LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
Operator : MD/SY
Sample : J4344-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

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\82U072718W.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	1.088	136	155	174	rBV	337187	389453	15.12%	1.618%
2	1.760	355	364	377	rBV	84534	111060	4.31%	0.461%
3	1.950	414	423	435	rBV	110396	145961	5.67%	0.606%
4	2.307	522	534	548	rVB4	34838	76133	2.96%	0.316%
5	2.445	565	577	595	rBV2	13662	27387	1.06%	0.114%
6	2.712	643	660	663	rBV2	51730	90634	3.52%	0.377%
7	2.747	664	671	678	rVV	121019	226427	8.79%	0.941%
8	2.796	679	686	708	rVB	108439	206374	8.01%	0.857%
9	3.001	733	750	772	rBV	155321	330444	12.83%	1.373%
10	3.307	825	845	870	rBV	69791	150592	5.85%	0.626%
11	3.863	997	1018	1033	rBV3	17669	51608	2.00%	0.214%
12	4.098	1073	1091	1113	rVV	226601	613068	23.80%	2.547%
13	4.217	1113	1128	1149	rVB7	7030	26276	1.02%	0.109%
14	4.741	1272	1291	1309	rBV3	30695	82791	3.21%	0.344%
15	4.889	1320	1337	1353	rBV	195125	442593	17.18%	1.839%
16	4.995	1353	1370	1386	rVV2	426311	1078246	41.86%	4.479%
17	5.085	1386	1398	1417	rVB2	97049	240919	9.35%	1.001%
18	5.226	1425	1442	1449	rBV2	31215	72674	2.82%	0.302%
19	5.313	1458	1469	1497	rVB	191265	414945	16.11%	1.724%
20	5.487	1504	1523	1535	rBV3	72285	172244	6.69%	0.716%
21	5.558	1535	1545	1557	rVV4	44012	106736	4.14%	0.443%
22	5.628	1557	1567	1582	rVB4	34034	74964	2.91%	0.311%
23	5.889	1635	1648	1673	rBV	373470	741898	28.80%	3.082%
24	6.419	1793	1813	1838	rBV2	160412	383314	14.88%	1.592%
25	6.648	1873	1884	1899	rVB4	21239	41843	1.62%	0.174%
26	6.741	1899	1913	1928	rBV4	27479	58722	2.28%	0.244%
27	6.905	1952	1964	1987	rVB4	33946	79637	3.09%	0.331%
28	7.053	1993	2010	2021	rBV5	16299	35393	1.37%	0.147%
29	7.262	2063	2075	2085	rBV5	16995	32211	1.25%	0.134%
30	7.570	2148	2171	2202	rBV	713129	1304803	50.66%	5.420%
31	7.715	2203	2216	2230	rBV2	38075	77634	3.01%	0.323%
32	7.921	2268	2280	2299	rVB2	78013	143796	5.58%	0.597%
33	8.050	2302	2320	2331	rBV2	26218	50456	1.96%	0.210%
34	8.493	2442	2458	2467	rBV4	21418	39128	1.52%	0.163%

LSC Area Percent Report

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Sample : J4344-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

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\82U072718W.M
Title : SW846 8260

35	8.609	2484	2494	2504	rBV3	22726	36041	1.40%	0.150%
36	9.094	2632	2645	2666	rBV	567914	950850	36.92%	3.950%
37	9.191	2667	2675	2682	rVV2	19687	31450	1.22%	0.131%
38	9.252	2682	2694	2720	rVV	1614996	2575748	100.00%	10.700%
39	9.381	2721	2734	2751	rVV	854514	1351241	52.46%	5.613%
40	10.172	2969	2980	2991	rBV	124188	194728	7.56%	0.809%
41	10.310	3011	3023	3038	rBV	593792	947981	36.80%	3.938%
42	10.590	3099	3110	3124	rVB	174797	271349	10.53%	1.127%
43	10.683	3128	3139	3144	rBV	344444	549561	21.34%	2.283%
44	10.776	3158	3168	3181	rVB	202010	310469	12.05%	1.290%
45	10.975	3219	3230	3249	rBV	315643	495472	19.24%	2.058%
46	11.152	3272	3285	3299	rVB	758523	1146155	44.50%	4.761%
47	11.294	3311	3329	3333	rBV	25480	42220	1.64%	0.175%
48	11.326	3333	3339	3355	rVB	71176	108542	4.21%	0.451%
49	11.429	3362	3371	3379	rBV2	55670	84341	3.27%	0.350%
50	11.487	3379	3389	3404	rVV	788547	1240257	48.15%	5.152%
51	11.573	3405	3416	3429	rVB	195818	299802	11.64%	1.245%
52	11.770	3464	3477	3485	rBV3	313702	580162	22.52%	2.410%
53	11.811	3486	3490	3499	rVV	155330	208232	8.08%	0.865%
54	11.876	3499	3510	3529	rVB4	100019	227415	8.83%	0.945%
55	11.982	3533	3543	3555	rBV2	28426	45713	1.77%	0.190%
56	12.059	3556	3567	3579	rVB	101493	152595	5.92%	0.634%
57	12.149	3584	3595	3600	rBV	83145	132345	5.14%	0.550%
58	12.178	3600	3604	3614	rVB	84260	113130	4.39%	0.470%
59	12.255	3614	3628	3634	rBV	107725	167777	6.51%	0.697%
60	12.287	3635	3638	3649	rVB2	38098	52956	2.06%	0.220%
61	12.358	3649	3660	3668	rBV2	107436	198585	7.71%	0.825%
62	12.496	3691	3703	3710	rBV3	20599	35457	1.38%	0.147%
63	12.551	3710	3720	3732	rVB3	60563	110671	4.30%	0.460%
64	12.651	3745	3751	3760	rVV	78363	114874	4.46%	0.477%
65	12.705	3760	3768	3780	rVB	124003	194858	7.57%	0.809%
66	12.792	3784	3795	3810	rBV3	52603	92591	3.59%	0.385%
67	12.882	3811	3823	3829	rBV2	23105	37042	1.44%	0.154%
68	12.930	3829	3838	3844	rBV3	44038	72151	2.80%	0.300%
69	12.966	3844	3849	3857	rVB2	50536	65730	2.55%	0.273%
70	13.033	3864	3870	3877	rVB	22277	30088	1.17%	0.125%
71	13.088	3877	3887	3890	rBV3	44900	69620	2.70%	0.289%

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ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Title : SW846 8260

72	13.123	3890	3898	3913	rVV3	197148	361955	14.05%	1.504%
73	13.188	3915	3918	3928	rVB3	22274	32071	1.25%	0.133%
74	13.239	3928	3934	3941	rBV	25872	34803	1.35%	0.145%
75	13.290	3941	3950	3974	rVB	135123	229606	8.91%	0.954%
76	13.409	3975	3987	3994	rBV3	42135	69621	2.70%	0.289%
77	13.461	3994	4003	4010	rBV3	52378	77565	3.01%	0.322%
78	13.512	4010	4019	4026	rBV4	105374	164420	6.38%	0.683%
79	13.612	4045	4050	4057	rVB	52280	60940	2.37%	0.253%
80	13.744	4075	4091	4100	rBV	190950	324365	12.59%	1.347%
81	13.789	4100	4105	4117	rVV2	34704	55121	2.14%	0.229%
82	13.860	4117	4127	4140	rVV	80832	127882	4.96%	0.531%
83	13.956	4140	4157	4169	rVV4	66374	143445	5.57%	0.596%
84	14.017	4169	4176	4185	rVB4	27715	44163	1.71%	0.183%
85	14.155	4208	4219	4236	rBV3	44543	84386	3.28%	0.351%
86	14.352	4265	4280	4293	rBV4	76689	172550	6.70%	0.717%
87	14.480	4312	4320	4329	rVB4	19258	29849	1.16%	0.124%
88	14.544	4330	4340	4352	rVB4	30443	58405	2.27%	0.243%
89	14.679	4370	4382	4396	rBV3	79492	144070	5.59%	0.598%
90	14.882	4431	4445	4455	rBV2	86422	174951	6.79%	0.727%
91	14.940	4456	4463	4479	rVB4	23890	45919	1.78%	0.191%
92	15.065	4493	4502	4516	rBV2	62345	114570	4.45%	0.476%
93	15.593	4655	4666	4684	rVB3	13118	32630	1.27%	0.136%
94	16.065	4805	4813	4820	rBV4	17766	30610	1.19%	0.127%

Sum of corrected areas: 24072460

LSC Report - Integrated Chromatogram

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Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
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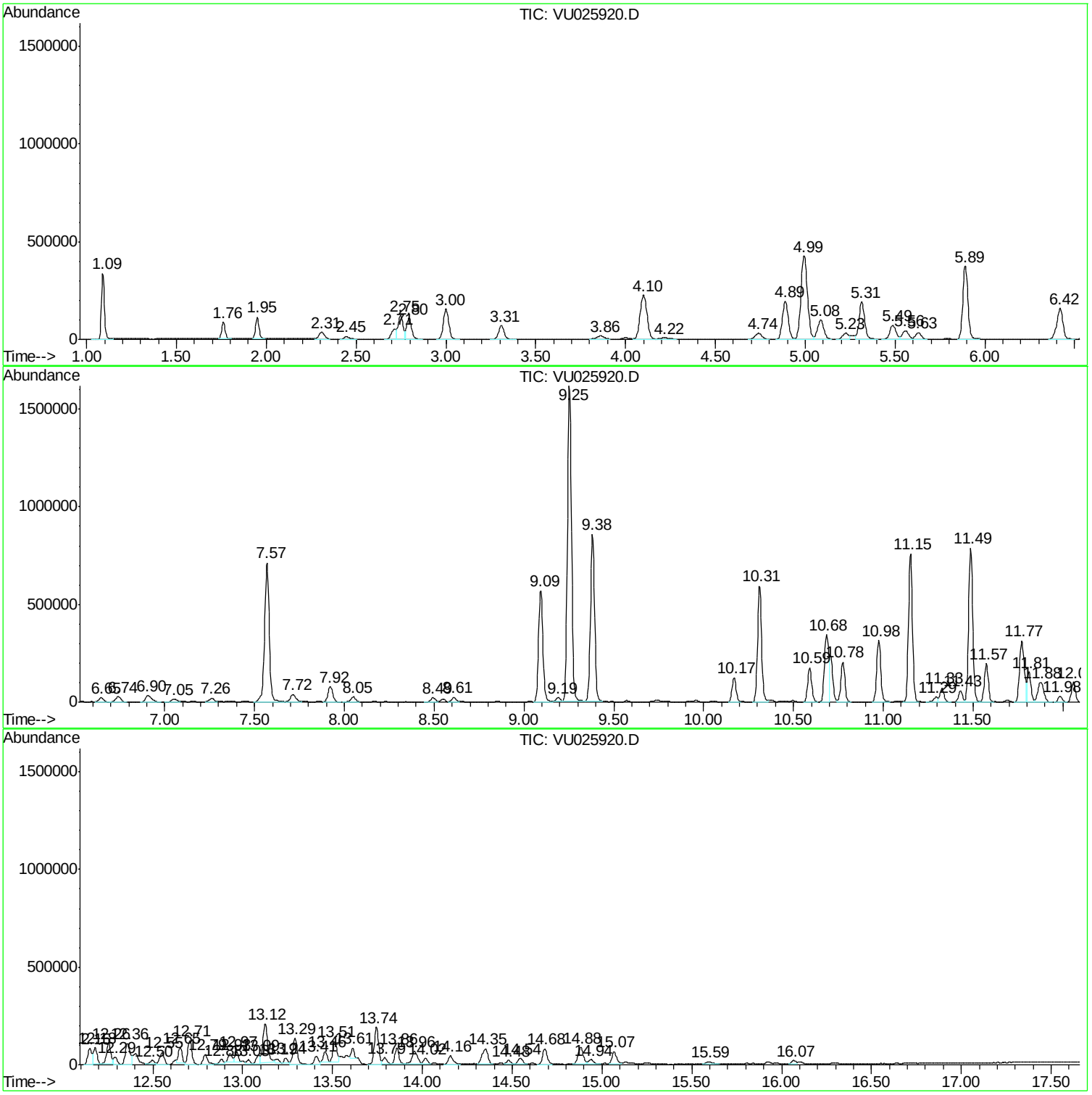
Sample : J4344-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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



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Quant Title : SW846 8260

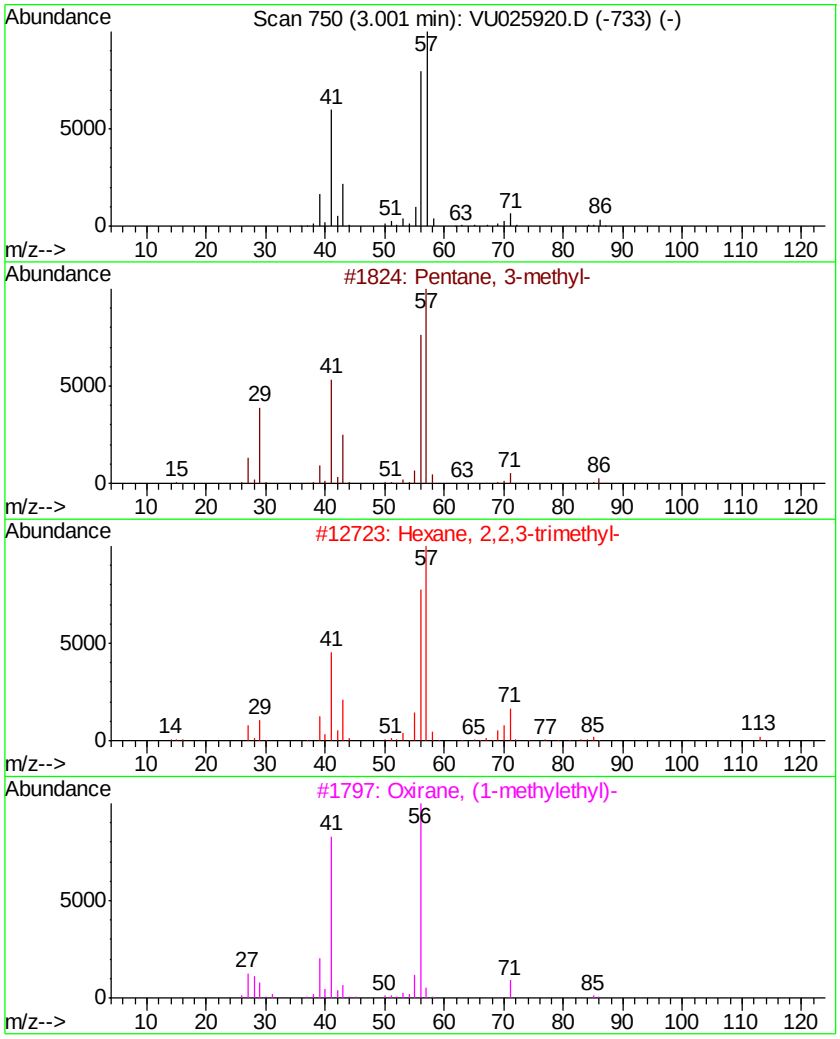
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

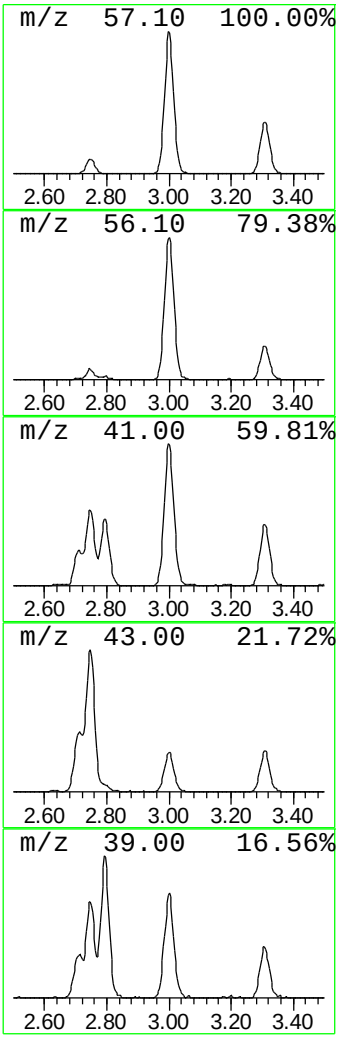
****

Peak Number 2 Pentane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	15.32 ug/l	330444	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	59
4	n-Hexane	86	C6H14	000110-54-3	56
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45





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Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
Operator : MD/SY
Sample : J4344-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

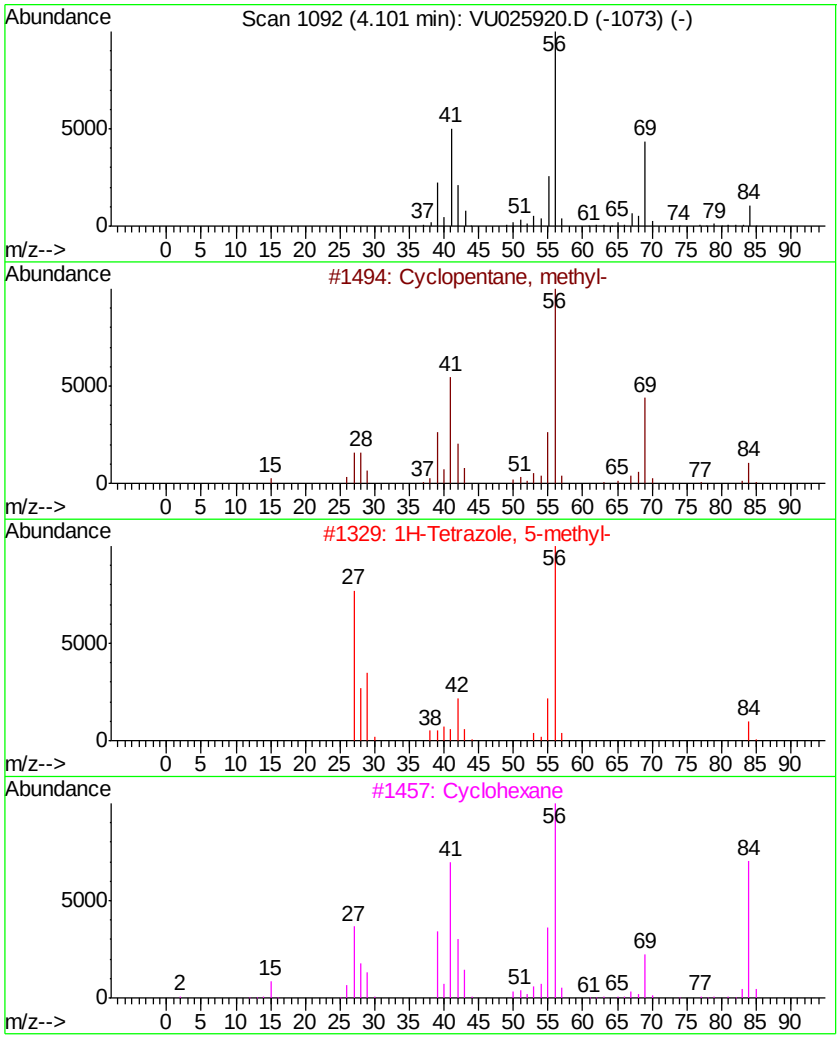
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

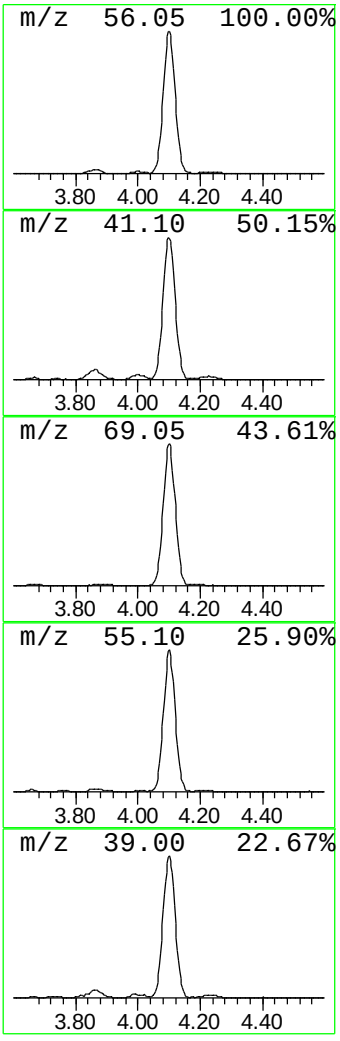
****

Peak Number 3 Cyclopentane, methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.10	28.43 ug/l	613068	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3	Cyclohexane	84	C6H12	000110-82-7	78
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5	Cyclobutane	56	C4H8	000287-23-0	53





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
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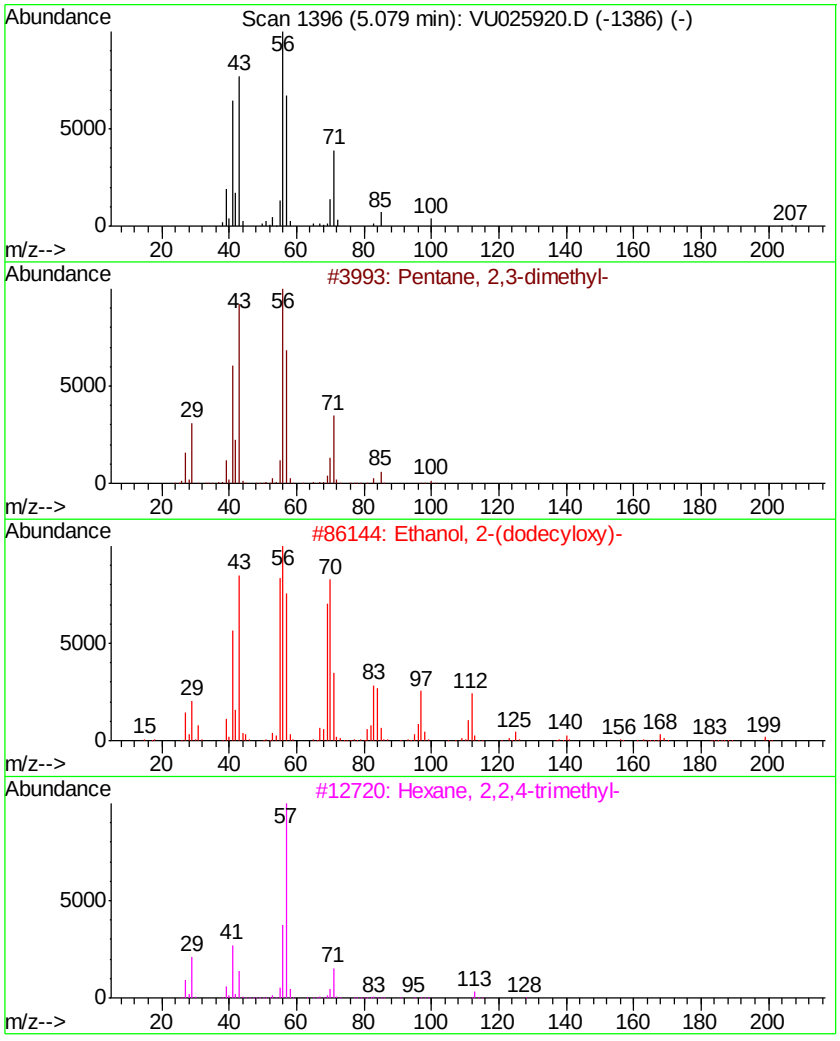
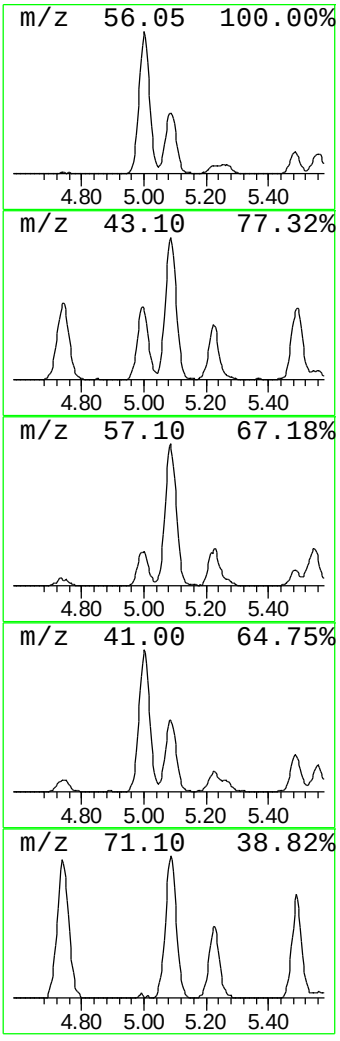
****

Peak Number 4 Pentane, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	11.17 ug/l	240919	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	64
3	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	43
5	Heptane	100	C7H16	000142-82-5	38





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ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

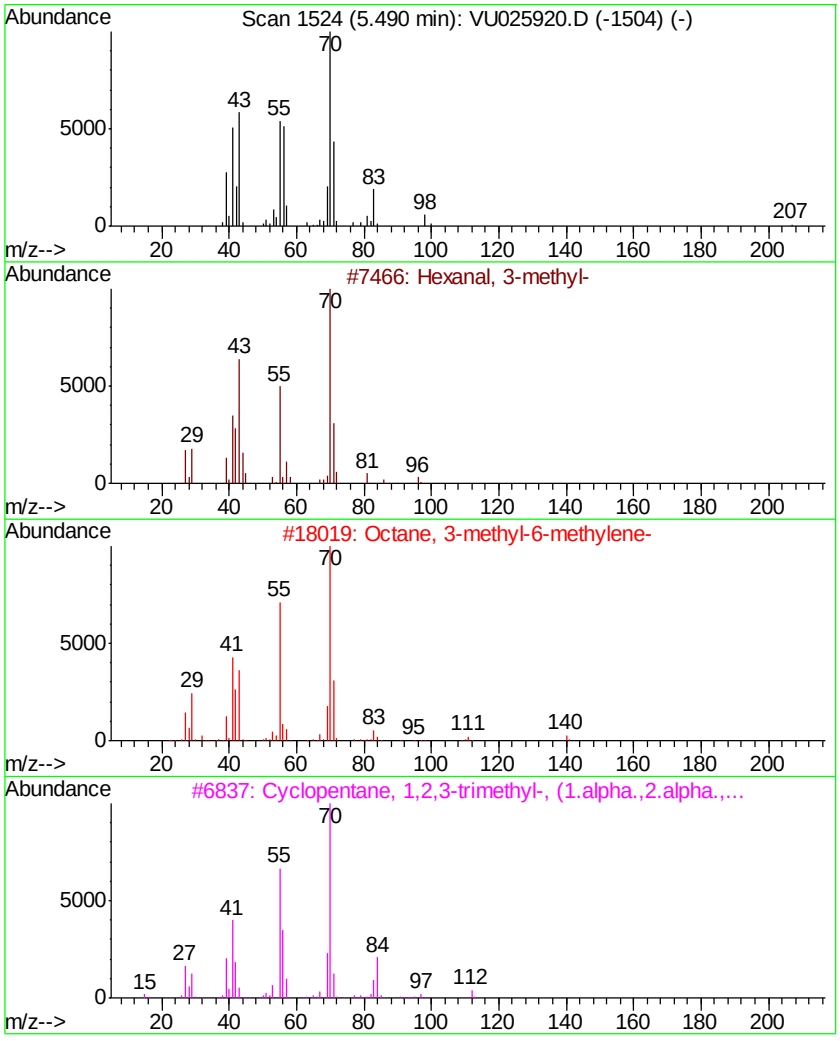
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

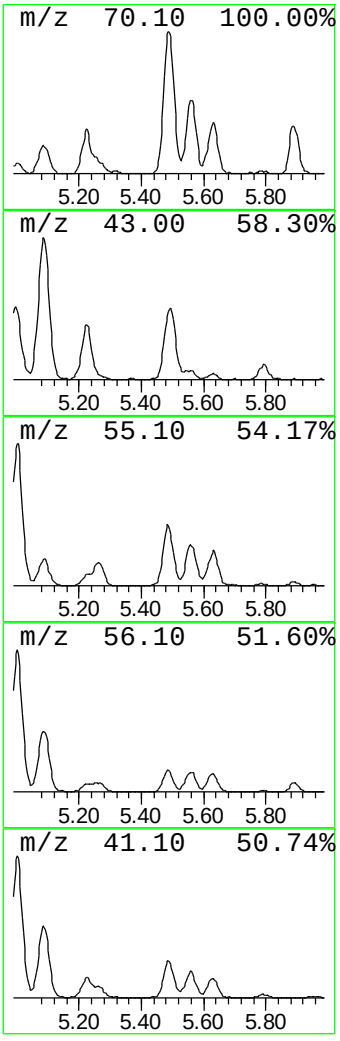
****

Peak Number 5 Hexanal, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.49	11.61 ug/l	172244	1,4-Difluorobenzene	5.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexanal, 3-methyl-	114	C7H14O	019269-28-4	59
2	Octane, 3-methyl-6-methylene-	140	C10H20	074630-07-2	59
3	Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	53
4	Cyclohexane, (1,1-dimethylpropyl)-	154	C11H22	031797-64-5	50
5	Undecane, 3-methylene-	168	C12H24	071138-64-2	50





Library Search Compound Report

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ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

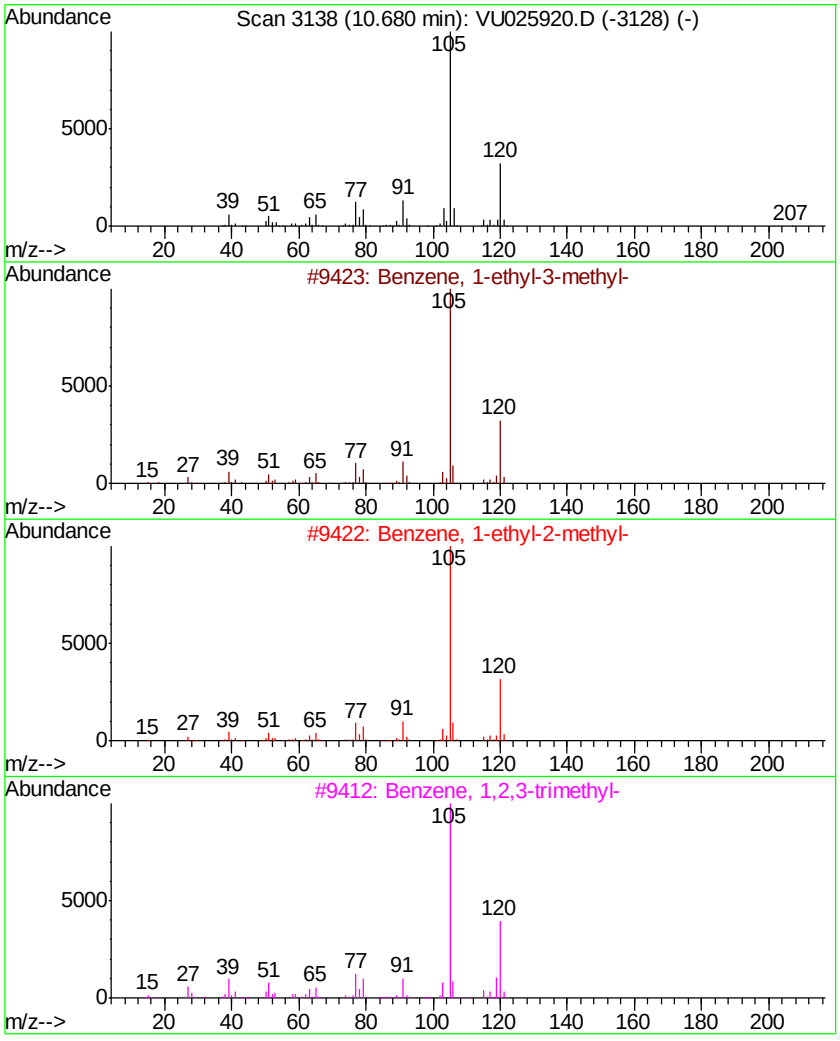
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

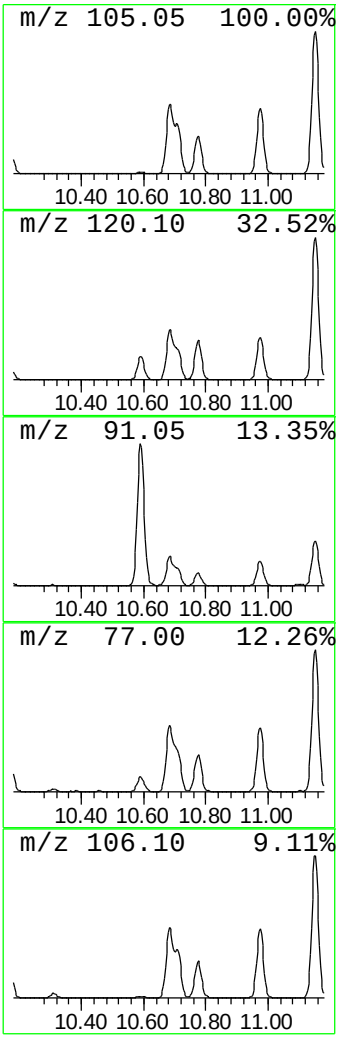
****

Peak Number 6 Benzene, 1-ethyl-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	22.16 ug/l	549561	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





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Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

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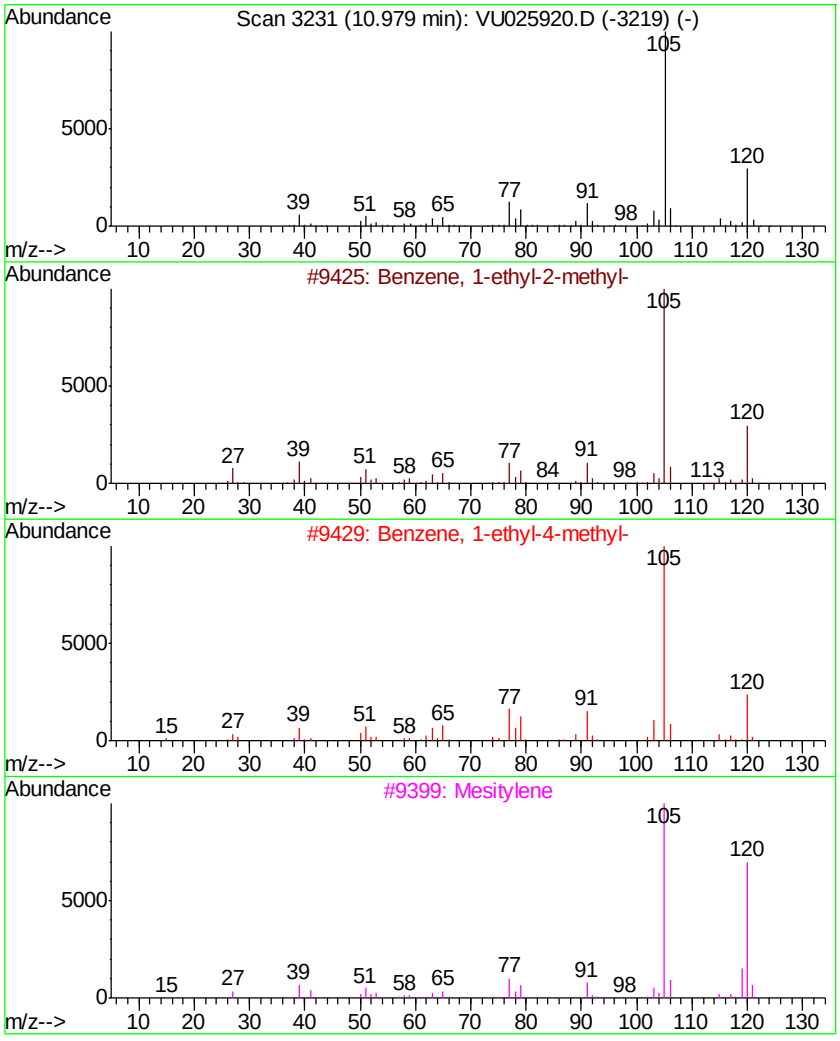
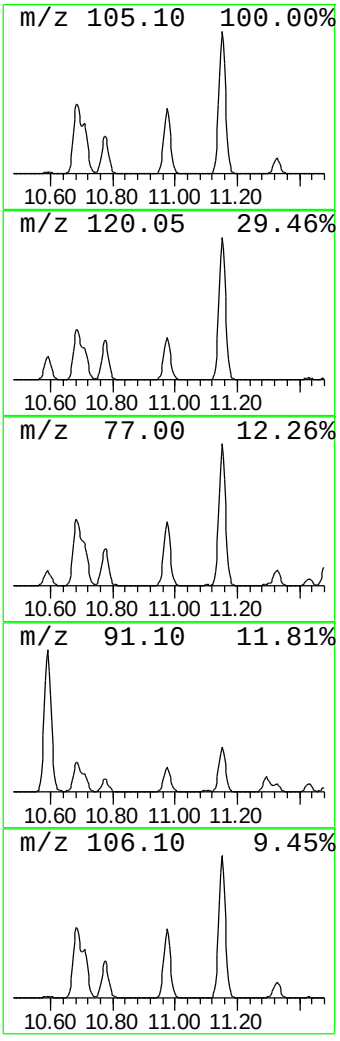
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.98	19.97 ug/l	495472	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
Operator : MD/SY
Sample : J4344-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

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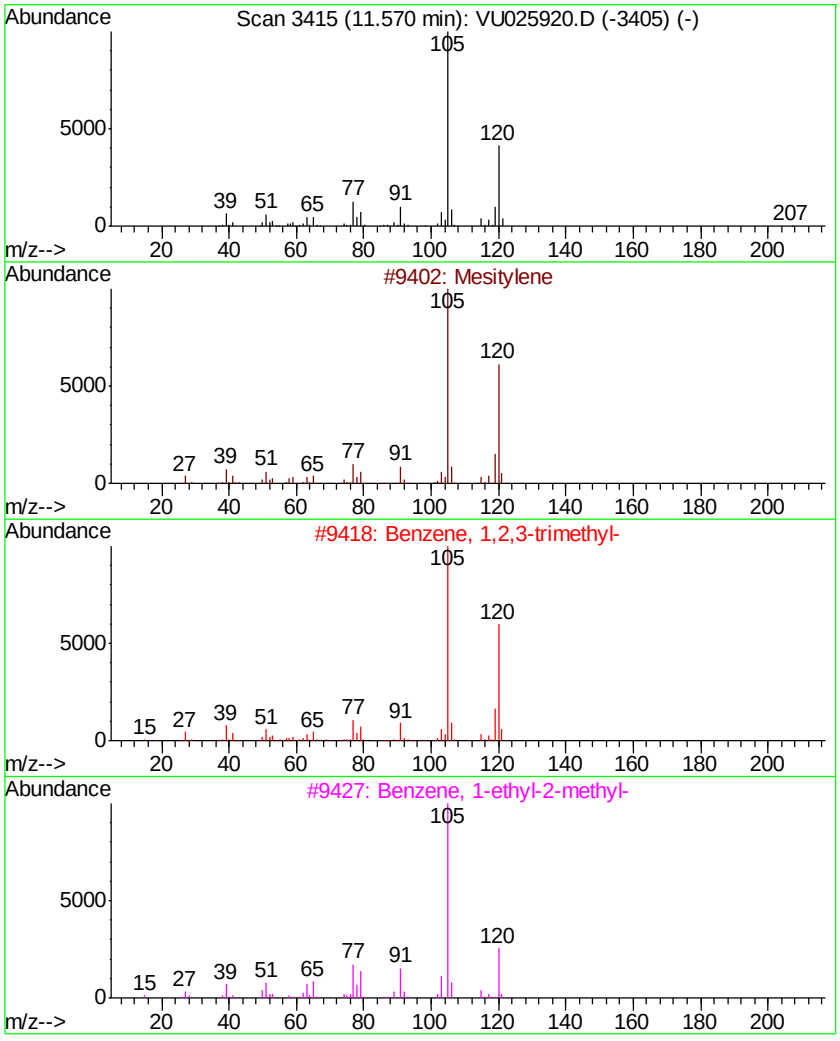
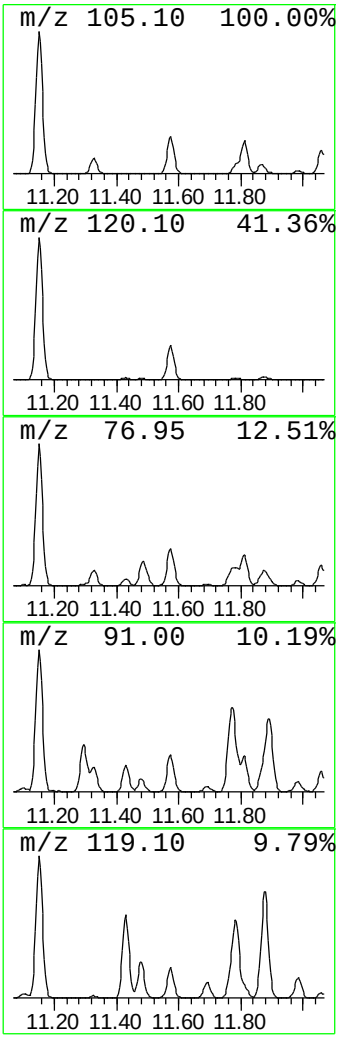
Peak Number 8 Mesitylene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	12.09 ug/l	299802	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
Operator : MD/SY
Sample : J4344-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

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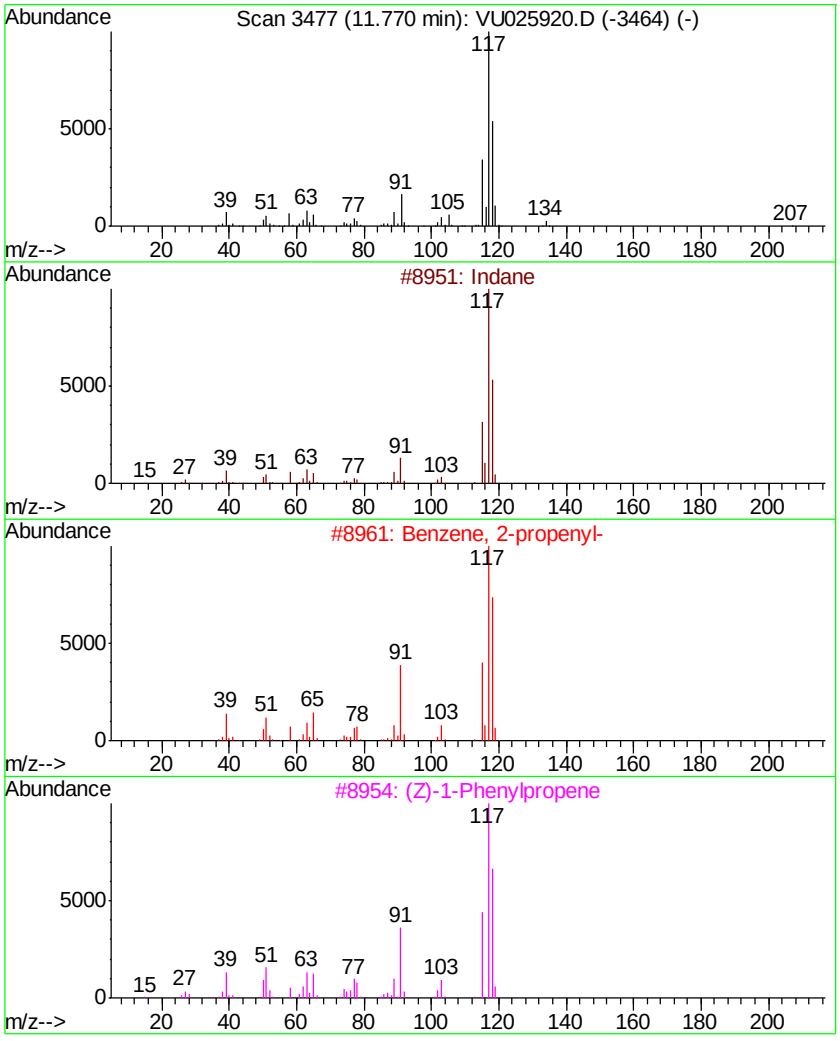
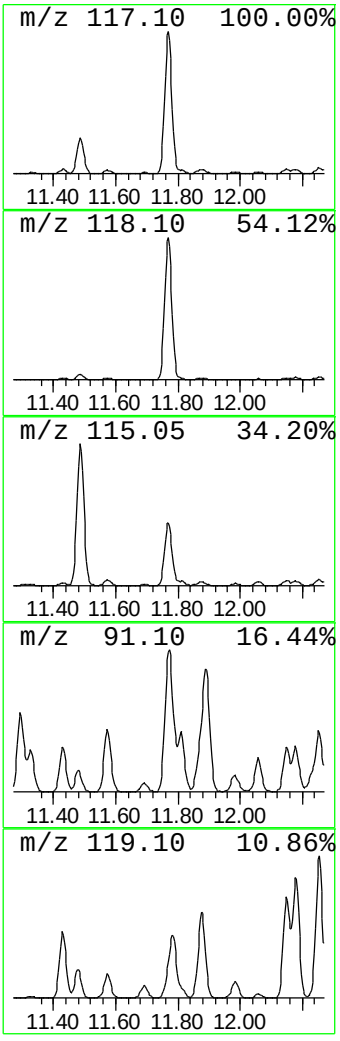
Peak Number 9 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	23.39 ug/l	580162	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
3	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	76
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
5	Deltacyclene	118	C9H10	007785-10-6	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU080618\
Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
Operator : MD/SY
Sample : J4344-09
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.M
Quant Title : SW846 8260

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

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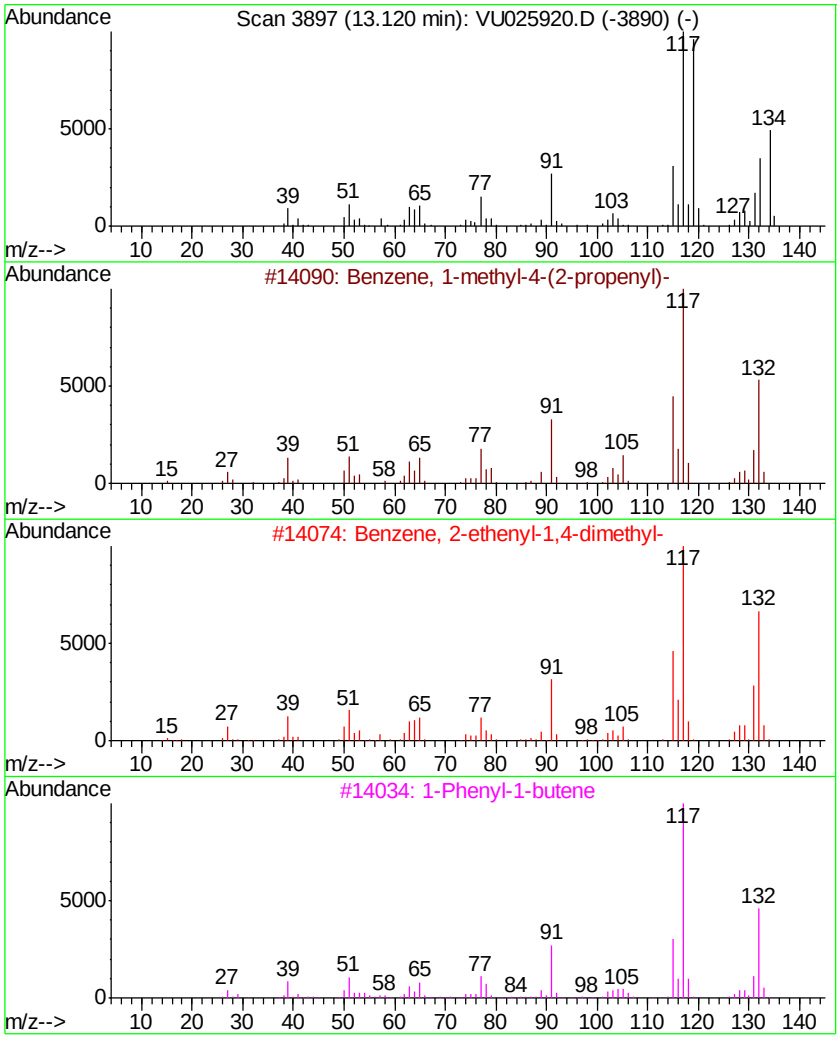
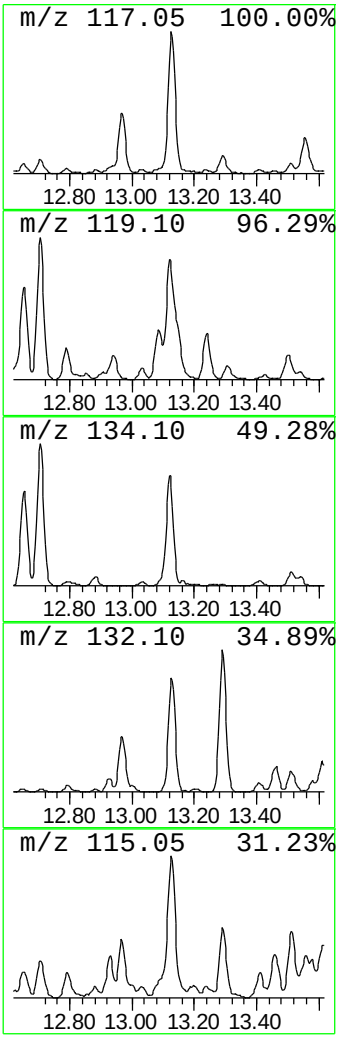
Peak Number 10 Benzene, 1-methyl-4-(2-prop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	14.59 ug/l	361955	1,4-Dichlorobenzene-d4	11.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	80
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	80
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	55
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	50
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU080618\
Data File : VU025920.D
Acq On : 07 Aug 2018 05:25
Operator : MD/SY
Sample : J4344-09
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : FL-0528

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U072718W.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-	3.00	15.3	ug/l	330444	1	4.99	1078250	50.0
Cyclopentane, met...	4.10	28.4	ug/l	613068	1	4.99	1078250	50.0
Pentane, 2,3-dime...	5.08	11.2	ug/l	240919	1	4.99	1078250	50.0
Hexanal, 3-methyl-	5.49	11.6	ug/l	172244	2	5.89	741898	50.0
Benzene, 1-ethyl-...	10.68	22.2	ug/l	549561	4	11.49	1240260	50.0
Benzene, 1-ethyl-...	10.98	20.0	ug/l	495472	4	11.49	1240260	50.0
Mesitylene	11.57	12.1	ug/l	299802	4	11.49	1240260	50.0
Indane	11.77	23.4	ug/l	580162	4	11.49	1240260	50.0
Benzene, 1-methyl...	13.12	14.6	ug/l	361955	4	11.49	1240260	50.0