LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

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\82U070918W.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	1.089	139	155	177	rBV	686603	822838	27.53%	2.233%
2	1.761	354	364	378	rVB	49222	65126	2.18%	0.177%
3	2.307	522	534	550	rVB2	20312	40230	1.35%	0.109%
4	2.709	643	659	681	rBV	206514	432781	14.48%	1.175%
5	3.002	736	750	767	rBV2	23927	51863	1.74%	0.141%
6	3.860	999	1017	1035	rVB3	16519	46472	1.55%	0.126%
7	4.002	1039	1061	1088	rBV2	158706	442614	14.81%	1.201%
8	4.230	1111	1132	1152	rVB4	29368	92078	3.08%	0.250%
9	4.738	1270	1290	1309	rBV2	30736	83875	2.81%	0.228%
10	4.889	1320	1337	1354	rBV	182710	410577	13.74%	1.114%
11	4.989	1354	1368	1385	rVV	273040	607306	20.32%	1.648%
12	5.085	1385	1398	1425	rVB	107550	262838	8.79%	0.713%
13	5.265	1434	1454	1461	rBV2	125599	298148	9.98%	0.809%
14	5.313	1461	1469	1489	rVB	191594	409170	13.69%	1.110%
15	5.542	1503	1540	1565	rBV	872685	2167781	72.53%	5.883%
16	5.892	1634	1649	1671	rBV	368463	730943	24.46%	1.984%
17	6.397	1789	1806	1820	rBV2	133271	285986	9.57%	0.776%
18	6.481	1820	1832	1843	rVV2	87909	162218	5.43%	0.440%
19	6.593	1855	1867	1882	rVB2	126863	261283	8.74%	0.709%
20	6.744	1897	1914	1932	rVB	129481	259381	8.68%	0.704%
21	6.931	1949	1972	1993	rBV	647153	1315783	44.02%	3.571%
22	7.053	1993	2010	2039	rVV	1274419	2610253	87.33%	7.084%
23	7.262	2063	2075	2084	rBV2	51761	107927	3.61%	0.293%
24	7.448	2117	2133	2144	rBV5	17042	38677	1.29%	0.105%
25	7.529	2144	2158	2160	rBV2	89032	134727	4.51%	0.366%
26	7.567	2161	2170	2189	rVB	756206	1344518	44.98%	3.649%
27	7.715	2203	2216	2231	rBV2	107657	221523	7.41%	0.601%
28	7.860	2249	2261	2271	rBV6	18633	36060	1.21%	0.098%
29	7.921	2271	2280	2288	rBV3	24759	41134	1.38%	0.112%
30	8.050	2302	2320	2331	rBV	31440	61422	2.06%	0.167%
31	8.136	2331	2347	2364	rVB9	11519	37158	1.24%	0.101%
32	8.493	2441	2458	2469	rBV5	22693	47904	1.60%	0.130%
33	8.609	2483	2494	2503	rBV2	69277	124384	4.16%	0.338%
34	8.850	2557	2569	2581	rBV2	35934	62487	2.09%	0.170%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

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\82U070918W.M
Title : SW846 8260

35	9.095	2632	2645	2651	rBV	528552	945366	31.63%	2.566%
36	9.191	2666	2675	2686	rVB	74802	133488	4.47%	0.362%
37	9.368	2719	2730	2747	rBV5	64031	147808	4.95%	0.401%
38	9.448	2747	2755	2763	rBV	27514	42144	1.41%	0.114%
39	9.574	2781	2794	2809	rBV2	49975	85109	2.85%	0.231%
40	9.747	2825	2848	2859	rBV6	85672	242430	8.11%	0.658%
41	9.808	2859	2867	2879	rVB3	27535	48698	1.63%	0.132%
42	9.924	2886	2903	2906	rBV4	62433	110647	3.70%	0.300%
43	9.956	2907	2913	2926	rVB	174078	283475	9.48%	0.769%
44	10.053	2932	2943	2958	rBV6	66777	141515	4.73%	0.384%
45	10.169	2971	2979	2987	rVV	64707	94277	3.15%	0.256%
46	10.223	2987	2996	3010	rVB6	44080	94425	3.16%	0.256%
47	10.310	3012	3023	3035	rBV	528238	825469	27.62%	2.240%
48	10.381	3035	3045	3053	rVV3	87521	139929	4.68%	0.380%
49	10.500	3073	3082	3092	rVB3	159808	251218	8.41%	0.682%
50	10.635	3116	3124	3131	rBV7	18450	37207	1.24%	0.101%
51	10.699	3137	3144	3156	rVB5	49575	100250	3.35%	0.272%
52	10.866	3188	3196	3209	rVB7	24667	49788	1.67%	0.135%
53	10.985	3217	3233	3248	rBV	150456	277831	9.30%	0.754%
54	11.104	3252	3270	3279	rBV	159866	264972	8.87%	0.719%
55	11.294	3311	3329	3335	rBV	346732	622006	20.81%	1.688%
56	11.326	3335	3339	3355	rVB	258267	364055	12.18%	0.988%
57	11.406	3356	3364	3378	rVB	15336	38965	1.30%	0.106%
58	11.487	3378	3389	3404	rBV	610856	1005606	33.64%	2.729%
59	11.693	3441	3453	3467	rVB	521051	829522	27.75%	2.251%
60	11.783	3467	3481	3494	rBV4	478369	799285	26.74%	2.169%
61	11.863	3495	3506	3529	rVB4	185025	380628	12.73%	1.033%
62	11.985	3532	3544	3557	rBV6	38255	82400	2.76%	0.224%
63	12.255	3614	3628	3634	rVV	1603399	2479014	82.94%	6.728%
64	12.291	3634	3639	3650	rVV2	720524	1105636	36.99%	3.001%
65	12.358	3650	3660	3679	rVV2	1279440	2988898	100.00%	8.112%
66	12.439	3679	3685	3695	rVB	116224	172785	5.78%	0.469%
67	12.503	3695	3705	3707	rBV	60850	83851	2.81%	0.228%
68	12.541	3708	3717	3732	rVB	507175	805217	26.94%	2.185%
69	12.651	3732	3751	3765	rBV2	768631	1436338	48.06%	3.898%
70	12.725	3765	3774	3784	rVB2	49466	76797	2.57%	0.208%
71	12.792	3784	3795	3811	rVB2	163011	262291	8.78%	0.712%

LSC Area Percent Report

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Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

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\82U070918W.M
Title : SW846 8260

72	12.882	3812	3823	3829	rBV	148427	219202	7.33%	0.595%
73	12.927	3829	3837	3844	rVB	140472	188826	6.32%	0.512%
74	12.969	3844	3850	3854	rBV	52036	61158	2.05%	0.166%
75	13.001	3854	3860	3871	rVB	105481	162507	5.44%	0.441%
76	13.123	3873	3898	3909	rBV2	1312255	2235581	74.80%	6.067%
77	13.194	3915	3920	3928	rVB3	54192	83781	2.80%	0.227%
78	13.242	3928	3935	3945	rVV	54132	98971	3.31%	0.269%
79	13.294	3947	3951	3975	rVB4	48086	91628	3.07%	0.249%
80	13.413	3979	3988	3994	rBV2	29751	49742	1.66%	0.135%
81	13.458	3994	4002	4010	rVB	58568	92610	3.10%	0.251%
82	13.512	4010	4019	4027	rBV2	227949	342720	11.47%	0.930%
83	13.580	4027	4040	4045	rVV4	111942	254473	8.51%	0.691%
84	13.612	4045	4050	4075	rVB2	161852	306220	10.25%	0.831%
85	13.734	4077	4088	4097	rBV2	54741	100293	3.36%	0.272%
86	13.789	4098	4105	4117	rVB3	40251	65091	2.18%	0.177%
87	13.860	4117	4127	4140	rVB2	44921	71677	2.40%	0.195%
88	13.956	4140	4157	4168	rBV5	47967	109601	3.67%	0.297%
89	14.017	4169	4176	4185	rVB4	25806	41715	1.40%	0.113%
90	14.342	4266	4277	4290	rBV6	36607	90647	3.03%	0.246%
91	14.683	4372	4383	4400	rBV3	46178	87126	2.91%	0.236%
92	15.069	4492	4503	4517	rBV2	68001	125680	4.20%	0.341%

Sum of corrected areas: 36846054

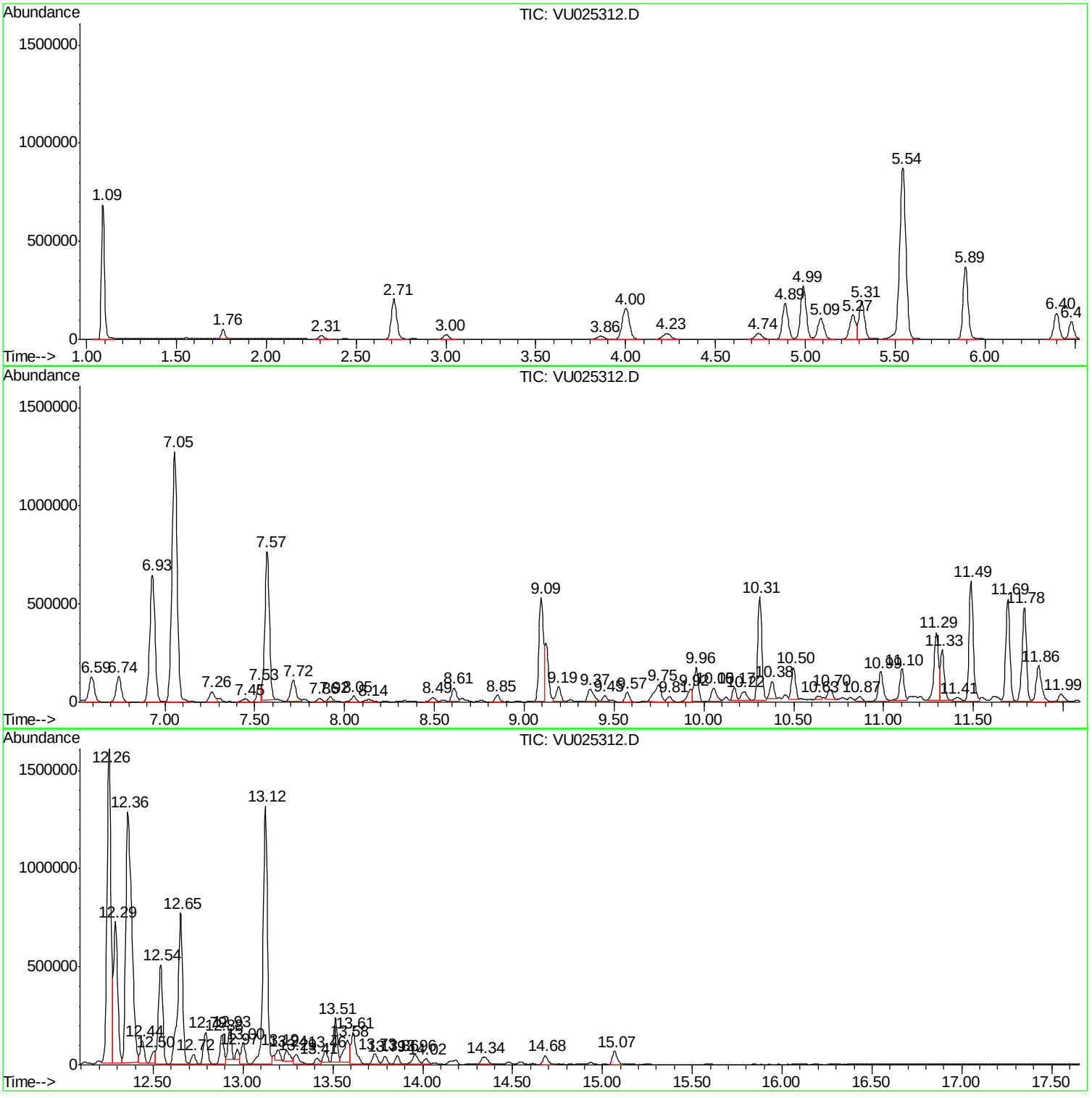
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

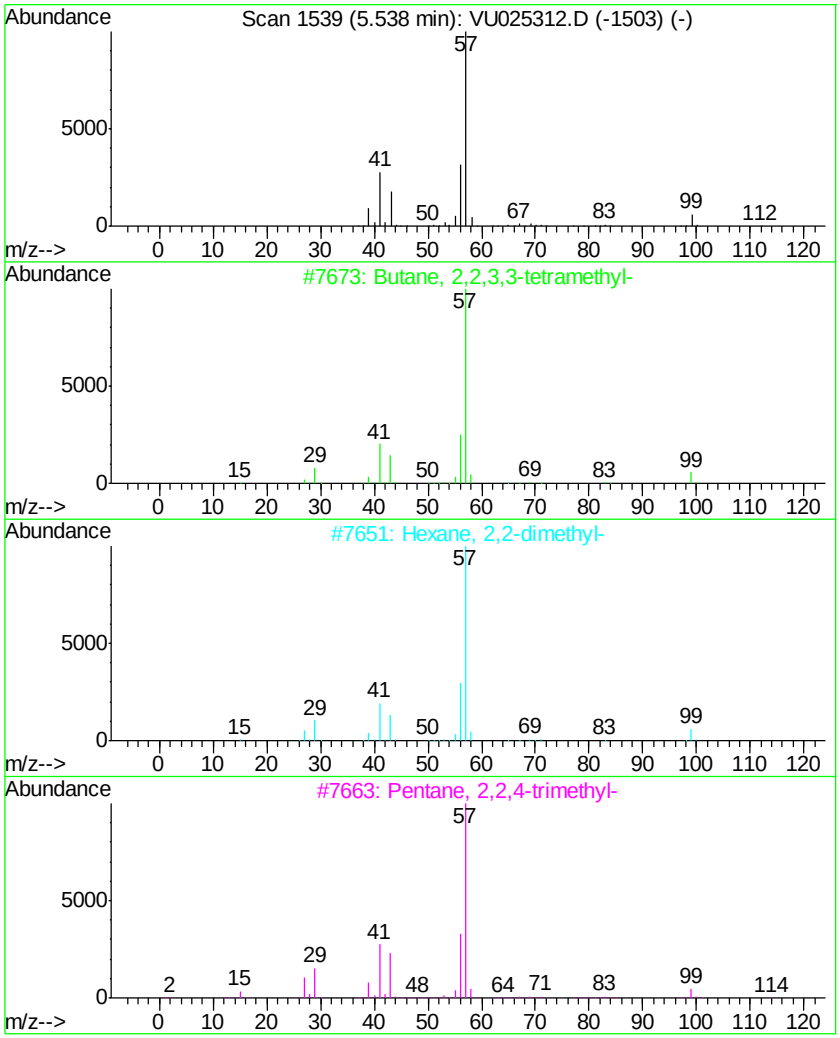
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

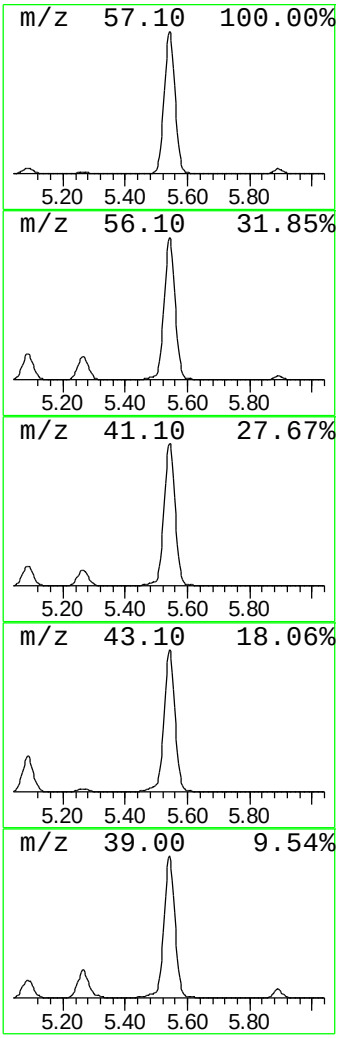
****

Peak Number 2 Butane, 2,2,3,3-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.54	148.29 ug/l	2167780	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
3	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	53





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

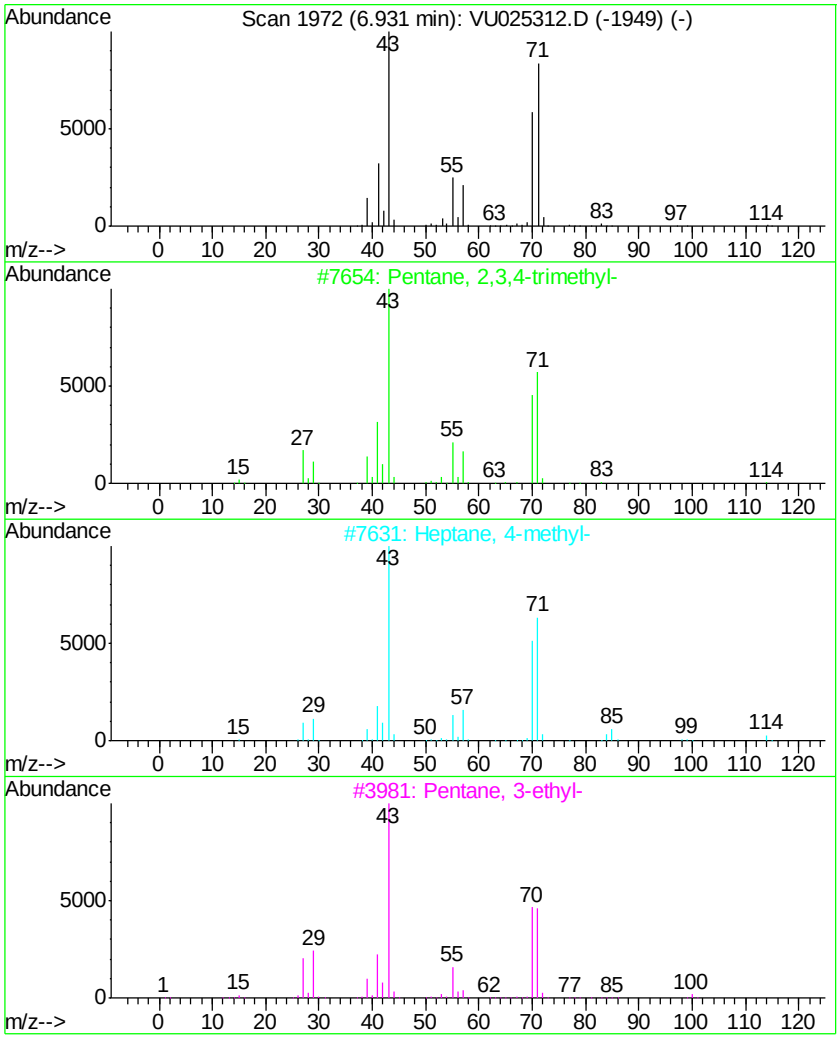
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

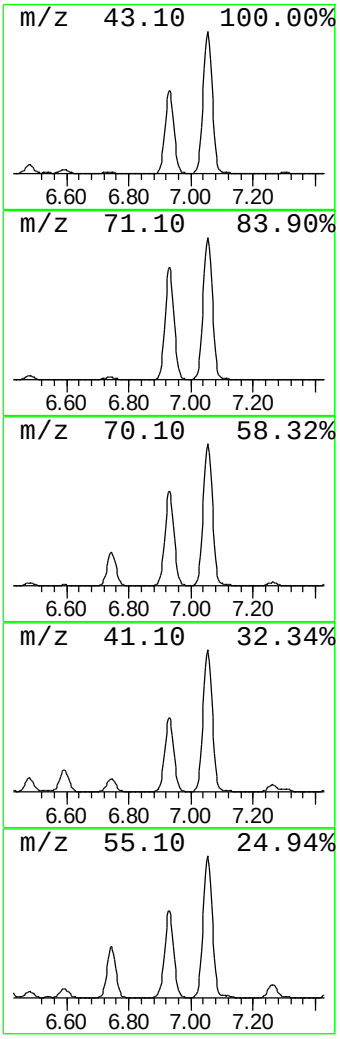
****

Peak Number 3 Pentane, 2,3,4-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	90.01 ug/l	1315780	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3	Pentane, 3-ethyl-	100	C7H16	000617-78-7	83
4	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	78
5	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

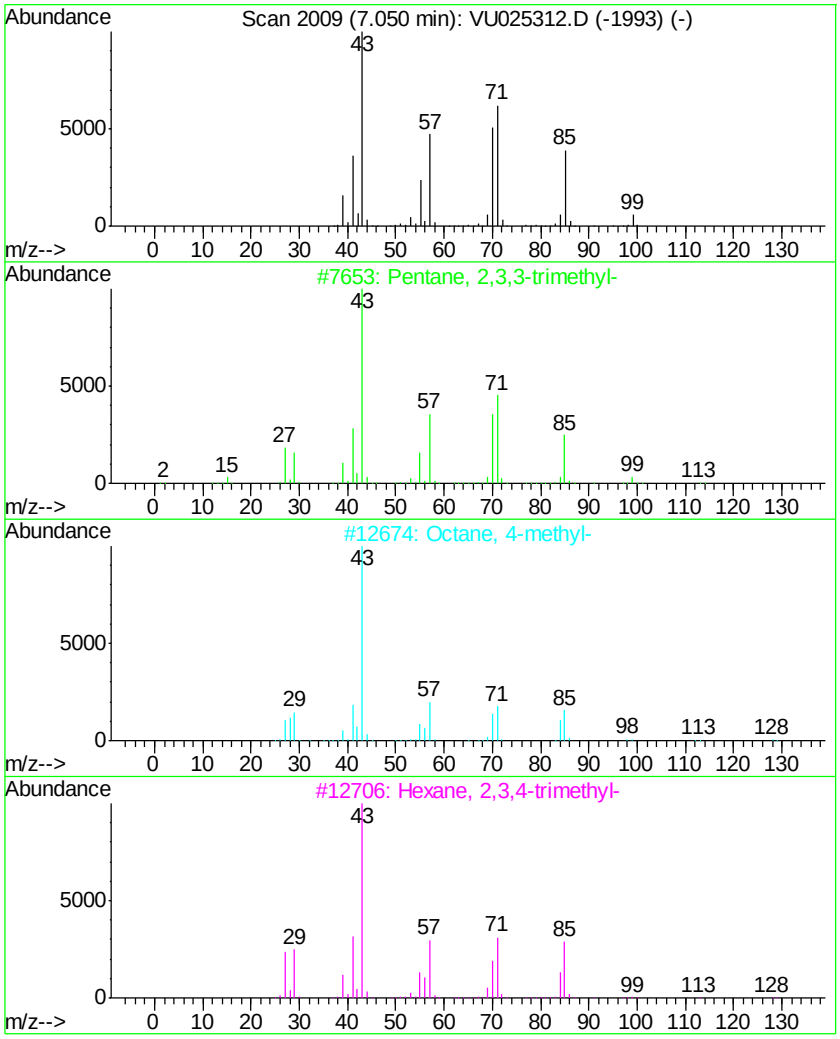
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

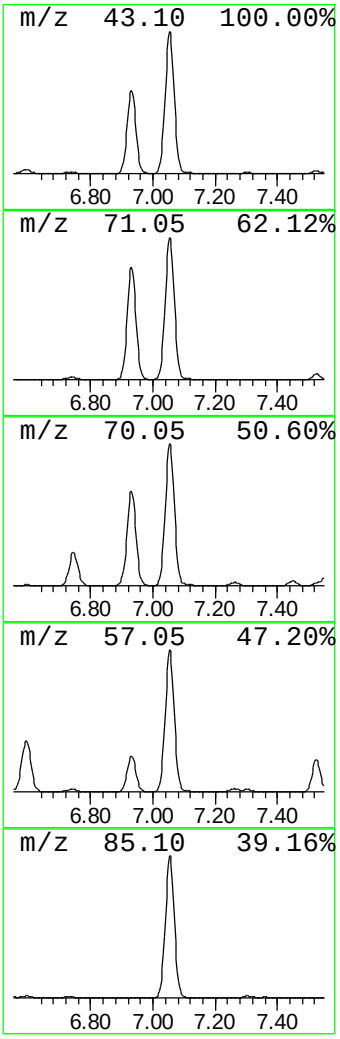
****

Peak Number 4 Pentane, 2,3,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.05	178.55 ug/l	2610250	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Octane, 4-methyl-	128	C9H20	002216-34-4	78
3	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	64
4	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
5	Heptane, 4-methyl-	114	C8H18	000589-53-7	53





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

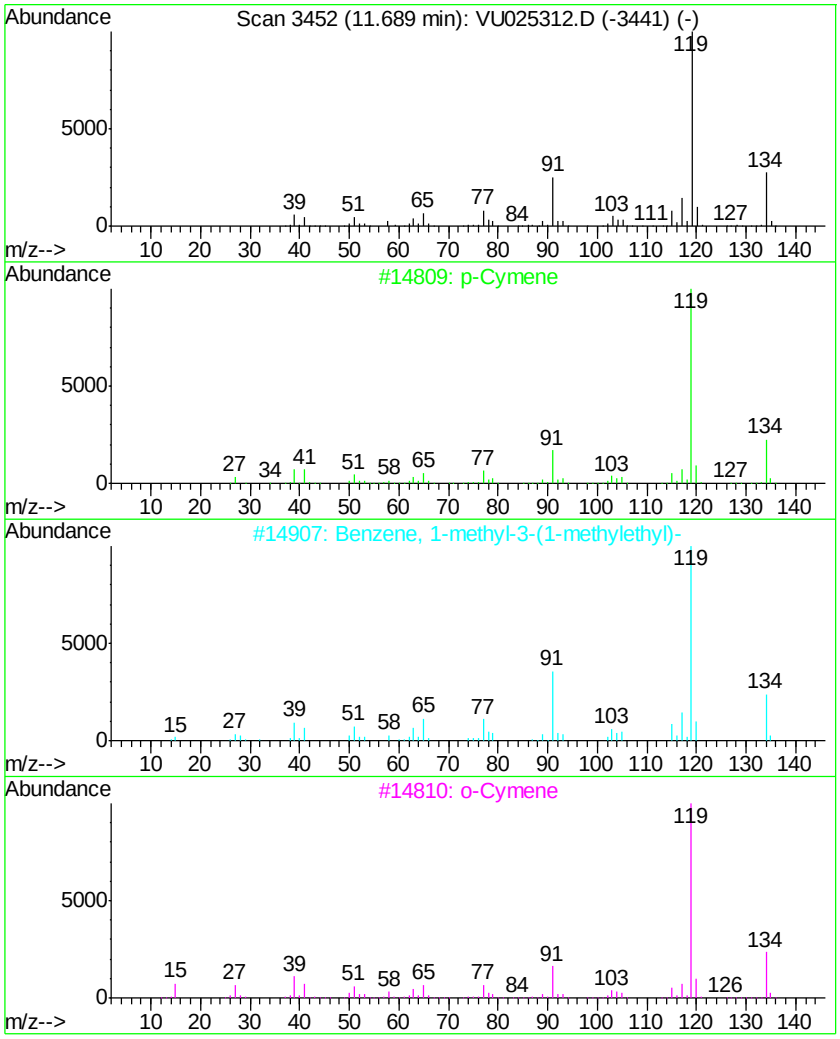
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

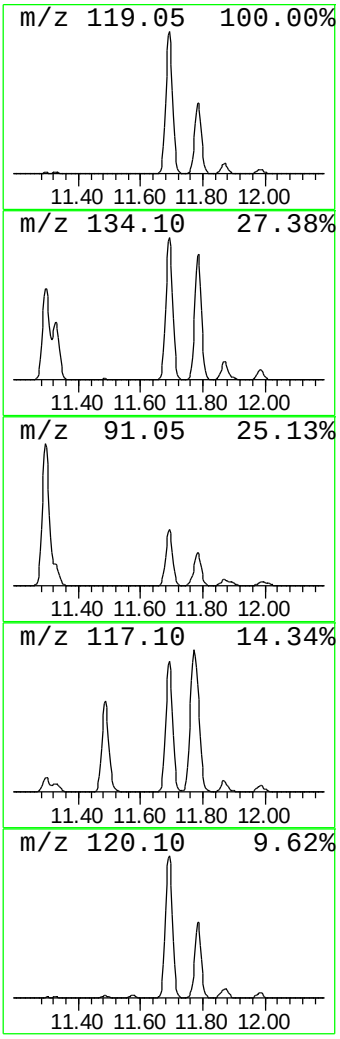
****

Peak Number 5 p-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	41.24 ug/l	829522	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
3	o-Cymene	134	C10H14	000527-84-4	94
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

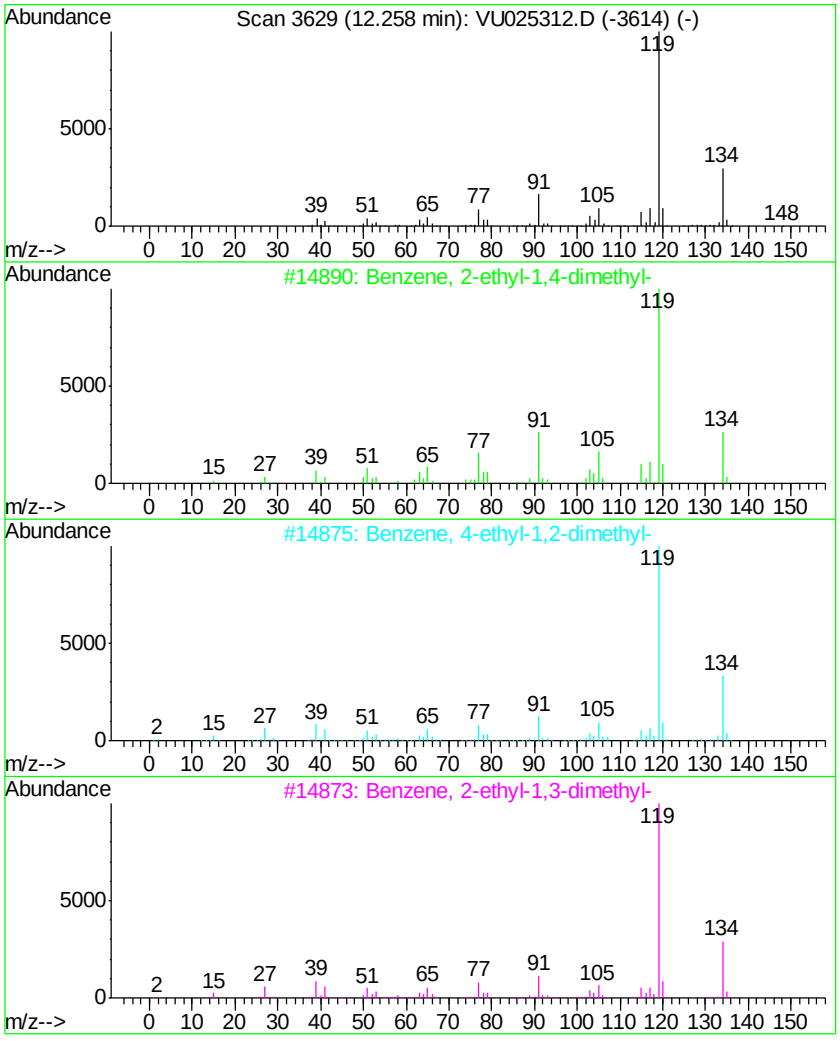
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

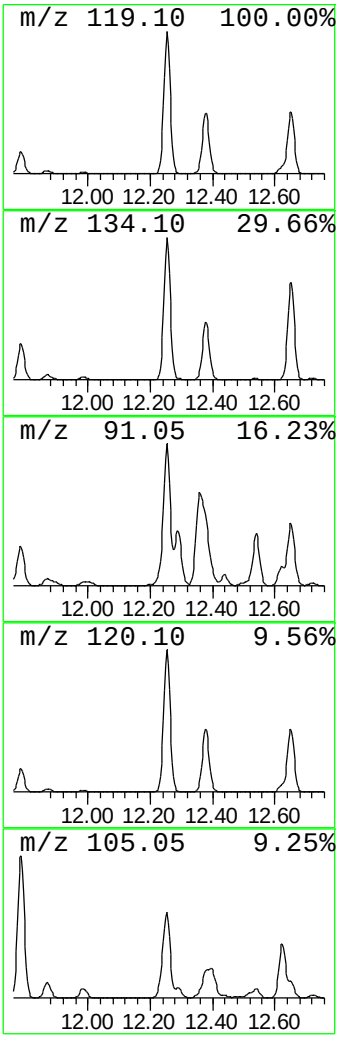
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	123.26 ug/l	2479010	1,4-Dichlorobenzene-d4	11.49

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	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	o-Cymene	134	C10H14	000527-84-4	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

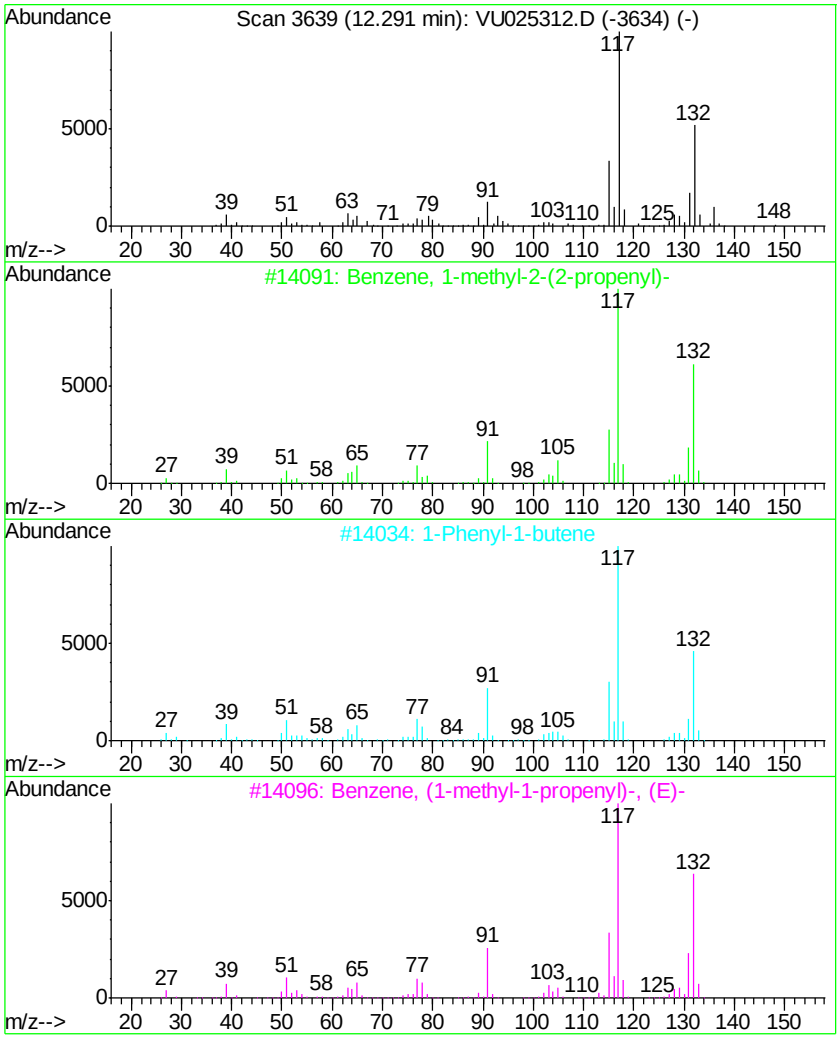
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

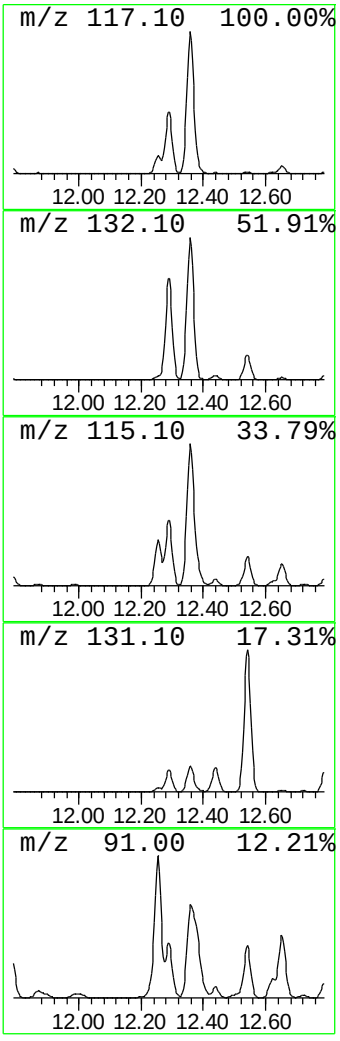
****

Peak Number 7 Benzene, 1-methyl-2-(2-prop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	54.97 ug/l	1105640	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	87
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

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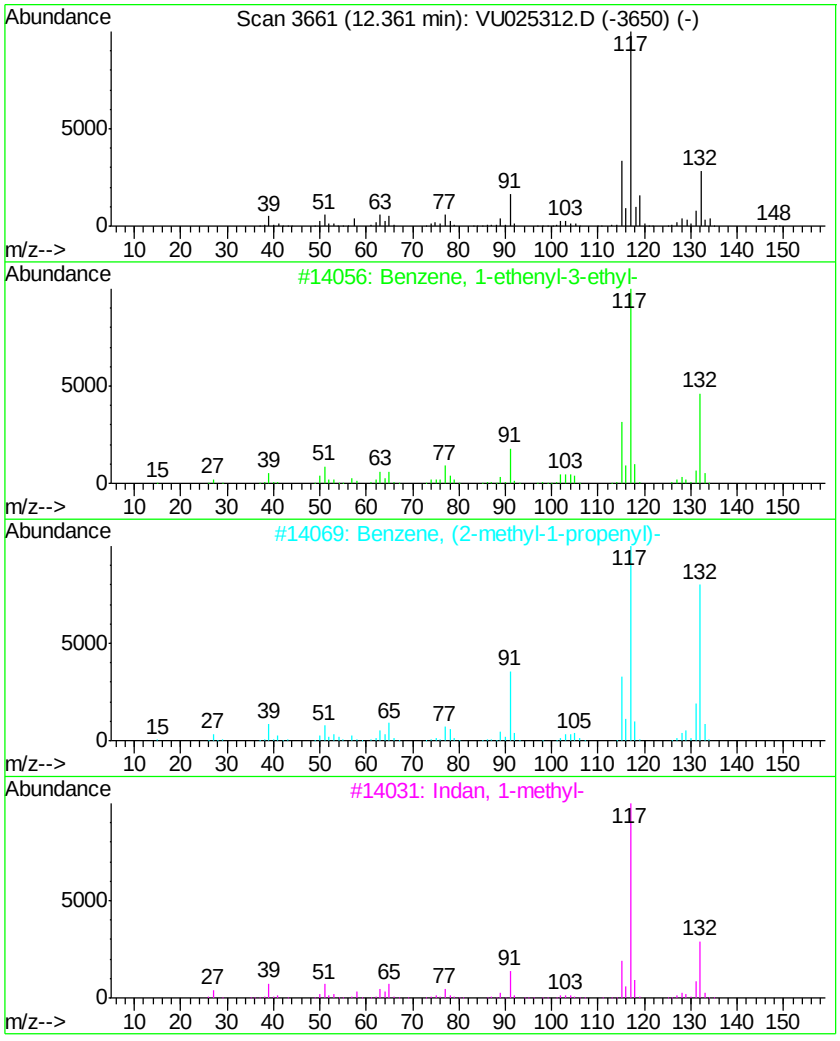
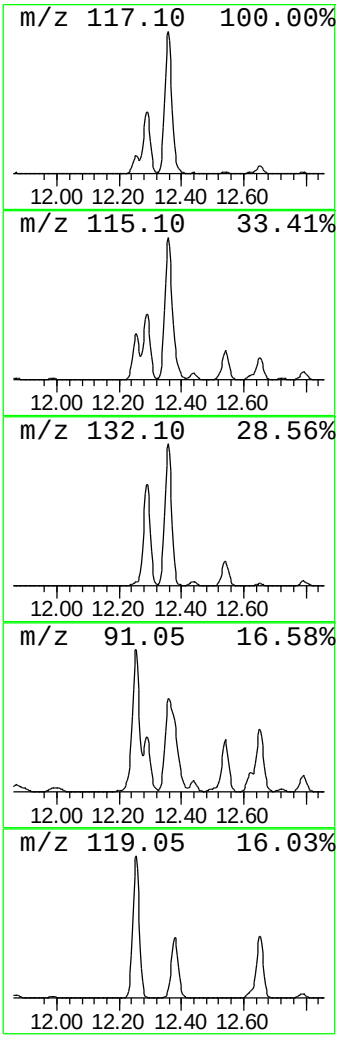
Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.36	148.61 ug/l	2988900	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

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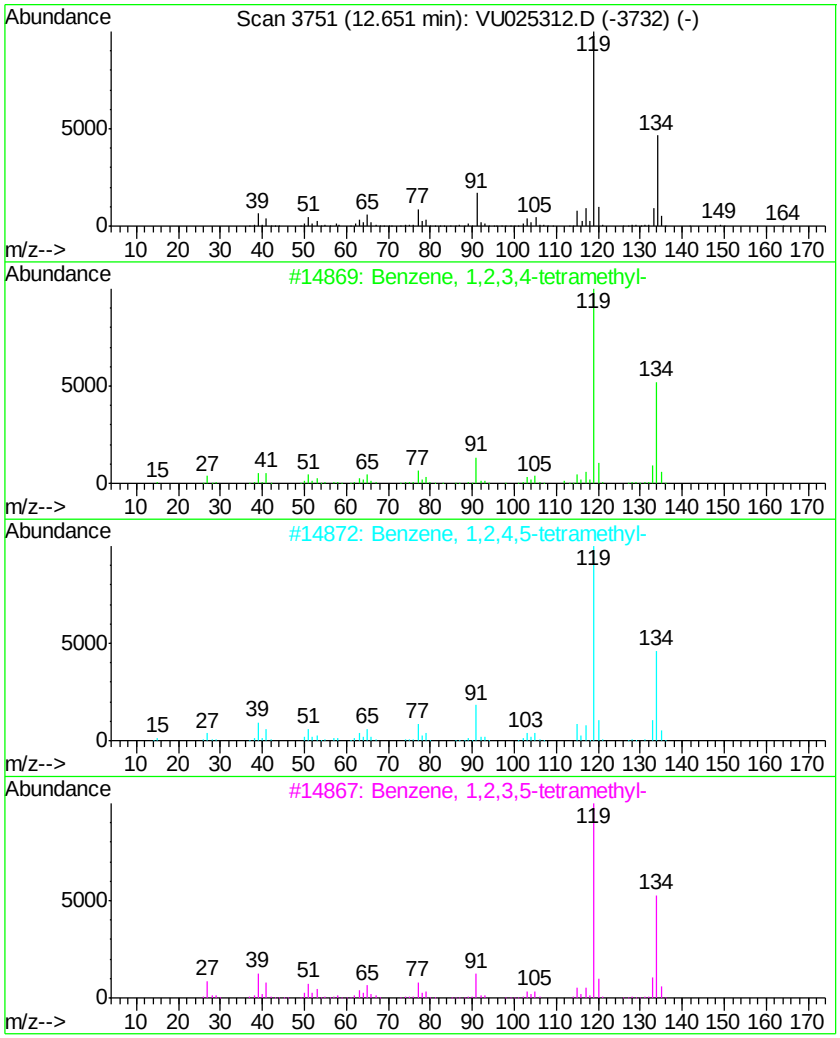
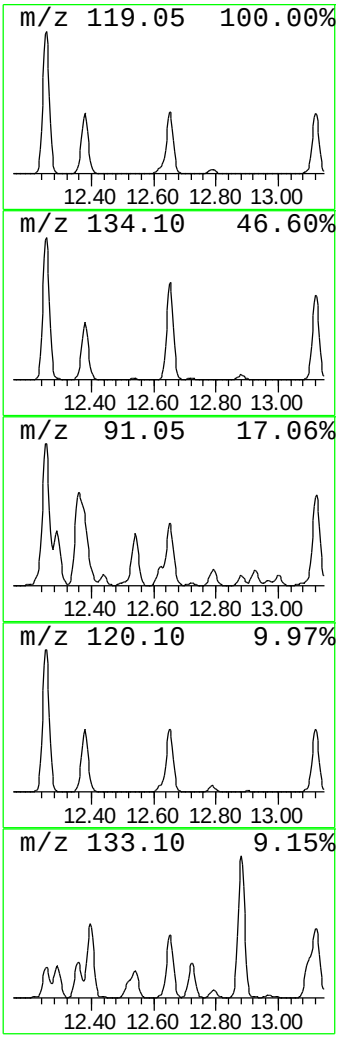
Peak Number 9 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	71.42 ug/l	1436340	1,4-Dichlorobenzene-d4	11.49

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	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.M
Quant Title : SW846 8260

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

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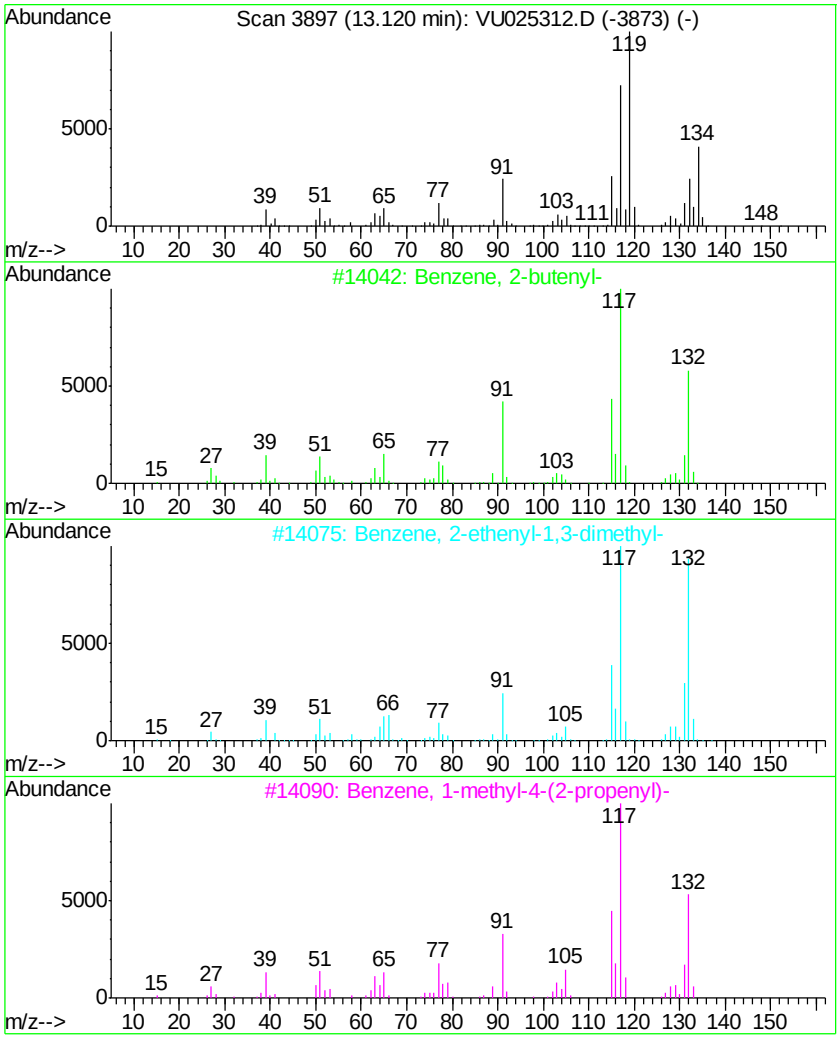
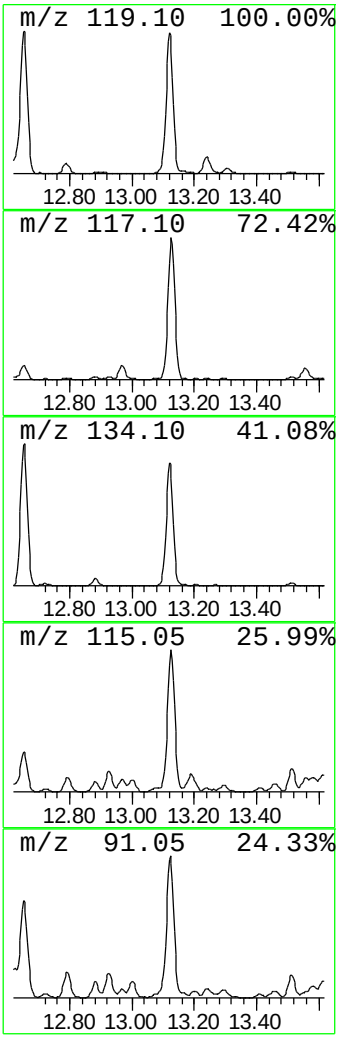
Peak Number 10 Benzene, 2-butenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	111.16 ug/l	2235580	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	64
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU071118\
Data File : VU025312.D
Acq On : 11 Jul 2018 14:14
Operator : MD/SY
Sample : J3898-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : DUP-0494-180704

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U070918W.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
Butane, 2,2,3,3-t...	5.54	148.3	ug/l	2167780	2	5.89	730943	50.0
Pentane, 2,3,4-tr...	6.93	90.0	ug/l	1315780	2	5.89	730943	50.0
Pentane, 2,3,3-tr...	7.05	178.6	ug/l	2610250	2	5.89	730943	50.0
p-Cymene	11.69	41.2	ug/l	829522	4	11.49	1005610	50.0
Benzene, 2-ethyl-...	12.26	123.3	ug/l	2479010	4	11.49	1005610	50.0
Benzene, 1-methyl...	12.29	55.0	ug/l	1105640	4	11.49	1005610	50.0
Benzene, 1-etheny...	12.36	148.6	ug/l	2988900	4	11.49	1005610	50.0
Benzene, 1,2,3,4-...	12.65	71.4	ug/l	1436340	4	11.49	1005610	50.0
Benzene, 2-butenyl-	13.12	111.2	ug/l	2235580	4	11.49	1005610	50.0