926453	63.24%	5.329%
47	12.721	3764	3773	3784	rVV3	42345	76380	2.51%	0.211%
48	12.789	3784	3794	3807	rVB3	149321	229749	7.54%	0.636%
49	12.882	3812	3823	3829	rBV	127430	202669	6.65%	0.561%
50	12.930	3830	3838	3842	rVV3	124413	184741	6.06%	0.511%
51	12.966	3843	3849	3858	rVB	389567	545390	17.90%	1.509%
52	13.126	3876	3899	3910	rBV2	1316822	2136532	70.14%	5.910%
53	13.194	3912	3920	3928	rVV4	90050	182031	5.98%	0.504%
54	13.239	3928	3934	3943	rVV	90741	163462	5.37%	0.452%
55	13.297	3944	3952	3971	rVB4	123466	243494	7.99%	0.674%
56	13.409	3978	3987	3994	rBV2	105112	176417	5.79%	0.488%
57	13.458	3995	4002	4010	rVB	161807	246600	8.10%	0.682%
58	13.512	4010	4019	4028	rBV2	364246	537296	17.64%	1.486%
59	13.580	4028	4040	4044	rBV3	159126	269520	8.85%	0.746%
60	13.612	4045	4050	4057	rVB	190371	252676	8.29%	0.699%
61	13.734	4076	4088	4097	rBV3	78952	143191	4.70%	0.396%
62	13.792	4098	4106	4116	rVV2	59888	97213	3.19%	0.269%
63	13.856	4116	4126	4139	rVV	323756	497343	16.33%	1.376%
64	13.953	4139	4156	4168	rVV2	235690	492976	16.18%	1.364%
65	14.017	4168	4176	4185	rVV3	92620	160986	5.28%	0.445%
66	14.065	4186	4191	4208	rVB5	27025	46092	1.51%	0.128%
67	14.155	4208	4219	4237	rBV2	112457	261520	8.59%	0.723%
68	14.335	4265	4275	4290	rBV2	224834	485962	15.95%	1.344%
69	14.480	4311	4320	4330	rVV2	94183	161706	5.31%	0.447%
70	14.544	4331	4340	4352	rVB3	110721	199962	6.56%	0.553%
71	14.609	4352	4360	4370	rVB4	43193	68665	2.25%	0.190%

LSC Area Percent Report

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Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

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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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Title : SW846 8260

72	14.679	4370	4382	4400	rBV2	495123	856690	28.12%	2.370%
73	14.824	4418	4427	4433	rBV2	47826	75775	2.49%	0.210%
74	14.866	4433	4440	4453	rVV4	92195	186532	6.12%	0.516%
75	14.937	4454	4462	4476	rVB4	78217	150071	4.93%	0.415%
76	15.014	4478	4486	4490	rBV3	22953	32133	1.05%	0.089%
77	15.065	4490	4502	4516	rBV	1060051	1756624	57.67%	4.859%
78	15.133	4516	4523	4528	rBV3	31974	45972	1.51%	0.127%
79	15.255	4551	4561	4573	rVB3	27810	64438	2.12%	0.178%
80	15.593	4653	4666	4681	rVB2	64941	141471	4.64%	0.391%
81	15.805	4723	4732	4736	rBV	29649	47189	1.55%	0.131%
82	15.837	4737	4742	4753	rVB2	50890	79933	2.62%	0.221%
83	15.917	4755	4767	4778	rBV	129558	241751	7.94%	0.669%
84	16.062	4800	4812	4820	rBV	234455	383294	12.58%	1.060%
85	16.265	4866	4875	4878	rVV2	49081	71900	2.36%	0.199%
86	16.293	4879	4884	4904	rVB2	78519	132080	4.34%	0.365%
87	16.435	4920	4928	4938	rVB2	33983	51106	1.68%	0.141%

Sum of corrected areas: 36148501

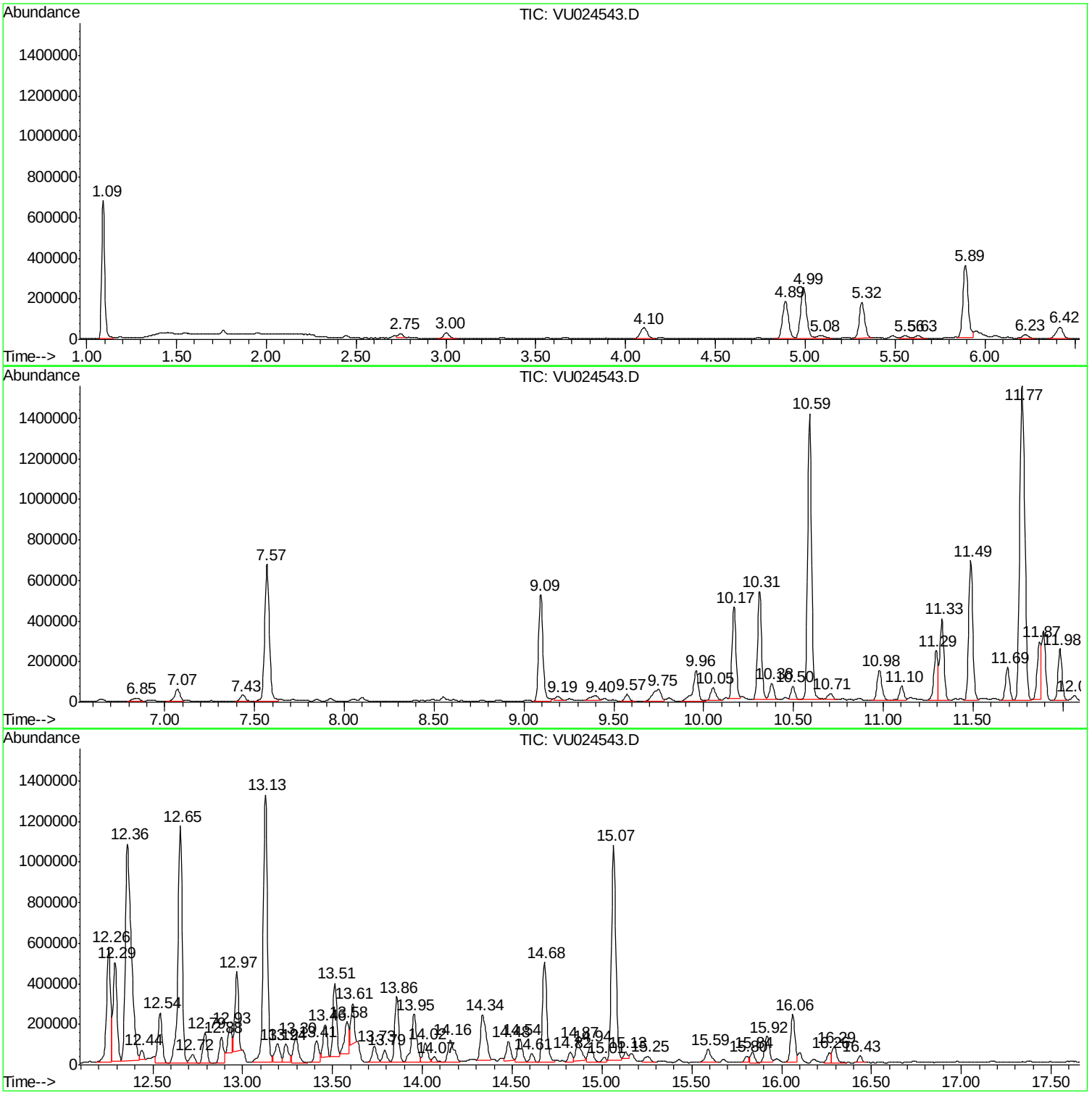
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.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_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

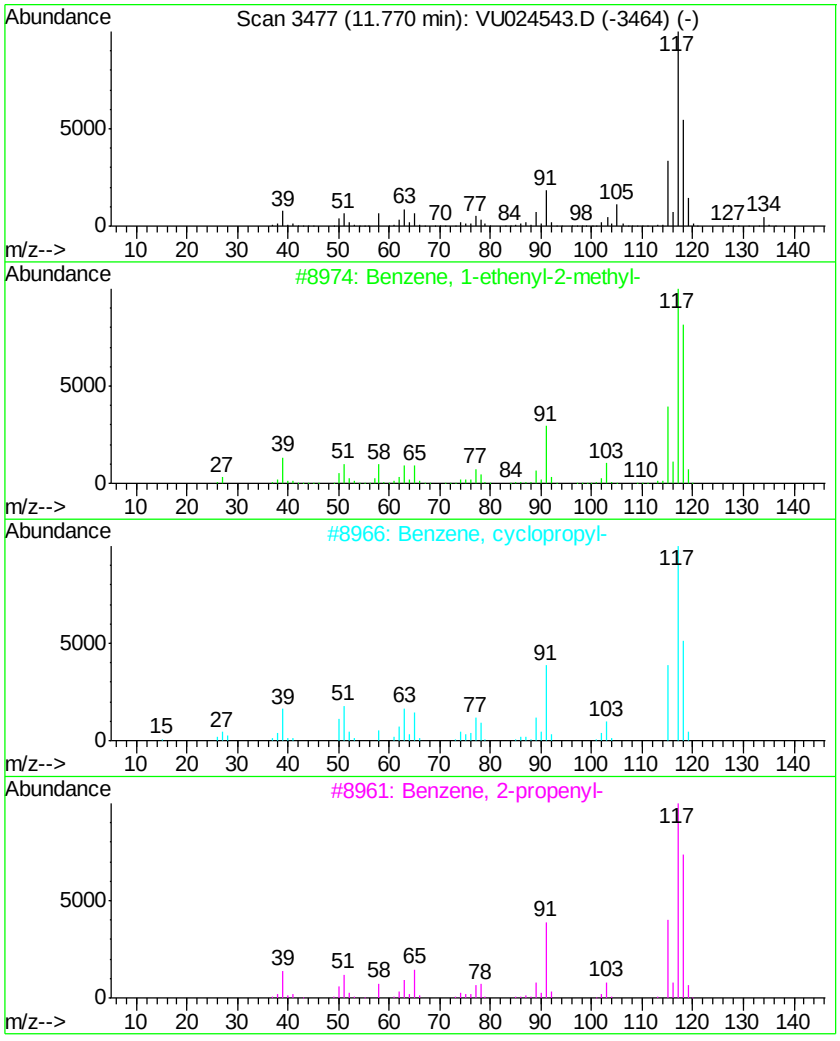
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

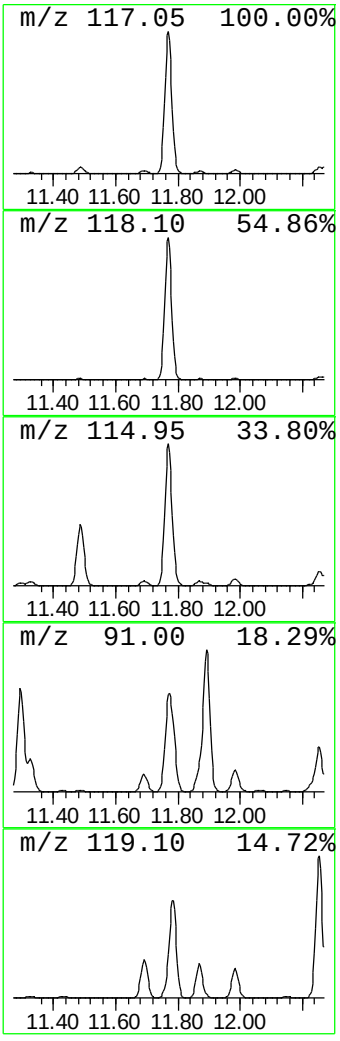
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	140.72 ug/l	3046150	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
4	Indane	118	C9H10	000496-11-7	70
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

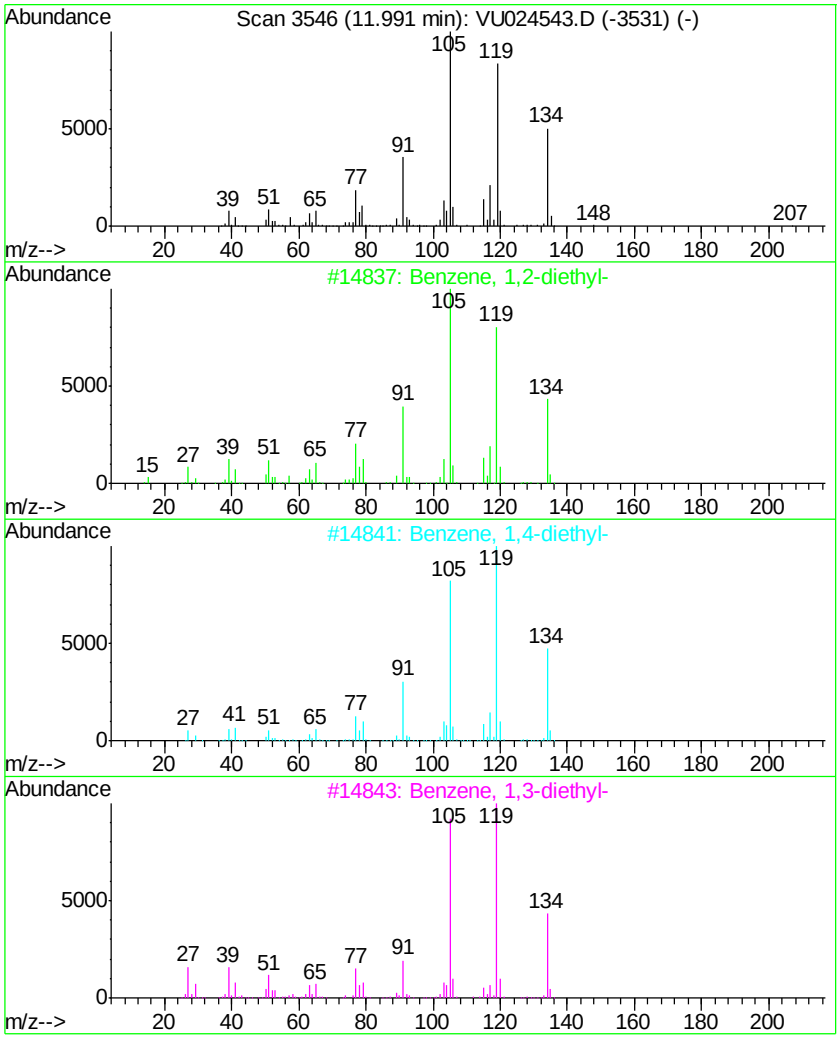
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

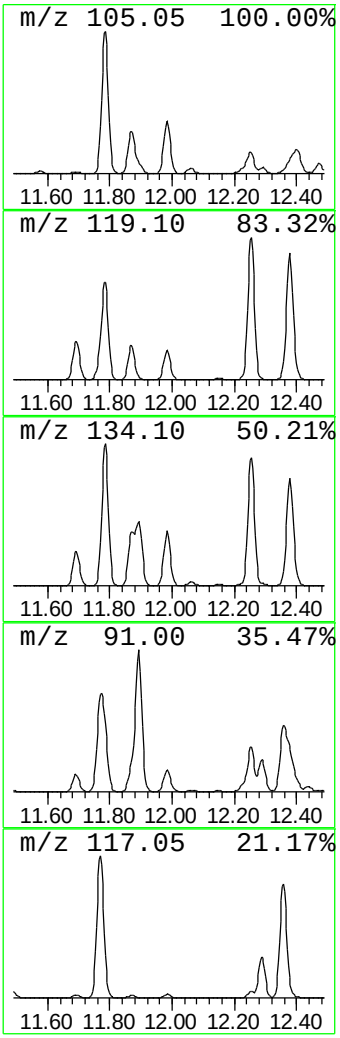
****

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.99	18.58 ug/l	402184	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	97
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	87
4	o-Cymene	134	C10H14	000527-84-4	76
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	74





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

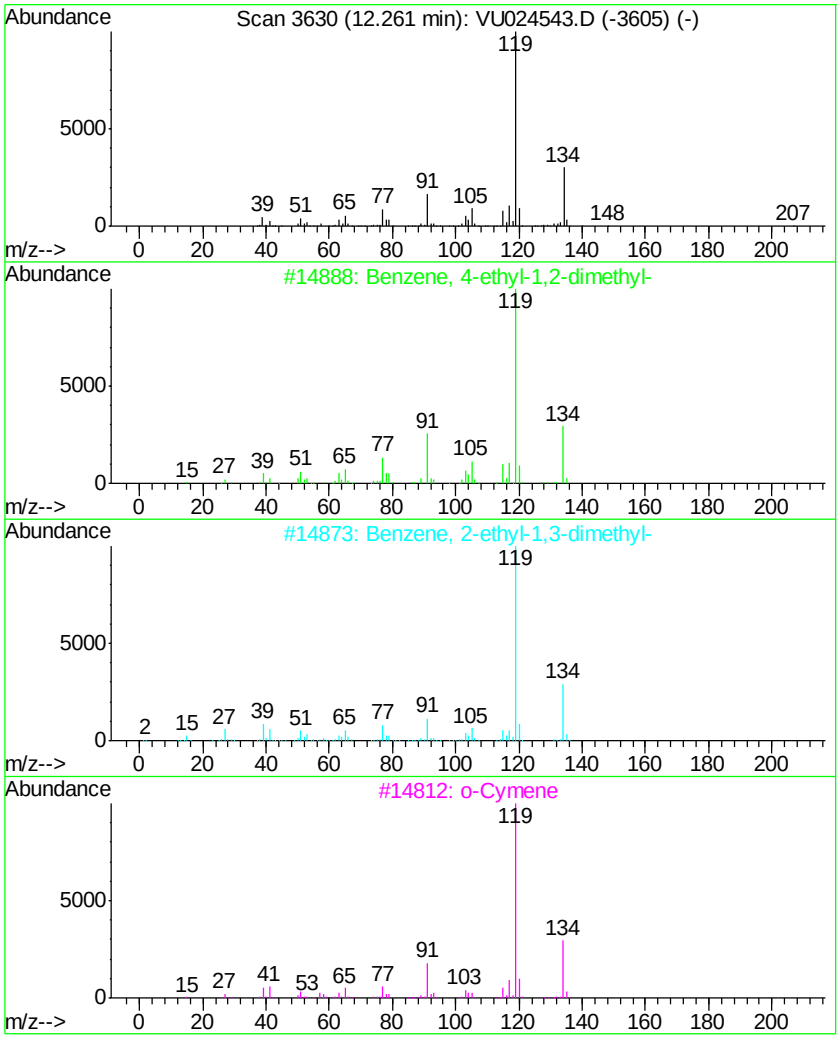
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

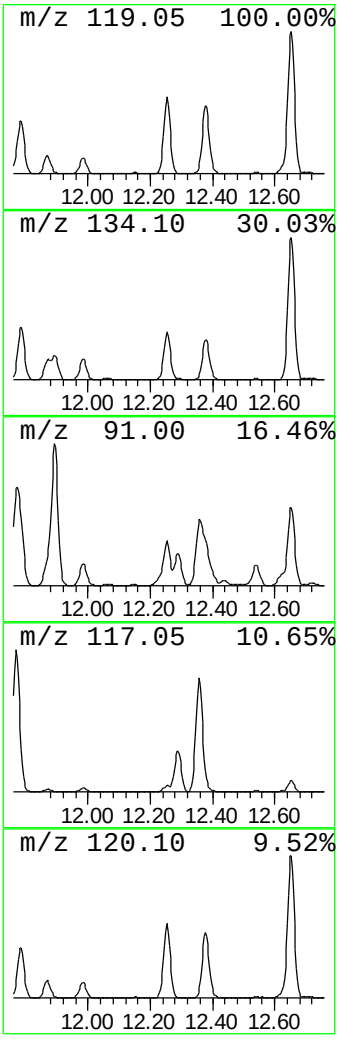
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	42.30 ug/l	915643	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	97
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

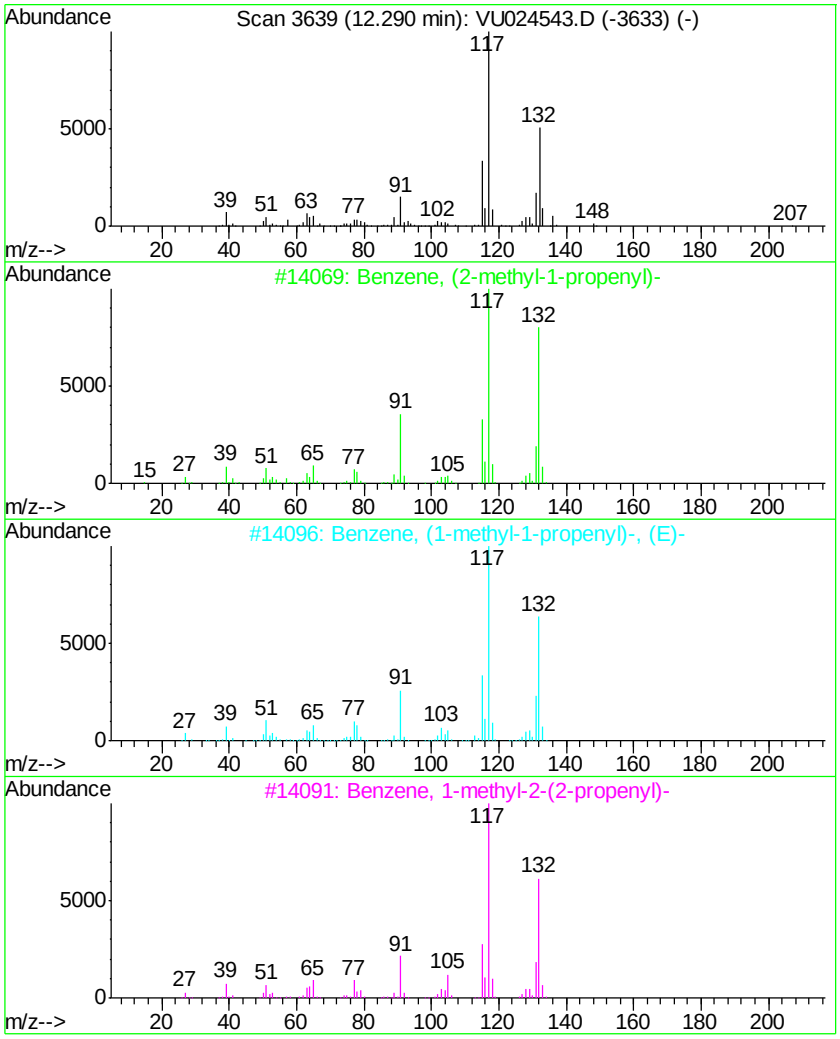
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

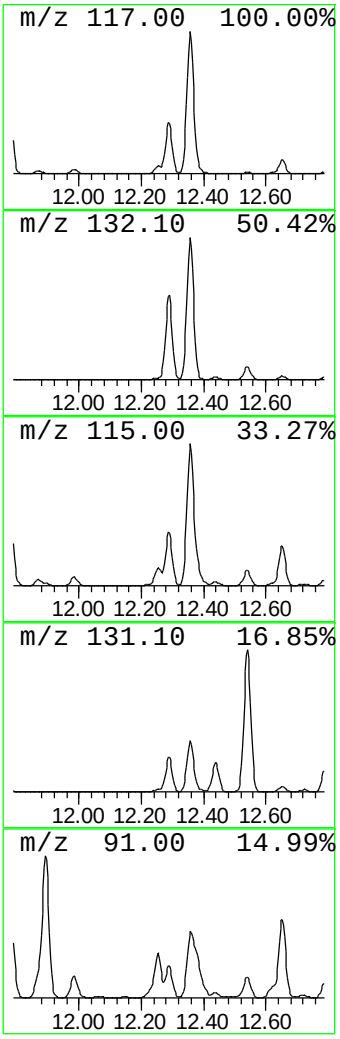
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	34.81 ug/l	753528	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	91
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	90





Library Search Compound Report

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Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
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Sample : J3443-07
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

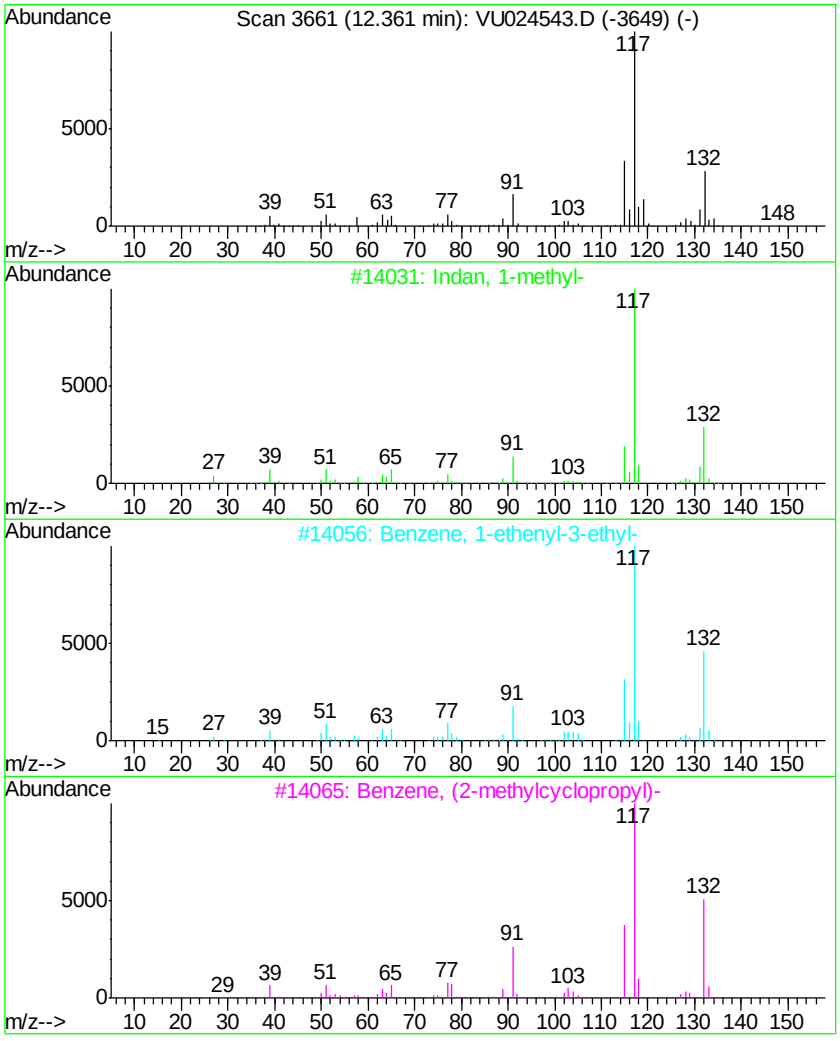
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

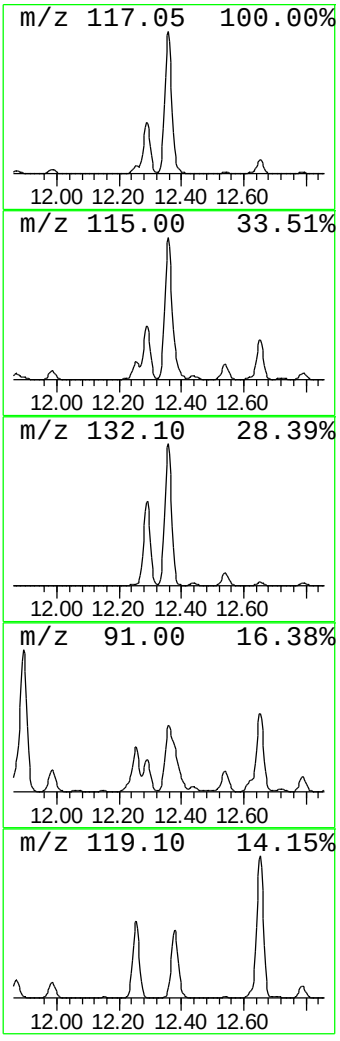
****

Peak Number 6 Indan, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	112.27 ug/l	2430300	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83





Library Search Compound Report

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Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

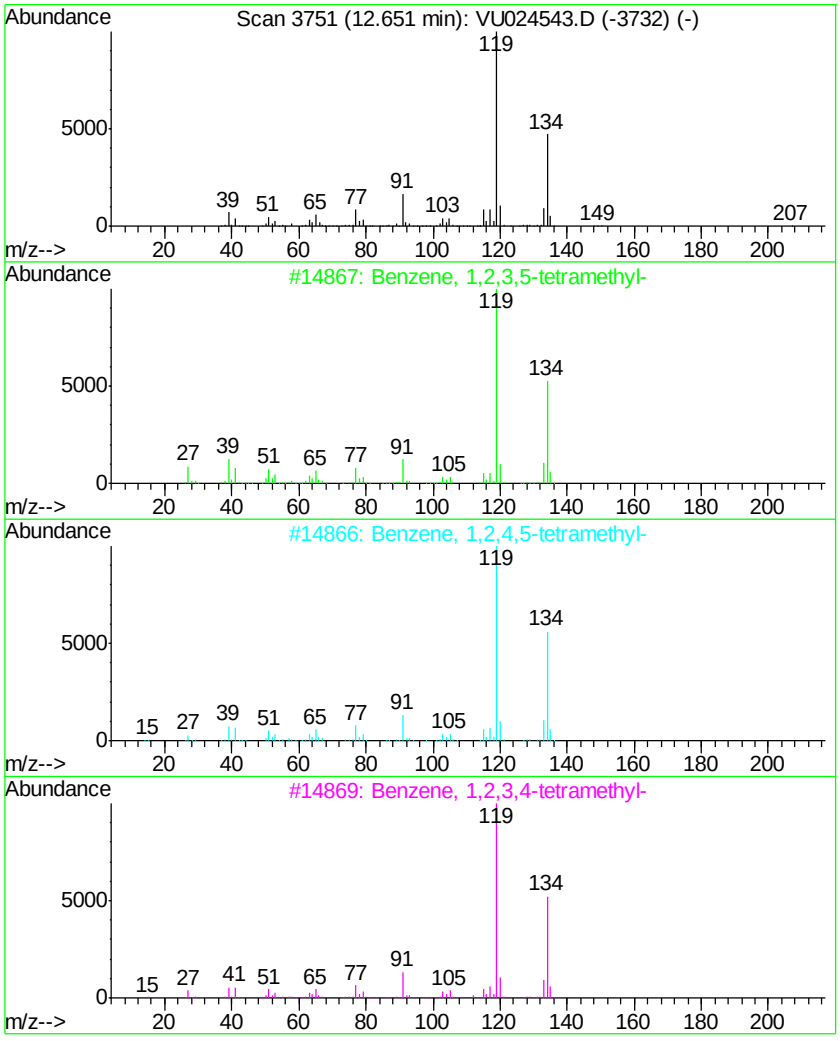
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

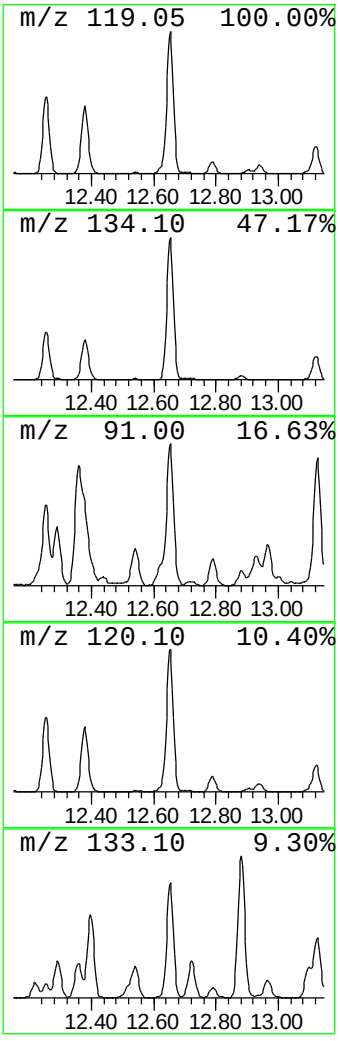
****

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	88.99 ug/l	1926450	1,4-Dichlorobenzene-d4	11.49

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





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

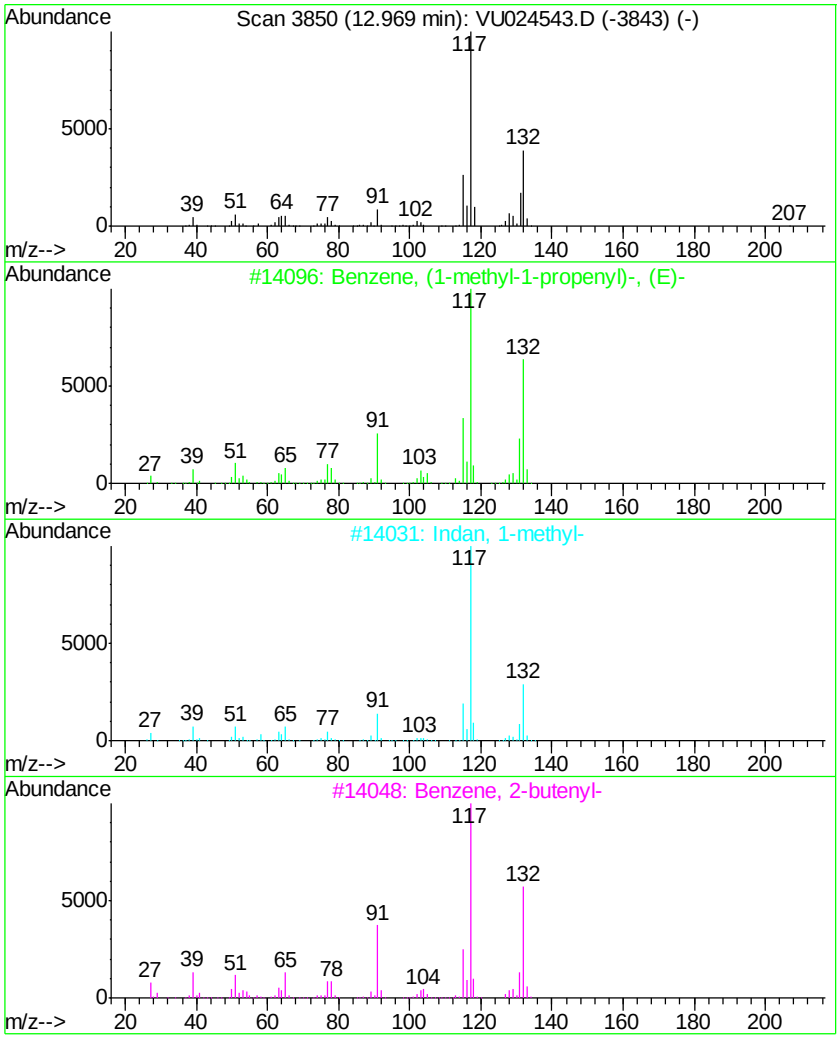
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

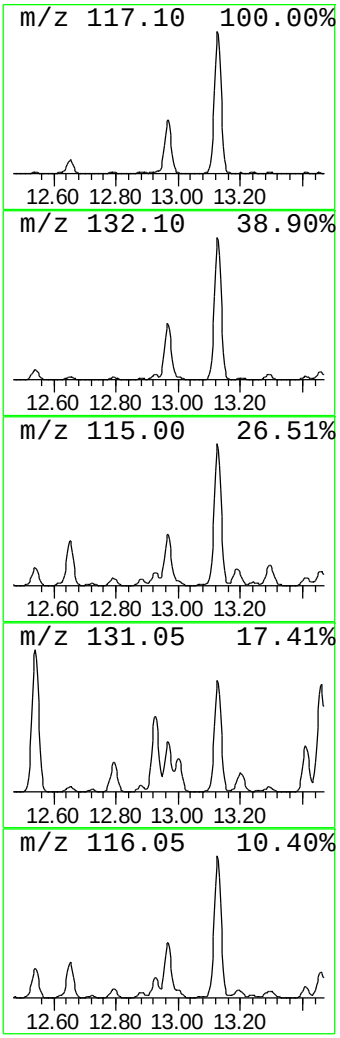
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, (1-methyl-1-propen... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.97	25.19 ug/l	545390	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	91
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	91
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

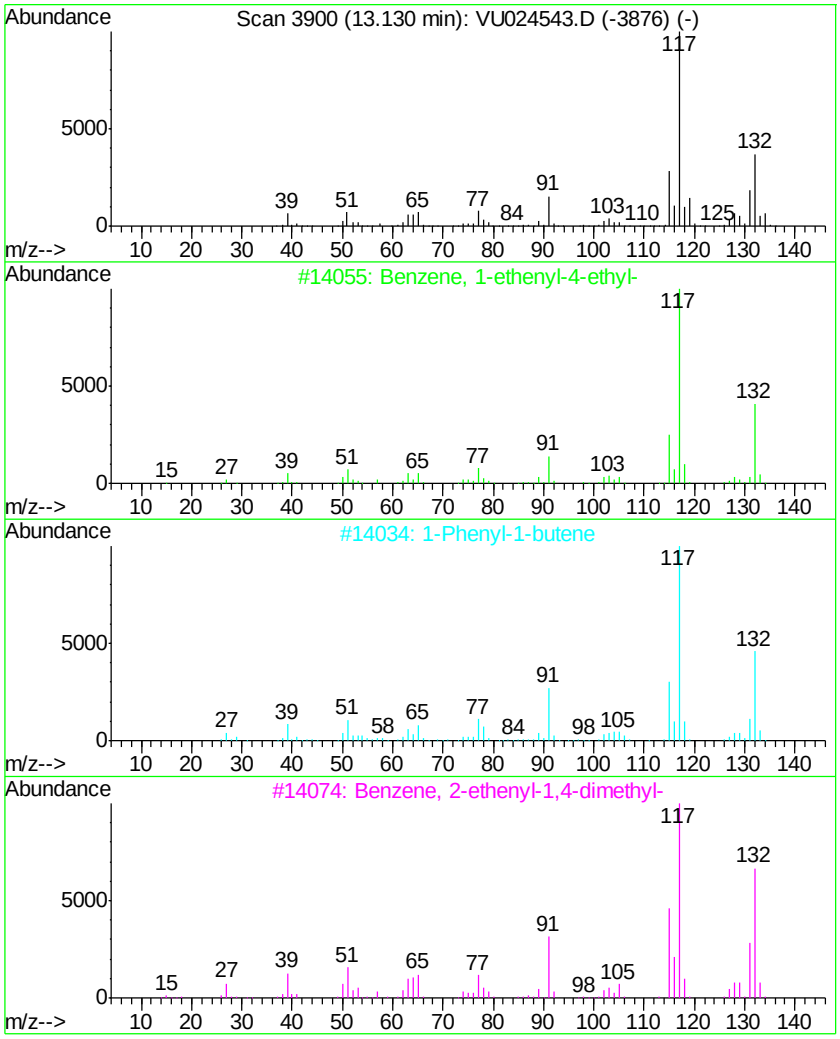
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

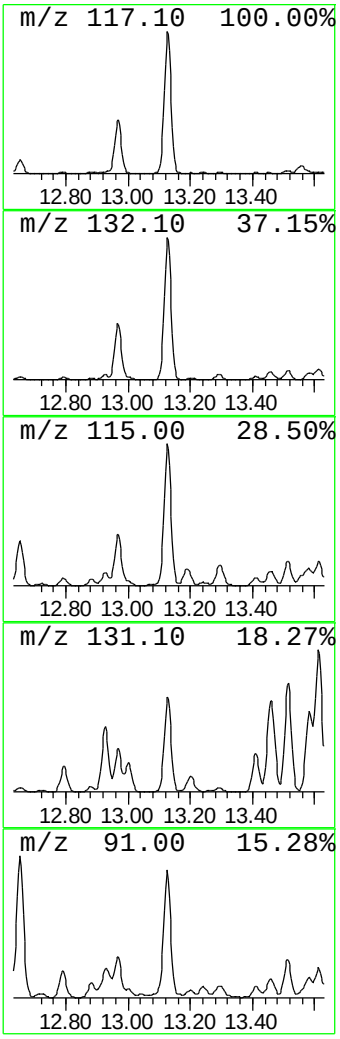
****

Peak Number 9 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	98.70 ug/l	2136530	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	93
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
5	Indan, 1-methyl-	132	C10H12	000767-58-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

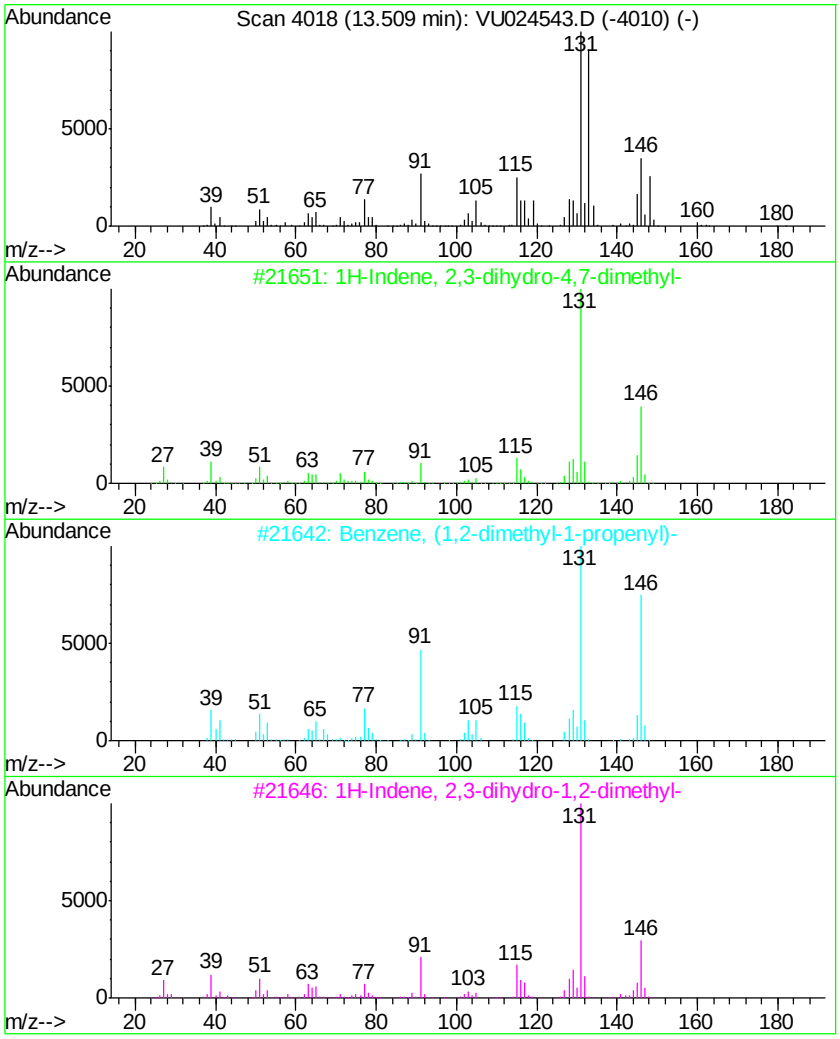
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.M
Quant Title : SW846 8260

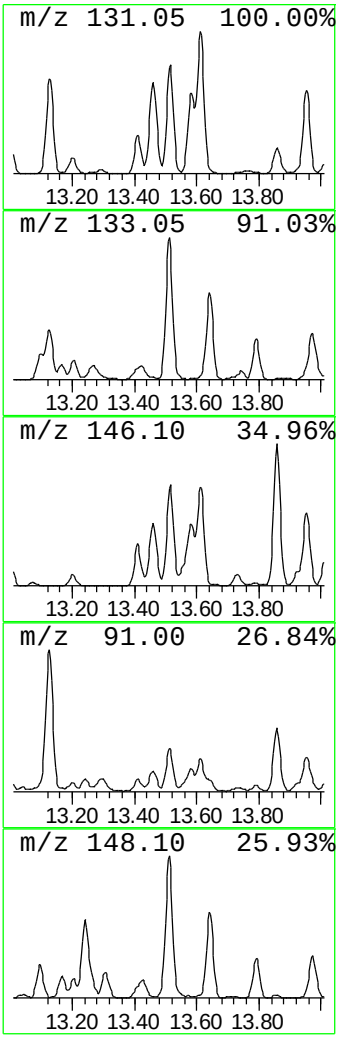
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	24.82 ug/l	537296	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	89
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	87
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU061418\
Data File : VU024543.D
Acq On : 14 Jun 2018 16:30
Operator : MD/SY
Sample : J3443-07
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : T3-MW-3-9.0-061118

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U061318W.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...	11.77	140.7	ug/l	3046150	4	11.49	1082380	50.0
Benzene, 1,2-diet...	11.99	18.6	ug/l	402184	4	11.49	1082380	50.0
Benzene, 4-ethyl-...	12.26	42.3	ug/l	915643	4	11.49	1082380	50.0
Benzene, (2-methy...	12.29	34.8	ug/l	753528	4	11.49	1082380	50.0
Indan, 1-methyl-	12.36	112.3	ug/l	2430300	4	11.49	1082380	50.0
Benzene, 1,2,3,5-...	12.65	89.0	ug/l	1926450	4	11.49	1082380	50.0
Benzene, (1-methy...	12.97	25.2	ug/l	545390	4	11.49	1082380	50.0
Benzene, 1-etheny...	13.13	98.7	ug/l	2136530	4	11.49	1082380	50.0
1H-Indene, 2,3-di...	13.51	24.8	ug/l	537296	4	11.49	1082380	50.0